## Data Abstraction \& Problem Solving with

2

## WALLS AND MIRRORS



## C++ Reserved Keywords

C++ reserves and gives predefined meanings to the following keywords. You may not redefine keywords or use them for other purposes. Keywords that appear in color are new since $\mathrm{C}++11$.

| alignas | decitype | namespace | struct |
| :---: | :---: | :---: | :---: |
| alignof | default | new | switch |
| and | delete | noexcept | template |
| and_eq | do | not | this |
| asm | doub7e | not_eq | thread_1ocal |
| auto | dynamic_cast | nul1ptr | throw |
| bitand | e1se | operator | true |
| bitor | enum | or | try |
| bool | explicit | or_eq | typedef |
| break | export | private | typeid |
| case | extern | protected | typename |
| catch | false | public | union |
| char | float | register | unsigned |
| char16_t | for | reinterpret_cast | using |
| char32_t | friend | return | virtual |
| class | goto | short | void |
| comp 1 | if | signed | volatile |
| const | inline | sizeof | wchar_t |
| const_cast | int | static | while |
| constexpr | 1ong | static_assert | xor |
| continue | mutable | static_cast | xor_eq |


| Operator | Meaning | Associativity | Usage |
| :---: | :---: | :---: | :---: |
| * | multiply | left | expr * expr |
| 1 | divide | left | expr / expr |
| \% | modulo | left | expr \% expr |
| + | add | left | expr + expr |
| - | subtract | left | expr - expr |
| << | bitwise shift left $\ddagger$ | left | expr $\ll$ expr |
| >> | bitwise shift right $\ddagger$ | left | expr >> expr |
| < | less than | left | expr $<$ expr |
| <= | less than or equal to | left | expr $<=$ expr |
| > | greater than | left | expr $>$ expr |
| >= | greater than or equal to | left | expr $>=$ expr |
| = $=$ | equal | left | expr $=$ expr |
| ! $=$ | not equal | left | expr ! = expr |
| \& | bitwise AND | left | expr \& expr |
| $\wedge$ | bitwise EXCLUSIVE OR | left | expr $\wedge$ expr |
| I | bitwise OR | left | expr \| expr |
| \& \& | logical AND | left | expr \& \& expr |
| 11 | logical OR | left | expr \|| expr |
| ? | conditional | left | expr ? expr : expr |
| = | assign | left | lvalue $=$ expr |
| *= | multiply and assign | left | lvalue = expr |
| /= | divide and assign | left | lvalue /= expr |
| \%= | modulo and assign | left | $l v a l u e \%=$ expr |
| += | add and assign | left | $l$ lvalue $+=$ expr |
| -= | subtract and assign | left | lvalue $=$ expr |
| <<= | shift left and assign | left | lvalue <<= expr |
| >>= | shift right and assign | left | lvalue >>= expr |
| \& $=$ | AND and assign | left | lvalue $\&=$ expr |
| \|= | OR and assign | left | lvalue \|= expr |
| $\wedge=$ | EXCLUSIVE OR and assign | left | lvalue $\wedge=$ expr |
|  |  |  |  |
| , | comma | left | expr, expr |

$\ddagger$ Typically overloaded for I/O

# Data Abstraction \& Problem Solving with C++ 

## WALLS AND MIRRORS

## SIXTH EDITION



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> University of Rhode Island

## Timothy Henry <br> University of Rhode Island

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Library of Congress Cataloging-in-Publication Data on file. the publication of the first edition, we all have gained experience with teaching data abstraction in an objectoriented way using $\mathrm{C}++$. This edition reflects that experience and the many comments and suggestions received from faculty and students alike.

I am happy to introduce Dr. Timothy Henry, my co-author and colleague at the University of Rhode Island. Together, we have given this book a much needed revision and a new look. However, our goal remains to give students a superior foundation in data abstraction, object-oriented programming, and other modern problemsolving techniques. All $\mathrm{C}++$ code has been rewritten with a focus on safe and secure programming practices. It also adheres to the $\mathrm{C}++11$ standard.

We hope that you enjoy reading this book. Like many others before you, you can learn-or teach-data structures in an effective and sustainable way.

## Talk to Us

Walls and Mirrors continues to evolve. Your comments, suggestions, and corrections will be greatly appreciated. Here are a few ways to reach us:

- E-mail: carrano@acm.org
- Facebook: www.facebook.com/makingitreal
- Twitter: twitter.com/Frank_M_Carrano
- Blog: frank-m-carrano.com/makingitreal

The topics that we cover in this book deal with the various ways of organizing data so that a given application can access and manipulate data in an efficient way. These topics are fundamental to your future study of computer science, as they provide you with the foundation of knowledge required to create complex and reliable software. Whether you are interested in designing video games or software for robotic-controlled surgery, the study of data structures is vital to your success. Even if you do not study all of the topics in this book now, you are likely to encounter them later. We hope that you will enjoy reading the book, and that it will serve as a useful reference tool for your future courses.

The walls and mirrors in the title represent two fundamental problem-solving techniques that appear throughout the presentation. Data abstraction isolates and hides the implementation details of a module from the rest of the program, much as a wall can isolate and hide you from your neighbor. Recursion is a repetitive technique that solves a problem by solving exactly the same but smaller problems, much as images in facing mirrors grow smaller with each reflection.

Please be sure to browse the rest of this preface to see the features that will help you in your studies. To help you learn and to review for exams, we have included such learning aids as video tutorials (VideoNotes), checkpoint questions with answers, margin notes, programming tips, chapter summaries, and a glossary. As a help during programming, you will find $\mathrm{C}++$ reference material in the appendices and inside the covers. You should review the list of this book's features given later in this preface in the section "Features to Enhance Learning."

The presentation makes some basic assumptions about your knowledge of $\mathrm{C}++$. Some of you may need to review this language or learn it for the first time by consulting the appendices of this book. This book covers C++ classes and other relevant aspects of the language in new C++ Interludes that occur throughout the book between certain chapters. These interludes do not assume that you already know their topics. We assume no experience with recursive functions, which are included in Chapters 2 and 5.

All of the $\mathrm{C}++$ source code that appears in this book is available for your use. Later in this preface, the description of supplementary materials tells you how to obtain these files, as well as the VideoNotes and other online documents.

## Organization

This book's organization, sequencing, and pace of topic coverage make learning and teaching easier by focusing your attention on one concept at a time, by providing flexibility in the order in which you can cover topics, and by clearly distinguishing between the specification and implementation of abstract data types, or ADTs. To accomplish these goals, we have organized the material into 21 chapters. Most chapters focus on either the specification and use of an ADT or its various implementations. You can choose to cover the specification of an ADT followed by its implementations, or you can treat the specification and use of several ADTs before you consider any implementation issues. The book's organization makes it easy for you to choose the topic order that you prefer.

## Table of Contents at a Glance

The following list shows the overall composition of the book. A further chapter-by-chapter description appears later. Note that gray highlighted sections are available online.

| Chapter 1 | Data Abstraction: The Walls |
| :--- | :--- |
| C++ Interlude 1 | C++ Classes |
| Chapter 2 | Recursion: The Mirrors |
| Chapter 3 | Array-Based Implementations |
| C++ Interlude 2 | Pointers, Polymorphism, and Memory Allocation |
| Chapter 4 | Link-Based Implementations |
| Chapter 5 | Recursion as a Problem-Solving Technique |
| Chapter 6 | Stacks |
| C++ Interlude 3 | Exceptions |
| Chapter 7 | Stack Implementations |
| Chapter 8 | Lists |
| Chapter 9 | List Implementations |
| Chapter 10 | Algorithm Efficiency |
| Chapter 11 | Sorting Algorithms and Their Efficiency |
| C++ Interlude 4 | Class Relationships and Reuse |
| Chapter 12 | Sorted Lists and Their Implementations |
| Chapter 13 | Queues and Priority Queues |
| Chapter 14 | Queue Implementations |
| C++ Interlude 5 | Overloaded Operators and Friend Access |
| Chapter 15 | Trees |
| Chapter 16 | Tree Implementations |
| C++ Interlude 6 | Iterators |
| Chapter 17 | Heaps |
| Chapter 18 | Dictionaries and Their Implementations |
| Chapter 19 | Balanced Search Trees |
| Chapter 20 | Graphs |
| Chapter 21 | Processing Data in External Storage |
| C++ Interlude 7 | The Standard Template Library |
| Appendix A | Review of C++ Fundamentals |
| Appendix B | Important Themes in Programming |
| Appendix C | The Unified Modeling Language |
| Appendix D | The Software Life Cycle |
| Appendix E | Mathematical Induction |
| Appendix F | Algorithm Verification |
| Appendix G | Files |
| Appendix H | C++ Header Files and Standard Functions |
| Appendix I | C++ Documentation Systems |
| Appendix J | ASCII Character Codes |
| Appendix K | C++ for Java Programmers |
| Appendix L | C++ for Python Programmers |
| Cer |  |

## What's New?

This edition of Walls and Mirrors is a significant revision of the previous edition, yet remains committed to a pedagogical approach that makes the material accessible to students at the introductory level. Although everything looks new, you will find the coverage that you enjoyed in previous editions is still here. At a glance, the book has more - but shorter - chapters, a second color, and new C++ Interludes. Let's examine the details.

Organization. The book begins with a shorter Chapter 1, so that it can focus on the specification of abstract data types (ADTs). After general discussions of problem solving, good programming practices, and ADTs, we specify a simple ADT - the bag. We define the bag's operations within a C++ template interface in a non-threatening way. We have moved some sections from the original first chapter to the appendices.

By introducing the bag as the first ADT we consider, we make the difficult topic of linked data more accessible to students. Adding or removing the first node in a chain of linked nodes is the easiest task, and these simple manipulations are the ones we need to use for a linked implementation of the bag. The next ADT that we consider is the stack, a more useful data container that has the same simple chain in one of its definitions. Moreover, many students are already familiar with stacks. Later, the treatment of lists looks at the more involved operations of adding and removing a node that lies between existing nodes of a chain.

The rest of the coverage will be familiar to previous users of Walls and Mirrors but often you will find ADTs presented in two chapters rather than one. These chapters separate the discussions of specification and implementation. We will describe each chapter later in this preface. To summarize, this organization

- Replaces Chapter 1 with an introduction to ADTs and template interfaces using the ADT bag.
- Provides a more focused introduction to array-based implementations and link-based implementations using the ADT bag.
- Makes the topic of linked data more accessible to students by discussing it progressively as we introduce the ADTs bag, stack, and list.
- Places greater emphasis on data abstraction as a problem solving tool.
- Enhances the separation between specification and implementation of basic ADTs by placing them in successive chapters.
- Specifies ADTs informally at first, then in UML, and ultimately in C++ template interfaces.
- Demonstrates safe and secure programming practices in completely revised code that adheres to the $\mathrm{C}++11$ standard.
- Covers the ADT heap in its own chapter.
- Reorganizes the coverage of the ADT dictionary (formerly called the ADT table) into two chapters.

C++ Interludes. The introduction of C++ classes and other aspects of C++ that we need for our presentation and were previously covered in Chapters 3,4 , and 8 of the previous edition are now featured in $\mathrm{C}++$ Interludes. Seven of these "mini-chapters" appear throughout the book to cover relevant C++ topics as we need them. Note that these interludes separate the details of $\mathrm{C}++$ from the discussion of data structures.

VideoNotes. Online tutorials are a Pearson feature that provides visual and audio support to the presentation given throughout the book. They offer students another way to recap and reinforce key concepts. VideoNotes allow for self-paced instruction with easy navigation, including the ability to select, play, rewind, fast-forward, and stop within each video. Unique VideoNote icons appear throughout this book whenever a video is available for a particular concept or problem. A detailed list of the 49 VideoNotes for this text and their associated locations in the book can be found on page xxiii. VideoNotes are free with the purchase of a new textbook. To purchase access to VideoNotes, please go to

Other features. Walls and Mirrors now has many new features that enhance its usefulness to both readers and instructors. This edition

- Adds a second color to enhance the effectiveness of the illustrations, distinguish pseudocode from C++ code, and provide visual interest.
- Includes Notes and Programming Tips to emphasize key material and offer programming advice.
- Distinguishes major pieces of code in a new Listing element.
- Replaces the self-test exercises at the ends of chapters with Checkpoint Questions placed throughout the chapters.
- Numbers sections and subsections for easy reference.
- Includes transition guides from Python to C++ and Java to C++.


## Features to Enhance Learning

The pedagogical features and organization of this book were carefully designed to facilitate learning and to allow instructors to tailor the material easily to a particular course. These features help students not only during their first reading of the material, but also during subsequent review.

Notes Important ideas are presented or summarized in highlighted paragraphs and are meant to be read in line with the surrounding text.


Programming Tips Suggestions to improve or facilitate programming are featured as soon as they become relevant.

Examples Numerous examples illuminate new concepts.

Checkpoint Questions Questions are posed throughout each chapter, integrated within the text, that reinforce the concept just presented. These "checkpoint" questions help readers to understand the material, since answering them requires pause and reflection. Solutions to these questions are provided online.

VideoNote
VideoNotes Online tutorials provide additional instruction in a more dynamic form than a static textbook.

Margin Notes Brief phrases in the margins help you review material or locate particular content.

Chapter Summaries Each chapter ends with a list of key ideas that summarize what was presented.

Glossary of Terms A glossary of all terms introduced in this book is available online.

Exercises and Programming Projects Further practice is available by solving the exercises and programming projects at the end of each chapter. Unfortunately, we cannot give readers the answers to these exercises and programming projects, even if they are not enrolled in a class. Only instructors who adopt the book can receive selected answers from the publisher. For help with these exercises and projects, you will have to contact your instructor.

## Accessing Instructor and Student Resource Materials

The following items are available on the publisher's website at

```
www.pearsonhighered.com/carrano
```

- $\mathrm{C}++$ code as it appears in the book
- A link to any misprints that have been discovered since the book was published
- Links to additional online content, which is described next


## Instructor Resources

The following protected material is available to instructors who adopt this book by logging onto Pearson's Instructor Resource Center, accessible from
www. pearsonhighered.com/carrano

- PowerPoint lecture slides
- Test bank
- Instructor solutions manual
- Figures from the book

Additionally, instructors can access the book's Companion Website for the following online premium content, also accessible from
www. pearsonhighered.com/carrano

- Instructional VideoNotes
- Answers to the Checkpoint Questions
- A glossary of terms

Please contact your Pearson sales representative for an instructor access code. Contact information is available at www. pearsonhighered.com/replocator.

## Student Resources

The following material is available to students by logging onto the book's Companion Website accessible from www. pearsonhighered.com/carrano:

- Instructional VideoNotes
- Answers to the Checkpoint Questions
- A glossary of terms

Students must use the access card located in the front of the book to register for and then enter the Companion Website. Students without an access code can purchase access from the Companion Website by following the instructions listed there.

## Chapter Overview

$\mathbf{R}_{\text {e }}$eaders of this book should have completed a programming course, preferably in C++. Appendix A covers the essentials of $\mathrm{C}++$ that we assume readers will know. You can use this appendix as a review or as the basis for making the transition to $\mathrm{C}++$ from another programming language. Note that Appendices K and L offer some help for those who are transitioning from Java or Python, respectively.

- Chapters 1 through 5: Chapter 1 introduces object-oriented concepts and focuses on the specification of abstract data types (ADTs). Our ADTs now store data whose data type is chosen by the client. To accomplish this, each specification of an ADT includes a $C++$ template interface. As an example, Chapter 1 introduces the ADT bag. Much of the software engineering material that was in Chapter 1 is now in the appendices.

Next is C++ Interlude 1, which presents C++ classes. It gives more details about template interfaces-like the one presented in Chapter 1-and shows how to use inheritance to define a class derived from an interface.

As it did in earlier editions, Chapter 2 introduces recursion, and Chapter 5 develops it as a problemsolving tool. Recursion is revisited often throughout the book.

We clearly separate the specification, use, and implementation of the bag by dividing the material across several chapters. For example, Chapter 1 specifies the bag and provides several examples of its use. Chapter 3 covers implementations that use arrays. Just before Chapter 4 introduces chains of linked nodes and uses one in the definition of a class of bags, C++ Interlude 2 covers pointers, polymorphism, and dynamic memory allocation. Both Chapters 3 and 4 include recursion in their presentation.

In a similar fashion, we separate specification from implementation throughout most of the book when we discuss various other ADTs. You can choose to cover the chapters that specify and use the ADTs and then later cover the chapters that implement them. Or you can cover the chapters as they appear, implementing each ADT right after studying its specification and use. A list of chapter prerequisites appears later in this preface to help you plan your path through the book.

Chapter 3 does more than simply implement the ADT bag. It shows how to approach the implementation of a class by initially focusing on core methods. When defining a class, it is often useful to implement and test these core methods first and to leave definitions of the other methods for later. Chapter 4 follows this approach in its development of a link-based implementation of the ADT bag.

- Chapters 6 and 7: Chapter 6 discusses stacks, giving examples of their use, and Chapter 7 implements the stack using an array and then again using a chain. Between these chapters is $\mathrm{C}++$ Interlude 3, which discusses C++ exceptions. Chapter 7 then shows how to use an exception in the implementation of the ADT stack, when a client violates a method's precondition.
- Chapters 8 and 9: The next two chapters introduce the ADT list. We discuss this container abstractly and then implement it by using an array and then a chain of linked nodes. Once again, we use exceptions to enforce method preconditions.
- Chapters 10 and 11: Chapter 10 introduces the complexity of algorithms, a topic that we integrate into future chapters. Chapter 11 discusses various sorting techniques and their relative complexities. We consider both iterative and recursive versions of these algorithms.
- Chapter 12: Chapter 12 introduces the sorted list, looking at a linked implementation and its efficiency. We then talk about the relationship between a list and a sorted list and show how to use the list as a base class for the sorted list. Note that Chapter 12 is preceded by C++ Interlude 4 that discusses class relationships and the various ways a class can be reused. Chapter 12 puts that discussion into immediate use.
- Chapters 13 and 14: Chapter 13 presents the ADTs queue and priority queue, along with some uses of these containers. In doing so, we give an example of simulation that uses both ADTs, and finally summarize the difference between position oriented and value oriented ADTs. Chapter 14 implements the queue, and introduces tail pointers, circularly linked chains, and circular arrays. We offer an implementation of a priority queue by using a sorted list, but note that a better approach will come later when we introduce the ADT heap.
- Chapters 15 through 17: Before we begin the next chapter, C++ Interlude 5 introduces overloaded operators and friend access. We overload operators when we define classes of trees in this group of chapters. Chapter 15 discusses trees-binary, binary search, and general-and their possible uses. Chapter 16 considers implementations of these trees, and briefly introduces the tree sort. C++ Interlude 6 presents iterators in the context of a list. Chapter 17 introduces the ADT heap and shows how to implement it by using an array. We then use a heap to implement the priority queue and to sort an array.
- Chapter 18: This chapter covers the specification and use of the ADT dictionary (formerly called the table in the previous edition). We look at implementations of the dictionary that use an array or a binary search tree. We then introduce hashing and use it as a dictionary implementation.
- Chapter 19: Chapter 19 introduces balanced search trees. Included in this chapter are the 2-3, 2-4, and redblack trees, as well as AVL trees. These trees are considered as implementations of the ADT dictionary.
- Chapter 20: Next, we discuss graphs, suggest two ways to implement them, and look at several applications.
- Chapter 21: This last chapter considers data storage in external direct access files. Merge sort is modified to sort such data, and external hashing and B-tree indexes are used to search it. These searching algorithms are generalizations of the internal hashing schemes and 2-3 trees already developed. Finally, C++ Interlude 7 ends the main presentation by discussing the containers and algorithms available in the C++ Standard Template Library (STL).
- Appendices A through L: The appendices provide supplemental information. As we mentioned earlier, Appendix A reviews C++ up to but not including classes. Appendices B, C, D, and F contain sections that were in Chapter 1 of the previous edition, namely important aspects of programming, the Unified Modeling Language (UML), the software life cycle, and algorithm verification. Appendix E covers mathematical induction, and Appendix G covers input and output with external files. Appendix H provides a list of C++ header files and standard functions, and Appendix I considers the javadoc commenting style and defines the tags that we use in this book. Appendix J is simply a chart of the ASCII character codes. Finally, Appendices K and L are brief transition guides to C++ for those who know Java or Python, respectively.


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T
his book evolved from the original Intermediate Problem Solving and Data Structures: Walls and Mirrors by Paul Helman and Robert Veroff ( $\bigcirc \dagger 1986$ by The Benjamin/Cummings Publishing Company, Inc.). Professors Helman and Veroff introduced two powerful analogies, walls and mirrors, that have made it easier for us to teach-and to learn-computer science. This work builds on their organizational framework and overall perspective and includes some technical and textual content, examples, figures, and exercises derived from the original work.

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Each chapter, interlude, and appendix assumes that the reader has studied certain previous material. This list indicates those prerequisites. Numbers represent chapter numbers, and letters reference appendices. The notation $\mathrm{I} x$ represents $\mathrm{C}++$ Interlude $x$. Underlined prerequisites indicate an incidental or a partial dependence. You can use this information to plan a path through the book.

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## Data Abstraction: The Walls

## Chapter

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## Prerequisites

Appendix A Review of C++ Fundamentals
Appendix B Important Themes in Programming
Appendix C The Unified Modeling Language

Coding without a solution design increases debugging time

A solution specifies a system of interacting objects

An object is an instance of a class

OOA explores a problem, not a solution

Specify what to do, not how to do it

This chapter summarizes several fundamental principles that serve as the basis for dealing with the complexities of large programs. The discussion both reinforces the basic principles of programming and demonstrates that writing well-designed and well-documented programs is essential. The chapter also introduces algorithms and data abstraction and indicates how these topics relate to the book's main theme of problem solving. Data abstraction is a technique for increasing the modularity of a program-for building "walls" between a program and its data structures. During the design of a solution, you will discover that you need to support several operations on the data and therefore need to define abstract data types (ADTs). Only after you have clearly specified the operations of an ADT should you consider data structures for implementing it. This chapter will introduce a simple ADT and use it to demonstrate the advantages of ADTs in general.

In subsequent chapters, we examine ways of organizing and using data. As we focus on these new ideas, pay attention to how all of the solutions we look at adhere to the basic principles discussed in this chapter.

### 1.1 Object-Oriented Concepts

Where did you begin when you wrote your last program? After reading the problem specifications and going through the requisite amount of procrastination, most novice programmers simply begin to write code. Obviously, their goal is to get their programs to execute, preferably with correct results. Therefore, they run their programs, examine error messages, insert semicolons, change the logic, delete semicolons, pray, and otherwise torture their programs until they work. Most of their time is probably spent checking both syntax and program logic. Certainly, your programming skills are better now than when you wrote your first program, but are you able to write a really large program by using the approach just described? Maybe, but there are better ways.

Whereas a first course in computer science typically emphasizes programming issues, the focus of this book is on the broader issues of problem solving. Here the term problem solving refers to the entire process of taking the statement of a problem and developing a computer program that solves that problem. This process requires you to pass through many phases, from gaining an understanding of the problem to be solved, through designing a conceptual solution, to implementing the solution as a computer program. In a large software project, problem solving requires not only the development of a computer program, but also the examination of user workflows and possible changes to how a user performs daily tasks.

Object-oriented analysis and design (OOAD) is a process for solving problems. From an object-oriented perspective, a solution is a computer program consisting of a system of interacting classes of objects. An object has a set of characteristics and behaviors related to the solution. Each object is responsible for some aspect of the solution. A set of objects having the same type is called a class. An object of a class is also known as an instance of the class. When you create an object of a class, you instantiate the object.

OOAD helps us to discover and describe these objects and classes. These techniques give us a starting place for moving from a problem statement to a solution.

### 1.1.1 Object-Oriented Analysis and Design

Object-oriented analysis (OOA) is the process of understanding what the problem is and what the requirements of a solution are. It is the initial stage in problem solving. During analysis you get an accurate perception of what end users expect the solution to be and do. The requirements of a solution give you a description of what a solution must be and what a solution must do-without imposing how to design or implement that solution.

During OOA, you express the problem and the requirements of a solution in terms of relevant objects. These objects may represent real-world objects, software systems, or ideas. Using OOA, you describe these objects and their interactions among one another. Analysis work involves the discovery associated with understanding the problem. It does not involve thinking about a solution to the problem, and instead focuses on thinking about the problem itself.

The results of analysis act as input to the process of design. During object-oriented design (OOD), you describe a solution to the problem, fulfilling the requirements you discovered during analysis. You express the solution in terms of software objects, and you note how those objects will collaborate. Objects collaborate when they send each other messages; that is, objects call on one another to perform operations. The interactions among objects are as important as the objects themselves and require careful planning. To solve the problem efficiently, the collaborations among objects should be meaningful and minimal.

During OOD, you typically create one or more models of a solution for the problem. Some of the models emphasize the interactions among objects; others show the relationships among the objects. Taken together, the models create a design that can be implemented in $\mathrm{C}++$ or any other object-oriented language.

### 1.1.2 Aspects of an Object-Oriented Solution

Unless otherwise stated, a solution to a problem in this book is a computer program. A program comprises modules working together. A module is a self-contained unit of code and could be a single, stand-alone function, a class method, a class itself, a group of several functions or classes that work closely together, or other blocks of code. Exactly how a module is defined depends on the type and size of the application. Functions and methods implement algorithms, which are step-by-step recipes for performing a task within a finite period of time. One action that an algorithm often performs is operating on a collection of data.

When designing a solution, your challenge is to create a good set of modules. These modules must store, move, and alter data. They also use methods to communicate with one another. When constructing a solution, you must organize your data collection so that you can operate on the data easily in the manner that an algorithm requires. In fact, most of this book describes ways of organizing data.

Object-oriented programming languages allow us to build classes of objects. A class combines the attributes-or characteristics-of objects of a single type together with the objects' operationsor behaviors-into a single unit. The individual data items specified in a class are called data members. The operations specified in the class are referred to as methods or member functions. Attributes are typically data, and the behaviors, or methods, often operate on that data. In programming languages such as $\mathrm{C}++$ and Java, classes specify the attributes and operations for the objects.

Encapsulation is a technique that hides inner details. Whereas functions encapsulate behavior, objects encapsulate data as well as behavior. For example, a clock encapsulates the time-an attribute-along with certain operations, such as setting or displaying the time. You can request that a clock perform those operations but you cannot see how they are done (unless you have a mechanical clock with see-through sides!).

Classes can inherit properties and operations from other classes. For example, once you have defined a base class of clocks, you can design a subclass of alarm clocks that inherits the properties of a clock but adds operations that provide the functionality of an alarm. You can produce an alarm clock quickly, because the clock portion is done. Thus, inheritance - another object-oriented conceptallows you to reuse classes you defined earlier for a related purpose by extending that implementation or making slight modifications.

Inheritance may make it impossible for the compiler to determine which operation you require in a particular situation. However, polymorphism-which literally means many forms-enables this

OOA explores a problem in terms of its objects

OOD explores a solution to a problem

OOD explores a solution's objects and their collaborations

## Modules implement

 algorithms, which often manipulate dataObjects encapsulate attributes (data) and behaviors
(operations)

Encapsulation hides inner details

Inheritance supports reusing software

Time devoted to analysis and design is time well spent

A highly cohesive module performs one well-defined task

A robust module performs well under unusual conditions
determination to be made at execution time. That is, the outcome of a particular operation depends upon the object that performs the operation. For example, you can create a pointer to a clock object, myClock, in your program in such a way that it could reference either a clock object or an alarm clock. When myClock is asked to display the time, the compiler cannot determine whether it should use the clock implementation to display the time or the alarm clock implementation, since it does not know to which class of clocks the object referenced by myClock belongs. Polymorphism allows the compiler to simply note that the meaning of an operation is unknown until execution time.


## Note: Three principles of object-oriented programming

1. Encapsulation: Objects combine data and operations.
2. Inheritance: Classes can inherit properties from other classes.
3. Polymorphism: Objects can determine appropriate operations at execution time.

### 1.2 Achieving a Better Solution

The last program you wrote most likely solved the given problem correctly. However, was it the best possible solution? If you spent little-if any-time doing analysis and design, the solution probably left something to be desired. If you were to code the same program again, you would doubtless produce a better solution than your first attempt. However, if you spent some extra time analyzing the problem and designing a solution, you would probably get your best solution.

Suppose that you generated three correct but different solutions. Can you identify aspects of each solution that makes it better than the other solutions? What are these aspects? What should you focus on to create better solutions?

Creating a good set of modules for a moderate-sized problem is more art than science. It requires experience on the part of the programmer. A given problem likely has no "best" set of modules. Some sets of modules - and their interactions-might be better than others in light of certain measures. Moreover, for a sufficiently large problem, several different sets of modules could be considered "best," depending upon the measure used. The "better" designs, however, do adhere to certain principles, which we examine next.

### 1.2.1 Cohesion

Each module should perform one well-defined task; that is, it should be highly cohesive. A highly cohesive module brings several immediate benefits to a design or solution.

First, the module, if well named, promotes self-documenting, easy-to-understand code. For example, a highly cohesive function called sort should do nothing but sort. What this function does is clear from its name: If this function also prints the sorted values, it is not cohesive.

Second, a highly cohesive module is easy to reuse in other software projects. If a solution for another problem is being developed, the highly cohesive sort function can be used without change. If this function also prints the sorted values, it is much less likely to be the right sorting routine for the job.

Third, a highly cohesive module is much easier to maintain. Because the highly cohesive sort function does nothing but sort, fixing a logical error is simpler. If this function prints the sorted values, too, the printing code will complicate the function's maintenance. A highly cohesive module has but one task that might need revision.

Fourth, a highly cohesive module is more robust; that is, it is less likely to be affected by change. The highly cohesive sort function will require change only if the system requires a different kind of sort. For example, you might need to sort data into descending, rather than ascending, order, or you might need a faster sort.

Like many object-oriented principles, cohesion can be described in human terms. A person with low cohesion has "too many irons in the fire." Such people tend to get bogged down in everything that they need to get done, and nothing they do gets done well. They could become more cohesive by delegating some of their responsibilities to others.

Note: A guiding principle of OOD is that each class should have a single, well-defined responsibility. The methods of a class should be highly cohesive and related directly to supporting the responsibility of the class. The responsibilities of a class are functionally equivalent to the tasks that the class needs to perform. If a class has too many responsibilities, it should be split into multiple classes, each with a single responsibility taken from the original class.

### 1.2.2 Coupling

Coupling is a measure of the dependence among modules. This dependence, for example, could involve sharing data structures or calling each other's methods. Ideally, the modules in a design should be independent of one another. However, some degree of coupling is necessary to get work done. That is, modules should be loosely coupled, and highly coupled modules should be avoided.

Loose coupling benefits a system in several ways. First, a module with loose coupling tends to create a system that is more adaptable to change. If class $A$ depends on - that is, is highly coupled to - a class $B$ and class $B$ is changed, it is very likely that these changes will affect class $A$ and break it.

Second, a module with loose coupling creates a system that is easier to understand. If class $A$ depends on a class $B$, understanding how class $A$ works requires an understanding of class $B$. Thus, class $A$ is difficult to understand in isolation. A solution with a high degree of coupling can become nearly impossible to understand.

Third, a module with loose coupling increases the reusability of that module. If class $A$ depends on a class $B$, reusing class $A$ in another program is complicated by the need to include class $B$ in that program as well. Reusing coupled modules requires reusing all of the modules together as a unit. Often this is not desirable, or possible.

Fourth, a module with loose coupling has increased cohesion. Moreover, highly cohesive modules tend to be loosely coupled. As the level of cohesion associated with a module goes down, that module does more unrelated work, as we saw in the previous section. This has the side effect of causing the module to be coupled with many other modules from other areas in the program.

Again, realize that some coupling is required; coupling cannot and should not be eliminated from designs and solutions. To get work done, objects must collaborate. But collaboration requires objects to depend on one another. Tasks that an object has delegated to other objects create coupling between these objects. This coupling is necessary, but it should be kept to a minimum. However, other factors may influence a design. Thus, some designs with more coupling are better than other designs with less coupling because of these other factors.

Note: If, in the past, you have spent little or no time on analysis and design for your programs, you must change this habit! The end result of OOD should be a modular solution that is easy to translate into the constructs of a particular programming language. By spending adequate time with analysis and design, you will spend less time writing and
debugging your program.


A loosely coupled module is independent

Write specifications for each module before implementing it

A module's contract specifies the module's purpose, assumptions, input, and output

Specify the data flow among modules

## FIGURE 1-1 The task sort is a module separate from the MyProgram module



### 1.3 Specifications

When you design a modular solution to a problem, each module simply states what it does but not how it does it. No one module may "know" how any other module performs its task-it may know only what that task is. For example, if one part of a solution is to sort some data, one of the modules may be a sorting algorithm, as Figure 1-1 illustrates. The other modules know that the sorting module sorts, but they do not know how it sorts. In this way, the various components of a solution are kept isolated from one another.

Using this idea, you can write the modules in relative isolation from one another, knowing what each one will do but not necessarily how each will eventually do it. It is essential that a module's specifications be written and understood.

### 1.3.1 Operation Contracts

An operation contract documents how a method can be used and what limitations it has. You should begin specifying this contract during analysis, finish the specification during design, and then document the contract in your code, particularly within the header files. In this way, programmers who use your code can understand what contract they need to honor for the method to generate correct results.

There is certain information you need to provide for each method. A method's interface will specify how to call the method and the number, order, and types of arguments it expects-as already discussed. You also need to specify what should be true before the method is called and what will be true after the method finishes execution.

During design, it is also important that you clearly specify not only the purpose of each module, but also the data flow among modules. For example, you should provide answers to these questions for each module: What data is available to the module before its execution? What does the module assume? What actions have taken place, and what does the data look like, after the module executes? Thus, you should specify in detail the assumptions, input, and output for each module.

For example, if you as program designer needed to sort an array of integers, you might write the following specifications for a sort function:

The function will receive an array of num integers, where num $>0$.
The function will return the array with the integers sorted.

You can view these specifications as the terms of a contract between your function and the module that calls it.

This contract helps programmers understand what responsibilities the module will have to the other modules in the solution. Whoever writes the sort function must live up to this contract. After the sort function has been written and tested, the contract tells the rest of the program how to call the sort function properly, as well as the result of doing so.

Notice, however, that a module's contract does not commit the module to a particular way of performing its task. If another part of the program assumes anything about the algorithm, it does so at its own risk. Thus, for example, if at some later date you rewrite your function to use a different sorting algorithm, you should not need to change the rest of the program at all. As long as the new function honors the terms of the original contract, the rest of the program should be oblivious to the change.

This should not be news to you. Although you might not have explicitly used the term contract before, the concept should be familiar. You write a contract when you write a function's precondition, which is a statement of the conditions that must exist at the beginning of a function, as well as when you write its postcondition, which is a statement of the conditions at the end of a function. For example, the sort function that adheres to the previous contract could appear in pseudocode ${ }^{1}$ as

```
// Sorts an array.
// Precondition: anArray is an array of num integers; num >0.
// Postcondition: The integers in anArray are sorted.
sort(anArray, num)
```

These particular pre- and postconditions actually are deficient, as can be the case in a first-draft contract. For example, does sorted mean ascending order or descending order? How large can num be? While implementing this function, you might assume that sorted means ascending order and that num will not exceed 100. Imagine the difficulties that can arise when another person tries to use sort to sort an array of 500 integers into descending order. This user will not know your assumptions unless you documented them by revising the contract as follows:

```
// Sorts an array into ascending order.
// Precondition: anArray is an array of num integers and l <= num <= MAX_ARRAY,
// where MAX_ARRAY is a global constant that specifies the maximum size of anArray.
// Postcondition: anArray[0]<= anArray[1]<= ... <= anArray[num - 1];
// num is unchanged.
sort(anArray, num)
```

When you write a precondition, begin by describing the method or function's input arguments, mention any global named constants that it uses, and finally list any assumptions that it makes. When you write a postcondition, describe what changes the module has made. Note that in the case of a method or function that returns a value-which is technically a part of the postcondition-the value should be described.

Novice programmers tend to dismiss the importance of precise documentation, particularly when they are simultaneously designer, programmer, and user of a small program. If you design sort but do not write down the terms of the contract, will you remember them when you later implement the function? Will you remember how to use sort weeks after you have written it? To refresh your memory, would you rather examine your program code or read a simple set of pre- and postconditions? As the size of a program increases, good documentation becomes even more important, regardless of whether you are sole author or part of a team.

Specifications are the terms of a contract

An operation contract should not describe how a module will perform its task

Operation contracts should include precise preconditions and postconditions

First-draft specifications

Revised specifications

Precise documentation is essential

[^0]

Note: An operation contract completely specifies a module's purpose, assumptions, input, and output.

Note: The program component that uses a module is the module's client. The user is a person who uses a program.

### 1.3.2 Unusual Conditions

You as a class designer need to make decisions about how to treat unusual conditions and include these decisions in your specifications. The documentation for methods and functions should reflect these decisions.

In general, you can address unusual situations in several ways. You can

- Assume that the invalid situations will not occur. This assumption is not as naive as it might sound. A method could state as an assumption-that is, a precondition-restrictions to which a client must adhere. It is then up to the client to check that the precondition is satisfied before invoking the method. As long as the client obeys the restriction, the invalid situation will not occur.
- Ignore the invalid situations. A method could simply do nothing when given invalid data. Doing absolutely nothing, however, leaves the client without knowledge of what happened.
- Guess at the client's intention. Like the previous option, this choice can cause problems for the client.
- Return a value that signals a problem. For example, a method can return a boolean value that indicates its success or failure.
- Throw an exception. Throwing an exception is often a desirable way for a $\mathrm{C}++$ method to react to unusual events that occur during its execution. The method can simply report a problem without deciding what to do about it. The exception enables each client to act as needed in its own particular situation. For simplicity right now, we will adopt the philosophy that methods should throw exceptions only in truly unusual circumstances, when no other reasonable solution exists. You can learn about exceptions in C++ Interlude 3.


Note: A first draft of a module's specifications often overlooks or ignores situations that you really need to consider. You might intentionally make these omissions to simplify this first draft. Once you have written the major portions of the specifications, you can concentrate on the details that make the specifications complete.

### 1.3.3 Abstraction

Specifications do not indicate how to implement a module

Abstraction separates the purpose of a module from its implementation. Modularity breaks a solution into modules; abstraction specifies each module clearly before you implement it in a programming language. For example, what does the module assume, and what action does it take? What task is this module responsible for when called on? Such specifications clarify the design of your solution, because you can focus on its high-level functionality without the distraction of implementation details. In addition, they help you modify one part of a solution without significantly affecting the other parts. For example, you should be able to change the sorting algorithm in the previous example without affecting the rest of the solution.

As the problem-solving process proceeds, you gradually refine the modules until eventually you implement their actions by writing code-typically, classes and their methods. Separating the purpose of a module from its implementation is known as functional (or procedural) abstraction. Once a module is written, you can use it without knowing the particulars of its algorithm as long as you have a statement of its purpose and a description of its arguments. Assuming that the module is documented properly, you can use it knowing only its specifications. You do not need to look at its implementation.

Functional abstraction is essential to team projects. After all, in a team situation, you have to use modules written by others, frequently without knowledge of their algorithms. Can you actually use such a module without studying its code? In fact, you do this each time you use a C++ Standard Library function, such as sqrt in the C++ math library cmath. Because sqrt is precompiled, you do not have access to its source statements. Furthermore, it may be that sqrt was written in a language other than $\mathrm{C}++$ ! There is so much about sqrt that you do not know; yet you can use it in your program without concern, as long as you know its specifications. If you pass sqrt a floating-point expression, it will return the floating-point square root of the value of that expression. You can use sqrt even though you do not know its implementation.

Consider now a collection of data and a set of operations on the data. The operations might include ones that add new data to the collection, remove data from the collection, or search for some data. Data abstraction focuses on what the operations do with the collection of data, instead of on how you implement them. The other modules of the solution "know" what operations they can perform, but they do not know how the data is stored or how the operations are performed.

For example, you have used an array, but have you ever stopped to think about what an array actually is? There are many pictures of arrays throughout this book. They might resemble the way a C++ array is implemented on a computer, and then again they might not. In either case, you can use an array without knowing what it "looks like"-that is, how it is implemented. Although different systems may implement arrays in different ways, the differences are transparent to the programmer.

For instance, regardless of how the array years is implemented, you can always store the value 1492 in location index of the array by using the statement

```
years[index] = 1492;
```

and later display the value by using the statement

```
cout << years[index] << endl;
```

Thus, you can use an array without knowing the details of its implementation, just as you can use the function sqrt without knowing the details of its implementation. Let's explore this idea in more detail.

### 1.3.4 Information Hiding

As you have seen, abstraction tells you to write functional specifications for each module that describe its outside, or public, view. However, abstraction also helps you identify details that you should hide from public view-details that should not be in the specifications but rather should be private. The principle of information hiding tells you not only to hide such details within a module, but also to ensure that no other module can tamper with these hidden details.

While writing a module's specifications, you must identify details that you can hide within the module. The principle of information hiding involves not only hiding these details, but also making them inaccessible from outside a module. One way to understand information hiding is to imagine walls around the various tasks a program performs. These walls prevent the tasks from becoming entangled. The wall around each task prevents the other tasks from "seeing" how that task is performed.

Complete specifications enable you to use a module without knowing its implementation

Both functional and data abstraction ask you to think "what," not "how"

All modules and ADTs should hide something

The isolation of the modules cannot be total, however. Although MyProgram does not know how the task sort is performed, it must know what the task sort is and how to initiate it. For example, suppose your program needs to operate on a sorted array of names. The program may, for instance, need to search the array for a given name or display the names in alphabetical order. The program thus needs a function sort that sorts an array of names. Although the rest of the program knows that sort can sort an array, it should not care how sort accomplishes its task.

Thus, imagine a tiny slit in the wall, as Figure 1-2 illustrates. The slit is not large enough to allow the outside world to see the function's inner workings, but items can pass through the slit into and out of the function. This slit is the prototype, declaration, or header of the function. The slit comprises the function or method's name, parameter list, and return type. For example, you can pass the array into the function sort, and the function can pass the sorted array out to you. What goes in and comes out is governed by the terms of the function's specifications, or contract: If you use the function in this way, this is exactly what it will do for you.

FIGURE 1-2 Tasks communicate through a slit in the wall


Suppose that a faster sort algorithm is developed. Since the function sort is isolated from the other modules in the program, the new algorithm can be implemented in the sort function without affecting those other modules. Thus, if MyProgram uses the task sort, and if the algorithm and implementation for performing the sort changes, MyProgram will not be affected. As Figure 1-3 illustrates, the wall prevents MyProgram's algorithm from depending on sort's algorithm.


Note: Information hiding limits the ways in which you need to deal with modules and data. As a user of a module, you do not worry about the details of its implementation. As an implementer of a module, you do not worry about its uses.

A module's interface is the only way to interact with that module

### 1.3.5 Minimal and Complete Interfaces

The interface for a class is made up of the publicly accessible methods and data. Typically, a class interface contains only methods, as you will see, because publicly accessible data fields generally

FIGURE 1-3 A revised implementation communicates through the same slit in the wall

cause problems. The interface for a class describes the only way for programmers to interact with that class. Thus, interfaces are where collaboration between objects takes place. It is through its interface that a class is coupled to other classes. Designing good class interfaces is an important skill.

Each class should be easy to understand. Thus, when designing a class, you should keep the number of its methods small. However, your classes should provide programmers the power to do what they need to do easily. These desires are at odds.

A complete interface for a class is one that will allow programmers to accomplish any reasonable task, given the responsibilities of that class. A minimal interface for a class is one that contains a method if and only if that method is essential to that class's responsibilities. Programmers can more easily learn how to interact with a class that has a minimal interface, as it has fewer methods to understand. Classes with minimal interfaces are also much easier to maintain. Changes to a class with an extensive interface could affect many of its methods.

You should evaluate classes in terms of how complete and minimal their interfaces are. It is important that interfaces be complete. Of somewhat less importance is having a minimal interface; sometimes a nonessential method is just too useful to omit.

The interface to a function or method is more accurately called its signature. The signature consists of a function's name; the number, order, and types of its arguments; and any qualifiers such as const that might apply. A signature looks very much like the function's prototype, but does not include its return type. You use the interface to a function or method when you call it, sending it the correct number, order, and type of arguments.

### 1.4 Abstract Data Types

Often the solution to a problem requires operations on data. Such operations are broadly described in one of three ways:

Typical operations on data

- Add data to a data collection.
- Remove data from a data collection.
- Ask questions about the data in a data collection.

An ADT is not a fancy name for a data structure

Specifications indicate what ADT operations do, but not how to implement them

## Carefully specify

 an ADT's operations before you implement themThe details of the operations, of course, vary from application to application, but the overall theme is the management of data. Realize, however, that not every problem uses or requires all of these operations.

Most of this book is about data abstraction. To enable you to think abstractly about data, you should define an abstract data type, or ADT. An ADT is a collection of data and a set of operations on the data. You can use an ADT's operations, if you know their specifications, without knowing how the operations are implemented or how the data is stored.

Ultimately, someone-perhaps you-will implement the ADT by using a data structure, which is a construct that you can define within a programming language to store a collection of data. For example, you might store the data in a C++ array of strings or in an array of objects or in an array of arrays.

For example, suppose that you need to store a collection of names in a manner that allows you to search rapidly for a given name. A collection of name items providing for rapid searches is the description of a simple ADT. The description of an ADT's operations must be rigorous enough to specify completely their effect on the data, yet must specify neither how to store the data nor how to carry out the operations. For example, the ADT operations should not specify whether to store the data in consecutive memory locations or in disjoint memory locations. You choose a particular data structure when you implement an ADT.

When a program must perform data operations that are not directly supported by the language, you should first design an ADT and carefully specify what the ADT operations are to do (the contract). Then-and only then-should you implement the operations with a data structure. If you implement the operations properly, the rest of the program will be able to assume that the operations perform as specified-that is, that the terms of the contract are honored. However, the program must not depend on a particular technique for supporting the operations.


## Note: ADTs versus data structures

- An abstract data type is a specification for a group of values and the operations on those values.
- A data structure is an implementation of an ADT within a programming language.

To give you a better idea of the conceptual difference between an ADT and a data structure, consider a refrigerator's ice dispenser, as Figure 1-4 illustrates. It has water as input and produces as output either chilled water, crushed ice, or ice cubes, according to which one of three buttons you push.

FIGURE 1-4 A dispenser of chilled water, crushed ice, and ice cubes


User view from specifications


Technician view

It also has an indicator that lights when no ice is available. The water is analogous to data; the operations are chill, crush, cube, and isEmpty (or noIce). At this level of design, the dispenser is analogous to an ADT; you are not concerned with how the dispenser will perform its operations, only that it performs them. If you want crushed ice, do you really care how the dispenser accomplishes its task, as long as it does so correctly? Thus, after you have specified the dispenser's operations, you can design many uses for crushed ice without knowing how the dispenser accomplishes its tasks and without the distraction of engineering details.

Eventually, however, someone must build the dispenser. Exactly how will this machine produce crushed ice, for example? It could first make ice cubes and then either crush them between two steel rollers or smash them into small pieces by using hammers. Many other techniques are possible. The internal structure of the dispenser corresponds to the implementation of the ADT in a programming language, that is, to a data structure.

Although the owner of the dispenser does not care about its inner workings, he or she does want a design that is as efficient in its operation as possible. Similarly, the dispenser's manufacturer wants a design that is as easy and inexpensive to build as possible. You should have these same concerns when you choose a data structure to implement an ADT. Even if you do not implement the ADT yourself, but instead use an already implemented ADT, you-like the person who buys a refrigerator-should care about at least the ADT's efficiency.

Notice that steel walls surround the dispenser. The only breaks in the walls accommodate the input (water) to the machine and its output (chilled water, crushed ice, or ice cubes). Thus, the machine's interior mechanisms are not only hidden from the user, but also inaccessible. In addition, the mechanism of one operation is hidden from and inaccessible to another operation.

This modular design has benefits. For example, you can improve the operation crush by modifying its implementation without affecting the other modules. You could also add an operation to the machine without affecting the original three operations. Thus, both abstraction and information hiding are at work here.

To summarize, data abstraction results in a wall of ADT operations between data structures and the program that accesses the data within these data structures, as Figure 1-5 illustrates. If you are on

A program should not depend on the details of an ADT's implementation

FIGURE 1-5 A wall of ADT operations isolates a data structure from the program that uses it


Using an ADT is like using a vending machine
the program's side of the wall, you will see an interface that enables you to communicate with the data structure. That is, you ask the ADT operations to manipulate the data in the data structure, and they pass the results of these manipulations back to you.

This process is analogous to using a vending machine. You press buttons to communicate with the machine and obtain something in return. The machine's external design dictates how you use it, much as an ADT's specifications govern what its operations are and what they do. As long as you use a vending machine according to its design, you can ignore its inner technology. As long as you agree to access data only by using ADT operations, your program can be oblivious to any change in the data structures that implement the ADT.

The following pages describe several abstract data types, focusing on the specifications of their operations, but not their implementations.

### 1.4.1 Designing an ADT

The design of an abstract data type should evolve naturally during the problem-solving process. As an example of how this process might occur, suppose that you want to determine the dates of all the holidays in a given year. One way to do this is to examine a calendar. That is, you could consider each day in the year and ascertain whether that day is a holiday. The following pseudocode is a possible solution to this problem:

```
// Displays the dates of all holidays in a given year.
1istHolidays(year)
    date = date offirst day of year
    while (date is before the first day of year + 1)
    {
        if (date is a holiday)
            write (date," is a holiday")
        date = date of next day
    }
```

What data is involved here? Clearly, this problem operates on dates, where a date consists of a month, day, and year. What operations will you need to solve the holiday problem? Your ADT must specify and restrict the legal operations on the dates, just as the fundamental data type int restricts you to operations such as addition and comparison. You can see from the previous pseudocode that you must

- Determine the date of the first day of a given year.
- Decide whether a date is before another date.
- Decide whether a date is a holiday.
- Determine the date of the day that follows a given date.

Thus, you could specify the following operations for an ADT date in pseudocode that uses the notation of the Unified Modeling Language (UML), as described in Appendix C:
// Returns the date of the first day of a given year.
+getFirstDay(year: integer): Date
// Returns true if this date is before the given date; otherwise returns false.
+isBefore(otherDate: Date): boolean
// Returns true if this date is a holiday; otherwise returns false.
+isHoliday(): boolean
// Returns the date of the day after this date.
+getNextDay(): Date

The 1istHolidays pseudocode now appears as follows:

```
// Displays the dates of all holidays in a given year.
listHolidays(year: integer): void
    date = getFirstDay (year)
    while (date.isBefore (getFirstDay (year + 1)))
    \{
        if (date.isHoliday())
            write (date," is a holiday ")
    date \(=\) date.getNextDay()
\}
```

Thus, you can design an ADT by identifying data and choosing operations that are suitable to your problem. After specifying the operations, you use them to solve your problem independently of the implementation details of the ADT.

An appointment book. As another example of an ADT design, imagine that you want to create a computerized appointment book that spans a one-year period. Suppose that you make appointments only on the hour and half hour between $8 \mathrm{a} . \mathrm{m}$. and $5 \mathrm{p} . \mathrm{m}$. You want your system to store a brief notation about the nature of each appointment along with the date and time.

To solve this problem, you can define an ADT appointment book. The data items in this ADT are the appointments, where an appointment consists of a date, time, and purpose. What are the operations? Two obvious operations are

- Make an appointment for a certain date, time, and purpose. (You will want to be careful that you do not make an appointment at an already occupied time.)
- Cancel the appointment for a certain date and time.

In addition to these operations, it is likely that you will want to

- Ask whether you have an appointment at a given time.
- Get the purpose of your appointment at a given time.

Finally, ADTs typically have initialization and destruction operations that we assume but do not specify at this stage.

Thus, the ADT appointment book can have the following operations:

```
// Returns true if an appointment exists for the date and time specified,
// false otherwise.
+isAppointment(apptDate: Date, apptTime: Time): boolean
// Inserts the appointment for the date, time, and purpose specified as long as
// it does not conflict with an existing appointment.
// Returns true if successful, false otherwise.
+makeAppointment(apptDate: Date, apptTime: Time,
    apptPurpose: string): boolean
// Deletes the appointment for the date and time specified.
// Returns true if successful, false otherwise.
+cancelAppointment(apptDate: Date, apptTime: Time): boolean
// Gets the purpose of the appointment at the given date and time, if one exists.
// Otherwise returns an empty string.
+getAppointmentPurpose(apptDate: Date, apptTime: Time): string
```

You can use these ADT operations to design other operations on the appointments. For example, suppose you want to change the date or time of a particular appointment within the existing

You can use an ADT without knowledge of its implementation

You can use an ADT to implement another ADT
appointment book apptBook. The following pseudocode indicates how to accomplish this task by using the previous ADT operations:

## // Change the date or time of an appointment.

```
Get the following data from the user: oldDate, oldTime, newDate, newTime
// Get purpose of appointment.
oldPurpose = apptBook.getAppointmentPurpose(oldDate, oldTime)
if (oldPurpose is not the empty string)
{
    // See whether a new date/time is available.
    if (apptBook.isAppointment(newDate, newTime))
        // New date/time is booked
        write("You already have an appointment at ", newTime,
            " on ", newDate)
    else
    {
        // New date/time is available; cancel old appointment; make new one.
        apptBook.cance1Appointment(oldDate, oldTime)
        if (apptBook.makeAppointment(newDate, newTime, oldPurpose))
            write("Your appointment has been rescheduled to ",
                newTime, " on ", newDate)
    }
}
e1se
    write("You do not have an appointment at ", oldTime,
        " on ", oldDate)
```

Again notice that you can design applications that use ADT operations without knowing how the ADT is implemented. The exercises at the end of this chapter provide examples of other tasks that you can perform with this ADT.

### 1.4.2 ADTs That Suggest Other ADTs

Both of the previous examples require you to represent a date; the appointment book example also requires you to represent the time. C++ has date-time objects specified in ctime that you can use to represent the date and the time. You can also design ADTs to represent these items in a more object-oriented way. It is not unusual for the design of one ADT to suggest other ADTs. In fact, you can use one ADT to implement another ADT.

This final example also describes an ADT that suggests other ADTs for its implementation. Suppose that you want to design a database of recipes. You could think of this database as an ADT. The recipes are the data items, and some typical operations on the recipes could include the following:

```
// Inserts a recipe into the database.
+insertRecipe(aRecipe: Recipe): boolean
// Deletes a recipe from the database.
+deleteRecipe(aRecipe: Recipe): boolean
// Gets the named recipe from the database.
+getRecipe(name: string): Recipe
```

This level of the design does not indicate such details as where insertRecipe will place a recipe into the database.

Now imagine that you want to design an operation that scales a recipe retrieved from the database: If the recipe is for $n$ people, you want to revise it so that it will serve $m$ people. Suppose that
the recipe contains measurements such as $21 / 2$ cups, 1 tablespoon, and $1 / 4$ teaspoon. That is, the quantities are given as mixed numbers-integers and fractions-in units of cups, tablespoons, and teaspoons.

This problem suggests another ADT—measurement—with the following operations:

```
// Returns this measure.
+getMeasure(): Measurement
// Sets this measure to another one.
+setMeasure(m: Measurement)
// Returns this measure multiplied by a fractional scale factor, which has no units.
+scaleMeasure(scaleFactor: float): Measurement
// Returns this measure converted from its old units to new units.
+convertMeasure(oldUnits: MeasureUnit,
    newUnits: MeasureUnit): Measurement
```

Suppose that you want the ADT measurement to perform exact fractional arithmetic. Because our planned implementation language $\mathrm{C}++$ does not have a data type for fractions, and floatingpoint arithmetic is not exact, another ADT called fraction is in order. Its operations could include addition, subtraction, multiplication, and division of fractions. For example, you could specify addition as

```
// Returns the sum, reduced to lowest terms, of this fraction and the given fraction.
+add(other: Fraction): Fraction
```

Moreover, you could include operations to convert a mixed number to a fraction and vice versa when feasible. When you finally implement the ADT measurement, you can use the ADT fraction. That is, you can use one ADT to implement another ADT.

### 1.5 The ADT Bag

Imagine a paper bag, a reusable cloth bag, or even a plastic bag. People use bags when they shop, pack a lunch, or eat potato chips. A bag contains things; it is a container of a collection of objects. We could consider the bag to be an abstract data type.

Let's specify and use the ADT bag. Knowing just its interface, you can use a bag in a program. You do not need to know how the entries in the bag are represented or how the bag operations are implemented. These specifics are hidden behind the bag's wall of abstraction, and your program will not depend on these specifics. As you will see, this important program characteristic is what data abstraction is all about.

What distinguishes a bag from other ADTs? A bag doesn't do much more than contain its items. It doesn't order them in a particular way, nor does it prevent duplicate items. While describing the behaviors for the ADT bag that we'll design in this chapter, let's keep in mind that we are specifying an abstraction inspired by an actual physical bag. For example, a paper bag holds things of various dimensions and shapes in no particular order and without regard for duplicates. Our abstract bag will hold unordered and possibly duplicate objects, but let's insist that these objects have the same or related types.

### 1.5.1 Identifying Behaviors

Since a bag contains a finite number of objects, reporting how many objects it contains could be one of a bag's behaviors:

- Get the number of items currently in the bag.

A related behavior detects if a bag is empty:

- See whether the bag is empty.

We should be able to add and remove objects:

- Add a given object to the bag.
- Remove an occurrence of a specific object from the bag, if possible.
- Remove all objects from the bag.

Note that the add operation does not indicate where in the bag an object should go. Remember that a bag does not order its contents. The remove operation looks for a particular item in the bag. If it finds it, it takes it out. If the bag contains several objects equal to the removed one, they remain in the bag. If the operation can't find the specific object in the bag, it can't remove it, and just says so. Finally, the second remove operation simply empties the bag of all objects.

What is in a particular bag? The answer to this question can be answered by the following operations:

- Count the number of times a certain object occurs in the bag.
- Test whether the bag contains a particular object.
- Look at all objects that are in the bag.

We have enough behaviors for now. At this point, we would have written all the behaviors on a piece of paper or, as Appendix C suggests, on the class-responsibility-collaboration (CRC) card pictured in Figure 1-6.

FIGURE 1-6 A CRC card for a class Bag

| Bag |
| :--- |
| Responsibilities |
| Get the number of items currently in the bag |
| See whether the bag is empty |
| Add a given object to the bag |
| Remove an occurrence of a specific object from |
| the bag, if possible |
| Remove all objects from the bag |
| Count the number of times a certain object occurs in the bag |
| Test whether the bag contains a particular object |
| Look at all objects that are in the bag |
| Collaborations |
| The class of objects that the bag can contain |

Since a bag is an abstract data type, we only describe its data and specify its operations. We do not indicate how to store the data or how to implement its operations. Don't think about arrays, for example. You first need to clearly know what the bag operations do: Focus on what the operations do, not on how they do them. That is, you need a detailed set of specifications before you can use a bag in a program. In fact, you should specify the bag operations before you even decide on a programming language.

### 1.5.2 Specifying Data and Operations

Before we can implement a bag in C++, we need to describe its data and specify in detail the methods that correspond to the bag's behaviors. We'll name the methods, choose their parameters, decide their return types, and write comments to fully describe their effect on the bag's data. Our eventual goal, of course, is to write a C++ header file and comments for each method. However, we first will express the methods in pseudocode and then in UML notation.

The first behavior on our CRC card gives rise to a method that returns a count of the current number of entries in the bag. The corresponding method has no parameters and returns an integer. In pseudocode, we have the following specification, again using UML notation:
// Returns the current number of entries in the bag.
+getCurrentSize(): integer
We can test whether the bag is empty by using a boolean-valued method, again without parameters. Its specification in pseudocode is

```
// Returns true if the bag is empty.
+isEmpty(): boolean
```

We now want to add a given object to the bag. We can name the method add and give it a parameter to represent the new entry. We could write the following pseudocode:
// Adds a new entry to the bag. add (newEntry)

We might be tempted to make add a void method, but if the bag is full, we cannot add a new entry to it. What should we do in this case?

Here are two options that we can take when add cannot complete its task:

- Do nothing. We cannot add another item, so we ignore it and leave the bag unchanged.
- Leave the bag unchanged, but signal the client that the addition is impossible.

The first option is easy, but it leaves the client wondering what happened. Of course, we could state as a precondition of add that the bag must not already be full. Then the client has the responsibility to avoid adding a new entry to a full bag.

The second option is the better one, and it is not too hard to specify or implement. How can we indicate to the client whether the addition was successful? We could throw an exception if the addition is not successful, but we will leave this approach for later and use another way. Displaying an error message is not a good choice, as you should let the client dictate all written output. Since the addition is either successful or not, we can simply have the method add return a boolean value.

Thus, we can specify the method add more completely as
+add(newEntry: ItemType): boolean
where newEntry's data type is ItemType. In C++ Interlude 1, we will discuss the implementation details of using templates and typedef statements to give us flexibility in the type of data our ADT can hold.

Use the UML to describe classes at various stages of their development

Two behaviors involve removing entries from a bag: remove a particular entry and remove all entries. Suppose we name the methods and any parameters and begin to specify them in pseudocode as follows:

```
// Removes one occurrence of a particular entry from the bag, if possible.
remove(anEntry)
// Removes all entries from the bag.
clear()
```

What return types are these methods?
The remove method won't be able to remove a given entry if it isn't in the bag. We will have the method return a boolean value, much as add does, so it can indicate success or not. Thus, the UML specification for the method is

```
+remove(anEntry: ItemType): boolean
```

The method clear can be a void method: We just want to empty the bag, not get any of its contents. Thus, we refine its specification as follows:

```
+clear(): void
```

The remaining behaviors do not change the contents of the bag. One of these behaviors counts the number of times a given object occurs within the bag. Its specification in pseudocode is

```
// Counts the number of times a given entry appears in the bag.
+getFrequencyOf(anEntry: ItemType): integer
```

Another method tests whether the bag contains a given object. We specify it in pseudocode as follows:

```
// Tests whether the bag contains a given entry.
+contains(anEntry: ItemType): boolean
```

Finally, we want to look at the contents of the bag. Rather than providing a method that displays the entries in the bag, we will define one that places these entries into a vector. The client is then free to display any or all of them in any way desired. Here is the specification for our last method:

```
// Gets all entries in the bag.
+toVector(): vector
```

As we developed the previous specifications for the bag's methods, we represented them using UML notation. Figure 1-7 shows the result of doing so.

FIGURE 1-7 UML notation for the class Bag

| Bag |
| :--- |
|  |
| +getCurrentSize(): integer |
| +isEmpty(): boolean |
| +add(newEntry: ItemType): boolean |
| +remove(anEntry: ItemType): boolean |
| +clear(): void |
| +getFrequencyOf(anEntry: ItemType): integer |
| +contains(anEntry: ItemType): boolean |
| +toVector(): vector |

Notice that the CRC card and UML diagram do not reflect all of the details, such as assumptions and unusual circumstances. However, after you have identified such conditions, you should specify how your methods will behave under each one. You should write down your decisions about how you want your methods to behave, as we have done in the following table. Later, you can incorporate these informal descriptions into the $\mathrm{C}++$ comments that document your methods.

## Abstract Data Type: Bag

## DATA

- A finite number of objects, not necessarily distinct, in no particular order, and having the same data type.
- The number of objects in this collection.


## Operations

| Pseudocode | Description |
| :---: | :---: |
| getCurrentSize() | Task: Reports the current number of objects in this bag. Input: None. <br> Output: The number of objects currently in the bag. |
| isEmpty () | Task: Sees whether this bag is empty. <br> Input: None. <br> Output: True or false according to whether the bag is empty. |
| add (newEntry) | Task: Adds a given object to this bag. <br> Input: newEntry is an object. <br> Output: True or false according to whether the addition succeeds. |
| remove (anEntry) | Task: Removes an occurrence of a particular object from this bag, if possible. Input: anEntry is an object. <br> Output: True or false according to whether the removal succeeds. |
| clear() | Task: Removes all objects from this bag. Input: None. <br> Output: None. |
| getFrequencyOf(anEntry) | Task: Counts the number of times an object occurs in this bag. Input: anEntry is an object. <br> Output: The integer number of times anEntry occurs in the bag. |
| contains(anEntry) | Task: Tests whether this bag contains a particular object. <br> Input: anEntry is an object. <br> Output: True or false according to whether anEntry occurs in the bag. |
| toVector() | Task: Gets all objects in this bag. <br> Input: None. <br> Output: A vector containing all entries currently in the bag. |

## Note: Dealing with unusual conditions

As mentioned earlier, you must make decisions about how to treat unusual conditions. The documentation for the ADT bag should reflect both these decisions and the details in the previous discussion. For example, our specification of the method remove indicates that it will return false if a client tries to remove an entry from an empty bag. Instead, we could have given the method the precondition that the bag not be empty. It is then up to the client to check that the precondition is satisfied before invoking the method. Notice that the client can use other methods of the ADT bag, such as isEmpty and getCurrentSize, to help with this task.

Note: A first draft of an ADT's specifications often overlooks or ignores situations that you really need to consider. You might intentionally make these omissions to simplify this first draft. Once you have written the major portions of the specifications, you can concentrate on the details that make the specifications complete.

Question 1 Write specifications using UML notation for a function that computes the sum of the first five positive integers in an array of $n$ arbitrary integers.

### 1.5.3 An Interface Template for the ADT

As your specifications become more detailed, they increasingly should reflect your choice of programming language. Ultimately, you can write C++ headers for the bag's methods and organize them into a header file for the class that will implement the ADT.

The C++ header file in Listing 1-1 contains the methods for an ADT bag and detailed comments that describe their behaviors. If you are unfamiliar with $\mathrm{C}++$ classes, templates, and virtual methods, $\mathrm{C}++$ Interlude 1 provides a refresher on $\mathrm{C}++$ classes and discusses $\mathrm{C}++$ templates, virtual methods, and abstract base classes. By using templates and abstract base classes, we can take advantage of the object-oriented design concepts of encapsulation, polymorphism, and inheritance beginning with our first ADT.

Note: To emphasize the distinction between the design of an ADT and its implementation with a data structure, this textbook uses templates and abstract base classes in the description of our ADTs. We will call the header file containing the abstract base class a client interface. ${ }^{2}$ A programmer should not need anything more than the client interface to use a class in your program, as it completely specifies the methods of the ADT.

As you examine the interface in Listing 1-1, notice the decisions that were made to address the unusual situations mentioned in the previous section. In particular, each of the methods add, remove, and contains returns a value. Note the commenting tags, which begin with @ and are described in Appendix I, that we use to specify the methods.

[^1]For now, the items in the bag will be objects of the same class. To accommodate entries of any single class type, the bag methods use a generic type ItemType for each entry. To give meaning to the identifier ItemType, we must write template<class ItemType> on the line before the class header. Once the actual data type is chosen by a client, the compiler will use that data type wherever ItemType appears.

The class BagInterface is a C++ abstract base class. An abstract base class, or simply an abstract class, in C++ contains at least one method that is declared as virtual and has no implementation. An abstract class cannot be instantiated; it can only be used as a base class. The subclass must then implement the methods specified but not defined in the abstract class.

## LISTING 1-1 A file containing a C++ interface for bags

```
/** @file BagInterface.h */
#ifndef _BAG_INTERFACE
#define _BAG_INTERFACE
#include <vector>
temp1ate<c1ass ItemType>
class BagInterface
{
public:
```

    /** Gets the current number of entries in this bag.
    @return The integer number of entries currently in the bag. */
    virtual int getCurrentSize() const \(=0\);
    /** Sees whether this bag is empty.
        @return True if the bag is empty, or false if not. */
    virtual bool isEmpty() const \(=0\);
    /** Adds a new entry to this bag.
        @post If successful, newEntry is stored in the bag and
        the count of items in the bag has increased by 1.
    @param newEntry The object to be added as a new entry.
    @return True if addition was successful, or false if not. */
    virtual bool add(const ItemType\& newEntry) \(=0\);
    /** Removes one occurrence of a given entry from this bag,
        if possible.
    @post If successful, anEntry has been removed from the bag
        and the count of items in the bag has decreased by 1.
    @param anEntry The entry to be removed.
    @return True if removal was successful, or false if not. */
    virtual bool remove(const ItemType\& anEntry) \(=0\);
    /** Removes al1 entries from this bag.
        @post Bag contains no items, and the count of items is 0. */
    virtual void clear() \(=0\);
    /** Counts the number of times a given entry appears in bag.
    @param anEntry The entry to be counted.
    @return The number of times anEntry appears in the bag. */
    virtual int getFrequencyOf(const ItemType\& anEntry) const \(=0\);
    ```
    /** Tests whether this bag contains a given entry.
    @param anEntry The entry to locate.
    @return True if bag contains anEntry, or false otherwise. */
    virtual bool contains(const ItemType& anEntry) const = 0;
    /** Empties and then fills a given vector with all entries that
    are in this bag.
    @return A vector containing al7 the entries in the bag. */
    virtual vector<ItemType> toVector() const = 0;
}; // end BagInterface
```

The specifications of these operations are the sole terms of the contract for the ADT bag: If you request that these operations be performed, this is what will happen. The specifications contain no mention of how to store the bag or how to perform the operations; they tell you only what you can do to the bag. It is of fundamental importance that the specification of an ADT not include implementation issues. This restriction on the specification of an ADT is what allows you to build a wall between an implementation of an ADT and the program that uses it - that is, the client. The behavior of the operations is the only thing on which a program should depend.

Note: Although writing an interface before implementing a class is certainly not required, doing so enables you to document your specifications in a concise way. You then can use the code in the interface as an outline for the actual class. Having an interface also provides a data type for an ADT that is independent of a particular class definition. Chapters 3 and 4 will develop different implementations of a class of bags. Code written with respect to an interface allows you to more easily replace one implementation of a bag with another.

### 1.5.4 Using the ADT Bag

Imagine that we hire a programmer to implement the ADT bag in $\mathrm{C}++$, given the interface and specifications that we have developed so far. If we assume that these specifications are clear enough for the programmer to complete the implementation, we can use the bag's operations in a program without knowing the details of the implementation. That is, we do not need to know how the programmer implemented the bag to be able to use it. We need to know only what the ADT bag does.

The following example demonstrates how we can use a bag and assumes that we have a C++ class, Bag, that implements the C++ abstract class BagInterface given in Listing 1-1.

Example. Suppose you invited three friends to your home to watch your favorite soccer team on television. To make the commercial breaks more interesting, you make up the following game. From a deck of cards, you remove the entire suit of clubs: the cards Ace, Two, Three, ..., Jack, Queen, and King of clubs. You randomly select six of these cards and place them into a bag. Your friends then guess which cards are in the bag. Each time they correctly guess a card, it is removed from the bag.

When the bag is empty, the friend with the most cards wins. Listing 1-2 shows a simple program that plays this game.

## LISTING 1-2 A program for a card guessing game

```
#include <iostream> // For cout and cin
#include <string> // For string objects
#include "Bag.h" // For ADT bag
using namespace std;
int main()
{
    string clubs[] = { "Joker", "Ace", "Two", "Three",
                    "Four", "Five", "Six", "Seven",
                    "Eight", "Nine", "Ten", "Jack",
                    "Queen", "King" };
```

    // Create our bag to hold cards.
    Bag<string> grabBag;
    // Place six cards in the bag.
    grabBag.add(clubs[1]);
    grabBag.add(clubs[2]);
    grabBag.add(clubs[4]);
    grabBag.add(clubs[8]);
    grabBag.add(clubs[10]);
    grabBag.add(clubs[12]);
    // Get friend's guess and check it.
    int guess \(=0\);
    while (!grabBag.isEmpty())
    \{
        cout << "What is your guess?"
            << " (1 for Ace to 13 for King):";
        cin >> guess;
        // Is card in the bag?
        if (grabBag.contains(clubs[guess]))
        \{
            // Good guess - remove card from the bag.
            cout << "You get the card!\n";
            grabBag. remove(clubs[guess]);
        \}
        e1se
        \{
            cout << "Sorry, card was not in the bag.\n";
        \} // end if
    \} // end while
    cout << "No more cards in the bag. Game over! \n";
    return 0;
    \}; // end main

Programming Tip: After you design a class, try writing some code that uses your class before you implement the class. You will not only see whether your design works for the problem at hand, but also test your understanding of your own design and check the comments that document your specifications. You might discover problems with either your class design or your specifications. If so change your design and specifications, and try using the class again.

## Question 2 What is an abstract data type?

## Question 3 What steps should you take when designing an ADT?

## Summary

1. Object-oriented analysis (OOA) is used during the initial stage in problem solving to understand what the problem is and what the requirements of a solution are.
2. During object-oriented design (OOD), you describe a solution to a problem whose requirements are discovered during OOA. You express the solution in terms of software objects.
3. Encapsulation is a principle of object-oriented programming whereby you hide the inner details of functions and objects. Functions encapsulate behavior, and objects-that is, instances of a class-encapsulate data as well as behavior.
4. Inheritance is another concept related to object-oriented programming, allowing you to reuse already defined classes by extending their definitions or making slight modifications.
5. Polymorphism is the third principle of object-oriented programming, whereby objects determine appropriate operations at execution time.
6. Each module should be highly cohesive; that is, it should perform one well-defined task.
7. Coupling is a measure of the dependence among modules. Module should be loosely coupled. A function or method should be as independent as possible and perform one well-defined task.
8. UML is a modeling language used to express object-oriented designs. It provides a notation to specify the data and operations and uses diagrams to show relationships among classes.
9. An operation contract documents how a module can be used and what limitations it has.
10. A function or method should always include an initial comment that states its purpose, its precondition-that is, the conditions that must exist at the beginning of a module-and its postcondition-the conditions at the end of a module's execution.
11. For problems that primarily involve data management, encapsulate data with operations on that data by designing classes. Practice abstraction - that is, focus on what a module does instead of how it does it.
12. Data abstraction is a technique for controlling the interaction between a program and its data structures. It builds walls around a program's data structures, just as other aspects of modularity build walls around a program's algorithms. Such walls make programs easier to design, implement, read, and modify.
13. The specification of a set of data-management operations, together with the data values on which they operate, defines an abstract data type (ADT).
14. Only after you have fully defined an ADT should you think about how to implement it. The proper choice of a data structure to implement an ADT depends both on the details of the ADT operations and on the context in which you will use the operations.

## Exercises

1. The price of an item you want to buy is given in dollars and cents. You pay for it in cash by giving the clerk $d$ dollars and $c$ cents. Write specifications for a function that computes the change, if any, that you should receive. Include a statement of purpose, the preconditions and postconditions, and a description of the arguments.
2. A date consists of a month, day, and year. Consider the class Date of such dates. Suppose that Date represents the month, day, and year as integers. For example, July 4, 1776, is month 7, day 4, and year 1776.
a. Write specifications for a method within Date that advances any given date by one day. Include a statement of purpose, the preconditions and postconditions, a description of the arguments, and a description of any return value.
b. Write a C++ implementation of this method. Design and specify any other method that you need. Include comments that will be helpful to someone who will maintain your implementation in the future.
3. Write a pseudocode function in terms of the ADT appointment book, described in Section 1.4.1, for each of the following tasks. Do you need to add operations to the ADT to perform these tasks?
a. Change the purpose of the appointment at a given date and time.
b. Display all the appointments for a given date.

Imagine that you have just left a store with a bag of groceries. You are concerned that the fragile items will not survive the trip home, so when you reach your car, you place those items into their own bag. If Bag is a class of bags, write C++ statements that remove all the items from storeBag and place them into one of two new bags, as follows: Place all occurrences of bread and eggs into fragileBag, and all other items into groceryBag. When you are done, storeBag should be empty. Assume that grocery items are represented by strings.
5. Suppose that a bag contains strings that represent various grocery items. Write a $\mathrm{C}++$ function that removes and counts all occurrences of a given string from such a bag. Your function should return this number. Use comments in javadoc style to fully specify your function. Accommodate the possibility that the given bag is either empty or does not contain any occurrences of the given string.
6. The union of two bags is a new bag containing the combined contents of the original two bags. Design and specify a method union for the ADT bag that returns as a new bag the union of the bag receiving the call to the method and the bag that is the method's one argument. Include sufficient comments to fully specify the method.

Note that the union of two bags might contain duplicate items. For example, if object x occurs five times in one bag and twice in another, the union of these bags contains $x$ seven times. Specifically, suppose that bag1 and bag2 are bags; bag1 contains the strings $\mathrm{a}, \mathrm{b}$, and c ; and bag2 contains the strings $\mathrm{b}, \mathrm{b}, \mathrm{d}$, and e . The expression bag1.union(bag2) returns a bag containing the strings $a, b, b, b, c, d$, and e. Note that union does not affect the contents of bag1 and bag2.
7. The intersection of two bags is a new bag containing the entries that occur in both of the original two bags. Design and specify a method intersection for the ADT bag that returns as a new bag the intersection of the bag receiving the call to the method and the bag that is the method's one argument. Include sufficient comments to fully specify the method.

Note that the intersection of two bags might contain duplicate items. For example, if object x occurs five times in one bag and twice in another, the intersection of these bags contains $x$ two times. Specifically, suppose that bag1 and bag2 are bags; bag1 contains the strings $a, b$, and $c$; and bag2 contains the strings $b, b, d$, and $e$. The expression bag1.intersection(bag2) returns a bag containing only the string b. Note that intersection does not affect the contents of bag1 and bag2.
8. The difference of two bags is a new bag containing the entries that would be left in one bag after removing those that also occur in the second. Design and specify a method difference for the ADT bag that returns as a new bag the difference of the bag receiving the call to the method and the bag that is the method's one argument. Include sufficient comments to fully specify the method.

Note that the difference of two bags might contain duplicate items. For example, if object x occurs five times in one bag and twice in another, the difference of these bags contains $x$ three times. Specifically, suppose that bag1 and bag2 are bags; bag1 contains the strings a, b, and c; and bag2 contains the strings b, b, d, and e. The expression bag1.difference(bag2) returns a bag containing only the strings a and c. Note that difference does not affect the contents of bag1 and bag2.
9. Consider the ADT polynomial-in a single variable $x$-whose operations include the following:
degree() // Returns the degree of a polynomial.
coefficient (power) // Returns the coefficient of the xpower term.
changeCoefficient(newCoefficient, power) // Replaces the coefficient of
// the $x^{\text {power }}$ term with newCoefficient.
For this problem, consider only polynomials whose exponents are nonnegative integers. For example,
$p=4 x^{5}+7 x^{3}-x^{2}+9$
The following examples demonstrate the ADT operations on this polynomial.
p. degree() is 5 (the highest power of a term with a nonzero coefficient)
p. coefficient (3) is 7 (the coefficient of the $x^{3}$ term)
p. coefficient (4) is 0 (the coefficient of a missing term is implicitly 0 )
p.changeCoefficient $(-3,7)$ changes the polynomial p to $-3 x^{7}+4 x^{5}+7 x^{3}-x^{2}+9$

Using these ADT operations, write statements to perform the following tasks:
a. Display the coefficient of the term that has the highest power.
b. Increase the coefficient of the $x^{3}$ term by 8 .
c. Compute the sum of two polynomials.

## Programming Problems

1. Add a Transaction class to the banking example in Figure C-4 of Appendix C. This class keeps track of the date, time, amount, and type of transaction (checking or savings).
2. Consider a program that will read employee data, sort the data by employee identification number, write out the sorted data, and compute various statistics about the data, such as the average age of an employee. Design a modular solution to this problem using UML notation. What classes and methods did you identify during the design of your solution? Write complete specifications, including preconditions and postconditions, for each module.
3. A matrix is a rectangular array of numerical values. You can add or multiply two matrices to form a third matrix. You can multiply a matrix by a scalar, and you can transpose a matrix. Design an ADT that represents a matrix that has these operations.

Specify each ADT operation by stating its purpose, by describing its parameters, and by writing a pseudocode version of its header. Then write a C++ interface for the methods of a matrix. Include javadoc-style comments in your code.
4. A ring is a collection of items that has a reference to a current item. An operation-let's call it advance - moves the reference to the next item in the collection. When the reference reaches the last item, the next advance operation moves the reference back to the first item. A ring also has operations to get the current item, add an item, and remove an item. The details of where an item is added and which item is removed are up to you.

Design an ADT to represent a ring of objects. Specify each operation by stating its purpose, by describing its parameters, and by writing a pseudocode version of its header. Then write a C++ interface for a ring's methods. Include javadoc-style comments in your code.
5. A set is a special bag that does not allow duplicates. Specify each operation for a set of objects by stating its purpose, by describing its parameters, and by writing preconditions, postconditions, and a pseudocode version of its header. Then write a C++ interface for the set. Include javadoc-style comments in your code.
6. Imagine a pile of books on your desk. Each book is so large and heavy that you can remove only the top one from the pile. You cannot remove a book from under another one. Likewise, you can add another book to the pile only by placing it on the top of the pile. You cannot add a book beneath another one.

If you represent books by their titles alone, design a class that you can use to track the books in the pile on your desk. Specify each operation by stating its purpose, by describing its parameters, and by writing a pseudocode version of its header. Then write a C++ interface for the pile's methods. Include javadoc-style comments in your code.

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## C++ Interlude

## C++ Classes

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## Prerequisites

Appendix A Review of C++ Fundamentals
Chapter 1 Data Abstraction: The Walls

Throughout this book we design abstract data types and implement them as data structures using C++. This C++ Interlude provides a refresher on C++ classes and then introduces additional C++ tools we can use to define our abstract data types in a flexible manner that maintains the wall between our design and the implementation.

After reviewing classes, we will look at class templates. This C++ construct gives us the power to specify the data type of the items contained in a data structure in a very generic way. For example, if you had spent two weeks developing a great class to represent a bag of integers, wouldn't it be great if you could easily use the same code for a bag that holds strings? Class templates allow you to define classes that are independent of the type of data stored in the data structure. When a client is ready to instantiate an object of the class, the client can specify the type of data the object holds.

Header (.h), or specification, files in C++ provide a mechanism to partially separate the design of a class from the implementation in the source, or implementation (.cpp), file. The header file must also contain a description of both the data fields for the class and any private methods used by the class. A client does not need to know about the private methods or data fields to use the class in a program. To provide a public interface for an ADT, you can write an abstract base class, thereby separating design from implementation. An abstract base class allows the client to take full advantage of polymorphism when using our class.

To introduce these concepts, let's look at a simple problem that illustrates all three of the fundamental concepts of object-oriented programming: encapsulation, inheritance, and polymorphism.

C++ classes

## C1.1 A Problem to Solve

Suppose that a friend who is creating a video game asked you to design and develop a group of classes to represent three types of boxes carried by the characters in the game. Each type of box can only hold one item. A character can put an item in the box or look at the item in the box. The three types of boxes are:

- Plain box-a plain old box that holds only one item.
- Toy box-a box that has color and holds only one item.
- Magic box-a box that holds only one item, but magically changes it to the first item that was ever stored in the box.

Let's begin by designing an ADT plain box. It is the simplest of the three boxes, and we may be able to use aspects of its implementation for the other two boxes. Since game characters can only place items in the box or look at them, the box needs only two public methods, setItem to place an item in the box and getItem to get the stored item. Here is the UML notation for those methods:

```
+setItem(theItem: ItemType)
+getItem(): ItemType
```

We can define ItemType as the type of item stored in the box by using a typedef statement. For example, to have the box hold a double, we can write

## typedef doub7e ItemType;

To have the box hold another type of data, we need only to replace the word double with the new data type. Listing C1-1 shows the declaration of our class PlainBox that would appear in a header file. To save space here, we show only comments that describe the class element type and omit the specifications, preconditions, and postconditions. We have used the object-oriented concept of encapsulation to group together data - the item stored in the box-with the methods that operate on that data: the two constructors, a method setItem to change the item's value, and a method getItem to return the item's value.

## LISTING C1-1 The header file for the class PlainBox

```
/** @file PlainBox.h */
#ifndef _PLAIN_BOX
#define _PLAIN_BOX
// Set the type of data stored in the box
typedef doub7e ItemType;
```

```
// Declaration for the class PlainBox
class PlainBox
{
private:
    // Data field
    ItemType item;
public:
    // Defau7t constructor
    PlainBox();
    // Parameterized constructor
    PlainBox(const ItemType& theItem);
    // Method to change the value of the data field
    void setItem(const ItemType& theItem);
    // Method to get the value of the data field
    ItemType getItem() const;
}; // end P7ainBox
#endif
```

Let's look in detail at this $\mathrm{C}++$ class declaration.

## C1.1.1 Private Data Fields

The data field item is declared in a private section of the class declaration. Every data field in every class presented in this textbook is in a private section. This restricts access to the data field to just the class in which it is defined. Typically, we provide methods-such as setItem and getItem-to access the data fields. In this way, the class controls how and whether other classes can access the data fields. This design principle should lead to programs that not only are easier to debug, but also have fewer logical errors from the beginning.

Clients and derived classes should not have direct access to the data fields of a class. If a class designer believes that a derived class might need to access or modify the data fields of the base class, you should still make the data fields private but provide protected methods so that any derived classes can access or modify the data. More information on access modifiers, such as private and protected, is presented in a later C++ Interlude.

As we build more complex data structures, we need to guarantee their integrity. This is simpler to do if we restrict access to the data fields to only our class. For example, suppose we have a class that stores items in an array and maintains a count of the number of items used in the array. Each time our class adds or removes an item from the array, the item count needs to be modified to reflect the change. If a client of our class had direct access to the array, the client could add and remove entries in the array and neglect to update the item counter. This counter would not accurately reflect the number of entries in the array, resulting in, for example, the loss of data or an abnormal termination of the program.

## C1.1.2 Constructors and Destructors

Classes have two types of special methods, called constructors and destructors. A constructor allocates memory for new instances of a class and can initialize the object's data to specified values. A destructor destroys an instance of a class when the object's lifetime ends. A typical class has several
constructors but only one destructor. For many classes, you can omit the destructor. In such cases, the compiler generates a destructor for you. For the classes in this C++ Interlude, the compiler-generated destructor is sufficient. C++ Interlude 2 discusses how and why you would write your own destructor.

In C++, a constructor has the same name as the class. Constructors have no return type-not even void-and cannot use return to return a value. A class can have more than one constructor, as is the case for the class P1ainBox. One of the constructors, the default constructor, has no parameters. Typically, a default constructor initializes data fields to values that the class implementation chooses. Other constructors have parameters. These parameterized constructors initialize data fields to values chosen by the client but approved by the constructor. The compiler decides which constructor to call by matching the argument list supplied by the client with the parameters of the available constructors. A match occurs when the arguments and parameters correspond in number, data type, and order.

()
Note: If you do not define any constructors for a class, the compiler creates a default constructor - one without parameters. Once you define a constructor, the compiler does not create any of its own. Therefore, if you define a parameterized constructor but not a default constructor, your class will not have a default constructor.

When you declare an instance of the class, a constructor is invoked implicitly. For example, the statement

```
PlainBox myBox;
```

invokes the default constructor, which creates the object myBox and initializes the data field item to a value given in the constructor's definition. Notice that you do not include parentheses after myBox when invoking the default constructor. The statement

```
PlainBox myBox(specialValue);
```

invokes the parameterized constructor, which, as you will see, initializes the data field item to specialValue. Thus, specialValue is stored in the box.

## C1.1.3 Methods

As discussed in Chapter 1, methods implement the algorithms that solve a problem. A method prototype conveys three important pieces of information to a client wishing to use our class: the method name, the number and types of its parameters, and the method's return type. The prototype describes the slit in the wall of abstraction. Prototypes for methods a client can use are in the public section of the class declaration.

An accessor method in a class accesses, or gets, the value of a data field. Its name often begins with the word get. Accessor methods do not change the data fields of an object. An important part of the PlainBox class declaration is the accessor method getItem, which is labeled with the keyword const:

ItemType getItem() const;
This keyword is a signal to both the compiler and other programmers that the method does not change the data fields of the object. Another way to think of a const method is that the object is the same after calling this kind of method as it was before calling the method. Using const in a method declaration protects you as a programmer when you implement the algorithm for the method. As you write the code for a method declared with const, the compiler can check your code to verify that you did not modify any data fields.

Programming Tip: Methods declared const can access and use data fields, so labeling all accessor methods with the const declaration is appropriate and a good idea.

A mutator method in a class changes the value of a data field. Often, the name of a mutator method begins with the word set. The method setItem is an example of a mutator method. Mutator methods cannot be declared as const. Likewise, constructors cannot be declared const, because they must initialize the data fields of an object.

Passing parameters by constant reference. The method setItem and the parameterized constructor both have a parameter theItem that is passed by constant reference, as you can see from their declarations:

```
void setItem(const ItemType& theItem);
P1ainBox(const ItemType& theItem);
```

Passing a parameter by constant reference provides several benefits to both the client of a class and the programmer implementing the class. Passing an argument by reference to a method, especially when the argument is a complex object, saves time and memory, since the method can access or modify the object without copying it. The risk with this technique is that the method has access to data declared outside of its class; the method has "broken through the wall" and can modify an item owned by the client that invoked the method. To keep the efficiency of passing by reference and still protect the data of our client, we use the keyword const before the declaration of that parameter. The method treats the parameter as a constant that cannot be modified. Using const with parameters passed by reference protects client objects and reduces the chance of side effects. Our method can still access and use an object passed by constant reference, but the compiler flags any modifications to the object as errors.

## Note: Guidelines for safe and secure programming

- Declare all data fields in the private section of the class declaration.
- Declare as const any method that does not change the object (accessor methods).
- Declare any parameter passed by reference as const, unless you are certain it must be modified by the method, in which case the method should be either protected or private.


## C1.1.4 Preventing Compiler Errors

You can declare C++ variables only once within each program block. A second declaration, or redefinition, of a variable in the same block results in a compiler error. For example, suppose you have the following loop:

```
for (int index = 0; index < count; index++)
{
    int index = count / 6; // index is redefined in loop - compiler error!
    tota1 += index;
} // end for
```

Redefining index inside the loop results in a compile-time error, since index was already defined in the for statement. The same restriction on redefinition applies to entire classes. Let's see how we could accidentally redefine our PlainBox class.

If we want to use a PlainBox object in our program, for instance in our main function, we must include PlainBox. h in the file containing the main function, main.cpp, so the compiler knows about the public interface for our class. We also need to include the header file for our PlainBox class in PlainBox.cpp so the compiler knows about the data fields and methods of the class when it compiles the class's implementation.

This creates a problem, since the compiler reads our PlainBox definition twice, even though the class needs to be defined only once. We don't know which file the compiler will try to compile first, so both files must include PlainBox.h. This duplication results in the class being redefined when the header file is read again. We need a way to prevent the compiler from reading the class definition a second time. The \#i fndef, \#define, and \#endif preprocessor directives provide a solution.

You are already familiar with the \#include directive, which includes the contents of another file in the current file. You use \#ifndef to conditionally include a class definition. For example, the header file for PlainBox in Listing C1-1 contains

```
#ifndef _PLAIN_BOX
#define _PLAIN_BOX
```

as its first two lines. The directive \#i fndef means "If _PLAIN_BOX is not defined, then ..." If the compiler had not defined the name _PLAIN_BOX, it would process the code that follows until it reached the \#endif directive at the end of the file.

The \#define directive defines the name _PLAIN_BOX. If another file includes the class definition of PlainBox, the name _PLAIN_BOX will already have been defined. That file's \#i fndef directive will cause the preprocessor to skip any of the code that follows; that is, the code will be hidden from the compiler, and so the compiler will not see our class definition more than once.

In this textbook, all header files use these preprocessor directives to protect against including class definitions multiple times. Though you can use any names in the directives, the name to test in \#i fndef and the name defined by \#define must be the same. We write such names in uppercase and begin them with an underscore character.

## C1.2 Implementing a Solution

After we have designed our solution, the next step is to implement each of the methods we declared in the header file. Listing C1-2 shows the implementation, or source code, file for the class PlainBox. The method implementations are simple, but there are a few syntax items we should discuss.

## LISTING C1-2 Implementation file for the PlainBox class

/** @file PlainBox.cpp */
\#include "PlainBox.h"
PlainBox: : PlainBox()
\{
\} // end default constructor
PlainBox::PlainBox (const ItemType\& theItem)
\{
item = theItem;
\} // end constructor
void P1ainBox::setItem(const ItemType\& theItem)

```
{
    item = theItem;
} // end setItem
ItemType PlainBox::getItem() const
{
    return item;
} // end getItem
```

The header file PlainBox.h contains our class declaration. We must use the preprocessor directive \#include to include the PlainBox class declaration so the compiler can validate our method headers and provide access to class data fields. The method headers must match those provided in the class declaration. Including the header file does not tell the compiler that the methods defined here are part of the PlainBox class. To do so, you must precede the constructor and method names with the class name followed by two colons-that is,

```
PlainBox::
```

The reason for this requirement is that $\mathrm{C}++$ allows a source file to contain the implementations of methods for several classes and stand-alone functions. The namespace indicator PlainBox: : is a prefix that indicates to the compiler that the method is a part of the PlainBox namespace. A C++ namespace is a syntax structure, such as a class, that allows you to group together declarations of data and methods under a common name, such as PlainBox. Once a method has been defined as part of a namespace, it has access to all the data and methods in that namespace. Thus, the constructor and methods in Listing C1-2 have access to PlainBox's data fields and methods.


Note: To better modularize our code, the implementation files in this book will contain only methods from a single class.

## C1.3 Templates

At this point, our PlainBox class looks similar to many of the classes you probably have implemented in a previous C++ course. The class works well, as long as the characters in the game want to store a double value in a box. But what do we do if one character wants to store a double and a second has a string or MagicWand object to store? Since our current PlainBox class can store only a double, we would need to create new classes-PlainBoxForStrings and PlainBoxForWands, for exampleto hold different object types. These classes would function in exactly the same way as the PlainBox class, so we could copy the code with only a few changes. For example, we have to change the typedef in each header file. For the PlainBoxForStrings class, the typedef would be
typedef string ItemType;
and for the PlainBoxForWands class, the typedef would be

## typedef MagicWand ItemType;

We also would need to change the names of the constructors and the namespace indicators to reflect the new class names. If we had a more complex class, these changes would require a substantial amount of effort and would be an error-prone process. Later, if our friend changed the
requirements for PlainBox, we would need to go through each of the PlainBox class variations and make the required changes.

The root of the problem in this scenario is that the programmer must know what types of objects will be stored in the box before the program is built. The programmer must then write a different class for each type of object that is to be stored. The functionality of each box is the same, but the type of data differs. Templates enable the programmer to separate the functionality of an implementation from the type of data used in the class. Listing C1-3 contains the header file for a template version of the class PlainBox.

## LISTING C1-3 Template header file for the PlainBox class

```
/** @file PTainBox.h */
#ifndef _PLAIN_BOX
#define _PLAIN_BOX
template<class ItemType>; // Indicates this is a template definition
// Declaration for the class PlainBox
class PlainBox
{
private:
    // Data field
    ItemType item;
public:
    // Defau7t constructor
    PlainBox();
    // Parameterized constructor
    P1ainBox(const ItemType& theItem);
    // Mutator method that can change the value of the data field
    void setItem(const ItemType& theItem);
    // Accessor method to get the value of the data field
    ItemType getItem() const;
}; // end P7ainBox
#include "PlainBox.cpp" // Include the implementation file
#endif
```

As you can see from Listing C1-3, changing our earlier definition of the class PlainBox in Listing C1-1 to be a template requires the change of only one line in our header file: The typedef was removed and replaced with a statement to indicate this class is a template:
template<class ItemType>;
However, we also must include the implementation file just prior to the \#endif directive by writing

```
#include "PlainBox.cpp"
```

This addition is necessary because the compiler does not compile a template class until it sees the client's instantiation of the template and knows the actual data type corresponding to the data-type parameter ItemType.

Programming Tip: In your development environment, do not add the implementation file-PlainBox.cpp, for example-to the project. It will automatically be included when it is needed.

The implementation file PlainBox.cpp requires a few changes, but these changes all follow a pattern. Prior to each method's definition, you write the same template statement:

```
template<class ItemType>;
```

to indicate that the method is a template. The namespace indicator, PlainBox<ItemType> : :, must also precede each method name to reflect that the method's definition is based on ItemType. These changes are shown in Listing C1-4.

## LISTING C1-4 Implementation file for the PlainBox template class

```
/** @file PTainBox.cpp */
```

template<c1ass ItemType>;
PlainBox<ItemType>: : PlainBox()
\{
\} // end default constructor
template<class ItemType>;
PlainBox<ItemType>: :PlainBox(const ItemType\& theItem)
\{
item = theItem;
\} // end constructor
template<c1ass ItemType>;
void PlainBox<ItemType>: :setItem(const ItemType\& theItem)
\{
item = theItem;
\} // end setItem
template<c1ass ItemType>;
ItemType PlainBox<ItemType>::getItem() const
\{
return item;
\} // end getItem

To instantiate an instance of PlainBox, you write the data type of the item to be placed in a box, surrounded by angle brackets:

```
P1ainBox<doub7e> numberBox; // A box to hold a doub7e
PlainBox<string> nameBox; // A box to hold a string object
PlainBox<MagicWand> wandBox; // A box to hold a MagicWand object
```

Methods of these box instances are invoked as before:

```
doub7e hea7th = 6.5;
numberBox.setItem(health);
```

```
string secretName = "Rumpelstiltskin";
nameBox.setItem(secretName);
MagicWand elfWand;
wandBox.setItem(e7fWand);
```

Note: By using templates, you can define a class that involves data of any type, even data types that are created after you designed and implemented your class.

## C1.4 Inheritance

Now that the class PlainBox has been written, we will look at the other two boxes. The toy box is very similar to our plain box, but it has an additional characteristic, color, which we can represent with an enumerated type Color. We set the color of our toy box when it is created. The characters in the game can ask about the box's color but cannot change it. Here is the UML notation for those methods:

```
+getColor(): string
+setItem(ItemType theItem)
+getItem(): ItemType
```

The methods setItem and getItem behave exactly the same as the similarly named methods implemented for the PlainBox class. We can reuse the code in PlainBox by using inheritance. Inheritance allows us to reuse and extend work we have already completed and tested.

## C1.4.1 Base Classes and Derived Classes

We can use PlainBox as a base class, or superclass, for our ToyBox class. You can think of a base class as a parent class. The ToyBox class is the derived class, or subclass, of the PlainBox class. To indicate that ToyBox is derived from PlainBox, we use the following syntax in the class header of ToyBox:

```
class ToyBox : public PlainBox<ItemType>
```

Any instance of the derived class is also considered to be an instance of the base class and can be used in a program anywhere that an instance of the base class can be used. Also, when the keyword public is used with the base class, any of the publicly defined methods or data fields in the base class can be used by instances of the derived class. In C++, a derived class inherits all the members of its base class, except the constructors and destructor. That is, a derived class has the data fields and methods of the base class in addition to the members it defines, though it can access only publicly defined members. A derived class can also revise any inherited public method.

As you can see in Listing C1-5, ToyBox defines only the constructors and new methods specific to it, if it has not needed to revise any methods of the base class. The methods setItem and getItem are inherited from the PlainBox class and can be used in our ToyBox class.

## LISTING C1-5 Template header file for the class ToyBox

/** @file ToyBox.h */
\#ifndef _TOY_BOX
\#define _TOY_BOX
\#include "PlainBox.h"

```
enum Color {BLACK, RED, BLUE, GREEN, YELLOW, WHITE};
temp1ate<c1ass ItemType>
class ToyBox : public PlainBox<ItemType>
{
private:
    Color boxColor;
public:
    ToyBox();
    ToyBox(const Color& theColor);
    ToyBox(const ItemType& theItem, const Color& theColor);
    Color getColor() const;
}; // end ToyBox
#include "ToyBox.cpp"
#endif
```

An instance of the class ToyBox has two data fields-item, which is inherited, and boxColor, which is new. Because an instance of a derived class can invoke any public method in the base class, an instance of ToyBox has all the methods that PlainBox defines; new constructors; a new, compilergenerated destructor; and a new method getColor. Although an instance of a derived class contains copies of inherited data fields, the code for inherited methods is not copied.

A derived class cannot access the private members of the base class directly, even though they are inherited. Inheritance does not imply access. After all, you can inherit a locked vault but be unable to open it. In the current example, the data field item of PlainBox is private, so you can reference it only within the definition of the class PlainBox and not within the definition of ToyBox. However, the class ToyBox can use PlainBox's public methods setItem and getItem to set or obtain the value of item indirectly.

## LISTING C1-6 Implementation file for the class ToyBox

```
/** @file ToyBox.cpp */
temp1ate<class ItemType>
ToyBox<ItemType>::ToyBox()
{
    PlainBox<ItemType>();
    boxColor = BLACK;
} // end default constructor
template<c1ass ItemType>
ToyBox<ItemType>::ToyBox(const Color& theColor)
{
    PlainBox<ItemType>();
    boxColor = theColor;
} // end constructor
temp1ate<c1ass ItemType>
ToyBox<ItemType>: :ToyBox(const ItemType& theItem,
const Color& theColor)
```

```
{
    P1ainBox<ItemType>();
    PlainBox<ItemType>::setItem(theItem);
    boxColor = theColor;
} // end constructor
temp1ate<class ItemType>
Color ToyBox<ItemType>::getColor() const
{
    return boxColor;
} // end getColor
```

Inheritance enables you to reuse software components when you define a new class. Our ToyBox class does not inherit the constructors from PlainBox, but it can use a PlainBox constructor or method to initialize the private data field in PlainBox. To use the default constructor of our base class, we place the statement

PlainBox<ItemType>();
as the first statement in each of the ToyBox constructors. This ensures that the data fields in the base class are correctly initialized. Only the default constructor can be called in this way. If we need more control over setting base-class data fields, a derived-class constructor must call the appropriate baseclass mutator methods, as shown in the third ToyBox constructor. In this event, the base-class namespace indicator should precede the method name, so the compiler knows where to find the method's code. Using the namespace indicator is necessary only within constructors, as you'll see when we implement the MagicBox class.

A client of our ToyBox class can create and use a ToyBox object in much the same way that we did for PlainBox:

```
string favoriteToy = "Jack-in-the-Box";
ToyBox<string> myToyCase(favoriteToy, RED); // A red toy box
string oldToy = myToyCase.getItem(); // oldToy is a Jack-in-the-Box
favoriteToy = "Spinning Top";
myToyCase.setItem(favoriteToy); // myToyCase now holds a Spinning Top
```


## C1.4.2 Overriding Base-Class Methods

You can add as many new members to a derived class as you like. Although you cannot revise a base class's private data fields and should not reuse their names, you can redefine inherited methods; this is called overriding a base-class method. A method in a derived class overrides, or redefines, a method in the base class if the two methods have the same name and parameter declarations-that is, if they have the same signatures.

To see an example of when we may need to redefine a base-class method, let's consider the implementation of the class MagicBox. This box magically changes any item placed inside to a copy of the first item it ever held. A simple way to do this is to store only the first item. Future calls to setItem will do nothing. Therefore, every call to getItem returns the first item stored in the box.

As with the toy box, we can derive our class from the PlainBox class, but we also need a data field to indicate if an initial item has been stored in our box. We can use a boolean field, firstItemStored, which is set to true once the first item has been stored in the box. We can then check this flag to see what setItem should do. If firstItemStored is false, we know that the argument passed to
setItem is the first item we are asked to store. If firstItemStored is true, we know this is not the first item, and we simply do nothing.

Here is the UML notation for our MagicBox methods:

```
+setItem(ItemType theItem)
+getItem(): ItemType
```

These methods are like those in the class PlainBox, but we need a different implementation of setItem to meet the special requirements of this box. Therefore, we need to override the PlainBox implementation of setItem. Listing C1-7 gives the header file for our class MagicBox. It has the private data field firstItemStored, a default constructor, a parameterized constructor, and the special version of setItem that we need.

## LISTING C1-7 Header file for the class MagicBox

```
/** @file MagicBox.h */
#ifndef _MAGIC_BOX
#define _MAGIC_BOX
#include "PlainBox.h"
template<class ItemType>
class MagicBox : public PlainBox<ItemType>
{
private:
    bool firstItemStored;
public:
    MagicBox();
    MagicBox(const ItemType& theItem);
    void setItem(const ItemType& theItem);
}; // end MagicBox
#include "MagicBox.cpp"
#endif
```

The implementation of the default constructor must call the base-class constructor and then initialize the data field firstItemStored to true to indicate that the first item has not been stored yet. The method setItem must check this data field and, if no item has been stored, call the base-class method setItem to store the parameter in the data field item. You can see these implementations in Listing C1-8.

## LISTING C1-8 Implementation file for the class MagicBox

```
/** @file MagicBox.cpp */
template<class ItemType>
MagicBox<ItemType>: :MagicBox()
{
    PlainBox<ItemType>();
```

```
    firstItemStored = false; // Box has no magic initially
} // end default constructor
temp1ate<c1ass ItemType>
MagicBox<ItemType>: :MagicBox(const ItemType& theItem)
{
    firstItemStored = false; // Box has no magic initially
    setItem(theItem);
    // Box has magic now
} // end constructor
temp7ate<class ItemType>
void MagicBox<ItemType>::setItem(const ItemType& theItem)
{
    if (!firstItemStored)
    {
        P1ainBox<ItemType>::setItem(theItem);
        firstItemStored = true; // Box now has magic
    } // end if
} // end setItem
```

To simplify the parameterized constructor, we can call the setItem method to set the first item stored in the box. However, the MagicBox class has two setItem methods available, its own and the base-class version. As discussed earlier, we would precede a method name with the base-class namespace indicator to tell the compiler to use the base-class version of the method. To use this class's setItem method, we can simply invoke the method as shown in the parameterized constructor. Without a namespace indicator in front of the method name, the compiler assumes that the programmer is referring to the current class and inserts code to call that version of the method.

?Question 1 Revise the parameterized constructor to call the base-class's constructor instead of MagicBox's constructor.

## C1.5 Virtual Methods and Abstract Classes

When we designed the interface for our ADT bag in Chapter 1, we created an abstract class that describes each of the public methods. Recall that an abstract class in $\mathrm{C}++$ contains at least one method declared as virtual that has no implementation. An abstract class can only serve as a base class; it cannot be instantiated. The derived class must define the methods specified but not already implemented by the abstract class.

## C1.5.1 Virtual Methods

Using the keyword virtual in front of the prototype, or header, of the method tells the $\mathrm{C}++$ compiler that the code this method executes is determined at runtime, not when the program is compiled. A method declared this way is referred to as a virtual method. Why should we declare a method as virtual?

The rules of inheritance allow us to use a derived class anywhere that its base class is used. For example, we could declare a variable that is a plain box:

```
PlainBox<string> cardBox;
```

Then we could create a MagicBox object and assign it to the variable cardBox:

```
cardBox = MagicBox<string>("Queen of Hearts");
```

Later, the player might store a different string in the box, and so we execute

```
cardBox.setItem("Jack of Spades");
```

Because the compiler considers the object cardBox to be of type PlainBox<string>, the PlainBox version of the method setItem is used in the previous statement. The PlainBox implementation of setItem stores the value of its parameter in item. Thus, a problem arises when our client calls the method getItem:

```
string myFirstCard = cardBox.getItem();
```

The method getItem returns the string "Jack of Spades" instead of "Queen of Hearts". This is not the behavior the client desires, because cardBox was assigned an instance of MagicBox.

If we declare P1ainBox's method setItem to be virtual in the P1ainBox header file by writing

```
virtual ItemType setItem() const;
```

the version of setItem invoked during execution is determined by the specific type of object stored in nameBox. In our example, the MagicBox version of setItem would be called, since that is the type of object most recently stored in cardBox. Having this decision delayed until the program executes is an example of polymorphism and is a key benefit of object-oriented programming.

To fully implement our box example as a demonstration of polymorphic code, and to thoroughly discuss virtual methods and abstract classes, we need to use pointers. Pointers are introduced in $\mathrm{C}++$ Interlude 2, and so we will delay further discussion and the completion of this example until then.

## C1.5.2 Abstract Classes

The previous example used a virtual method so that a choice could be made during program execution between two implementations of the setItem method. We can use this same idea to assist with our goal of designing each ADT as a public interface that describes its functionality and is implementation independent. Declaring our ADT methods as virtual allows an application using our class to take advantage of polymorphism when the ADT's methods are invoked.

One significant difference between such an ADT and the previous example that uses the PlainBox class is that we do not want to provide an implementation of the ADT methods. Instead, we want to force the classes derived from the ADT to provide the implementations. Yet if we do not provide an implementation of the ADT, the compiler or linker will issue an error message when we try to build a program using our class.

We can avoid this error by writing our methods as pure virtual methods. A pure virtual method is a virtual method that has no implementation. We can now say that an abstract class is one that has at least one pure virtual method. To tell the compiler that a virtual method is a pure virtual method, you write $=0$ before the semicolon at the end of the method prototype, as in the following example:

```
virtual void setItem(const ItemType& theItem) = 0;
```

Listing C1-9 defines the template class BoxInterface that provides a public interface for the box classes described earlier in this interlude. BoxInterface is an abstract class, because it contains at least one pure virtual method. Remember that abstract classes cannot be directly instantiated.

```
LISTING C1-9 An abstract class that is an interface for the ADT box
/** @file BoxInterface.h */
#ifndef _BOX_INTERFACE
#define _BOX_INTERFACE
template <class ItemType>
class BoxInterface
{
public:
    virtual void setItem(const ItemType& theItem) = 0;
    virtual ItemType getItem() const = 0;
}; // end BoxInterface
#endif
```

We can indicate that our class PlainBox is derived from BoxInterface by changing its class header to
class PlainBox : public BoxInterface<ItemType>
In the future when we design the public interface for an ADT, we will express that interface as an abstract class. When we introduce pointers in $\mathrm{C}++$ Interlude 2, we will explore the importance and use of abstract classes in depth. Until then, we will use only abstract classes as a design tool for our ADTs.

## Recursion: The Mirrors

## Chapter

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The goal of this chapter is to ensure that you have a basic understanding of recursion, which is one of the most powerful techniques available to the computer scientist. This chapter assumes that you have had little or no previous introduction to recursion. If, however, you have already studied recursion, you can review this chapter as necessary.

By presenting several relatively simple problems, the chapter demonstrates the thought processes that lead to recursive solutions. These problems are diverse and include examples of counting, searching, and organizing data. In addition to presenting recursion from a conceptual viewpoint, this chapter discusses techniques that will help you understand the mechanics of recursion. These techniques are particularly useful for tracing and debugging recursive functions.

Some recursive solutions are far more elegant and concise than the best of their nonrecursive counterparts. For example, the classic Towers of Hanoi problem appears to be quite difficult, yet it has an extremely simple recursive solution. On the other hand, some recursive solutions are terribly inefficient, as you will see, and should not be used.

Chapter 5 continues the formal discussion of recursion by examining more difficult problems. Recursion will play a major role in many of the solutions that appear throughout the remainder of this book.

### 2.1 Recursive Solutions

Recursion breaks a problem into smaller identical problems

## Some recursive

 solutions are inefficient and impractical
## Complex problems

 can have simple recursive solutionsRecursion is an extremely powerful problem-solving technique. Problems that at first appear to be quite difficult often have simple recursive solutions. Like other problem-solving techniques, recursion breaks a problem into several smaller problems. What is striking about recursion is that these smaller problems are of exactly the same type as the original problem-mirror images, so to speak.

Did you ever hold a mirror in front of another mirror so that the two mirrors face each other? You will see many images of yourself, each behind and slightly smaller than the other. Recursion is like these mirror images. That is, a recursive solution solves a problem by solving a smaller instance of the same problem! It then solves this new problem by solving an even smaller instance of the same problem. Eventually, the new problem will be so small that its solution will be either obvious or known. This solution will lead to the solution of the original problem.

For example, suppose that you could solve problem $P_{1}$ if you had the solution to problem $P_{2}$, which is a smaller instance of $P_{1}$. Suppose further that you could solve problem $P_{2}$ if you had the solution to problem $P_{3}$, which is a smaller instance of $P_{2}$. If you knew the solution to $P_{3}$ because it was small enough to be trivial, you would be able to solve $P_{2}$. You could then use the solution to $P_{2}$ to solve the original problem $P_{1}$.

Recursion can seem like magic, especially at first, but as you will see, it is a very real and important problem-solving approach that is an alternative to iteration. An iterative solution involves loops. You should know at the outset that not all recursive solutions are better than iterative solutions. In fact, some recursive solutions are impractical because they are so inefficient. Recursion, however, can provide elegantly simple solutions to problems of great complexity.

As an illustration of the elements in a recursive solution, consider the problem of looking up a word in a dictionary. Suppose you wanted to look up the word "vademecum." Imagine starting at the beginning of the dictionary and looking at every word in order until you found "vademecum." That is precisely what a sequential search does, and for obvious reasons, you want a faster way to perform the search.

One such approach is the binary search, which in spirit is similar to the way in which you actually use a dictionary. You open the dictionary - maybe to a point near its middle - and by glancing at
the page, determine which "half" of the dictionary contains the desired word. The following pseudocode is a first attempt to formalize this process:

```
/ Search a dictionary for a word by using a recursive binary search
    if (the dictionary contains only one page)
        Scan the page for the word
    e1se
    \{
        Open the dictionary to a point near the middle
        Determine which half of the dictionary contains the word
        if (the word is in the first half of the dictionary)
            Search the first half of the dictionary for the word
        e1se
            Search the second half of the dictionary for the word
    \}
```

Parts of this solution are intentionally vague: How do you scan a single page? How do you find the middle of the dictionary? Once the middle is found, how do you determine which half contains the word? The answers to these questions are not difficult, but they would only obscure the solution strategy right now.

The previous search strategy reduces the problem of searching the dictionary for a word to a problem of searching half of the dictionary for the word, as Figure 2-1 illustrates. Notice two important points. First, once you have divided the dictionary in half, you already know how to search the appropriate half: You can use exactly the same strategy that you employed to search the original dictionary. Second, note that there is a special case that is different from all the other cases: After you have divided the dictionary so many times that you are left with only a single page, the halving ceases. At this point, the problem is sufficiently small that you can solve it directly by scanning the single page that remains for the word. This special case is called the base case (or basis or degenerate case).

## FIGURE 2-1 A recursive solution



This strategy is called divide and conquer. You solve the dictionary search problem by first dividing the dictionary into two halves and then conquering the appropriate half. You solve the smaller problem by using the same divide-and-conquer strategy. The dividing continues until you reach the base case. As you will see, this strategy is inherent in many recursive solutions.

To further explore the nature of the solution to the dictionary problem, consider a slightly more rigorous formulation.

```
search(aDictionary: Dictionary, word: string)
    if (aDictionary is one page in size)
        Scan the page for word
    else
    {
Open aDictionary to a point near the middle Determine which half of aDictionary contains word
```

A binary search of a dictionary

A base case is a special case whose solution you know

A binary search uses a divide-andconquer strategy

A recursive function calls itself

Each recursive call solves an identical, but smaller, problem

A test for the base case enables the recursive calls to stop

Eventually, one of the smaller problems must be the base case

```
if (word is in the first half of aDictionary)
    search ( first half of aDictionary, word)
else
    search(second half of aDictionary, word)
}
```

Writing the solution as a function allows several important observations:

1. One of the actions of the function is to call itself; that is, the function search calls the function search. This action is what makes the solution recursive. The solution strategy is to split aDictionary in half, determine which half contains word, and apply the same strategy to the appropriate half.
2. Each call to the function search made from within the function search passes a dictionary that is one-half the size of the previous dictionary. That is, at each successive call to search (aDictionary, word), the size of aDictionary is cut in half. The function solves the search problem by solving another search problem that is identical in nature but smaller in size.
3. There is one search problem that you handle differently from all of the others. When aDictionary contains only a single page, you use another approach: You scan the page directly. Searching a one-page dictionary is the base case of the search problem. When you reach the base case, the recursive calls stop and you solve the problem directly.
4. The manner in which the size of the problem diminishes ensures that you will eventually reach the base case.

These facts describe the general form of a recursive solution. Though not all recursive solutions fit these criteria as nicely as this solution does, the similarities are far greater than the differences. As you attempt to construct a new recursive solution, you should keep in mind the following four questions.

## Note: Four questions for constructing recursive solutions

1. How can you define the problem in terms of a smaller problem of the same type?
2. How does each recursive call diminish the size of the problem?
3. What instance of the problem can serve as the base case?
4. As the problem size diminishes, will you reach this base case?

Now consider two relatively simple problems: computing the factorial of a number and writing a string backward. Their recursive solutions further illustrate the points raised by the solution to the dictionary search problem. These examples also illustrate the difference between a recursive valued function-which returns a value - and a recursive void function.

### 2.2 Recursion That Returns a Value

The mechanics of recursion are clearer when the recursive function returns a value instead of just performing an action, and so we examine one in detail.

### 2.2.1 A Recursive Valued Function: The Factorial of $\boldsymbol{n}$

Computing the factorial of an integer $n$ is a good first example because its recursive solution is easy to understand and neatly fits the mold described earlier. However, because the problem has a simple and efficient iterative solution, you should not use the recursive solution in practice.

To begin, consider the familiar iterative definition of factorial( $n$ ) (more commonly written as $n!$ ):

$$
\begin{aligned}
& \operatorname{factorial}(n)=n \times(n-1) \times(n-2) \times \cdots \times 1 \quad \text { for an integer } n>0 \\
& \operatorname{factorial}(0)=1
\end{aligned}
$$

The factorial of a negative integer is undefined. You should have no trouble writing an iterative factorial function based on this definition.

To define factorial( $n$ ) recursively, you first need to define factorial( $(n)$ in terms of the factorial of a smaller number. To do so, simply observe that the factorial of $n$ is equal to the factorial of $(n-1)$ multiplied by $n$; that is,

$$
\begin{aligned}
\operatorname{factorial}(n) & =n \times[(n-1) \times(n-2) \times \cdots \times 1] \\
& =n \times \operatorname{factorial}(n-1)
\end{aligned}
$$

The definition of factorial ( $n$ ) in terms of factorial ( $n-1$ ), which is an example of a recurrence relation, implies that you can also define factorial ( $n-1$ ) in terms of factorial( $n-2$ ), and so on. This process is analogous to the dictionary search solution, in which you search a dictionary by searching a smaller dictionary in exactly the same way.

The definition of factorial ( $n$ ) lacks one key element: the base case. As was done in the dictionary search solution, here you must define one case differently from all the others, or else the recursion will never stop. The base case for the factorial function is factorial $(0)$, which you know is 1 . Because $n$ originally is greater than or equal to zero and each call to factorial decrements $n$ by 1 , you will always reach the base case. With the addition of the base case, the complete recursive definition of the factorial function is

$$
\operatorname{factorial}(n)= \begin{cases}1 & \text { if } n=0 \\ n \times \operatorname{factorial}(n-1) & \text { if } n>0\end{cases}
$$

To be sure that you understand this recursive definition, apply it to the computation of facto$\operatorname{rial}(4)$. Because $4>0$, the recursive definition states that

$$
\text { factorial }(4)=4 \times \text { factorial }(3)
$$

Similarly,

$$
\begin{aligned}
\text { factorial }(3) & =3 \times \text { factorial }(2) \\
\text { factorial }(2) & =2 \times \text { factorial }(1) \\
\text { factorial }(1) & =1 \times \text { factorial }(0)
\end{aligned}
$$

You have reached the base case, and the definition directly states that

## factorial $(0)=1$

At this point, the application of the recursive definition stops and you still do not know the answer to the original question: What is factorial(4)? However, the information to answer this question is now available:

Because factorial $(0)=1$, factorial $(1)=1 \times 1=1$
Because factorial $(1)=1$, factorial $(2)=2 \times 1=2$
Because $\operatorname{factorial}(2)=2$, factorial $(3)=3 \times 2=6$
Because factorial $(3)=6$, factorial $(4)=4 \times 6=24$
You can think of recursion as a process that divides a problem into a task that you can do and a task that a friend can do for you. For example, if I ask you to compute factorial(4), you could first determine whether you know the answer immediately. You know immediately that factorial $(0)$ is 1 -that is, you know the base case-but you do not know the value of factorial(4) immediately. However, if your friend computes factorial(3) for you, you could compute factorial(4) by multiplying factorial(3) and 4. Thus, your task will be to do this multiplication, and your friend's task will be to compute factorial(3).

Your friend now uses the same process to compute factorial( 3 ) as you are using to compute factorial(4). Thus, your friend determines that factorial(3) is not the base case, and so asks another friend to compute factorial(2). Knowing factorial(2) enables your friend to compute factorial(3), and when you learn the value of factorial(3) from your friend, you can compute factorial(4).

A recurrence relation

A recursive definition of factorial

Notice that the recursive definition of factorial(4) yields the same result as the iterative definition, which gives $4 \times 3 \times 2 \times 1=24$. To prove that the two definitions of factorial are equivalent for all nonnegative integers, you would use mathematical induction. (See Appendix E.) Chapter 5 discusses the close tie between recursion and mathematical induction.

The recursive definition of the factorial function has illustrated two points. (1) Intuitively, you can define factorial $(n)$ in terms of factorial $(n-1)$. (2) Mechanically, you can apply the definition to determine the value of a given factorial. Even in this simple example, applying the recursive definition required quite a bit of work. That, of course, is where the computer comes in.

Once you have a recursive definition of factorial( $n$ ), it is easy to construct a $\mathrm{C}++$ function that implements the definition:

```
/** Computes the factorial of the nonnegative integer n.
    @pre n must be greater than or equal to 0.
    @post None.
    @return The factorial of n; n is unchanged. */
int fact(int n)
{
    if (n == 0)
        return 1;
    e1se // n>0, so n-1>= 0. Thus, fact(n-1) returns ( }n-1\mathrm{ )!
        return n * fact(n-1); // n* (n-1)! is n!
} // end fact
```

Suppose that you use the statement

```
cout << fact(3);
```

to call the function. Figure 2-2 depicts the sequence of computations that this call would require.

## FIGURE 2-2 fact(3)



This function fits the model of a recursive solution given earlier in this chapter as follows:

1. One action of fact is to call itself.
2. At each recursive call to fact, the integer whose factorial you need to compute is diminished by 1 .
3. The function handles the factorial of 0 differently from all the other factorials: It does not generate a recursive call. Rather, you know that fact ( 0 ) is 1 . Thus, the base case occurs when $n$ is 0 .
4. Given that $n$ is nonnegative, item 2 of this list assures you that you will always reach the base case.


Note: A recursive algorithm must have a base case, whose solution you know directly without making any recursive calls. Without a base case, a recursive function will generate an infinite sequence of calls.

The function fact requires as its precondition a nonnegative value of $n$. At the time of the recursive call fact $(n-1), n$ is positive, so $n-1$ is nonnegative. Because the recursive call satisfies fact's precondition, you can expect that fact $(n-1)$ will return the factorial of $n-1$. Therefore, $\mathrm{n} *$ fact $(\mathrm{n}-1)$ is the factorial of $n$. Chapter 5 uses mathematical induction to prove formally that fact $(\mathrm{n})$ returns the factorial of $n$.

If you ever violated fact's precondition, the function would not behave correctly. That is, if the calling program ever passed a negative value to fact, an infinite sequence of recursive callsterminated only by a system-defined limit-would occur, because the function would never reach the base case. For example, fact ( -4 ) would call fact ( -5 ), which would call fact ( -6 ), and so on.

The function ideally should protect itself by testing for a negative $n$. If $n<0$, the function could, for example, either return 0 or set an error flag. Appendix B discusses error checking in Sections B. 2 "Style" and B. 5 "Fail-Safe Programming"; you might want to review that discussion at this time.

At an intuitive level, it should be clear that the function fact implements the recursive definition of factorial. Now consider the mechanics of executing this recursive function. The logic is straightforward, except perhaps for the expression in the e1 se clause. This expression can be explained as follows:

1. Each operand of the product $n *$ fact $(n-1)$ is evaluated.
2. The second operand-fact $(n-1)$-is a call to the function fact. Although this is a recursive call (the function fact calls the function fact), there really is nothing special about it. Imagine substituting a call to another function-the standard function abs, for example-for the recursive call to fact. The principle is the same: Simply evaluate the function.

In theory, evaluating a recursive function is no more difficult than evaluating a nonrecursive function. In practice, however, manual bookkeeping can quickly get out of hand. The next section introduces a systematic way to trace the actions of a recursive function. For a computer, the bookkeeping is simple but can use more memory than is allocated for such a task.

Question 1 The following function computes the sum of the first $n \geq 1$ integers. Show how this function satisfies the properties of a recursive function.

```
/** Computes the sum of the integers from 1 through n.
```

/** Computes the sum of the integers from 1 through n.
apre n > 0.
apre n > 0.
@post None.
@post None.
@param n A positive integer.
@param n A positive integer.
@return The sum 1 + 2 + . . + n. */

```
@return The sum 1 + 2 + . . + n. */
```

fact satisfies the four criteria of a recursive solution

Violating fact's precondition causes "infinite" recursion

An activation record is created for each function call

Label each recursive call in the function

Each time a function is called, a new box represents its local environment

```
int sumUpTo(int n)
{
    int sum = 0;
    if (n == 1)
        sum = 1;
    else // n > 1
        sum = n + sumUpTo(n - 1);
    return sum;
} // end sumUpTo
```


### 2.2.2 The Box Trace

You can use a box trace both to help you understand recursion and to debug recursive functions. However, such a mechanical device is no substitute for an intuitive understanding of recursion. The box trace illustrates how compilers frequently implement recursion. As you read the following description of the technique, realize that each box roughly corresponds to an activation record, which a compiler typically uses in its implementation of a function call. C++ Interlude 2 will discuss activation records further.

The box trace is illustrated here for the recursive function fact, which returns a value.

1. Label each recursive call in the body of the recursive function. Several recursive calls might occur within a function, and it will be important to distinguish among them. These labels help you keep track of the correct place to which you must return after a function call completes. For example, mark the expression fact $(n-1)$ within the body of the function with the letter A:
```
if (n == 0)
    return 1;
else
    return n * fact(n-1);
```

                                    (A)
    You return to point A after each recursive call, substitute the computed value for the term fact ( $n-1$ ), and continue execution by evaluating the expression $n *$ fact $(n-1)$.
2. Represent each call to the function during the course of execution by a new box in which you note the local environment of the function. More specifically, each box will contain

- The values of the arguments of the argument list.
- The function's local variables.
- A placeholder for the value returned by each recursive call from the current box. Label this placeholder to correspond to the labeling in step 1.
- The value of the function itself.

When you first create a box, you will know only the values of the input arguments. You fill in the other values as you determine them from the function's execution. For example, you would create the box in Figure 2-3 for the call fact (3). (You will see in later examples that you must handle reference arguments somewhat differently than value arguments and local variables.)
3. Draw an arrow from the statement that initiates the recursive process to the first box. Then, when you create a new box after a recursive call, as described in step 2, you draw an arrow from the box that makes the call to the newly created box. Label each arrow to correspond to

FIGURE 2-4 The beginning of the box trace

the label (from step 1) of the recursive call; this label indicates exactly where to return after the call completes. For example, Figure 2-4 shows the first two boxes generated by the call to fact in the statement cout << fact (3).
4. After you create the new box and arrow as described in steps 2 and 3 , start executing the body of the function. Each reference to a value in the function's local environment references the corresponding value in the current box, regardless of how you generated the current box.
5. On exiting the function, cross off the current box and follow its arrow back to the box that called the function. This box now becomes the current box, and the label on the arrow specifies the exact location at which execution of the function should continue. Substitute the value returned by the just-terminated function call for the appropriate placeholder in the current box.

Figure 2-5 is a complete box trace for the call fact (3). In the sequence of diagrams in this figure, the current box is the deepest along the path of arrows and is highlighted in blue, whereas crossed-off boxes are dashed and shaded gray.

## FIGURE 2-5 Box trace of fact (3)

The initial call is made, and method fact begins execution:

```
n=3
A: fact(n-1)=?
return ?
```

At point A a recursive call is made, and the new invocation of the method fact begins execution:
\(\left.$$
\begin{array}{|l|}\hline \begin{array}{l}n=3 \\
\text { A: fact }(n-1)=? \\
\text { return? }\end{array}
$$ <br>

\hline\end{array}\right]\)| $\mathrm{n}=2$ |
| :--- |
| A: fact( $\mathrm{n}-1)=?$ |
| return ? |

At point $A$ a recursive call is made, and the new invocation of the method fact begins execution:


At point A a recursive call is made, and the new invocation of the method fact begins execution:


## FIGURE 2-5 Box trace of fact (3) (continued)

This is the base case, so this invocation of fact completes and returns a value to the caller:


The method value is returned to the calling box, which continues execution:


The current invocation of fact completes and returns a value to the caller:


The method value is returned to the calling box, which continues execution:


The current invocation of fact completes and returns a value to the caller:


The method value is returned to the calling box, which continues execution:


The current invocation of fact completes and returns a value to the caller:


The value 6 is returned to the initial call.

Note: The box trace, in conjunction with well-placed cout statements, can be a good aid in debugging recursive functions. Such statements should report the point in the program from which each recursive call occurs, as well as the values of input arguments and local variables at both entry to and exit from the function. Be sure to remove these cout statements from the final version of the function

Question 2 Write a box trace of the function given in Checkpoint Question 1.

### 2.3 Recursion That Performs an Action

Recursive functions need not return a value; they can be void functions.

### 2.3.1 A Recursive Void Function: Writing a String Backward

Now consider a problem that is slightly more difficult: Given a string of characters, write it in reverse order. For example, write the string "cat" as "tac". To construct a recursive solution, you should ask the four questions in the Note at the end of Section 2.1.

You can construct a solution to the problem of writing a string of length $n$ backward in terms of the problem of writing a string of length $n-1$ backward. That is, each recursive step of the solution diminishes by 1 the length of the string to be written backward. The fact that the strings get shorter and shorter suggests that the problem of writing some very short strings backward can serve as the base case. One very short string is the empty string, the string of length zero. Thus, you can choose for the base case the problem

## Write the empty string backward

The solution to this problem is to do nothing at all-a very straightforward solution indeed! (Alternatively, you could use the string of length 1 as the base case.)

Exactly how can you use the solution to the problem of writing a string of length $n-1$ backward to solve the problem of writing a string of length $n$ backward? This approach is analogous to the one used to construct the solution to the factorial problem, where you specified how to use factorial( $n-1$ ) in the computation of factorial ( $n$ ). Unlike the factorial problem, however, the string problem does not suggest an immediately clear way to proceed. Obviously, not any string of length $n-1$ will do. For example, there is no relation between writing the string "apple" (of length 5) backward and writing "pear" (a string of length 4) backward. You must choose the smaller problem carefully so that you can use its solution in the solution to the original problem.

The string of length $n-1$ that you choose must be a substring (part) of the original string. Suppose that you strip away one character from the original string, leaving a substring of length $n-1$. For the recursive solution to be valid, the ability to write the substring backward, combined with the ability to perform some minor task, must result in the ability to write the original string backward. Compare this approach with the way you computed factorial recursively: The ability to compute factorial ( $n-1$ ), combined with the ability to multiply this value by $n$, resulted in the ability to compute factorial( $n$ ).

You need to decide which character to strip away and which minor task to perform. Consider the minor task first. Because you are writing characters, a likely candidate for the minor task is writing a single character. As for the character that you should strip away from the string, there are several possible alternatives. Two of the more intuitive alternatives are

The base case

How can you write an n-character string backward, if you can write an ( $n-1$ )-character string backward?

Strip away the last character
or
Strip away the first character
Consider the first of these alternatives, stripping away the last character, as Figure 2-6 illustrates.

FIGURE 2-6 A recursive solution


For the solution to be valid, you must write the last character in the string first. Therefore, you must write the last character before you write the remainder of the string backward. A high-level recursive solution, given the string $s$, is
writeBackward writes a string backward

```
writeBackward(s: string)
    if (the string is empty)
        Do nothing-this is the base case
    e1se
    \{
        Write the last character of s
        writeBackward(s minus its last character)
    \}
```

Note: A recursive solution must involve one or more smaller problems that are each closer to a base case than is the original problem. You must be sure that these smaller problems eventually reach the base case. Failure to do so could result in an algorithm that does not terminate.

This solution to the problem is conceptual. To obtain a C++ function, you must resolve a few implementation issues. Suppose that the function will receive one argument: a string s to be written backward. All characters, including blanks, are part of the string. The C++ function writeBackward appears as follows:

```
/** Writes a character string backward.
@pre The string s to write backward.
@post None.
```

```
@param s The string to write backward. */
void writeBackward(string s)
{
    int length = s.size(); // Length of string
    if (length > 0)
    {
            // Write the last character
            cout << s.substr(length - 1, 1);
            // Write the rest of the string backward
            writeBackward(s.substr(0, length - 1)); // Point A
    } // end if
    // length == 0 is the base case - do nothing
} // end writeBackward
```

Notice that the recursive calls to writeBackward use successively shorter versions of the string s, ensuring that the base case will be reached. Because the function does nothing when it reaches the base case, it does not deal with the base case explicitly. The base case is implicit.

You can trace the execution of writeBackward by using the box trace. As was true for the function fact, each box contains the local environment of the recursive call-in this case, the input argument s and the local variable length. The trace will differ somewhat from the trace of
writeBackward does not return a computed value

## FIGURE 2-7 Box trace of writeBackward("cat")

The initial call is made, and the function begins execution:

```
s = "cat"
length = 3
```

Output line: t
Point A (writeBackward(s)) is reached, and the recursive call is made.
The new invocation begins execution:


## Output line: ta

Point A is reached, and the recursive call is made
The new invocation begins execution:


FIGURE 2-7 Box trace of writeBackward("cat") (continued)
Output line: tac
Point A is reached, and the recursive call is made.
The new invocation begins execution:


This is the base case, so this invocation completes.
Control returns to the calling box, which continues execution:


This invocation completes. Control returns to the calling box, which continues execution:


This invocation completes. Control returns to the calling box, which continues execution:


This invocation completes. Control returns to the statement following the initial call.

Another solution. Now consider a slightly different approach to the problem. Recall the two alternatives for the character that you could strip away from the string: the last character or the first character. The solution just given strips away the last character of the string. It will now be interesting to construct a solution based on the second alternative:

## Strip away the first character

To begin, consider a simple modification of the previous pseudocode solution that replaces each occurrence of last with first. Thus, the function writes the first character rather than the last and then recursively writes the remainder of the string backward.

```
writeBackward1(s: string)
    if (the string s is empty)
    Do nothing - this is the base case
    e1se
    {
        Write the first character of s
        writeBackward1(s minus its first character)
    }
```

Does this solution do what you want it to? If you think about this function, you will realize that it writes the string in its normal left-to-right direction instead of backward. After all, the steps in the pseudocode are

## Write the first character of s <br> Write the rest of s

These steps simply write the string s. Naming the function writeBackward does not guarantee that it will actually write the string backward-recursion really is not magic!

You can write s backward correctly by using the following recursive formulation:
Write s minus its first character backward
Write the first character of s
In other words, you write the first character of s only after you have written the rest of s backward. This approach leads to the following pseudocode solution:

```
writeBackward2(s: string)
    if (the string s is empty)
        Do nothing - this is the base case
    else
    {
        writeBackward2(s minus its first character)
        Write the first character of s
    }
```

The translation of writeBackward2 into C++ is similar to that of the original writeBackward function and is left as an exercise.

It is instructive to carefully trace the actions of the two pseudocode functions writeBackward and writeBackward2. First, add statements to each function to provide output that is useful to the trace, as follows:

```
writeBackward(s: string)
    cout << "Enter writeBackward with string: " << s << endl;
    if (the string is empty)
        Do nothing-this is the base case
    else
    {
        cout << "About to write last character of string: "
                << s << endl;
        Write the last character of s
        writeBackward(s minus its last character) // Point A
    }
    cout << "Leave writeBackward with string: " << s << end7;
writeBackward2(s: string)
    cout << "Enter writeBackward2 with string: "
        << s << endl;
    if (the string is empty)
        Do nothing -this is the base case
    e7se
    {
        writeBackward2(s minus its first character) // Point A
        cout << "About to write first character of string: "
            << S << endl;
        Write the first character of s
    }
    cout << "Leave writeBackward2 with string: " << s << endl;
```

Figures 2-8 and 2-9 show the output of the revised pseudocode functions writeBackward and writeBackward2, when initially given the string "cat".
cout statements can help you trace the logic of a recursive function

FIGURE 2-8 Box trace of writeBackward("cat") in pseudocode

The initial call is made, and the function begins execution:
s = "cat"

Output stream:

Enter writeBackward with string: cat
About to write last character of string: cat
t
Point A is reached, and the recursive call is made. The new invocation begins execution:


Output stream:

> Enter writeBackward with string: cat About to write last character of string: cat t
> Enter writeBackward with string: ca
> About to write last character of string: ca
> a

Point $A$ is reached, and the recursive call is made. The new invocation begins execution:


Output stream:

> Enter writeBackward with string: cat About to write last character of string: cat t
> Enter writeBackward with string: ca
> About to write last character of string: ca
> a
> Enter writeBackward with string: c
> About to write last character of string: c
> c

Point A is reached, and the recursive call is made. The new invocation begins execution:


This invocation completes execution, and a return is made.

Output stream：

Enter writeBackward with string：cat
About to write last character of string：cat
t
Enter writeBackward with string：ca
About to write last character of string：ca
a
Enter writeBackward with string：c
About to write last character of string：c
c
Enter writeBackward with string：
Leave writeBackward with string：


This invocation completes execution，and a return is made．
Output stream：

Enter writeBackward with string：cat
About to write last character of string：cat
t
Enter writeBackward with string：ca
About to write last character of string：ca
a
Enter writeBackward with string：c
About to write last character of string：c

## c

Enter writeBackward with string：
Leave writeBackward with string：
Leave writeBackward with string：c


This invocation completes execution，and a return is made．
Output stream：
Enter writeBackward with string：cat
About to write last character of string：cat

## t

Enter writeBackward with string：ca
About to write last character of string：ca
a
Enter writeBackward with string：c
About to write last character of string：c

## FIGURE 2-8 Box trace of writeBackward("cat") in pseudocode (continued)

c
Enter writeBackward with string:
Leave writeBackward with string:
Leave writeBackward with string: c
Leave writeBackward with string: ca


This invocation completes execution, and a return is made.
Output stream:

$$
\begin{aligned}
& \text { Enter writeBackward with string: cat } \\
& \text { About to write last character of string: cat } \\
& \mathrm{t} \\
& \text { Enter writeBackward with string: ca } \\
& \text { About to write last character of string: ca } \\
& \text { a } \\
& \text { Enter writeBackward with string: c } \\
& \text { About to write last character of string: c } \\
& \text { c } \\
& \text { Enter writeBackward with string: } \\
& \text { Leave writeBackward with string: } \\
& \text { Leave writeBackward with string: c } \\
& \text { Leave writeBackward with string: ca } \\
& \text { Leave writeBackward with string: cat }
\end{aligned}
$$

Well-placed but temporary cout statements can help you debug a recursive function

Remove cout statements after you have debugged the function

You need to be comfortable with the differences between these two functions. The recursive calls that the two functions make generate a different sequence of values for the argument $s$. Despite this fact, both functions correctly write the string argument backward. They compensate for the difference in the sequence of values for $s$ by writing different characters in the string at different times relative to the recursive calls. In terms of the box traces in Figures 2-8 and 2-9, writeBackward writes a character just before generating a new box (just before a new recursive call), whereas writeBackward2 writes a character just after crossing off a box (just after returning from a recursive call). When these differences are put together, the result is two functions that employ different strategies to accomplish the same task.

This example also illustrates the value of the box trace, combined with well-placed cout statements, in debugging recursive functions. The cout statements at the beginning, interior, and end of the recursive functions report the value of the argument $s$. In general, when debugging a recursive function, you should also report both the values of local variables and the point in the function where each recursive call occurred, as in this example:

```
abc(...);
    cout << "Calling function abc from point A.\n";
    abc(...) // This is point A
    cout << "Calling function abc from point B.\n";
    abc(...); // This is point B
```

Realize that the cout statements do not belong in the final version of the function.

## FIGURE 2-9 Box trace of writeBackward2("cat") in pseudocode

The initial call is made, and the function begins execution:

$$
s=\text { "cat" }
$$

## Output stream:

Enter writeBackward2 with string: cat
Point $A$ is reached, and the recursive call is made. The new invocation begins execution:


Output stream:

Enter writeBackward2 with string: cat
Enter writeBackward2 with string: at
Point $A$ is reached, and the recursive call is made. The new invocation begins execution:


Output stream:
Enter writeBackward2 with string: cat
Enter writeBackward2 with string: at
Enter writeBackward2 with string: t
Point $A$ is reached, and the recursive call is made. The new invocation begins execution:


This invocation completes execution, and a return is made.

Output stream:
Enter writeBackward2 with string: cat Enter writeBackward2 with string: at Enter writeBackward2 with string: t
Enter writeBackward2 with string:
Leave writeBackward2 with string:


This invocation completes execution, and a return is made.

Output stream:
Enter writeBackward2 with string: cat
Enter writeBackward2 with string: at
Enter writeBackward2 with string: t
(continues)

FIGURE 2-9 Box trace of writeBackward2 ("cat") in pseudocode (continued)

Enter writeBackward2 with string:
Leave writeBackward2 with string:
About to write first character of string: $t$
t
Leave writeBackward2 with string: t


This invocation completes execution, and a return is made.

Output stream:
Enter writeBackward2 with string: cat Enter writeBackward2 with string: at Enter writeBackward2 with string: t Enter writeBackward2 with string: Leave writeBackward2 with string: About to write first character of string: $t$ t

Leave writeBackward2 with string: t
About to write first character of string: at a
Leave writeBackward2, string: at


This invocation completes execution, and a return is made.
Output stream:
Enter writeBackward2 with string: cat
Enter writeBackward2 with string: at
Enter writeBackward2 with string: t
Enter writeBackward2 with string:
Leave writeBackward2 with string:
About to write first character of string: $t$
t
Leave writeBackward2 with string: t
About to write first character of string: at
a
Leave writeBackward2 with string: at
About to write first character of string: cat
c
Leave writeBackward2 with string: cat

Question 3 Given an integer $n>0$, write a recursive function countDown that writes the integers $n, n-1, \ldots, 1$. Hint: What task can you do and what task can you ask a friend to do for you?

### 2.4 Recursion with Arrays

When you are working with an array, recursion can be a practical and powerful tool. For our first example, we write the contents of an array backward. We then look at several problems that involve searching an array.

### 2.4.1 Writing an Array's Entries in Backward Order

The solution to this problem is much like our first solution in Section 2.3.1 when we wrote a string backward. We can write the following pseudocode:

```
writeArrayBackward(anArray: char[])
    if (the array is empty)
        Do nothing-this is the base case
    else
    {
        Write the last character in anArray
        writeArrayBackward(anArray minus its last character)
    }
```

How will we pass anArray minus its last character to writeArrayBackward? We could pass the number of characters left in the array. At each recursive call, we would decrease this number by 1 . Alternatively, we could pass the index of the last character. That is, if last is this index, writeArrayBackward would act on the array anArray[0..1ast], that is, the portion anArray[0] through anArray[last]. The recursive call would then act on the subarray anArray[0.. last - 1]. A more general variation of this idea also passes the index of the first array character. So instead of assuming that this index is 0 , we would pass anArray[first..last] to writeArrayBackward.

We can write the function writeArrayBackward as follows:

```
/**Writes the characters in an array backward.
    @pre The array anArray contains size characters, where size >= 0.
    @post None.
    @param anArray The array to write backward.
    @param first The index of the first character in the array.
    @param last The index of the last character in the array. */
void writeArrayBackward(const char anArray[], int first, int last)
{
    if (first <= last)
    {
            // Write the 1ast character
            cout << anArray[last];
            // Write the rest of the array backward
            writeArrayBackward(anArray, first, last - 1);
        } // end if
    // first > last is the base case - do nothing
} // end writeArrayBackward
```

[^2]A binary search conquers one of its subproblems at each step


Question 4 In the previous definition of writeArrayBackward, why does the base case occur when the value of first exceeds the value of last?

Question 5 Write a recursive function that computes and returns the product of the first $n \geq 1$ real numbers in an array.

Question 6 Show how the function that you wrote for the previous question satisfies the properties of a recursive function.

Question 7 Write a recursive function that computes and returns the product of the integers in the array anArray[first..7ast].

### 2.4.2 The Binary Search

Searching is an important task that occurs frequently. Often, searches are for a particular entry in an array. We now will examine a few searching problems that have recursive solutions. Our goal is to develop further your understanding of recursion.

This chapter began with an intuitive approach to a binary search algorithm by presenting-at a high level-a way to find a word in a dictionary. We now develop this algorithm fully and illustrate some important programming issues.

Recall the earlier solution to the dictionary problem:

```
search(aDictionary: Dictionary, word: string)
    if (aDictionary is one page in size)
        Scan the page for word
    e7se
    {
        Open aDictionary to a point near the middle
        Determine which half of aDictionary contains word
        if (word is in the first half of aDictionary)
            search(first half of aDictionary, word)
        else
            search(second half of aDictionary, word)
    }
```

Now alter the problem slightly by searching an array anArray of integers for a given value, the target. The array, like the dictionary, must be sorted, or else a binary search is not applicable. Hence, assume that

```
anArray[0] anArray[1] anArray[2] ... anArray[size - 1]
```

where size is the size of the array. A high-level binary search for the array problem is

```
binarySearch(anArray: ArrayType, target: ValueType)
    if (anArray is of size 1)
    Determine if anArray's value is equal to target
    else
    {
        Find the midpoint of anArray
        Determine which half of anArray contains target
        if (target is in the first half of anArray)
            binarySearch(first half of anArray, target)
            else
            binarySearch(second half of anArray, target)
    }
```

Although the solution is conceptually sound, you must consider several details before you can implement the algorithm:

1. How will you pass half of anArray to the recursive calls to binarySearch? You can pass the entire array at each call but have binarySearch search only anArray [first.. 7ast], that is, the portion anArray [first] through anArray[last]. Thus, you would also pass the integers first and last to binarySearch:
```
binarySearch(anArray, first, last, target)
```

With this convention, the new midpoint is given by

```
mid = (first + last) / 2
```

Then binarySearch (first half of anArray, target) becomes
binarySearch(anArray, first, mid - 1, target)
and binarySearch (second half of anArray, target) becomes
binarySearch(anArray, mid + 1, 1ast, target)
2. How do you determine which half of the array contains target? One possible implementation of
if (target is in the first half of anArray)
is
if (target < anArray[mid])
However, there is no test for equality between target and anArray [mid]. This omission can cause the algorithm to miss target. After the previous halving algorithm splits anArray into halves, anArray [mid] is not in either half of the array. (In this case, two halves do not make a whole!) Therefore, you must determine whether anArray [mid] is the value you seek now, because later it will not be in the remaining half of the array. The interaction between the halving criterion and the termination condition (the base case) is subtle and is often a source of error. We need to rethink the base case.
3. What should the base case(s) be?As it is written, binarySearch terminates only when an array of size 1 occurs; this is the only base case. By changing the halving process so that anArray [mid] remains in one of the halves, it is possible to implement the binary search correctly so that it has only this single base case. However, it can be clearer to have two distinct base cases as follows:

- first $>$ last. You will reach this base case when target is not in the original array.
- target $==$ anArray $[\mathrm{mid}]$. You will reach this base case when target is in the original array.

These base cases are a bit different from any you have encountered previously. In a sense, the algorithm determines the answer to the problem from the base case it reaches. Many search problems have this flavor.
4. How will binarySearch indicate the result of the search?If binarySearch successfully locates target in the array, it could return the index of the array value that is equal to target. Because this index would never be negative, binarySearch could return a negative value if it does not find target in the array.
The C++ function binarySearch that follows implements these ideas. The two recursive calls to binarySearch are labeled as $X$ and $Y$ for use in a later box trace of this function.

The array halves are anArray[first. . mid-1] and anArray [mid+1..last]; neither half contains anArray[mid]

Determine whether anArray[mid] is the target you seek

```
/** Searches the array anArray[first] through anArray[last]
    for a given value by using a binary search.
@pre 0<= first, 7ast <= SIZE - 1, where SIZE is the
        maximum size of the array, and anArray[first] <=
        anArray[first + 1] <= ... <= anArray[last].
@post anArray is unchanged and either anArray[index] contains
            the given value or index == -1.
    @param anArray The array to search.
@param first The low index to start searching from.
@param last The high index to stop searching at.
@param target The search key.
@return Either index, such that anArray[index] == target, or -1.
*/
int binarySearch(const int anArray[], int first, int last, int target)
{
    int index;
    if (first > 1ast)
        index = -1; // target not in original array
    else
    {
        // If target is in anArray,
        // anArray[first] <= target <= anArray[last]
        int mid = first + (last - first) / 2;
        if (target == anArray[mid])
            index = mid; // target found at anArray[mid]
        else if (target < anArray[mid])
                    // Point X
            index = binarySearch(anArray, first, mid - 1, target);
        else
            // Point Y
            index = binarySearch(anArray, mid + 1, last, target);
    } // end if
    return index;
} // end binarySearch
```

Notice that if target occurs in the array, it must be in the segment of the array delineated by first and last. That is, the following is true:

```
anArray[first] target anArray[last]
```

Figure 2-10 shows box traces of binarySearch when it searches the array containing 1, 5, $9,12,15$, 21, 29, and 31. Notice how the labels $X$ and $Y$ of the two recursive calls to binarySearch appear in the diagram. Exercise 16 at the end of this chapter asks you to perform other box traces with this function.

Note: When developing a recursive solution, you must be sure that the solutions to the smaller problems really do give you a solution to the original problem. For example, binarySearch works because each smaller array is sorted and the value sought is between its first and last values.

There is another implementation issue-one that deals specifically with $\mathrm{C}++$ - to consider. Recall that an array is never passed to a function by value and is therefore not copied. This aspect of $\mathrm{C}++$ is particularly useful in a recursive function such as binarySearch. If the array anArray is large, many recursive calls to binarySearch might be necessary. If each call copied anArray, much
memory and time would be wasted. On the other hand, because anArray is not copied, the function can alter the array's values unless you specify anArray as const, as was done for binarySearch.

FIGURE 2-10 Box traces of binarySearch with anArray $=<1,5,9,12,15,21,29,31>$ : (a) a successful search for 9 ; (b) an unsuccessful search for 6


A box trace of a recursive function that has an array argument requires a new consideration. Because the array anArray is neither a value argument nor a local variable, it is not a part of the function's local environment, and so the entire array anArray should not appear within each box. Therefore, as Figure 2-11 shows, you represent anArray outside the boxes, and all references to anArray affect this single representation.


Note: Notice that the $\mathrm{C}++$ computation of the midpoint mid is

$$
\text { int mid }=\text { first }+(7 \text { ast }- \text { first }) / 2 \text {; }
$$

instead of

```
int mid=(first + last) / 2;
```

as the pseudocode would suggest. If you were to search an array of at least $2^{30}$, or about 1 billion, elements, the sum of first and last could exceed the largest possible int value of $2^{30}-1$. Thus, the computation first + last would overflow to a negative integer and result in a negative value for mid. If this negative value of mid was used as an array index, it would be out of bounds and cause incorrect results. The computation first + (7ast first) / 2 is algebraically equivalent to (first + 1ast) / 2 and avoids this error.

Because an array argument is always passed by reference, a function can alter it unless you specify the array as const

Represent reference arguments outside of the boxes in a box trace
maxArray conquers both of its subproblems at each step

FIGURE 2-11 Box trace with a reference argument

### 2.4.3 Finding the Largest Value in an Array

Suppose that you have an array anArray of integers and you want to find the largest value. You could construct an iterative solution without too much difficulty, but instead let's consider a recursive formulation:

```
```

if (anArray has only one entry)

```
```

if (anArray has only one entry)
maxArray (anArray) is the entry in anArray
maxArray (anArray) is the entry in anArray
else if (anArray has more than one entry)
else if (anArray has more than one entry)
maxArray(anArray) is the maximum of
maxArray(anArray) is the maximum of
maxArray(left half of anArray) and maxArray(right half of anArray)

```
```

        maxArray(left half of anArray) and maxArray(right half of anArray)
    ```
```



Notice that this strategy fits the divide-and-conquer model that the previous binary search algorithm used. That is, we proceed by dividing the problem and conquering the subproblems, as Figure 2-12 illustrates. However, there is a difference between this algorithm and the binary search algorithm. Although the binary search algorithm conquers only one of its subproblems at each step, maxArray conquers both. Because both subproblems are solved recursively, this approach is called multipath recursion. After maxArray conquers the subproblems, it must reconcile the two solutions-that is, it must find the maximum of the two maximums. Figure 2-13 illustrates the computations that are necessary to find the largest integer in the array that contains $1,6,8$, and 3 (denoted here by $\langle 1,6,8,3>$ ).

Question 7 Define the recursive C++ function maxArray that returns the largest value in an array and adheres to the pseudocode just given.

### 2.4.4 Finding the $\boldsymbol{k} t h$ Smallest Value of an Array

Our discussion of searching concludes with a more difficult problem. Although you could skip this example now, Chapter 11 uses aspects of it in a sorting algorithm.

FIGURE 2-12 Recursive solution to the largest-value problem


FIGURE 2-13 The recursive calls that maxArray $(<1,6,8,3>)$ generates


The previous two examples presented recursive techniques for finding the largest value in an arbitrary array and for finding an arbitrary value in a sorted array. This example describes a recursive solution for finding the $k^{\text {th }}$ smallest value in an arbitrary array anArray. Would you ever be interested in such a value? Statisticians often want the median value in a collection of data. The median value in an ordered collection of data occurs in the middle of the collection. In an unordered collection of data, the number of values that are smaller than the median value is about the same as the number of values that are larger. Thus, if you have 49 values, the 25 th smallest value is the median value.

Obviously, you could solve this problem by sorting the array. Then the $k^{\text {th }}$ smallest value would be anArray $[k-1]$. Although this approach is a legitimate solution, it does more than the problem requires; a more efficient solution is possible. The solution outlined here finds the $k^{\text {th }}$ smallest value without completely sorting the array.

By now, you know that you solve a problem recursively by writing its solution in terms of one or more smaller problems of the same type in such a way that this notion of smaller ensures that you will always reach a base case. For all of the earlier recursive solutions, the reduction in problem size

For all previous examples, you know the amount of reduction made in the problem size by each recursive call

You cannot predict in advance the size of either the smaller problems or the base case in the recursive solution to the $k^{\text {th}}$-smallestvalue problem

## Partition anArray

 into three parts: values $p, p$, and values $\geq p$between recursive calls is predictable. For example, the factorial function always decreases the problem size by 1 ; the binary search always halves the problem size. In addition, the base cases for all the previous problems except the binary search have a static, predefined size. Thus, by knowing only the size of the original problem, you can determine the number of recursive calls that are necessary before you reach the base case.

The solution that you are about to see for finding the $k^{\text {th }}$ smallest value departs from these techniques. Although you solve the problem in terms of a smaller problem, just how much smaller this problem is depends on the values in the array and cannot be predicted in advance. Also, the size of the base case depends on the values in the array, as it did for the binary search. (Recall that you reach one of the base cases for a binary search when the middle value is the one sought.)

The unpredictable nature of this solution is caused by the problem itself: The relationship between the rankings of the values in any predetermined parts of the array and the ranking of the values in the entire array is not strong enough to determine the $k^{\text {th }}$ smallest value. For example, suppose that anArray contains the values shown in Figure 2-14. Notice that 6, which is in anArray [3], is the third-smallest value in the first half of anArray and that 8, which is in anArray [4], is the third-smallest value in the second half of anArray. Can you conclude from these observations anything about the location of the third-smallest value in all of anArray? The answer is no; these facts about parts of the array do not allow you to draw any useful conclusions about the entire array. You should experiment with other fixed splitting schemes as well.

FIGURE 2-14 A sample array


The recursive solution proceeds by

1. Selecting a pivot value in the array
2. Cleverly arranging, or partitioning, the values in the array about this pivot value
3. Recursively applying the strategy to one of the partitions

Suppose that you want to find the $k^{\text {th }}$ smallest value in the array segment anArray [first. . 1 ast]. Let the pivot $p$ be any value of the array segment. (For now, ignore how to choose $p$.) You can partition the values of anArray [first. . 7ast] into three regions: $S_{1}$, which contains the values less than or equal to $p$; the pivot $p$ itself; and $S_{2}$, which contains the values greater than or equal to $p$. This partition implies that all of the values in $S_{1}$ are no larger than all of the values in $S_{2}$. Figure 2-15 illustrates this partition.

All values in anArray[first..pivotIndex-1] are less than or equal to $p$, and all values in anArray[pivotIndex+1..7ast] are greater than or equal to $p$. Notice that the sizes of the regions $S_{1}$ and $S_{2}$ depend on both $p$ and the other values of anArray[first. . 7ast].

This partition induces three smaller problems, such that the solution to one of the problems will solve the original problem:

1. If $S_{1}$ contains $k$ or more values, $S_{1}$ contains the $k$ smallest values of the array segment anArray[first. . 7ast]. In this case, the $k^{\text {th }}$ smallest value must be in $S_{1}$. Since $S_{1}$ is the array segment anArray[first. . pivotIndex-1], this case occurs if $k<$ pivotIndex - first +1 .

FIGURE 2-15 A partition about a pivot

2. If $S_{1}$ contains $k-1$ values, the $k^{\text {th }}$ smallest value must be the pivot $p$. This is the base case; it occurs if $k=$ pivotIndex - first +1 .
3. If $S_{1}$ contains fewer than $k-1$ values, the $k^{\text {th }}$ smallest value in anArray [first. . 1 ast] must be in $S_{2}$. Because $S_{1}$ contains pivotIndex - first values, the $k^{\text {th }}$ smallest value in anArray [first. . 7ast] is the $(k-(\text { pivotIndex }- \text { first }+1))^{\text {st }}$ smallest value in $S_{2}$. This case occurs if $k>$ pivotIndex - first +1 .

A recursive definition can summarize this discussion. Let
$\mathrm{kSma11}\left(\mathrm{k}\right.$, anArray, first, last) $=k^{\text {th }}$ smallest value in anArray[first..1ast]
After you select the pivot value $p$ and partition anArray[first. . 1ast] into $S_{1}$ and $S_{2}$, you have that kSmall(k, anArray, first, last) equals

- kSmal1(k, anArray, first, pivotIndex - 1) if $k<$ pivotIndex-first + 1
- p if $\mathrm{k}=$ pivotIndex - first +1
- kSmall(k - (pivotIndex - first + 1), anArray, pivotIndex + 1, last)
if $\mathrm{k}>$ pivotIndex - first +1
There is always a pivot, and because it is not part of either $S_{1}$ or $S_{2}$, the size of the array segment to be searched decreases by at least 1 at each step. Thus, you will eventually reach the base case: The desired value is a pivot. A high-level pseudocode solution is as follows.

The $k^{\text {th }}$ smallest value in anArray [first..7ast]

```
// Returns the kth smallest value in anArray[first. .1ast].
kSma11(k: integer, anArray: ArrayType,
        first: integer, last: integer): ValueType
    Choose a pivot value p from anArray[first. . 1ast]
    Partition the values of anArray[first. . 1ast] about p
    if (k < pivotIndex - first + 1)
        return \(\mathrm{kSmall}(\mathrm{k}\), anArray, first, pivotIndex - 1)
    else if (k == pivotIndex - first + 1)
        return \(p\)
    e1se
        return kSmal1 (k - (pivotIndex - first + 1) , anArray,
                        pivotIndex + 1, last)
```

This pseudocode is not far from a C++ function. The only questions that remain are how to choose the pivot value $p$ and how to partition the array about the chosen $p$. The choice of $p$ is arbitrary. Any $p$ in the array will work, although the sequence of choices will affect how soon you reach the base case. Chapter 11 gives an algorithm for partitioning the values about $p$. There you will see how to turn the function kSma 11 into a sorting algorithm.

### 2.5 Organizing Data

Given some data organized in one way, you might need to organize it in another way. Thus, you will actually change some aspect of the data and not, for example, simply search it. The problem in this section is called the Towers of Hanoi. Although this classic problem probably has no direct realworld application, we consider it because its solution so well illustrates the use of recursion.

### 2.5.1 The Towers of Hanoi

Many, many years ago, in a distant part of the Orient-in the Vietnamese city of Hanoi-the emperor's wiseperson passed on to join his ancestors. The emperor needed a replacement wiseperson. Being a rather wise person himself, the emperor devised a puzzle, declaring that its solver could have the job of wiseperson.

FIGURE 2-16 (a) The initial state; (b) move $n-1$ disks from A to C ; (c) move 1 disk from A to B ; (d) move $n-1$ disks from C to B
(a)

A

A

B
C
(c)
A

B


C
C

A
B
(d)
)

The emperor's puzzle consisted of $n$ disks (he didn't say exactly how many) and three poles: A (the source), B (the destination), and C (the spare). The disks were of different sizes and had holes in the middle so that they could fit on the poles. Because of their great weight, the disks could be placed only on top of disks larger than themselves. Initially, all the disks were on pole A, as shown in Figure 2-16a. The puzzle was to move the disks, one by one, from pole A to pole B. A person could also use pole C in the course of the transfer, but again a disk could be placed only on top of a disk larger than itself.

As the position of wiseperson was generally known to be a soft job, there were many applicants. Scholars and peasants alike brought the emperor their solutions. Many solutions were thousands of steps long, and many contained deeply nested loops and control structures. "I can't understand these solutions," bellowed the emperor. "There must be an easy way to solve this puzzle."

And indeed there was. A great Buddhist monk came out of the mountains to see the emperor. "My son," he said, "the puzzle is so easy, it almost solves itself." The emperor's security chief wanted to throw this strange person out, but the emperor let him continue.
"If you have only one disk (that is, $n=1$ ), move it from pole A to pole B." So far, so good, but even the village idiot could get that part right. "If you have more than one disk (that is, $n>1$ ), simply

1. Ignore the bottom disk and solve the problem for $n-1$ disks, with the small modification that pole $C$ is the destination and pole $B$ is the spare. (See Figure 2-16b.)
2. After you have done this, $n-1$ disks will be on pole C , and the largest disk will remain on pole A. So solve the problem for $n=1$ (recall that even the village idiot could do this) by moving the large disk from A to B . (See Figure 2-16c.)
3. Now all you have to do is move the $n-1$ disks from pole $C$ to pole $B$; that is, solve the problem with pole $C$ as the source, pole $B$ as the destination, and pole $A$ as the spare." (See Figure 2-16d.)

There was silence for a few moments, and finally the emperor said impatiently, "Well, are you going to tell us your solution or not?" The monk simply gave an all-knowing smile and vanished.

The emperor obviously was not a recursive thinker, but you should realize that the monk's solution is perfectly correct. The key to the solution is the observation that you can solve the Towers problem of $n$ disks by solving three smaller-in the sense of number of disks - Towers problems. Let towers (count, source, destination, spare) denote the problem of moving count disks from pole source to pole destination, using pole spare as a spare. Notice that this definition makes sense even if there are more than count disks on pole source; in this case, you concern yourself with only the top count disks and ignore the others. Similarly, the poles destination and spare might have disks on them before you begin; you ignore these, too, except that you may place only smaller disks on top of them.

You can restate the emperor's problem as follows: Beginning with $n$ disks on pole A and zero disks on poles B and C, solve towers (n, A, B, C). You can state the monk's solution as follows:

Step 1. Starting in the initial state-with all the disks on pole $A$-solve the problem

```
towers(n - 1, A, C, B)
```

That is, ignore the bottom (largest) disk and move the top $n-1$ disks from pole A to pole C , using pole B as a spare. When you are finished, the largest disk will remain on pole A , and all the other disks will be on pole C .

Step 2. Now, with the largest disk on pole $A$ and all others on pole $C$, solve the problem
towers (1, A, B, C)
That is, move the largest disk from pole A to pole B. Because this disk is larger than the disks already on the spare pole $C$, you really could not use the spare. However, fortunately-and obviously-you do not need to use the spare in this base case. When you are done, the largest disk will be on pole B , and all other disks will remain on pole C .

Step 3. Finally, with the largest disk on pole $B$ and all the other disks on pole C, solve the problem

```
towers(n - 1, C, B, A)
```

That is, move the $n-1$ disks from pole C to pole B , using A as a spare. Notice that the destination pole B already has the largest disk, which you ignore. When you are done, you will have solved the original problem: All the disks will be on pole $B$.

The problem statement

The solution

The solution to the Towers problem satisfies the four criteria of a recursive solution

The solution for three disks

The problem towers(count, source, destination, spare) has the following pseudocode solution:

```
solveTowers(count, source, destination, spare)
    if (count is l)
        Move a disk directly from source to destination
    e7se
    {
    solveTowers(count - 1, source, spare, destination)
    solveTowers(1, source, destination, spare)
    solveTowers(count - 1, spare, destination, source)
    }
```

This recursive solution follows the same basic pattern as the recursive solutions you saw earlier in this chapter:

1. You solve a Towers problem by solving other Towers problems.
2. These other Towers problems are smaller than the original problem; they have fewer disks to move. In particular, the number of disks decreases by 1 at each recursive call.
3. When a problem has only one disk - the base case - the solution is easy to solve directly.
4. The way that the problems become smaller ensures that you will reach a base case.

Solving the Towers problem requires you to solve many smaller Towers problems recursively. Figure 2-17 illustrates the resulting recursive calls and their order when you solve the problem for three disks.

Now consider a C++ implementation of this algorithm. Notice that since most computers do not have arms (at the time of this writing), the function moves a disk by giving directions to a human. Thus, the parameters that represent the poles are of type char, and the corresponding arguments could be 'A', 'B', and 'C'. The call solveTowers(3, 'A', 'B', 'C') produces this output:

```
Move top disk from pole A to pole B
Move top disk from pole A to pole C
Move top disk from pole B to pole C
Move top disk from pole A to pole B
Move top disk from pole C to pole A
Move top disk from pole C to pole B
Move top disk from pole A to pole B
```

FIGURE 2-17 The order of recursive calls that results from solveTowers (3, A, B , C)


```
The C++ function follows:
void solveTowers(int count, char source, char destination, char spare)
\{
    if (count == 1)
    \{
            cout << "Move top disk from pole " << source
                << " to pole " << destination << end1;
    \}
    else
    \{
        solveTowers(count - 1, source, spare, destination); // X
        solveTowers(1, source, destination, spare); // Y
        solveTowers(count - 1, spare, destination, source); // Z
    \} // end if
\} // end solveTowers
```

Question 8 Trace the execution of the function solveTowers to solve the Towers of Hanoi problem for two disks.

### 2.6 More Examples

The next three problems require you to count certain events or combinations of events or things. They are good examples of recursive solutions with more than one base case. However, these solutions are tremendously inefficient, and so are not practical. Do not let this inefficiency discourage you. Recursion can be useful and efficient, even though it is not always so. Your goal right now is to understand recursion by examining simple problems.

### 2.6.1 The Fibonacci Sequence (Multiplying Rabbits)

Rabbits are very prolific breeders. If rabbits did not die, their population would quickly get out of hand. Suppose we assume the following "facts," which were obtained in a recent survey of randomly selected rabbits:

- Rabbits never die.
- A rabbit reaches sexual maturity exactly two months after birth; that is, at the beginning of its third month of life.
- Rabbits are always born in male-female pairs. At the beginning of every month, each sexually mature male-female pair gives birth to exactly one male-female pair.

Suppose that you started with a single newborn male-female pair. How many pairs would there be in month 6 , counting the births that took place at the beginning of month 6 ? As 6 is a relatively small number, you can figure out the solution easily:

Month 1: 1 pair, the original rabbits.
Month 2: 1 pair still, because the rabbits are not yet sexually mature.
Month 3: 2 pairs; the original pair has reached sexual maturity and has given birth to a second pair.
Month 4: 3 pairs; the original pair has given birth again, but the pair born at the beginning of month 3 are not yet sexually mature.
Month 5: 5 pairs; all rabbits alive in month 3 ( 2 pairs) are now sexually mature. Add their offspring to those pairs alive in month 4 ( 3 pairs) to yield 5 pairs.
Month 6: 8 pairs; 3 newborn pairs from the pairs alive in month 4 plus 5 pairs alive in month 5 .

The number of pairs in month $n$

Two base cases are necessary because there are two smaller problems
rabbit computes the Fibonacci sequence but does so inefficiently

You can now construct a recursive solution for computing $\operatorname{rabbit}(n)$, the number of pairs alive in month $n$. You must determine how you can use $\operatorname{rabbit}(n-1)$ to compute $\operatorname{rabbit}(n)$. Observe that $\operatorname{rabbit}(n)$ is the sum of the number of pairs alive just prior to the start of month $n$ and the number of pairs born at the start of month $n$. Just prior to the start of month $n$, there are rabbit $(n-1)$ pairs of rabbits. Not all of these rabbits are sexually mature at the start of month $n$. Only those that were alive in month $n-2$ are ready to reproduce at the start of month $n$. That is, the number of pairs born at the start of month $n$ is rabbit $(n-2)$. Therefore, you have the recurrence relation

$$
\operatorname{rabbit}(n)=\operatorname{rabbit}(n-1)+\operatorname{rabbit}(n-2)
$$

Figure 2-18 illustrates this relationship.
FIGURE 2-18 Recursive solution to the rabbit problem


This recurrence relation-like some previous examples-solves a problem by solving more than one smaller problem of the same type. This does not add much conceptual difficulty, but you must be very careful when selecting the base case. The temptation is simply to say that rabbit(1) should be the base case because its value is 1 according to the problem's statement. But what about rabbit(2)? Applying the recursive definition to $\operatorname{rabbit}(2)$ would yield

$$
\operatorname{rabbit}(2)=\operatorname{rabbit}(1)+\operatorname{rabbit}(0)
$$

Thus, the recursive definition would need to specify the number of pairs alive in month 0 - an undefined quantity.

One possible solution is to define $\operatorname{rabbit}(0)$ to be 0 , but this approach seems artificial. A slightly more attractive alternative is to treat rabbit(2) itself as a special case with the value of 1 . Thus, the recursive definition has two base cases, rabbit(2) and rabbit(1). The recursive definition becomes

$$
\operatorname{rabbit}(n)= \begin{cases}1 & \text { if } n \text { is } 1 \text { or } 2 \\ \operatorname{rabbit}(n-1)+\operatorname{rabbit}(n-2) & \text { if } n>2\end{cases}
$$

Incidentally, the series of numbers $\operatorname{rabbit}(1), \operatorname{rabbit}(2), \operatorname{rabbit}(3)$, and so on is known as the Fibonacci sequence, which models many naturally occurring phenomena.

A C++ function to compute $\operatorname{rabbit}(n)$ is easy to write from the previous definition:

```
/** Computes a term in the Fibonacci sequence.
    @pre n is a positive integer.
    @post None.
    @param n The given integer.
    @return The nth Fibonacci number. */
int rabbit(int n)
{
        if (n <= 2)
            return 1;
        else // n > 2, so n - 1>0 and n-2 >0
            return rabbit(n-1) + rabbit(n-2);
} // end rabbit
```



The number of acceptable parades of length $n$ that end with a float

The number of acceptable parades of length $n$ that end with a band

The number of acceptable parades of length $n$

Should you actually use this function? Figure 2-19 illustrates the recursive calls that rabbit (7) generates. Think about the number of recursive calls that rabbit(10) generates. At best, the function rabbit is inefficient. Thus, its use is not feasible for large values of $n$. This problem is discussed in more detail at the end of this chapter, at which time you will see some techniques for generating a more efficient solution from this same recursive relationship.


Note: A recursive solution that recomputes certain values frequently can be quite inefficient. In such cases, iteration may be preferable to recursion.

### 2.6.2 Organizing a Parade

You have been asked to organize the Fourth of July parade, which will consist of bands and floats in a single line. Last year, adjacent bands tried to outplay each other. To avoid this problem, the sponsors have asked you never to place one band immediately after another. In how many ways can you organize a parade of length $n$ ?

Assume that you have at least $n$ marching bands and $n$ floats from which to choose. When counting the number of ways to organize the parade, assume that the parades band-float and float-band, for example, are different parades and count as two ways.

The parade can end with either a float or a band. The number of ways to organize the parade is simply the sum of the number of parades of each type. That is, let
$P(n)$ be the number of ways to organize a parade of length $n$
$F(n)$ be the number of parades of length $n$ that end with a float
$B(n)$ be the number of parades of length $n$ that end with a band
Then

$$
P(n)=F(n)+B(n)
$$

First, consider $F(n)$. You will have a parade of length $n$ that ends with a float simply by placing a float at the end of any acceptable parade of length $n-1$. Hence, the number of acceptable parades of length $n$ that end with a float is precisely equal to the total number of acceptable parades of length $n-1$; that is

$$
F(n)=P(n-1)
$$

Next, consider $B(n)$. The only way a parade can end with a band is if the unit just before the end is a float. (If it is a band, you will have two adjacent bands.) Thus, the only way to organize an acceptable parade of length $n$ that ends with a band is first to organize a parade of length $n-1$ that ends with a float and then add a band to the end. Therefore, the number of acceptable parades of length $n$ that end with a band is precisely equal to the number of acceptable parades of length $n-1$ that end with a float:

$$
B(n)=F(n-1)
$$

You use the earlier fact that $F(n)=P(n-1)$ to obtain

$$
B(n)=P(n-2)
$$

Thus, you have solved $F(n)$ and $B(n)$ in terms of the smaller problems $P(n-1)$ and $P(n-2)$, respectively. You then use

$$
P(n)=F(n)+B(n)
$$

to obtain

$$
P(n)=P(n-1)+P(n-2)
$$

The form of this recurrence relation is identical to the solution for the multiplying rabbits problem.

As you saw in the rabbit problem, two base cases are necessary, because the recurrence relation defines a problem in terms of two smaller problems. As you did for the rabbit problem, you can choose $n=1$ and $n=2$ for the base cases. Although both problems use the same values of $n$ for their base cases, there is no reason to expect that they use the same values for these base cases. That is, there is no reason to expect that rabbit(1) is equal to $P(1)$ and that $\operatorname{rabbit}(2)$ is equal to $P(2)$.

A little thought reveals that for the parade problem,

```
\(P(1)=2\) (The parades of length 1 are float and band.)
\(P(2)=3\) (The parades of length 2 are float-float, band-float, and float-band.)
```

In summary, the solution to this problem is

$$
\begin{aligned}
& P(1)=2 \\
& P(2)=3 \\
& P(n)=P(n-1)+P(n-2) \text { for } n>2
\end{aligned}
$$

This example demonstrates the following points about recursion:

- Sometimes you can solve a problem by breaking it up into cases-for example, parades that end with a float and parades that end with a band.
- The values that you use for the base cases are extremely important. Although the recurrence relations for $P$ and rabbit are the same, their base cases (when $n=1$ or 2 ) are different. This difference causes $\operatorname{rabbit}(n)$ and $P(n)$ to differ when $n$ is greater than 2 . For example, $\operatorname{rabbit}(20)=6,765$, while $P(20)=17,711$. The larger the value of $n$, the larger the discrepancy. You should think about why this is so.


### 2.6.3 Choosing $\boldsymbol{k}$ Out of $\boldsymbol{n}$ Things

A rock band would like to tour $n$ cities. Unfortunately, time will allow for visits to only $k$ cities. The band's agent considers the different choices for visiting $k$ cities out of the $n$ possibilities. Because time is short, the band members are not concerned about the order in which they visit the same $k$ cities.

Let $g(n, k)$ be the number of groups of $k$ cities chosen from $n$. If we consider city $C$, either we visit $C$ or we do not. If we do visit city $C$, we will have to choose $k-1$ other cities to visit from the $n-1$ remaining cities. Thus, the number of groups of cities that include $C$ is $g(n-1, k-1)$. On the other hand, if we do not visit city $C$, we will have to choose $k$ cities to visit from the remaining $n-1$ cities. The number of groups of cities that do not include $C$ is $g(n-1, k)$. Thus, we can compute $g(n, k)$ by solving two smaller counting problems of the same type; that is,

$$
g(n, k)=g(n-1, k-1)+g(n-1, k)
$$

We still need to find the base case(s) and demonstrate that each of the two smaller problems eventually reaches a base case. First, if the band had time to visit all $n$ cities-that is, if $k$ equals $n$-there is only one group of all the cities. Thus, the first base case is

$$
g(k, k)=1
$$

If $k<n$, the second term $g(n-1, k)$ in the recursive definition will reach this base case, because at each successive stage of the recursion, $n-1$ decreases until it reaches $k$. However, the first term, $g(n-1$, $k-1$ ), does not reach this base case. Since $n-1$ and $k-1$ both decrease at the same rate, they will never become equal. The first term does, in fact, approach another trivial selection problem. Just as there is only one group of all the cities $(k=n)$, there is also only one group of zero cities $(k=0)$. Thus, the second base case is

$$
g(n, 0)=1
$$

Two base cases are necessary because there are two smaller problems

The number of ways to choose $k$ out of $n$ things is the sum of the number of ways to choose $k-1$ out of $n-1$ things and the number of ways to choose kout of $n-1$ things

The number of groups of $k$ things recursively chosen out of $n$ things
and will be reached by the first term, $g(n-1, k-1)$. (Alternatively, you could define the second base case to be $g(n, 1)=n$.)

T
Note: When you solve a problem by solving two (or more) smaller problems, each of the smaller problems must be closer to a base case than the original problem.

For completeness, we add one final part to this recursive solution:

$$
g(n, k)=0 \quad \text { if } k>n
$$

Although $k$ could not be greater than $n$ in the context of this problem, the addition of this case makes the recursive solution more generally applicable.

To summarize, the following recursive solution solves the problem of choosing $k$ out of $n$ things:

$$
g(n, k)=\left\{\begin{array}{cl}
1 & \text { if } k=0 \\
1 & \text { if } k=n \\
0 & \text { if } k>n \\
g(n-1, k-1)+g(n-1, k) & \text { if } 0<k<n
\end{array}\right.
$$

You can easily derive the following function from this recursive definition:

```
/** Computes the number of groups of k out of n things.
    @pre n and k are nonnegative integers.
    @post None.
    @param n The given number of things.
    @param k The given number to choose.
    @return g(n, k). */
int getNumberOfGroups(int n, int k)
{
    if ( (k == 0) || (k == n) )
        return 1;
    else if (k > n)
        return > 0;
    else
        return g(n - 1, k - 1) + g(n - 1, k);
} // end getNumberOfGroups
```

Like the rabbit function, this function is inefficient and not practical to use. Figure 2-20 shows the number of recursive calls that the computation of $g(4,2)$ requires.


Note: When a recursive function contains more than one recursive call, you often will need more than one base case.

Question 9 Compute $g(4,2)$.

FIGURE 2-20 The recursive calls that $g(4,2)$ generates


### 2.7 Recursion and Efficiency

Recursion is a powerful problem-solving technique that often produces very clean solutions to even the most complex problems. Recursive solutions can be easier to understand and to describe than iterative solutions. By using recursion, you can often write simple, short implementations of your solution.

The overriding concern of this chapter has been to give you a solid understanding of recursion so that you will be able to construct recursive solutions on your own. Most of our examples, therefore, have been simple. Unfortunately, many of the recursive solutions in this chapter are so inefficient that you should not use them. The recursive functions binarySearch and solveTowers are the notable exceptions, as they are quite efficient. ${ }^{2}$

Two factors contribute to the inefficiency of some recursive solutions:

- The overhead associated with function calls
- The inherent inefficiency of some recursive algorithms

The first of these factors does not pertain specifically to recursive functions but is true of functions in general. In most implementations of $\mathrm{C}++$ and other high-level programming languages, a function call incurs a certain amount of bookkeeping overhead. As was mentioned earlier, each function call produces an activation record, which is analogous to a box in the box trace. Recursive functions magnify this overhead because a single initial call to the function can generate a large number of recursive calls. For example, the call factorial( $n$ ) generates $n$ recursive calls. On the other hand, the use of recursion, as is true of modularity in general, can greatly clarify complex programs. This clarification frequently more than compensates for the additional overhead.

Factors that contribute to the inefficiency of some recursive solutions

Recursion can clarify complex solutions

[^3]Do not use a recursive solution if it is inefficient and you have a clear, efficient iterative solution

The recursive version of rabbit is inherently inefficient

You can use
rabbit's
recurrence relation to construct an efficient iterative solution

Convert from recursion to iteration if it is easier to discover a recursive solution but more efficient to use an iterative solution

However, you should not use recursion just for the sake of using recursion. For example, you probably should not use the recursive factorial function in practice. You can easily write an iterative factorial function, given the iterative definition that was stated earlier in this chapter. The iterative function is almost as clear as the recursive one and is more efficient. There is no reason to incur the overhead of recursion when its use does not gain anything. Recursion is truly valuable when a problem has no simple iterative solutions.

The second point about recursion and efficiency is that some recursive algorithms are inherently inefficient. This inefficiency is a very different issue than that of overhead. It has nothing to do with how a compiler happens to implement a recursive function, but rather is related to the technique that the algorithm employs.

As an example, recall the recursive solution for the multiplying rabbits problem that you saw earlier in this chapter:

$$
\operatorname{rabbit}(n)= \begin{cases}1 & \text { if } n \text { is } 1 \text { or } 2 \\ \operatorname{rabbit}(n-1)+\operatorname{rabbit}(n-2) & \text { if } n>2\end{cases}
$$

The diagram in Figure 2-19 illustrated the computation of $\operatorname{rabbit}(7)$. Earlier, you were asked to think about what the diagram would look like for rabbit(10). If you thought about this question, you may have come to the conclusion that such a diagram would fill up most of this chapter. The diagram for $\operatorname{rabbit}(100)$ would fill up most of this universe!

The fundamental problem with rabbit is that it computes the same values over and over again. For example, in the diagram for $\operatorname{rabbit}(7)$, you can see that $\operatorname{rabbit}(3)$ is computed five times. When $n$ is moderately large, many of the values are recomputed literally trillions of times. This enormous number of computations makes the solution infeasible, even if each computation requires only a trivial amount of work.

However, do not conclude that the recurrence relation is of no use. One way to solve the rabbit problem is to construct an iterative solution based on this same recurrence relation. The iterative solution goes forward instead of backward and computes each value only once. You can use the following iterative function to compute $\operatorname{rabbit}(n)$ even for very large values of $n$.

```
/** Iterative solution to the rabbit prob7em. */
int iterativeRabbit(int n)
{
    // Initialize base cases:
    int previous = 1; // Initially rabbit(1)
    int current = 1; // Initially rabbit(2)
    int next = 1; // Result when n is 1 or 2
    // Compute next rabbit values when n >= 3
    for (int i = 3; i <= n; i++)
    {
        // current is rabbit(i - 1), previous is rabbit(i - 2)
        next = current + previous; // rabbit(i)
        previous = current; // Get ready for next iteration
        current = next;
    } // end for
    return next;
} // end iterativeRabbit
```

Thus, an iterative solution can be more efficient than a recursive solution. In certain cases, however, it may be easier to discover a recursive solution than an iterative solution. Therefore, you may need to convert a recursive solution to an iterative solution. This conversion process is easier if your recursive function calls itself once, instead of several times. Be careful when deciding whether your function calls itself more than once. Although the recursive function rabbit calls itself twice, the
function binarySearch calls itself once, even though you see two calls in the C++ code. Those two calls appear within an if statement; only one of them will be executed.

Converting a recursive solution to an iterative solution is even easier when the solitary recursive call is the last action that the function takes. This situation is called tail recursion. For example, the function writeBackward exhibits tail recursion because its recursive call is the last action that the function takes. Before you conclude that this is obvious, consider the function fact. Although its recursive call appears last in the function definition, fact's last action is the multiplication. Thus, fact is not tail-recursive.

Recall the definition of writeBackward:

```
void writeBackward(string s)
{
    int length = s.size();
    if (length > 0)
    {
        // Write last character
        cout << s.substr(length - 1, 1);
        writeBackward(s.substr(0, length - 1)); // Write rest
    } // end if
} // end writeBackward
```

Because this function is tail-recursive, its last recursive call simply repeats the function's action with altered arguments. You can perform this repetitive action by using an iteration that will be straightforward and often more efficient. For example, the following definition of writeBackward is iterative:

Because tail-recursive functions are often less efficient than their iterative counterparts, and because the conversion of a tail-recursive function to an equivalent iterative function is rather mechanical, some compilers automatically replace tail recursion with iteration. Eliminating
other forms of recursion is usually more complex and is a task that you would need to undertake, if mechanical, some compilers automatically replace tail recursion with iteration. Eliminating
other forms of recursion is usually more complex and is a task that you would need to undertake, if necessary.

Some recursive algorithms, such as rabbit, are inherently inefficient, while other recursive algorithms, such as the binary search, ${ }^{3}$ are extremely efficient. You will learn how to determine the relative efficiency of a recursive algorithm in more advanced courses concerned with the analysis of algorithms. Chapter 10 introduces some of these techniques briefly. Chapter 5 continues the discussion of recursion by examining several difficult problems that have straightforward recursive solutions. Other chapters in this book use recursion as a matter of course.


Question 10 Of the following recursive functions that you saw in this chapter, identify those that exhibit tail recursion: fact, writeBackward, writeBackward2, rabbit, $P$ in the parade problem, getNumberOfGroups, maxArray, binarySearch, and kSma11.

[^4]```
/** Iterative version. */
```

/** Iterative version. */
void writeBackward(string s)
void writeBackward(string s)
{
{
int length = s.size();
int length = s.size();
while (length > 0)
while (length > 0)
{
{
cout << s.substr(length - 1, 1);
cout << s.substr(length - 1, 1);
1ength--;
1ength--;
} // end while
} // end while
} // end writeBackward

```
} // end writeBackward
```

A tail-recursive function

## Removing tail

 recursion is often straightforward
## Summary

1. Recursion is a technique that solves a problem by solving a smaller problem of the same type.
2. When constructing a recursive solution, keep the following four questions in mind:
a. How can you define the problem in terms of a smaller problem of the same type?
b. How does each recursive call diminish the size of the problem?
c. What instance of the problem can serve as the base case?
d. As the problem size diminishes, will you reach this base case?
3. When constructing a recursive solution, you should assume that a recursive call's result is correct if its precondition has been met.
4. You can use the box trace to trace the actions of a recursive function. These boxes resemble activation records, which many compilers use to implement recursion. Although the box trace is useful, it cannot replace an intuitive understanding of recursion.
5. Recursion allows you to solve problems-such as the Towers of Hanoi-whose iterative solutions are difficult to conceptualize. Even the most complex problems often have straightforward recursive solutions. Such solutions can be easier to understand, describe, and implement than iterative solutions.
6. Some recursive solutions are much less efficient than a corresponding iterative solution due to their inherently inefficient algorithms and the overhead of function calls. In such cases, the iterative solution can be preferable. You can use the recursive solution, however, to derive the iterative solution.
7. If you can easily, clearly, and efficiently solve a problem by using iteration, you should do so.

## Exercises

1. The following recursive function getNumberEqual searches the array $x$ of $n$ integers for occurrences of the integer desiredValue. It returns the number of integers in x that are equal to desiredValue. For example, if x contains the ten integers $1,2,4,4,5,6,7,8,9$, and 12 , then getNumberEqual ( $x, 10,4$ ) returns the value 2 , because 4 occurs twice in x .
```
int getNumberEqual(const int x[], int n, int desiredValue)
{
    int count = 0;
    if (n <= 0)
        return 0;
    else
    {
        if (x[n - 1] == desiredValue)
            count = 1;
        return getNumberEqua1(x, n - 1, desiredValue) + count;
    } // end else
} // end getNumberEqual
```

Demonstrate that this function is recursive by listing the criteria of a recursive solution and stating how the function meets each criterion.
2. Perform a box trace of the following calls to recursive functions that appear in this chapter. Clearly indicate each subsequent recursive call.
a. rabbit(5)
b. countDown (5) (You wrote countDown in Checkpoint Question 3.)
3. Write a recursive function that will compute the sum of the first $n$ integers in an array of at least $n$ integers. Hint: Begin with the $n$th integer.
4. Given two integers, start and end, where end is greater than start, write a recursive $\mathrm{C}++$ function that returns the sum of the integers from start through end, inclusive.
5. a. Revise the function writeBackward, discussed in Section 2.3.1, so that its base case is a string of length 1.
b. Write a C++ function that implements the pseudocode function writeBackward2, as given in Section 2.3.1.
6. Describe the problem with the following recursive function:

```
void printNum(int n)
{
    cout << n << endl;
    printNum(n - 1);
} // end printNum
```

7. Given an integer $n>0$, write a recursive $\mathrm{C}++$ function that writes the integers $1,2, \ldots, n$.
8. Given an integer $n>0$, write a recursive $\mathrm{C}++$ function that returns the sum of the squares of 1 through $n$.
9. Write a recursive $\mathrm{C}++$ function that writes the digits of a positive decimal integer in reverse order.
10. a. Write a recursive $\mathrm{C}++$ function writeLine that writes a character repeatedly to form a line of $n$ characters. For example, writeLine('*', 5) produces the line $\%=* * *$.
b. Now write a recursive function writeBlock that uses writeLine to write $m$ lines of $n$ characters each. For example, writeBlock('*', 5, 3) produces the output
```
*****
*****
*****
```

11. What output does the following program produce?
```
int getValue(int a, int b, int n);
int main()
{
    cout << getValue(1, 7, 7) << end1;
    return 0;
} // end main
int getValue(int a, int b, int n)
{
    int returnValue = 0;
    cout << "Enter: a = " << a << " b = " << b << endl;
    int c = (a + b)/2;
    if (c * c <= n)
        returnValue = c;
    else
        returnValue = getValue(a, c-1, n);
    cout << "Leave: a = " << a << " b = " << b << endl;
    return returnValue;
} // end getValue
```

12. What output does the following program produce?
```
int search(int first, int last, int n);
int mystery(int n);
int main()
{
    cout << mystery(30) << endl;
    return 0;
} // end main
int search(int first, int last, int n)
{
    int returnValue = 0;
    cout << "Enter: first = " << first << " last = "
            << last << end7;
    int mid = (first + last)/2;
    if ( (mid * mid <= n) && (n < (mid+1) * (mid+1)) )
        returnValue = mid;
    else if (mid * mid > n)
        returnValue = search(first, mid-1, n);
    else
        returnValue = search(mid+1, last, n);
    cout << "Leave: first = " << first << " 1ast = "
            << last << endl;
    return returnValue;
} // end search
int mystery(int n)
{
    return search(1, n, n);
} // end mystery
```

13. Consider the following function that converts a positive decimal number to base 8 and displays the result.
```
void displayOctal(int n)
{
    if (n>0)
    {
        if (n / 8 > 0)
            displayOctal(n / 8);
        cout << n % 8;
    } // end if
} // end disp7ay0cta`
```

Describe how the algorithm works. Trace the function with $n=100$.
14. Consider the following program:

```
int f(int n);
int main()
{
    cout << "The value of f(8) is " << f(8) << endl;
    return 0;
} // end main
/** @pre n >= 0. */
int f(int n)
{
    cout << "Function entered with n=" << n << endl;
        switch (n)
```

```
    {
        case 0: case 1: case 2:
        return n + 1;
        default:
            return f(n-2) * f(n-4);
    } // end switch
} // end f
```

Show the exact output of the program. What argument values, if any, could you pass to the function $f$ to cause the program to run forever?
15. Consider the following function:

```
void recurse(int x, int y)
{
    if (y > 0)
    {
        x++;
        y--;
        cout << x << " " << y << end7;
        recurse(x, y);
        cout << x << " " << y << endl;
    } // end if
} // end recurse
```

Execute the function with $x=5$ and $y=3$. How is the output affected if x is a reference argument instead of a value argument?
16. Perform a box trace of the recursive function binarySearch, which appears in Section 2.4.2, with the array $1,5,9,12,15,21,29,31$ for each of the following search values:
a. 5
b. 13
c. 16

Imagine that you have 101 Dalmatians; no two Dalmatians have the same number of spots. Suppose that you create an array of 101 integers: The first integer is the number of spots on the first Dalmatian, the second integer is the number of spots on the second Dalmatian, and so on. Your friend wants to know whether you have a Dalmatian with 99 spots. Thus, you need to determine whether the array contains the integer 99.
a. If you plan to use a binary search to look for the 99 , what, if anything, would you do to the array before searching it?
b. What is the index of the integer in the array that a binary search would examine first?
c. If all of your Dalmatians have more than 99 spots, exactly how many comparisons will a binary search require to determine that 99 is not in the array?

This problem considers several ways to compute $x^{n}$ for some $n \geq 0$.
a. Write an iterative function power1 to compute $x^{n}$ for $n \geq 0$.
b. Write a recursive function power2 to compute $x^{n}$ by using the following recursive formulation:

$$
\begin{aligned}
& x^{0}=1 \\
& x^{n}=x \times x^{n-1} \text { if } n>0
\end{aligned}
$$

c. Write a recursive function power3 to compute $x^{n}$ by using the following recursive formulation:

$$
\begin{aligned}
& x^{0}=1 \\
& x^{n}=\left(x^{n / 2}\right)^{2} \text { if } n>0 \text { and } n \text { is even } \\
& x^{n}=x \times\left(x^{n / 2}\right)^{2} \text { if } n>0 \text { and } n \text { is odd }
\end{aligned}
$$

d. How many multiplications will each of the functions power1, power2, and power3 perform when computing $3^{32} ? 3^{19}$ ?
e. How many recursive calls will power2 and power3 make when computing $3^{32}$ ? $3^{19}$ ?
19. Modify the recursive rabbit function so that it is visually easy to follow the flow of execution. Instead of just adding "Enter" and "Leave" messages, indent the trace messages according to how "deep" the current recursive call is. For example, the call rabbit(4) should produce the output

```
Enter rabbit: n = 4
    Enter rabbit: n = 3
        Enter rabbit: n = 2
        Leave rabbit: n=2 value = 1
        Enter rabbit: n = 1
        Leave rabbit: n = 1 value = 1
    Leave rabbit: n = 3 value = 2
    Enter rabbit: n = 2
    Leave rabbit: n = 2 value = 1
Leave rabbit: n = 4 value = 3
```

Note how this output corresponds to Figure 2-19.
Consider the following recurrence relation:
$f(1)=1 ; f(2)=1 ; f(3)=1 ; f(4)=3 ; f(5)=5$;
$f(n)=f(n-1)+3 \times f(n-5)$ for all $n>5$.
a. Compute $f(n)$ for the following values of $n: 6,7,12,15$.
b. If you were careful, rather than computing $f(15)$ from scratch (the way a recursive $\mathrm{C}++$ function would compute it), you would have computed $f(6)$, then $f(7)$, then $f(8)$, and so on up to $f(15)$, recording the values as you computed them. This ordering would have saved you the effort of ever computing the same value more than once. (Recall the iterative version of the rabbit function discussed at the end of this chapter.)

Note that during the computation, you never need to remember all of the previously computed val-ues-only the last five. Taking advantage of these observations, write a C++ function that computes $f(n)$ for arbitrary values of $n$.

Write iterative versions of the following recursive functions: fact, writeBackward, binarySearch, and kSmall.

Prove that the function iterativeRabbit, which appears in Section 2.7, is correct by using invariants. (See Appendix F for a discussion of invariants.)
23. Consider the problem of finding the greatest common divisor (gcd) of two positive integers $a$ and $b$. The algorithm presented here is a variation of Euclid's algorithm, which is based on the following theorem: ${ }^{4}$

Theorem. If $a$ and $b$ are positive integers with $a>b$ such that $b$ is not a divisor of $a$, then $\operatorname{gcd}(a, b)=\operatorname{gcd}(b, a$ $\bmod b)$.

This relationship between $\operatorname{gcd}(a, b)$ and $\operatorname{gcd}(b, a \bmod b)$ is the heart of the recursive solution. It specifies how you can solve the problem of computing $\operatorname{gcd}(a, b)$ in terms of another problem of the same type. Also, if $b$ does divide $a$, then $b=\operatorname{gcd}(a, b)$, so an appropriate choice for the base case is $(a \bmod b)=0$.

This theorem leads to the following recursive definition:

$$
\operatorname{gcd}(a, b)= \begin{cases}b & \text { if }(a \bmod b)=0 \\ \operatorname{gcd}(b, a \bmod b) & \text { otherwise }\end{cases}
$$

The following function implements this recursive algorithm:

```
int gcd(int a, int b)
{
    if (a % b == 0) // Base case
        return b;
    else
        return gcd(b, a % b);
} // end gcd
```

a. Prove the theorem.
b. What happens if $b>a$ ?
c. How is the problem getting smaller? (That is, do you always approach a base case?) Why is the base case appropriate?
24. Let $c(n)$ be the number of different groups of integers that can be chosen from the integers 1 through $n-1$ so that the integers in each group add up to $n$ (for example, $n=4=[1+1+1+1]=[1+1+2]=[2+2]$ ). Write recursive definitions for $c(n)$ under the following variations:
a. You count permutations. For example, 1,2,1 and 1,1,2 are two groups that each add up to 4 .
b. You ignore permutations.

Consider the following recursive definition:

$$
\operatorname{Acker}(m, n)= \begin{cases}n+1 & \text { if } m=0 \\ \operatorname{Acker}(m-1,1) & \text { if } n=0 \\ \operatorname{Acker}(m-1, \operatorname{Acker}(m, n-1)) & \text { otherwise }\end{cases}
$$

This function, called Ackermann's function, is of interest because it grows rapidly with respect to the sizes of $m$ and $n$. What is $\operatorname{Acker}(1,2)$ ? Implement the function in $\mathrm{C}++$ and do a box trace of Acker $(1,2)$. (Caution: Even for modest values of $m$ and $n$, Ackermann's function requires many recursive calls.)

## Programming Problems

1. Implement a recursive function that computes $a^{n}$, where $a$ is a real number and $n$ is a nonnegative integer.
2. Implement the algorithm maxArray, discussed in Section 2.4.3, as a $\mathrm{C}++$ function. What other recursive definitions of maxArray can you describe?
3. Implement the binarySearch algorithm presented in this chapter for an array of strings.
4. Implement the algorithm kSma11, discussed in Section 2.4.4, as a C++ function. Use the first value of the array as the pivot.

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## Array-Based Implementations

## Chapter

 3
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C++ Interlude 1 C++ Classes
Chapter 2 Recursion: The Mirrors (for Section 3.3)You have seen that during the design of a solution, you must support severaloperations on data and therefore need to define abstract data types (ADTs). Only afteryou have clearly specified the operations of an ADT should you consider data structuresfor implementing it. This chapter explores implementation issues that involve arrays asthe underlying data structures.

An ADT is a collection of data and a set of operations on that data

Specifications indicate what ADT operations do, but not how to implement them

### 3.1 The Approach

We have said that a collection of data, together with a set of operations on that data, are called an abstract data type, or ADT. For example, suppose that you want to store a collection of names that you can search rapidly. The collection of names, together with operations that add a name, remove a name, and search for a name, can form an ADT that solves this problem.

Previously, we emphasized the specification of an abstract data type. When you design an ADT, you concentrate on what its operations do, but you ignore how you will implement them. That is, the description of an ADT's operations must be rigorous enough to specify completely their effect on the data, yet it must not specify how to store the data nor how to carry out the operations. For example, the operations of an ADT should not specify whether to store the data in consecutive memory locations or in disjoint memory locations. The result should be a set of clearly specified ADT operations.

How do you implement an ADT once its operations are clearly specified? That is, how do you store the ADT's data and carry out its operations? You begin by choosing particular data structures to store the data. Recall that a data structure is a construct that you can define within a programming language to store a collection of data. For example, C++ arrays are data structures.

Your first reaction to the implementation question might be to choose a data structure and then to write functions that access it in accordance with the ADT's operations. Although this point of view is not incorrect, hopefully you have learned not to jump right into code. In general, you should refine an ADT through successive levels of abstraction yielding successively more concrete descriptions of the ADT.

The choices that you make during the implementation process can affect the execution time of your code. For now, our analyses will be intuitive, but Chapter 10 will introduce you to quantitative techniques that you can use to weigh the trade-offs involved.

Recall that the client - that is, the program that uses the ADT - should see only a wall of available operations that act on data, as Figure 1-5 in Chapter 1 illustrates. Both the data structure that you choose to contain the data and the implementations of the ADT's operations are hidden behind the wall. By now, you should realize the advantage of this wall and take steps to prevent a client's direct access-either intentional or accidental-to the data structure, as shown in Figure 3-1. Why is such access undesirable? Public data fields allow the client to go around the wall of abstraction and access the ADT's data directly. For example, suppose that you use an array items to store an ADT's data. In a program that uses the ADT, you might, for example, accidentally access the first element in the array by writing

```
firstItem = items[0];
```

instead of by invoking an ADT operation. If you changed to another implementation of the ADT, your program would be incorrect. To correct your program, you would need to locate and change all occurrences of $i$ tems [0] -but first you would have to realize that items [0] is in error! Moreover, by violating the wall, a client could damage the ADT's data. If the data were ordered in a certain way, for example, the ADT's operations would ensure that the order was maintained. But if a client could alter the data directly, that order could be destroyed.

Note: Implementing an ADT as a C++ class provides a way for you to enforce the wall of an ADT, thereby preventing access of the data structure in any way other than by using the ADT's operations. A client then cannot damage the ADT's data. Moreover, the client is independent of the details of the ADT's implementation, because it adheres only to the ADT's specifications.

FIGURE 3-1 Violating the wall of ADT operations


### 3.1.1 Core Methods

The definition of a class that implements an ADT could be fairly involved. In general, you should not define the entire class and then attempt to test it. Instead, you should identify a group of core methods to both implement and test before continuing with the rest of the class definition. By leaving the definitions of the other methods for later, you can focus your attention and simplify your task. But what methods should be part of this group? In general, such methods should be central to the purpose of the class and allow reasonable testing. We sometimes will call a group of core methods a core group.

When dealing with a container such as a bag, for example, you cannot test most methods until you have created the container and placed items into it. Thus, adding objects to the container is a fundamental operation. If the method add does not work correctly, testing other methods such as remove would be pointless. Thus, the add method would be part of the group of core methods that we implement first.

To test whether add works correctly, we need a method that allows us to see the container's data. For the ADT bag, the method toVector serves this purpose, and so it is a core method. Any constructors are also fundamental and are in the core group. Similarly, other methods that a core method might call are part of the core group as well. Such methods might be public methods within the ADT or private "helper" methods.


Note: Methods such as add and remove that can alter the underlying data structure of a container are likely to have the most involved implementations. In general, you should define such methods before the others in the class. But since you can't test remove before add is correct, you should delay implementing it until after add is completed and thoroughly tested.

### 3.1.2 Using Fixed-Size Arrays

When implementing an ADT that represents a data collection, you need to store the data items and track their number. In an array-based implementation, you store the items in an array. How much of the array will the items occupy? Possibly all of it, but this is not likely. That is, you need to keep track of the array elements that you have assigned to contain the data and those that are available for use in the future. The maximum length of the array-its physical size-is a known, fixed value such as MAX_SIZE. You can keep track of the current number of items in the collection-that is, the collection's logical size - in a variable. An obvious benefit of this approach is that implementing an operation that reports this value will be easy.

When you add a first entry to an array, you typically place it in the array's first element-that is, the element whose index is 0 . Doing so, however, is not a requirement, especially for arrays that implement containers. For example, some container implementations can benefit by ignoring the array element whose index is 0 and using index 1 as the first element in the array. Sometimes you might want to use the elements at the end of the array before the ones at its beginning. For the ADT bag, we have no reason to be atypical, and so the objects in a bag can begin at index 0 of the array.

Another consideration is whether the container's entries should occupy consecutive elements of the array. Requiring the add method to place objects into an array consecutively is certainly reasonable, but why should we care, and is this really a concern? It is a concern, because we need to establish certain truths, or assertions, about our planned implementation so that the action of each method is not detrimental to other methods. For example, the method toVector must "know" where add has placed the entries. Our decision now also will affect what must happen later when we remove an entry from the container. If we insist that the bag's entries occupy consecutive array elements, will the method remove ensure that they remain so?

What happens after add places a new entry into the last available array element? A subsequent addition will be impossible unless an entry is first removed from the array. An add method that encounters a full array should either signal its client or allocate a larger array. In this chapter, add will return a boolean value to indicate whether it was successful. In C++ Interlude 2, which follows this chapter, you will learn how to resize the array. Rather than returning a value, a method can signal its client by throwing an exception. C++ Interlude 3 will discuss this approach.

### 3.2 An Array-Based Implementation of the ADT Bag

We will now make the previous discussion concrete by implementing the ADT bag as a class using an array to store its entries. Recall from Chapter 1 that the ADT bag's operations are

```
+getCurrentSize(): integer
+isEmpty(): boolean
+add(newEntry: ItemType): boolean
+remove(anEntry: ItemType): boolean
+clear(): void
+getFrequencyOf(anEntry: ItemType): integer
+contains(anEntry: ItemType): boolean
+toVector(): vector
```

These operations will become public methods in our class.
Each method will require access to both the array of bag entries and the current number of entries in the bag. Thus, we make the array and a counter data members of the class. Since the length of the
array defines the bag's capacity-that is, its maximum size-we will define a constant whose value is this size. To hide these data members from the clients of the class, we make them private. Thus, we define the following private data members within the header file for our class:

```
static const int DEFAULT_CAPACITY = 50;
ItemType items[DEFAULT_CAPACITY]; // Array of bag items
int itemCount; // Current count of bag items
int maxItems; // Max capacity of the bag
```

Figure 3-2 illustrates these data members, assuming a bag of integers.

FIGURE 3-2
An array-based implementation of the ADT bag


## Note: Data members should be private

By making a class's data members private, you control how a client can access or change their values. Doing so also makes debugging a program's logic easier. If the client could manipulate this data directly because the data was public, you would not know where to look for errors.

### 3.2.1 The Header File

Having chosen the data members, and knowing the details of BagInterface, which we developed in Chapter 1, we can write the header file shown in Listing 3-1 for our class of bags. We name the class ArrayBag and give it a default constructor, which is sufficient for this example. Notice the small value given to the constant DEFAULT_CAPACITY. It enables us to fill a bag completely and easily to verify that our class behaves gracefully in this situation. After you study C++ Interlude 2, you will be able to add another constructor that enables the client to set the capacity of the bag during program execution. We included the data field maxItems in anticipation of this enhancement.

Also notice that we declare a private method getIndexOf. This method returns the index of the element within the array items that contains a given entry. As you will see later, we can use getIndexOf to make the methods contains and remove more time efficient. You most likely would not decide to add getIndexOf to the header file until after you began work on the implementation of the public methods.

## LISTING 3-1 The header file for the class ArrayBag

```
/** Header file for an array-based implementation of the ADT bag.
@file ArrayBag.h */
#ifndef _ARRAY_BAG
#define _ARRAY_BAG
#include "BagInterface.h"
```

```
template<class ItemType>
class ArrayBag : public BagInterface<ItemType>
{
private:
    static const int DEFAULT_CAPACITY = 6;// Small size to test for a full bag
    ItemType items[DEFAULT_CAPACITY]; // Array of bag items
    int itemCount; // Current count of bag items
    int maxItems; // Max capacity of the bag
    // Returns either the index of the element in the array items that
    // contains the given target or -1, if the array does not contain
    // the target.
    int getIndexOf(const ItemType& target) const;
public:
    ArrayBag();
    int getCurrentSize() const;
    bool isEmpty() const;
    bool add(const ItemType& newEntry);
    bool remove(const ItemType& anEntry);
    void clear();
    bool contains(const ItemType& anEntry) const;
    int getFrequencyOf(const ItemType& anEntry) const;
    vector<ItemType> toVector() const;
}; // end ArrayBag
#include "ArrayBag.cpp"
#endif
``` a const method as a safeguard against an implementation error.

\subsection*{3.2.2 Defining the Core Methods}

Section 3.1.1 suggests that we consider the methods add and toVector as core methods and define them first. Since we will want to check that add correctly increments the count of the bag's entries, we should also implement getCurrentSize and isEmpty. Thus, we add these two methods to our core group, along with appropriate constructors.

As we define the methods for the class ArrayBag, we place them into the file ArrayBag.cpp, which begins as follows:
```

/** Implementation file for the class ArrayBag.
@file ArrayBag.cpp */
\#include "ArrayBag.h"

```

The constructor. The following default constructor initializes the current number of items in the bag to zero, as well as the bag's capacity:
```

template<class ItemType>
ArrayBag<ItemType>::ArrayBag()

```
```

{
itemCount = 0;
maxItems = DEFAULT_CAPACITY;
} // end default constructor

```

Although you can simply use an assignment statement to assign a value to a data member, as we have done here, it is preferable to use an initializer. Each initializer uses a functional notation that consists of a data member name followed by its initial value enclosed in parentheses-for example, itemCount (0). If you write more than one initializer, you separate them with commas. A colon precedes the first (or only) initializer. Thus, a better definition of ArrayBag's default constructor is
```

template<class ItemType>
ArrayBag<ItemType>::ArrayBag(): itemCount(0), maxItems(DEFAULT_CAPACITY)
{
} // end default constructor

```

Often the implementation of a constructor consists only of initializers, so its body is empty, as is the case here. Note that you can use these initializers with constructors but not with other methods.

Programming Tip: When a class has several data members, the constructor initializes them in the order in which they appear in the class definition instead of the order in which the initializers appear in the constructor definition. You should use the same order in both cases to avoid confusion, even if the initialization order does not make a difference.

The method add. To add a new item to a bag that is not full, we can place it right after the last item in the array by writing the following statement:
```

items[itemCount] = newEntry;

```

If we are adding to an empty bag, itemCount will be zero, and the assignment will be to items [0]. If the bag contains one entry, an additional entry will be assigned to items [1], and so on. In this way, no other items in the array need to move. After each addition to the bag, we increase the counter itemCount. Figure 3-3 depicts this insertion. If add is successful, it returns true. Otherwise-if the bag is full, for example-add returns false, as you can see from its definition:
```

temp7ate<class ItemType>
bool ArrayBag<ItemType>::add(const ItemType\& newEntry)
{
bool hasRoomToAdd = (itemCount < maxItems);
if (hasRoomToAdd)
{
items[itemCount] = newEntry;
itemCount++;
} // end if
return hasRoomToAdd;
} // end add

```

Note: The entries in a bag have no particular order. Thus, the method add can place a new entry into any convenient element of the array items. In the previous definition of add, that element is the one immediately after the last element used.

FIGURE 3-3 Inserting a new entry into an array-based bag


Note: An array-based implementation of an ADT restricts the number of items that you can store. Thus, the implementation should check whether the array has space available before inserting a new item, and the client should take appropriate action if the insertion is impossible.


Programming Tip: Variables that are local to a method's implementation should not be data members of the class.

The method toVector. The method toVector in our initial core group gets the entries that are in a bag and returns them to the client within a vector. A simple loop within toVector adds the bag's entries to this vector.
```

template<class ItemType>
vector<ItemType> ArrayBag<ItemType>:: toVector() const
{
vector<ItemType> bagContents;
for (int i = 0; i < itemCount; i++)
bagContents.push_back(items[i]);
return bagContents;
} // end toVector

```

Note: The class vector is in the Standard Template Library (STL) and is described in Appendix A.

The methods getCurrentSize and isEmpty. The last two methods in our core group have the following straightforward definitions:
```

template<class ItemType>
int ArrayBag<ItemType>::getCurrentSize() const
{
return itemCount;
} // end getCurrentSize
template<class ItemType>
bool ArrayBag<ItemType>::isEmpty() const
{
return itemCount == 0;
} // end isEmpty

```

Question 1 What happens to the array items when the method add cannot add another entry to it, because it is already full?
Question 2 If a client of ArrayBag creates a vector containing five items, what happens to those items after the vector is passed as an argument to the method toVector?

\subsection*{3.2.3 Testing the Core Methods}

Getting ready. After defining the core methods, you should test them. But what about the other methods in BagInterface? Since ArrayBag-as given in Listing 3-1-adheres to the specifications in BagInterface, the \(\mathrm{C}++\) compiler will look for a definition of each method declared in this interface. Should we wait until we complete their definitions to begin testing? Absolutely not! Testing methods as you write them makes finding logical errors easier. However, instead of writing a complete implementation of each method in BagInterface, we can provide incomplete definitions of the methods we choose to temporarily ignore.

An incomplete definition of a method is called a stub. The stub needs only to keep the syntax checker happy. For example, for each method that returns a value, you can avoid syntax errors by adding a return statement that returns a dummy value. Methods that return a boolean value should return false, for example. This result is reasonable, since the method does not perform its specified behavior. On the other hand, void methods can simply have an empty body.

For instance, the method remove ultimately will return true or false, so its stub must contain a return statement and could appear as follows:
```

temp7ate<class ItemType>
boo1 ArrayBag<ItemType>::remove(const ItemType\& anEntry)
{
return fa1se; // STUB
} // end remove

```

A stub for the void method clear could be
```

template<class ItemType>
void ArrayBag<ItemType>::clear()
{
// STUB
} // end clear

```

Note that if you plan to call a stub within your test program, the stub should report that it was invoked by displaying a message.

Programming Tip: Do not wait until you complete the implementation of an ADT before testing it. By writing stubs, which are incomplete definitions of required methods, you can begin testing early in the process.

A test program. Listing 3-2 contains a program that tests the core methods add, toVector, getCurrentSize, and isEmpty of the class ArrayBag at this stage of its development. The main function creates an empty bag by using the default constructor and passes it to the function bagTester. This function calls another function, displayBag, that tests the methods toVector and getCurrentSize.

\section*{LISTING 3-2 A program that tests the core methods of the class ArrayBag}
```

\#include <iostream>
\#include <string>
\#include "ArrayBag.h"
using namespace std;
void disp1ayBag(ArrayBag<string>\& bag)
{
cout << "The bag contains " << bag.getCurrentSize()
<< " items:" << endl;
vector<string> bagItems = bag.toVector();
int numberOfEntries = (int)bagItems.size();
for (int i = 0; i < numberOfEntries; i++)
{
cout << bagItems[i] << " ";
} // end for
cout << endl << end1;
} // end displayBag
void bagTester(ArrayBag<string>\& bag)
{
cout << "isEmpty: returns " << bag.isEmpty()
<< "; should be 1 (true)" << endl;
displayBag(bag);
string items[] = {"one", "two", "three", "four", "five", "one"};
cout << "Add 6 items to the bag: " << endl;
for (int i = 0; i < 6; i++)
{
bag.add(items[i]);
} // end for
disp1ayBag(bag);
cout << "isEmpty: returns " << bag.isEmpty()
<< "; should be 0 (false)" << endl;
cout << "getCurrentSize: returns " << bag.getCurrentSize()
<< "; should be 6" << endl;
cout << "Try to add another entry: add(\"extra\") returns "
<< bag.add("extra") << endl;
} // end bagTester
int main()
{
ArrayBag<string> bag;
cout << "Testing the Array-Based Bag:" << endl;
cout << "The initial bag is empty." << endl;
bagTester(bag);
cout << "A11 done!" << endl;
return 0;
} // end main

```

\section*{Output}

Testing the Array-Based Bag:
The initial bag is empty.
isEmpty: returns 1 ; should be 1 (true)
The bag contains 0 items:

Add 6 items to the bag:
The bag contains 6 items:
one two three four five one
isEmpty: returns 0 ; should be 0 (false)
getCurrentSize: returns 6; should be 6
Try to add another entry: add("extra") returns 0
Al1 done!

Note: If you define a constructor for a class but do not also define a default constructor, the compiler will not generate one for you. In this case, a statement such as
```

    ArrayBag<string> aBag;
    ```
is illegal.

\subsection*{3.2.4 Implementing More Methods}

Now that we can add items to a bag successfully, we can define the other methods. Although we will not do so here, you should test each new definition as you complete it.

Programming Tip: Testing each method as you define it is especially important when its logic is not obvious. However, even the simplest method is susceptible to a careless mistake, which can lead to hours of frustrating debugging of other methods that depend on the supposedly correct method.

The method getFrequencyOf. To count the number of times a given object occurs in a bag, we count the number of times the object occurs in the array items. Using a while loop to cycle through the array's indices from 0 to itemCount - 1 , we compare the given object to every object in the array. Each time we find a match, we increment a counter. When the loop ends, we simply return the value of the counter.

The method definition follows:
```

temp7ate<c1ass ItemType>
int ArrayBag<ItemType>::getFrequencyOf(const ItemType\& anEntry) const

```
```

{
int frequency = 0;
int curIndex = 0; // Current array index
while (curIndex < itemCount)
{
if (items[curIndex] == anEntry)
{
frequency++;
} // end if
curIndex++; // Increment to next entry
} // end while
return frequency;
} // end getFrequency0f

```

The method contains. One easy way to define the method contains is to have it call the method getFrequency0f, which we just defined. If the frequency of occurrence of a given entry is greater than zero, the bag must contain that entry. If the frequency is zero, the bag does not contain the entry. Thus, you could define contains as follows:
```

template<class ItemType>
bool ArrayBag<ItemType>::contains(const ItemType\& target) const
{
return getFrequency0f(target) > 0;
} // end contains

```

Although this method will work, it usually will do more work than necessary. After all, getFrequencyOf must check every entry in the bag, whereas contains could quit looking at entries as soon as it finds one that is the same as the given entry. That is, its logic can be described by the following pseudocode:
```

contains(anEntry)
while (anEntry is not found and we have more array elements to check)
{
if (anEntry equals the next array entry)
anEntry is found in the array
}

```

This loop terminates under one of two conditions: Either anEntry has been found in the array or the entire array has been searched without success.

Based on this pseudocode, a definition of the method contains follows:
```

template<class ItemType>
boo1 ArrayBag<ItemType>::contains(const ItemType\& anEntry) const
{
bool found = false;
int curIndex = 0; // Current array index
while (!found \&\& (curIndex < itemCount))
{
if (anEntry == items[curIndex])
{
found = true;
} // end if
curIndex++; // Increment to next entry
} // end while
return found;
} // end contains

```

Question 3 What is an advantage and a disadvantage of calling the method getFrequencyOf from contains?

\section*{Note: Testing the additional methods}

As you define additional methods for the class ArrayBag, you should test them. While you could focus only on these additional methods, you really should form a test program incrementally. In this way, you test all of the methods you have defined so far. By using a small value for the constant DEFAULT_CAPACITY, you easily can test what happens when a bag becomes full.

\section*{Note: Two kinds of loops}

To count how many times an entry occurs in an array, the method getFrequencyOf uses a loop that cycles through all of the array's entries. In fact, the body of the loop executes itemCount times. In contrast, to indicate whether a given entry occurs in an array, the loop in the method contains ends as soon as the desired entry is discovered. The body of this loop executes between one and itemCount times. You should be comfortable writing loops that execute either a definitive or a variable number of times.

\subsection*{3.2.5 Methods That Remove Entries}

We have postponed the definition of the method remove until now because it involves more thought than the other methods. The method clear, however, is relatively simple, even though we are leaving it until last.

The method remove. We now want to remove a given entry-call it anEntry-from a bag. If the entry occurs more than once in the bag, we will remove only one occurrence. Exactly which occurrence is removed is unspecified. We will simply remove the first occurrence of anEntry that we encounter while searching for it. Adhering to our specifications in BagInterface, we will return either true or false to indicate whether the removal was successful.

Assuming that the bag is not empty, we search the array items until either we find anEntry in the array or we note where it occurs within the array. Figure 3-4 illustrates the array after a successful search.

FIGURE 3-4 The array items after a successful search for the string "Alice"


FIGURE 3-5 (a) A gap in the array items after deleting the entry in items [index] and decrementing itemCount; (b) shifting subsequent entries to avoid a gap; (c) the array after shifting


Now consider how to remove the located item from the array. You could blank it out, but this strategy can lead to gaps in the array, as Figure 3-5a illustrates. An array that is full of gaps has three significant problems:
- itemCount - 1 is no longer the index of the last item in the array. You would need another variable, lastPosition, to contain this index.
- Because the items are spread out, the method contains might have to look at every element of the array, even when only a few items are present.
- When items [maxItems - 1] is occupied, the bag could appear full, even when fewer than maxItems items are present.

We could get rid of that gap by copying each successive entry to the previous element in the array, beginning with the entry after the removed item and ending with the last entry, as shown in Figure 3-5b. The result of shifting these entries is shown in Figure 3-5c. This time-consuming approach is not necessary, however.

Remember that we are not required to maintain any particular order for a bag's entries. So instead of shifting array entries after removing an entry, we can replace the entry being removed with the last entry in the array, as follows. After locating anEntry in items [index], as Figure 3-6a indicates, we copy the entry in items [itemCount - 1] to items [index] (Figure 3-6b). We then ignore the entry in items[itemCount - 1] by decrementing itemCount.

The following pseudocode details the steps needed to remove a given entry from a bag. Note that by decrementing itemCount before using it in an index, we compute itemCount -1 once instead of twice.

Search the array items for anEntry
if (anEntry is in the bag at items [index])
\{
Decrement the counter itemCount
```

    items[index] = items[itemCount]
    return true
    }
e1se
return false

```

The initial search for anEntry is the same search as done in the method contains. Calling contains here is not much help, as we want to know not only whether the bag contains anEntry but also, if it does, where it occurs in the array. Clearly, we could modify the loop in contains slightly, but a better way places the loop in a private method that both contains and remove can call. This method is getIndexOf, which we declared earlier in the header file and whose definition follows:
```

temp1ate<c1ass ItemType>
int ArrayBag<ItemType>::getIndexOf(const ItemType\& target) const
{
bool found = false;
int result = -1;
int searchIndex = 0;
// If the bag is empty, itemCount is zero, so loop is skipped
while (!found \&\& (searchIndex < itemCount))
{
if (items[searchIndex] == target)
{
found = true;
result = searchIndex;
}
e1se
{
searchIndex++;
} // end if
} // end while
return result;
} // end get IndexOf

```

FIGURE 3-6 Avoiding a gap in the array while removing an entry


The method remove now has the following definition:
```

template<class ItemType>
boo1 ArrayBag<ItemType>::remove(const ItemType\& anEntry)
{
int locatedIndex = getIndexOf(anEntry);
bool canRemoveItem = !isEmpty() \&\& (locatedIndex > -1);
if (canRemoveItem)
{
itemCount--;
items[locatedIndex] = items[itemCount];
} // end if
return canRemoveItem;
} // end remove

```

\section*{Programming Tip: Should the method getIndexOf be public or private?}

You might think that the method getIndexOf, since it returns the index of an entry in the array items, would be useful to a client. While that might be true, there are important reasons why the method should be private. We declared the array items as private because it is detail of our array-based implementation. The client cannot access the entries in this array without invoking a public method of our class. The indices of these entries are also implementation details that are of no use to the client because it cannot access the array items by name. Any method that returns an index to a private array should be declared as private.

Question 4 Revise the definition of the method contains so that it calls the method getIndexOf.
Question 5 Should we revise the specification of the method contains so that if it locates a given entry within the bag, it returns the index of that entry?
Question 6 Revise the definition of the method getIndexOf so that it does not use a boolean variable.

The method clear. The method clear can make a bag appear empty simply by setting itemCount to zero. Thus, we have the following definition for this method:
```

template<class ItemType>
void ArrayBag<ItemType>::clear()
{
itemCount = 0;
} // end clear

```

\subsection*{3.2.6 Testing}

To test our completed class, we can add the following statements to the method bagTester that appears in Listing 3-2:
```

cout << "contains(\"three\"): returns " << bag.contains("three")
<< "; should be 1 (true)" << endl;
cout << "contains(\"ten\"): returns " << bag.contains("ten")
<< "; should be 0 (false)" << endl;

```
```

cout << "getFrequencyOf(\"one\"): returns "
<< bag.getFrequencyOf("one") << " should be 2" << end7;
cout << "remove(\"one\"): returns " << bag.remove("one")
<< "; should be 1 (true)" << endl;
cout << "getFrequencyOf(\"one\"): returns "
<< bag.getFrequencyOf("one") << " should be 1" << end7;
cout << "remove(\"one\"): returns " << bag.remove("one")
<< "; should be 1 (true)" << endl;
cout << "remove(\"one\"): returns " << bag.remove("one")
<< '; should be 0 (false)" << endl;
cout << endl;
displayBag(bag);
cout << "After clearing the bag, ";
bag.clear();
cout << "isEmpty: returns " << bag.isEmpty()
<< "; should be 1 (true)" << endl;

```

Note: References such as bag.itemCount and bag. items[4] would be illegal within the client, because \(i\) temCount and \(i\) tems are within the private portion of the class.

Note: To implement an ADT, given implementation-independent specifications of the ADT's operations, you first must choose a data structure to contain the data. Next, you declare a class within a header file. The ADT's operations are public methods within the class, and the ADT's data consists of class members that are typically private. You then implement the class's methods within an implementation file. The program that uses the class will be able to access the data only by using the ADT's operations.

Note: In practice, when choosing among implementations of an ADT , you must ask whether the fixed-size restriction of an array-based implementation presents a problem in the context of a particular application. The answer to this question depends on two factors. The obvious factor is whether, for a given application, you can predict in advance the maximum number of items in the ADT at any one time. If you cannot, it is quite possible that an operation - and hence the entire program-will fail because the ADT in the context of a particular application requires more storage than the array can provide.

On the other hand, if for a given application you can predict in advance the maximum number of items in an ADT at any one time, you must explore a more subtle factor: Would you waste storage by declaring an array to be large enough to accommodate this maximum number of items? Consider a case in which the maximum number of items is large, but you suspect that this number rarely will be reached. For example, suppose that a bag could contain as many as 10,000 items, but the actual number of items in the bag rarely exceeds 50 . If you declare 10,000 array locations at compilation time, at least 9,950 array locations will be wasted most of the time.

We can avoid both of the previous cases by allocating a larger-or smallerarray during execution. The next C++ Interlude will show you how.

\subsection*{3.3 Using Recursion in the Implementation}

The previous chapter gave several examples of how to use recursion to process an array. If you examine the definition of ArrayBag that we have just developed, you will see two loops that can be replaced by recursion: One is in the public method getFrequencyOf, and the other is in the private method getIndexOf. Let's look at the details.

\subsection*{3.3.1 The Method getIndexOf}

We begin with the private method that searches the array for a given entry. The iterative definition examines each entry in the array one at a time in sequential order until it either finds the desired one or reaches the end of the array without success. Suppose that searchIndex is the index of the next array element to examine. We have two base cases:
- If items [searchIndex] is the entry we seek, we are done.
- If searchIndex equals itemCount, we are done, because the entry is not in the array.

The recursive step is to search the rest of the array beginning with items [searchIndex +1 ].
To accomplish this recursion, getIndexOf needs searchIndex as a parameter. Since the method is private, we need not define yet another private method for it to call. However, we must revise its declaration in the header file as follows:
```

int getIndexOf(const ItemType\&target, int searchIndex) const;

```

We then can write the following recursive definition for this method:
```

temp1ate<class ItemType>
int ArrayBag<ItemType>::getIndexOf(const ItemType\& target, int searchIndex) const
{
int result = -1;
if (searchIndex < itemCount)
{
if (items[searchIndex] == target)
{
result = searchIndex;
}
else
{
result = getIndex0f(target, searchIndex + 1);
} // end if
} // end if
return result;
} // end getIndexOf

```

Because we have changed the header of the method, we need to locate and change the calls to getIndexOf within the other methods of ArrayBag. Thus, we will change the calls within the methods remove and contains from
```

getIndexOf(anEntry);

```
to
```

getIndexOf(anEntry, 0);

```

\subsection*{3.3.2 The Method getFrequencyOf}

The method getFrequencyOf also searches an array, much as getIndexOf does, but unlike getIndexOf, getFrequencyOf must examine every entry in the array. That is, the search does not end if an array entry matches the entry we seek.

Suppose that searchIndex is the index of the next array element to examine. A recursive definition for this search has only one base case: If searchIndex equals itemCount, we are done. However, we have two recursive steps:
- If items [searchIndex] is the entry we seek, the frequency of occurrence of this entry is one more than its frequency of occurrence in the rest of the array.
- If items [searchIndex] is not the entry we seek, the frequency of occurrence of the entry is the same as its frequency of occurrence in the rest of the array.

Just as we had to add another parameter to getIndexOf, the recursive method here needs searchIndex as a parameter. However, getFrequencyOf is a public method, so we cannot change its signature. Instead, getFrequencyOf must call a private method that accomplishes the recursive search. Thus, let's add the following declaration to the private portion of the header file:
```

int countFrequency(const ItemType\& target, int searchIndex) const;

```

The method of getFrequencyOf simply calls this private method, so its definition is as follows:
```

template<class ItemType>
int ArrayBag<ItemType>::getFrequencyOf(const ItemType\& anEntry) const
{
return countFrequency(anEntry, 0);
} // end getFrequencyOf

```

The recursive definition of the private method countFrequency is
```

template<class ItemType>
int ArrayBag<ItemType>::countFrequency(const ItemType\& target,
int searchIndex) const
{
if (searchIndex < itemCount)
{
if (items[searchIndex] == target)
{
return 1 + countFrequency(target, searchIndex + 1);
}
else
{
return countFrequency(target, searchIndex + 1);
} // end if
}
else
return 0; // Base case
} // end countFrequency

```

Stylistically, we prefer to have only one exit from a method. To revise countFrequency so that it has one return statement instead of three, we define a local variable, frequency, to contain the return value. The method then has the following definition. Note that the base case is implicit.
```

template<c1ass ItemType>
int ArrayBag<ItemType>: :countFrequency (const ItemType\& target,
int searchIndex) const

```
```

{
int frequency = 0;
if (searchIndex < itemCount)
{
if (items[searchIndex] == target)
{
frequency = 1 + countFrequency(target, searchIndex + 1);
}
else
{
frequency = countFrequency(target, searchIndex + 1);
} // end if
} // end if
return frequency;
} // end countFrequency

```

\section*{Summary}
1. By using a class to implement an ADT, you encapsulate the ADT's data and operations. In this way, you can hide implementation details from the program that uses the ADT. In particular, by making the class's data members private, you can change the class's implementation without affecting the client.
2. Given an interface that specifies an ADT in an implementation-independent way, derive a class from the interface and declare the class within a header file. Choose a data structure to contain the ADT's data. Then implement the class's methods within an implementation file.
3. You should make a class's data members private so that you can control how a client can access or change the data.
4. An array-based implementation of an ADT stores the ADT's data in an array.
5. Generally, you should not define an entire class and then attempt to test it. Instead, you should identify a group of core methods to both implement and test before continuing with the rest of the class definition.
6. Stubs are incomplete definitions of a class's methods. By using stubs for some methods, you can begin testing before the class is completely defined.
7. A client must use the operations of an ADT to manipulate the ADT's data.

\section*{Exercises}
1. Consider a bag of integers. Write a client function that computes the sum of the integers in the bag aBag.
2. Write a client function replace that replaces a given item in a given bag with another given item. The function should return a boolean value to indicate whether the replacement was successful.
3. The previous exercise describes the function replace. This operation exists outside of the ADT bag; that is, it is not an ADT bag operation. Instead, its implementation is written in terms of the ADT bag's operations.
a. What is an advantage and a disadvantage of the way that replace is implemented?
b. What is an advantage and a disadvantage of adding the operation replace to the ADT bag?
4. Design and implement an ADT that represents a rectangle. Include typical operations, such as setting and retrieving the dimensions of the rectangle, and finding the area and the perimeter of the rectangle.
5. Design and implement an ADT that represents a triangle. The data for the ADT should include the three sides of the triangle but could also include the triangle's three angles. This data should be in the private section of the class that implements the ADT.

Include at least two initialization operations: one that provides default values for the ADT's data, and another that sets this data to client-supplied values. These operations are the class's constructors. The ADT also should include operations that look at the values of the ADT's data; change the values of the ADT's data; compute the triangle's area; and determine whether the triangle is a right triangle, an equilateral triangle, or an isosceles triangle.
6. Write a recursive array-based implementation of the method toVector for the class ArrayBag.
7. Write a client function that merges two bags into a new third bag. Do not destroy the original two bags.
8. Specify and define a method for ArrayBag that removes a random entry from the bag.
9. Add a constructor to the class ArrayBag that creates a bag from a given array of entries.

\section*{Programming Problems}
1. Design and implement an ADT that represents the time of day. Represent the time as hours and minutes on a 24 -hour clock. The hours and minutes are the private data members of the class that implements the ADT. Include at least two initialization operations: one that provides a default value for the time, and another that sets the time to a client-supplied value. These operations are the class's constructors. Also include operations that set the time, increase the present time by a number of minutes, and display the time in 12 -hour and 24 -hour notations.
2. Design and implement an ADT that represents a calendar date. You can represent a date's month, day, and year as integers (for example, 4/1/2014). Include operations that advance the date by one day and display the date by using either numbers or words for the months. As an enhancement, include the name of the day.
3. Design and implement an ADT that represents a price in U.S. currency as dollars and cents. After you complete the implementation, write a client function that computes the change due a customer who pays \(x\) for an item whose price is \(y\).
4. Design and implement an ADT that represents a bank account. The data of the ADT should include the customer name, the account number, and the account balance. The initialization operation should set the data to client-supplied values. Include operations for a deposit and a withdrawal, the addition of interest to the balance, and the display of the statistics of the account.
5. Add the methods union, intersection, and difference to the class ArrayBag. Exercises 6, 7, and 8 of Chapter 1 describe these operations.
6. Implement the ADT set that you specified in Programming Problem 5 of Chapter 1 by using an array.
7. Implement the ADT pile that you specified in Programming Problem 6 of Chapter 1 by using an array.
8. Implement the ADT polynomial that Exercise 9 in Chapter 1 describes by using an array.
9. Implement the ADT appointment book, described in Section 1.4.1 of Chapter 1. Write a program that demonstrates your new class.
10. Specify and implement an ADT for fractions. Provide operations that add, subtract, multiply, and divide these numbers. The results of all arithmetic operations should be in lowest terms, so include a private method reduceToLowestTerms. Exercise 23 in Chapter 2 will help you with the details of this method. To simplify the determination of a fraction's sign, you can assume that the denominator of the fraction is positive.
11. Specify and implement an ADT for mixed numbers, each of which contains an integer portion and a fractional portion in lowest terms. Assume the existence of the ADT fraction, as described in the previous problem. Provide operations that add, subtract, multiply, and divide mixed numbers. The results of all arithmetic operations should have fractional portions that are in lowest terms. Also include an operation that converts a fraction to a mixed number.
12. Implement the ADT recipe book as described in Section 1.4.2 of Chapter 1. In doing so, implement the ADT measurement. Add operations as necessary. For example, you should add an operation to the recipe book to scale a recipe.
13. You can use either a set or a bag to create a spell checker. The set or bag serves as a dictionary and contains a collection of correctly spelled words. To see whether a word is spelled correctly, you see whether it is contained in the dictionary. Use this scheme to create a spell checker for the words in an external file. To simplify your task, restrict your dictionary to a manageable size.
14. Repeat the previous project to create a spell checker, but instead place the words whose spelling you want to check into a bag. The difference between the dictionary (the set or bag containing the correctly spelled words) and the bag of words to be checked is a bag of incorrectly spelled words.

\section*{Pointers, \\ Polymorphism, and Memory Allocation}

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\section*{Prerequisites}

Chapter 3 Array-Based Implementations

This C++ Interlude discusses memory allocation for variables and arrays and introduces you to pointers to complete the discussion on polymorphism begun in the first C++ Interlude. Pointers are a powerful tool for the programmer, but you must be careful to use them correctly. This interlude explains the need for pointers and shows when and how to safely use them. Certain aspects of pointers are not covered in this interlude, but those features either are unnecessary in this textbook or introduce unsafe practices.

We will introduce memory allocation during execution instead of compilation and show how pointers are involved with this process. We will also discuss how to create a new and larger array during execution, if an array becomes full. You will see examples of this process at the end of this interlude.

Throughout our discussion here, and throughout the book, we will stress safe and secure programming practices. To begin our discussion, let's consider how memory is allocated for local variables and parameters.

\section*{C2.1 Memory Allocation for Variables and Early Binding of Methods}

When you declare an ordinary variable x to have the data type int , the \(\mathrm{C}++\) compiler allocates a memory cell that can hold an integer. You use the identifier \(x\) to refer to this cell. To put the value 5 in the cell, you could write
```

int x = 5;

```

To display the value that is in the cell, you could write
```

cout << "The value of x is " << x << endl;

```

As discussed in Chapter 2, a function's locally declared variables such as x are placed into an activation record with its parameters and some bookkeeping data. These activation records are stored in an area of your application's memory called the run-time stack. Each time a function is called, an activation record is automatically created on the run-time stack. When the function ends, the activation record is destroyed, freeing the memory used for the local variables and parameters. At that point, the function's local variables and their values are no longer accessible to your program. Your program then returns to where the function was invoked and executes the statement following the function call. This behavior is the same for methods.

When you create an object, the storage for the data members of that object are also placed into an activation record for the currently executing function or method. The statements
```

PlainBox<string> myPlainBox;
MagicBox<string> myMagicBox = MagicBox<string>();

```
show two different ways of invoking default constructors to create instances of PlainBox and MagicBox, respectively. When these objects are instantiated, their data fields are placed on the runtime stack just as primitive data types are. Recall from C++ Interlude 1 that the names of those data fields are item and firstItemStored.

The compiler also knows that if you invoke the setItem method on these objects by writing
```

myP1ainBox.setItem("Fun Item");
myMagicBox.setItem("Secret Item");

```
the PlainBox version of the setItem method should be called for myPlainBox, and the MagicBox version of the setItem method should be called for myMagicBox. This choice, which is an example of early binding, is made during compilation and cannot be altered during execution.

Most of the time, this automatic memory management and early binding are all you need in your

Two situations when automatic memory management and early binding are insufficient
program. However, two situations can arise when they are not:
- You want to take advantage of polymorphism.
- You must access an object outside of the function or method that creates it.

\section*{C2.2 A Problem to Solve}

Suppose your friend was so impressed with your work on the three boxes discussed in C++ Interlude 1 that she asked you to continue helping with the video game. She would like you to write a function that takes two arguments: an object of any of the three types of boxes and an item of type string. The function should place the item in the box by invoking the box's setItem method.


Since ToyBox and MagicBox are derived from PlainBox, as Figure C2-1 illustrates, you may think that the following function definition would suffice:
```

void placeInBox(PlainBox<string>\& theBox, string theItem)
{
theBox.setItem(theItem);
} // end placeInBox

```

The parameter theBox can accept as an argument any PlainBox object or an object of a class derived from PlainBox. This function could then be used in the following sequence of statements:
```

string specialItem = "Riches beyond compare!";
string otherItem = "Hammer";
PlainBox<string> myPlainBox;
placeInBox(myPlainBox, specialItem);
MagicBox<string> myMagicBox;
placeInBox(myMagicBox, otherItem);
placeInBox(myMagicBox, specialItem); // specialItem is stored!
cout << myMagicBox.getItem() << end1; // "Riches beyond compare!"

```

Although this code compiles, it does not perform as you would expect. Since otherItem has already been stored in the magic box, specialItem should not replace that item. Unfortunately, when the item stored in the magic box is displayed, the output is Riches beyond compare. The reason is that, in our function, the statement
```

theBox.setItem(theItem);

```
invokes the PlainBox version of the setItem method instead of the MagicBox version. In this case, the compiler determined the version of the method to invoke from the type of the parameter theBox instead of from the type of its corresponding argument. The following set of statements has a similar result:
```

string specialItem = "Riches beyond compare!";
string otherItem = "Hammer";

```

An example of early binding

Another example of early binding
```

PlainBox<string> mySpecialBox = MagicBox<string>();
mySpecia1Box.setItem(otherItem);
mySpecialBox.setItem(specialItem); // specialItem is stored!
cout << mySpecialBox.getItem() << endl; // "Riches beyond compare!"

```

In both situations, the version of setItem that will be called is determined when the program is compiled. In this second case, even though we instantiated a MagicBox object for mySpecialBox, the variable mySpecia1Box is of type PlainBox, so the P1ainBox version of setItem is called. The same decision logic applies to our function placeInBox.

This code is correct from the compiler's perspective. The compiler assumes it is our intent to have the parameter theBox and the variable mySpecialBox behave as PlainBox objects. We need a way to communicate to the compiler that the code to execute should not be determined until the program is running. This is called late binding, which is an aspect of polymorphism. To solve this problem and have both our function and simple-code examples execute as we intend, we need two tools: pointer variables and virtual methods.

\section*{C2.3 Pointers and the Program's Free Store}

The new operator allocates memory on the free store

To take advantage of late binding, we do not want our objects to be in an activation record on the runtime stack. We need another location. When a C++ program begins execution, in addition to the runtime stack, the operating system sets aside memory for the code-called code storage or text storage - and for any global variables and static variables-called static storage. Your program is also given extra memory, called the heap, or free store, which a programmer can use to store data. Figure C2-2 illustrates these portions of memory.

We allocate memory for a variable on the free store by using the new operator. After allocating memory for the variable, the new operator returns the memory address of the variable in the free store so the program can use it. This memory address must be placed in a special type of variable called a pointer variable, or simply a pointer. A pointer variable contains the location, or address in memory, of a memory cell.

FIGURE C2-2 Sample program memory layout


FIGURE C2-3 Run-time stack and free store after myboxPtr points to a MagicBox object and its


Unlike variables on the run-time stack, which have their memory allocated and deallocated automatically, variables placed in the free store persist in memory even when the function or method that created them ends. This means that when a programmer allocates memory from the free store for a variable, the programmer has the responsibility to deallocate that memory when the variable is no longer needed. Failure to do so often results in a memory leak. A memory leak is memory that has been allocated for use but is no longer needed and cannot be accessed or deallocated.

To indicate that a variable is a pointer, we place the character * after the type of the data the pointer references. \({ }^{1}\) For example, the statement
MagicBox<string>* myBoxPtr = new MagicBox<string>();
creates a MagicBox object in the free store and places the address of the object in the local variable myBoxPtr. We say that myBoxPtr points to a MagicBox object. To call a method of an object that is in the free store, we use the notation ->:
```

string someItem = "Something Free";
myBoxPtr->setItem(someItem);

```

Figure C2-3 shows the state of memory and the local variables after the previous statements execute.

Observe that this newly created object has no programmer-defined name. The only way to access its methods is indirectly via the pointer that new creates, that is, by using myBoxPtr-> as in the previous example. Usually we will simplify the diagrams so that only the object in the free store and its pointer are drawn, as in Figure C2-4.

\footnotetext{
1 Some programmers place the * next to the name of the variable. The authors of this textbook believe placing the * next to the data type more clearly shows that the variable is of a pointer type.
}

If you allocate memory on the free store, you eventually must deallocate it

Indicate a pointer type by writing an asterisk after the data type

FIGURE C2-4 myBoxPtr and the object to which it points


Pointer variables are simply another type of variable and follow the same rules as other variable types. For example, if you declare a pointer variable that points to a ToyBox<string> object, you can have it point only to ToyBox<string> objects, not MagicBox<string> objects or ToyBox<double> objects, as the following statements illustrate:
```

ToyBox<string>* toyPtr = new ToyBox <string>(); // OK
ToyBox<string>* boxPtr = new MagicBox<string>(); // Error!
ToyBox<string>* somePtr = new ToyBox <double>(); // Error!

```

If you have two pointer variables that can point to the same type of object, you can make them point to the same object by using the assignment operator \(=\). For example, the statement
```

somePtr = toyPtr;

```
makes the variable somePtr point to the same ToyBox object to which toyPtr points. In such cases, it is helpful to think of the assignment operator as copying the value stored within the pointer variable on its right into the pointer variable on its left. Since pointers store addresses of objects, the value copied is the location of an object in the free store. The object itself is not copied. The result is that both pointer variables point to the same object, as shown in Figure C2-5.

If you declare a pointer variable but do not immediately create an object for it to reference, you should set the pointer to nul1ptr. An assignment like the following one is necessary because \(\mathrm{C}++\) does not initialize the pointer for you:
```

ToyBox<int>* myToyPtr = nu11ptr;

```

\section*{FIGURE C2-5 Two pointer variables that point to the same object}


\section*{C2.3.1 Deallocating Memory}

When the memory to which a pointer variable points is no longer needed, you deallocate it by using the delete operator. You should then set the pointer variable to nullptr as a signal that it no longer references or points to an object:
```

delete somePtr;
somePtr = nul1ptr;

```

If we did not set somePtr to nul1ptr in this example, somePtr would be an example of a dangling pointer, because it would still contain the address of an object that was deallocated. Dangling pointers can be the source of serious errors.

Always remember that if a method of a class uses new to create an object, some method of the class must use delete eventually to free the object's memory. It is a characteristic of safe and secure programming that any object allocating memory from the free store also assumes responsibility to deallocate that memory when it is no longer needed.

Pointers give the programmer a great amount of control over how memory is used and when binding occurs, as we will discuss shortly. They also can create severe problems and hard-to-find bugs if not properly used. The two most common errors involving pointers are the memory leak and the dangling pointer, as presented in this section. Let's look at how these problems can occur in a program so that you can prevent them in your code.

\section*{C2.3.2 Avoiding Memory Leaks}

Memory leaks occur when an object has been created in the free store, but the program no longer has a way to access it. Because the object cannot be accessed, it cannot be deleted and so takes up memory, even though it cannot be used. A program can quickly run out of memory and crash if it has several memory leaks. Even a single statement that is called multiple times can cause a memory leak.

Consider the following statements:
MagicBox<string>* myBoxPtr = new MagicBox<string>();
MagicBox<string>* yourBoxPtr = new MagicBox<string>();
yourBoxPtr = myBoxPtr; // Results in inaccessible object
Figure C2-6 traces the execution of these statements. Eventually, yourBoxPtr points to the same object as myBoxPtr, and so the object to which yourBoxPtr originally pointed is no longer accessible, resulting in the memory leak shown in part \(c\) of the figure.

To prevent this leak, we should not create the second object. Instead, either initialize yourBoxPtr to null ptr or simply set it to myBoxPtr:
```

MagicBox<string>* myBoxPtr = new MagicBox<string>();
MagicBox<string>* yourBoxPtr = myBoxPtr;

```

A more subtle memory leak occurs when a function or method creates an object in the free store and loses the pointer to it by either not returning that pointer to the caller or not storing it in a class data member. That is what happens in the function given in Listing C2-1.

\section*{LISTING C2-1 Poorly written function that allocates memory in the free store}
void myLeakyFunction(const double\& someItem)
\{
ToyBox<double>* someBoxPtr = new ToyBox<doub1e>();
someBoxPtr->setItem(someItem);
\} // end myLeakyFunction

The parameter someItem and the pointer someBoxPtr are both stored in an activation record on the run-time stack and so are automatically destroyed after the function ends. The object created in the free store by new ToyBox<double> (RED) is still in the free store. Since the only reference we had to it was someBoxPtr and that has been destroyed, we no longer have a way to get to that object, and so we have a memory leak.

If you use the new operator to allocate memory, you must use the delete operator to deallocate it

FIGURE C2-6 (a) Creating the first object; (b) creating the second object; (c) assignment causes an inaccessible object
(a) MagicBox<string>* myBoxPtr = new MagicBox<string>();

(b) MagicBox<string>* yourBoxPtr = new MagicBox<string>();

(c) yourBoxPtr \(=\) myBoxPtr;


Inaccessible object in free store (memory leak)

We have several options to fix this function. The first is to delete the object before the function terminates. We can do this by adding the lines
```

de1ete someBoxPtr;
someBoxPtr = nul1ptr;

```
at the end of the function. If the ToyBox object pointed to by someBoxPtr is needed only by this function, a better implementation choice would have been to not allocate memory from the free store, but to use a local variable such as is done in the following statements:

ToyBox<double> someBox(); // someBox is not a pointer variable someBox.setItem(someItem);

If the object created in this function is required outside the function, the function can return a pointer to it so that the caller can access the object. To implement this option, we would change the function's return type from void to ToyBox<double>* and return someBoxPtr, which is a pointer to a ToyBox<double> object. The resulting function definition is
```

ToyBox<doub7e>* pluggedLeakyFunction(const doub7e\& someItem)
{
ToyBox<doub7e>* someBoxPtr = new ToyBox<doub7e>();
someBoxPtr->setItem(someItem);
return someBoxPtr;
} // end p7uggedLeakyFunction

```

You could call this function by writing statements such as
```

doub7e boxValue = 4.321;
ToyBox<double>* toyPtr = pluggedLeakyFunction(boxValue);

```

Now toyPtr points to the object created by the ToyBox constructor in pluggedLeakyFunction. When a function returns a pointer to an object that it created in the free store, the segment of the program using that pointer must take responsibility for deleting the object. Otherwise, a memory leak could still occur. When documenting such a function, you should add comments to indicate this responsibility. For example, you could precede the definition of pluggedLeakyFunction with these comments:
```

/** Creates an object in the free store and returns a pointer
to it. Caller must delete the object when it is no longer
needed. */

```

Of course, this comment does not prevent a memory leak, but it alerts users of your function about the potential for a leak. The function could still be misused and called with a statement such as
```

pluggedLeakyFunction(boxValue); // Misused; returned pointer is lost

```
in which case the returned pointer to the object in the free store is lost and a memory leak occurs.
The best option for preventing a memory leak is to not use a function to return a pointer to a newly created object. Instead, you should define a class that has a method for this task. The class takes responsibility for deleting the object in the free store and ensures that there is no memory leak. At a minimum, such a class will have three parts: a method that creates the object in the free store, a data field that points to the object, and a method-the destructor-that deletes the object when the class instance is no longer needed. Although \(\mathrm{C}++\) Interlude 1 introduced class destructors, we must say more about them now.

Every \(\mathrm{C}++\) class has a destructor that has the same name as the class, but is preceded by the tilde \((\sim)\) character. For objects that are local variables, the destructor is called when the activation record containing the object is removed from the run-time stack. If the object was created using new and is stored in the free store, the destructor is called when the client uses the operator delete to free the memory allocated to the object.

Often, the compiler-generated destructor is sufficient for a class, but if the class itself creates an object on the free store by using the new operator, it is a safe and secure programming practice to implement a destructor to ensure that the memory for that object is freed.

Listing C2-2 gives the header file for a GoodMemory class that demonstrates how to avoid a memory leak. The pointer variable in our previous function pluggedLeakyFunction is now a data field in GoodMemory. As long as this field points to an existing object in the free store, we will have access to the object.

\section*{LISTING C2-2 Header file for the class GoodMemory}
```

/** @file GoodMemory.h */
\#include <string>
\#include "ToyBox.h"
using namespace std;
class GoodMemory
{

```

To prevent a memory leak, do not use a function to return a pointer to a newly created object
```

private:
ToyBox<string>* someBoxPtr;
public:
GoodMemory(); // Defau7t constructor
~GoodMemory(); // Destructor
void fixedLeak(const doub7e\& someItem);
}; // end GoodMemory

```

Because the class GoodMemory has a pointer variable for a data field, we should define a default constructor to initialize the pointer to nul1ptr. The destructor for this class simply needs to delete the object to which the pointer points. Listing C2-3 shows the implementation file for this class.
```

LISTING C2-3 Implementation file for the class GoodMemory
/** @file GoodMemory.cpp */
\#include "GoodMemory.h"
GoodMemory::GoodMemory() : someBoxPtr(nu11ptr)
{
} // end default constructor
GoodMemory::~GoodMemory()
{
de1ete someBoxPtr;
} // end destructor
void GoodMemory::fixedLeak(const double\& someItem)
{
someBoxPtr = new ToyBox<doub1e>();
someBoxPtr->setItem(someItem);
} // end fixedLeak

```

Unlike the original function myLeakyFunction, as presented in Listing C2-1, the following client function uses the class GoodMemory to guarantee that no memory leak occurs:
```

void goodFunction()
{
double boxValue = 4.321; // Original statement
GoodMemory gmObject; // Create a safe memory object
gmObject.fixedLeak(boxValue); // Perform the task
} // end goodFunction

```

The compiler greatly helps us here. First, the compiler can check to ensure that a GoodMemory object has been created before calling the method fixedLeak. This guarantees a safe memory allocation and deallocation. Then, since the variable gmObject is a local variable, the compiler automatically calls the GoodMemory destructor when execution of the function goodFunction ends. Thus, the memory allocated in the free store is freed.

\section*{C2.3.3 Avoiding Dangling Pointers}

As mentioned earlier, a dangling pointer is a pointer variable that no longer references a valid object. Four situations can cause dangling pointers, but we will cover only three of them because we consider the fourth one to be an unsafe practice. \({ }^{2}\)
- As Section C2.3.1 discusses, if you do not set a pointer variable to nul1 ptr after using delete, you leave the pointer dangling.
- Similarly, if you declare a pointer variable but do not assign it a value, the result is a dangling pointer. As with any C++ variable, when C++ creates a pointer variable, its value is undefined. That is, C++ does not automatically initialize or clear the memory that a variable represents. The compiler and the programmer have no way to check that the value in the pointer variable actually points to an object in the free store. That value might simply be whatever was in the memory when the pointer was created. For example, suppose that you write statements such as

MagicBox<int>* myMagicBoxPtr;
myMagicBoxPtr->getItem();
Here, the pointer variable myMagicBoxPtr is not assigned an object to point to. When the method getItem is called, the program will abort-usually with a segment fault errorbecause the program treats the value in myMagicBoxPtr as the address of a MagicBox<int> object and tries to find the getItem method at that address. Since it is very unlikely that a MagicBox<int> object happens to be there, the program ends abnormally.

If you need to create a pointer variable but do not have an object for it to point to, you should always set it to nul1ptr:
```

MagicBox<int>* myMagicBoxPtr = nu11ptr;

```

Then your code can compare the pointer variable to nullptr to see whether it points to a valid object on the free store, as in the following example:
```

if (myMagicBoxPtr != nul1ptr)
myMagicBoxPtr->getItem();

```

If the pointer is not nul1ptr, you can use the pointer to call a method.
- The third situation that can cause a dangling pointer is subtle, and the best way to guard against it is careful programming. Consider the following statements that we saw earlier when discussing memory leaks:
```

MagicBox<string>* myBoxPtr = new MagicBox<string>();
MagicBox<string>* yourBoxPtr = myBoxPtr;

```

Executing this code results in the memory configuration shown in Figure C2-7, where both yourBoxPtr and myBoxPtr point to the same object. We say that yourBoxPtr is an alias of myBoxPtr since they both refer to the same object. There is nothing wrong with this code so far, and it does not have any memory leaks.

Suppose we then execute the following statements:
```

de1ete myBoxPtr;
myBoxPtr = nul1ptr;
yourBoxPtr->getItem();

```

Situations that can cause a dangling pointer

FIGURE C2-7 Two pointers referencing (pointing to) the same object


Here we try to practice safe and secure programming by setting myBoxPtr to nul1 ptr after deleting the object that it points to. But the call yourBoxPtr->getItem() results in the program aborting. What happened? Figure C2-8 shows the state of memory just prior to this call. As you can see, the object pointed to by myBoxPtr was deleted, as it should have been. The problem arises because yourBoxPtr still references the object's location in the free store, even though the object no longer exists. Since the object no longer exists, yourBoxPtr is a dangling pointer and the program aborts when we try to call a method on that object.

FIGURE C2-8 Example of a dangling pointer


\section*{Programming Tip: How to avoid dangling pointers}
- Set pointer variables to nul1ptr either initially or when you no longer need them. If a class has a pointer variable as a data field, the constructor should always initialize that data field, to point either to an object or to nul1ptr. The class GoodMemory in Listing C2-3 demonstrates this safeguard.
- Test whether a pointer variable contains nul1ptr before using it to call a method.
- Try to reduce the use of aliases in your program. As you will see, that is not always possible or desirable in certain situations.
- Do not delete an object in the free store until you are certain that no other alias needs to use it.
- Set all aliases that reference a deleted object to nul1ptr when the object is deleted.

\section*{C2.4 Virtual Methods and Polymorphism}

Now that we have some of the basics of pointers behind us, we can dive into the implementation of polymorphism. To allow the compiler to perform the late binding necessary for polymorphism, you must declare the methods in the base class as virtual. In C++ Interlude 1, we began a discussion of
virtual methods-methods that use the keyword virtual to indicate that they can be overridden. In the example in Section C2.2, PlainBox is the base class, which we defined in Listing C1-3 of C++ Interlude 1. The code that we wrote then did not behave as we desired. To correct that problem, we must declare as virtual the methods of P7ainBox that we want other classes to override, as Listing C2-4 shows. Notice that only the header file of the base class (PlainBox) needs to be revised. Declaring the methods setItem and getItem as virtual makes it possible for the method code to be bound late.

\section*{LISTING C2-4 Revised header file for the class PlainBox}
```

/** @file PlainBox.h */
\#ifndef _PLAIN_BOX
\#define _PLAIN_BOX
template<class ItemType>; // Indicates this is a template
// Declaration for the class PlainBox
class PlainBox
{
private:
// Data field
ItemType item;
public:
// Default constructor
PlainBox();
// Parameterized constructor
PlainBox(const ItemType\& theItem);
// Mutator method that can change the value of the data field
virtual void setItem(const ItemType\& theItem);
// Accessor method to get the value of the data field
virtual ItemType getItem() const;
}; // end PTainBox
\#include "PlainBox.cpp" // Include the implementation file
\#endif

```

To fully implement late binding, we must create the variables in the free store and use pointers to reference them. Thus, we must also change the code from our first example of early binding in Section C2.2 to
```

string specialItem = "Riches beyond compare!";
string otherItem = "Hammer";

```
```

PlainBox<string>* myPlainBoxPtr = new PlainBox<string>();
placeInBox(myP7ainBoxPtr, specialItem);
MagicBox<string>* myMagicBoxPtr = new MagicBox<string>();
placeInBox(myMagicBoxPtr, otherItem);
placeInBox(myMagicBoxPtr, specialItem);
cout << myMagicBoxPtr->getItem() << endl;

```

Next, we must change the function \(p\) laceInBox to accept a pointer to a PlainBox<string> object:
```

void placeInBox(P1ainBox<string>* theBox, string theItem)
{
theBox->setItem(theItem);
} // end placeInBox

```

The last change is to free the memory used by the variables in the free store by adding the following statements to the program:
```

delete myP1ainBoxPtr;
myPlainBoxPtr = nul1ptr;
delete myMagicBoxPtr;
myMagicBoxPtr = nul1ptr;

```

Our function placeInBox now will call the correct version of setItem according to the type of box pointed to by its first argument.

The use of virtual methods has a significant impact on the future use of a class. Imagine that you had compiled the class PlainBox and its implementation before you wrote the derived class MagicBox. If you then wrote MagicBox, assuming access to the compiled class PlainBox, you could override getItem because it is virtual in PlainBox. As a result, you would change the behavior of getItem for instances of MagicBox, even though PlainBox was already compiled. That is, classes that define virtual methods are extensible: You can add capabilities to a derived class without having access to the ancestor's source statements.

\section*{Note: Key points about virtual methods}
- A virtual method is one that a derived class can override.
- You must implement a class's virtual methods. (Pure virtual methods are not included in this requirement.)
- A derived class does not need to override an existing implementation of an inherited virtual method.
- Any of a class's methods may be virtual. However, if you do not want a derived class to override a particular method, the method should not be virtual.
- Constructors cannot be virtual.
- Destructors can and should be virtual. Virtual destructors ensure that future descendants of the object can deallocate themselves correctly.
- A virtual method's return type cannot be overridden.

An ordinary C++ array is statically allocated

\section*{C2.5 Dynamic Allocation of Arrays}

When you declare an array in C++ by using statements such as
```

const int MAX_SIZE = 50;
doub7e myArray[MAX_SIZE];

```
the compiler reserves a specific number-MAX_SIZE, in this case-of memory cells for the array. This memory allocation occurs before your program executes, so it is not possible to wait until execution to give MAX_SIZE a value. We have already discussed the problem this fixed-size data structure causes when your program has more than MAX_SIZE items to place into the array.

You just learned how to use the new operator to allocate memory dynamically-that is, during program execution. Although Section C2.3 showed you how to allocate memory for a single variable or object, you actually can allocate memory for many at one time. If you write
```

int arraySize = 50;
doub7e* anArray = new double[arraySize];

```
the pointer variable anArray will point to the first item in an array of 50 items. Unlike MAX_SIZE, arraySize can change during program execution. You can assign a value to arraySize at execution time and thus determine how large your array is. That is good, but how do you use this array?

Regardless of how you allocate an array-statically, as in the first example, or dynamically, as in the second-you can use an index and the familiar array notation to access its elements. For example, anArray [0] and anArray [1] are the first two items in the array anArray.

When you allocate an array dynamically, you need to return its memory cells to the system when you no longer need them. As described earlier, you use the de1ete operator to perform this task. To deallocate the array anArray, you write
```

delete [ ] anArray;

```

Note that you include brackets when you apply de1ete to an array.
Now suppose that your program uses all of the array anArray, despite having chosen its size during execution. You can allocate a new and larger array, copy the old array into the new array, and finally deallocate the old array. Doubling the size of the array each time it becomes full is a reasonable approach. The following statements double the size of anArray:
```

doub7e* oldArray = anArray;
anArray = new double[2 * arraySize];
for (int index = 0; index < arraySize; index++)
anArray[index] = oldArray[index];
delete [ ] oldArray;
// Copy pointer to array
// Double array size
// Copy old array
// Deallocate old array

```

Subsequent discussions in this book will refer to both statically allocated and dynamically allocated arrays. Our array-based ADT implementations will use statically allocated arrays for simplicity. The programming problems will ask you to create array-based implementations that use dynamically allocated arrays. We will refer to such arrays as resizable.

\subsection*{2.5.1 A Resizable Array-Based Bag}

We can use a resizable array to implement the ADT bag so that the bag never becomes full—within the bounds of the particular computer, of course. If we look at the header file for the class ArrayBag, as given in Listing 3-1 of Chapter 3, and the implementation of the various methods, we see that only the add method needs to be changed. When the array-and hence the bag-becomes full, add can double the capacity of the array instead of failing to add the item and returning false. The revised method follows. Since the bag is never full, this version of the add method can always return true.
```

temp7ate<c1ass ItemType>
bool ArrayBag<ItemType> : :add(const ItemType\& newEntry)
{
bool hasRoomToAdd = (itemCount < maxItems);
if (!hasRoomToAdd)
{
ItemType* oldArray = items;

```

Use the new operator to allocate an array dynamically
delete returns a dynamically allocated array to the system for reuse

You can increase the size of a dynamically allocated array
```

    items = new ItemType[2 * maxItems];
    for (int index = 0; index < maxItems; index++)
        items[index] = oldArray[index];
    delete [ ] oldArray;
    maxItems = 2 * maxItems;
    } // end if
// We can always add the item
items[itemCount] = newEntry;
itemCount++;
return true;
} // end Resizab7eArrayBag add

```

Doubling the array items each time the bag is full is not as attractive as it might first seem. Each time you expand the size of an array, you must copy its contents. When the array is a small 50 -element array, you copy the 50 -element array to a 100 -element array before completing the addition. The next 49 additions then can be made quickly without copying the array.

If you have a 50,000 -element array that is full, you must allocate memory for 100,000 entries before copying the 50,000 entries in the current array. This takes considerably longer to do and a significantly greater amount of memory. It is possible that you needed storage for only 50,001 entries and will have 49,999 unused array elements wasting memory. Once you realize that your array is too large, you could reduce its size by allocating a smaller array and copying the entries to this new array. However, there is no way to predict whether or not you should do so. The application might need the extra array elements in the future.

In Chapter 4, we consider another implementation of the ADT bag that does not have this memory-allocation problem and that never becomes full.

\section*{Link-Based Implementations}

\section*{Chapter}

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C++ Interlude 2 C++ Pointers, Polymorphism, and Memory Allocation

T
his chapter introduces you to a link-based data structure using C++ pointers, which were described in C++ Interlude 2. You will learn how to work with such structures and how to use one to implement a class of bags. The material in this chapter is essential to much of the presentation in the chapters that follow.

Defining a pointer to a node

Dynamically allocating a node

Referencing a node member

\subsection*{4.1 Preliminaries}

Let's consider components that can be linked to one another. Each component-usually called a node - contains both a data item and a "pointer" to the next item. Because each node must contain two pieces of information-the data item and a pointer to the next node-it is natural to conclude that each node should be an object. One member of the object is the data item and the other is a pointer. Typically, such pointers are \(\mathrm{C}++\) pointer variables.

\section*{FIGURE 4-1 A node}


Figure 4-1 illustrates a single node, and Figure 4-2 shows several nodes linked together. Here, the data portion of each node contains a string. We use arrows to represent the pointers, but what type of pointer should you use within a node, and to what will it point? You might guess that the pointer should point to a string, but actually it must point to a node that contains the string. Because pointers can point to any data type except files-and since nodes are objects-pointers can, in fact, point to nodes. Thus, a node of type Node, for example, will have as one of its members a pointer to another node of type Node. For example, the statement
```

Node<string>* nodePtr;

```
defines a pointer variable nodePtr that can point to a node of type Node that contains a string as its data item. Nodes should be dynamically allocated. For example,
```

nodePtr = new Node<string>();

```
allocates a node to which nodePtr points. To access the members of a node, you will need the -> notation introduced in C++ Interlude 2, because the node does not have a user-defined name. Suppose our class of nodes has an accessor method, getItem, that returns the value of the data member item. To reference item in the node to which nodePtr points, you write nodePtr->getItem(). \({ }^{1}\) Likewise, if our class of nodes has the method getNext to access the data member next, the expression nodePtr \(->\) getNext () represents the pointer next. For example, if nodePtr points to the second node in Figure 4-2,

FIGURE 4-2 Several nodes linked together


1 An alternative notation, (*nodePtr). getItem(), is possible, but this book will not use it. Although this notation is analogous to the use of the dot operator to reference the members of a named structure, the \(->\) operator is more suggestive of pointers.
nodePtr->getItem() returns the string "cd" and nodePtr->getNext() returns a pointer to the third node. We can save this new pointer in nodePtr itself by writing
```

nodePtr = nodePtr->getNext();

```

If we execute the previous statement again, nodePtr will point to the fourth node in Figure 4-2. Executing the statement one more time makes nodePtr point to the last node shown in Figure 4-2. What is the value of the member next in this last node? That is, what value does nodePtr->getNext()return? We want this value to be a signal that we have reached the last of the linked nodes. If this value is nu11ptr, we easily can detect when we have reached the end of the linked nodes.

We have one more detail to consider: Nothing so far points to the first node. If you cannot get to the first node, you cannot get to the second node; and if you cannot get to the second node, you cannot get to the third node; and so on. The solution is to have an additional pointer whose sole purpose is to point to the first of several linked nodes. Such a pointer is called the head pointer or simply the head of the data structure.

Figure 4-3 illustrates a linked chain of nodes with a head pointer. Observe that the head pointer headPtr is different from the other pointers in the diagram in that it is not within one of the nodes. The variable headPtr simply enables you to access the chain's beginning. What value should headPtr contain if there are no nodes for it to point to? Assigning headPtr the value nu11ptr is a logical choice in this case.

FIGURE 4-3 A head pointer to the first of several linked nodes


Note: Although all of the pointers shown in Figure 4-3 point to nodes, the head pointer is a simple pointer variable, whereas the next members are within nodes.

Programming Tip: It is a common mistake to think that before you can assign headPtr a value, you first must create a new Node object. This misconception is rooted in the belief that the variable headPtr does not exist if it does not point to a node. This is not at all true; headPtr is a pointer variable waiting to be assigned a value. Thus, for example, you can assign nu11ptr to headPtr without first using new. In fact, the sequence
```

headPtr = new Node; // An incorrect use of new
headPtr = nul1ptr;

```
destroys the content of the only pointer-headPtr-to the newly created node, as Figure 4-4 illustrates. Thus, you have needlessly created a new node and then made it inaccessible. This action creates a memory leak in your program and should be avoided!

The head pointer points to the first node

If headPtr is nul7ptr, it points to nothing

FIGURE 4-4 A lost node
headPtr \(=\) new Node<string>(); headPtr \(=\) nul7ptr;


\subsection*{4.1.1 The Class Node}

Before we continue, let's examine the C++ code for the class Node. Listing 4-1 gives the header file for the class. In addition to the data members item and next that we introduced earlier, the class declares three constructors, two accessor methods-one each for item and next-and two mutator methods-again one each for item and next. Notice that within this class, a pointer to a node has the data type Node<ItemType>*.

\section*{LISTING 4-1 The header file for the template class Node}
```

/** @file Node.h */
\#ifndef _NODE
\#define _NODE
template<class ItemType>
class Node
{
private:
ItemType item; // A data item
Node<ItemType>* next; // Pointer to next node
public:
Node();
Node(const ItemType\& anItem);
Node(const ItemType\& anItem, Node<ItemType>* nextNodePtr);
void setItem(const ItemType\& anItem);
void setNext(Node<ItemType>* nextNodePtr);
ItemType getItem() const ;
Node<ItemType>* getNext() const ;
}; // end Node
\#include "Node.cpp"
\#endif

```

Listing 4-2 shows the implementation file for Node.

\section*{LISTING 4-2 The implementation file for the class Node}
```

/** @file Node.cpp */
\#include "Node.h"
\#include <cstddef>

```
```

template<class ItemType>
Node<ItemType>::Node() : next(nu11ptr)
{
} // end defau7t constructor
template<class ItemType>
Node<ItemType>::Node(const ItemType\& anItem) : item(anItem), next(nu11ptr)
{
} // end constructor
temp1ate<c1ass ItemType>
Node<ItemType>: :Node(const ItemType\& anItem, Node<ItemType>* nextNodePtr) :
item(anItem), next(nextNodePtr)
{
} // end constructor
temp1ate<c1ass ItemType>
void Node<ItemType>::setItem(const ItemType\& anItem)
{
item = anItem;
} // end setItem
template<class ItemType>
void Node<ItemType>::setNext(Node<ItemType>* nextNodePtr)
{
next = nextNodePtr;
} // end setNext
temp1ate<c1ass ItemType>
ItemType Node<ItemType>::getItem() const
{
return item;
} // end getItem
temp1ate<c1ass ItemType>
Node<ItemType>* Node<ItemType>::getNext() const
{
return next;
} // end getNext

```

\subsection*{4.2 A Link-Based Implementation of the ADT Bag}

Now that we have a class of nodes to use when defining link-based data structures, we will use it to define a link-based implementation of the ADT bag. Doing so will allow us to examine some of the basic practices for creating such implementations. Subsequent chapters will expand on these ideas.

Unlike array-based implementations, a link-based implementation does not impose a fixed maximum size on the data structure-except, of course, as imposed by the storage limits of the system. Additionally, data items do not move during an ADT's insertion and removal operations.

You need to represent the items in the ADT bag and its size. Figure 4-5 indicates one possible way to represent this data by using pointers. Here headPtr points to a linked chain of the nodes

VideoNote Core LinkedBag methods

FIGURE 4-5 A link-based implementation of the ADT bag

containing the items in the bag. The integer itemCount is the current number of items in the bag. Both headPtr and itemCount will be private data members of our class.

From Chapter 1, recall the interface BagInterface that describes the operations of the ADT bag. Also remember that we can use this interface when developing any implementation of this ADT. The bag operations in this interface, given in UML notation, are as follows:
```

+getCurrentSize(): integer
+isEmpty(): boolean
+add(newEntry: ItemType): boolean
+remove(anEntry: ItemType): boolean
+clear(): void
+getFrequencyOf(anEntry: ItemType): integer
+contains(anEntry: ItemType): boolean
+toVector(): vector

```

These operations correspond to the public methods in our class.

\subsection*{4.2.1 The Header File}

The header file given in Listing 4-3 declares the class LinkedBag. The private data members are headPtr, which points to the first node in the linked chain that contains the items in the bag, and itemCount, which counts these items. The private section of the class also declares a private method, whose purpose will become clear shortly. The public section declares a default constructor, a copy constructor-whose purpose is discussed near the end of Section 4.2.2-a destructor, and the public methods of the class. The public method declarations are identical to those in the class ArrayBag as given in Chapter 3. As you will see, the compiler-generated copy constructor and destructor are not sufficient for link-based implementations.

Recall from C++ Interlude 2 that destructors should be virtual. We have tagged LinkedBag's destructor with the keyword virtual in the header file. Failure to do so will result in a warning from the compiler.

\section*{LISTING 4-3 The header file for the class LinkedBag}
```

/** ADT bag: Link-based implementation.
@file LinkedBag.h */
\#ifndef _LINKED_BAG
\#define _LINKED_BAG
\#include "BagInterface.h"
\#include "Node.h"
template<class ItemType>
class LinkedBag : public BagInterface<ItemType>

```
```

{
private:
Node<ItemType>* headPtr; // Pointer to first node
int itemCount; // Current count of bag items
// Returns either a pointer to the node containing a given entry
// or the null pointer if the entry is not in the bag.
Node<ItemType>* getPointerTo(const ItemType\& target) const;
public:
LinkedBag();
LinkedBag(const LinkedBag<ItemType>\& aBag); // Copy constructor
virtua1 ~LinkedBag(); // Destructor should be virtual
int getCurrentSize() const;
bool isEmpty() const;
boo1 add(const ItemType\& newEntry);
bool remove(const ItemType\& anEntry);
void clear();
bool contains(const ItemType\& anEntry) const;
int getFrequencyOf(const ItemType\& anEntry) const;
vector<ItemType> toVector() const;
}; // end LinkedBag
\#include "LinkedBag.cpp"
\#endif

```

\subsection*{4.2.2 Defining the Core Methods}

Just as we did when we defined an array-based implementation of the ADT bag in Chapter 3, we begin our new implementation by defining the same core methods: the constructor, add, toVector, getCurrentSize, and isEmpty. We will write stubs for the remaining methods, including the copy constructor and destructor.

The constructor. The following default constructor initializes the head pointer and the current number of items in the bag:
```

template<class ItemType>
LinkedBag<ItemType>::LinkedBag() : headPtr(nu17ptr), itemCount(0)
{
} // end default constructor

```

The method add. A bag does not order its entries, so the method add can insert a new item at any convenient location within the linked chain of nodes that contains the bag's entries. The most convenient place for us to make this insertion is at the beginning of the chain, because the first node is the only one that we can access directly. Figure 4-6 illustrates how we add a new node to the beginning of an existing chain of nodes.

You must make headPtr point to the new node, and the new node must point to the node that had been at the beginning of the chain. Note how the following definition accomplishes this:
```

template<class ItemType>
boo1 LinkedBag<ItemType>::add(const ItemType\& newEntry)
{
// Add to beginning of chain: new node references rest of chain;

```

Inserting a node at the beginning of a linked chain

Insertion into an empty chain is really an insertion at the beginning of the chain

A traverse operation visits each node in the linked chain
// (headPtr is nul7ptr if chain is empty)
Node<ItemType>* newNodePtr = new Node<ItemType>();
newNodePtr->setItem(newEntry);
newNodePtr->setNext(headPtr); // New node points to chain
headPtr \(=\) newNodePtr; // New node is now first node
itemCount++;
return true;
\} // end add

FIGURE 4-6 Inserting at the beginning of a linked chain


Observe that if the bag is empty before the insertion headPtr is nul1ptr, so the next pointer of the new node is set to nullptr. This step is correct because the new item is the last item-as well as the first item-in the chain.


Programming Tip: Our method add uses the operator new to create a new node object and place it on the heap. Because C++ does not have garbage collection as Java does, it is our class's responsibility to ensure that the object is removed from the heap using delete. As discussed in the second C++ Interlude, the class destructor is one method where we do this. Since we created this node to store an item we are adding to the bag, it follows that another opportunity to delete the node is when we remove the item from the bag.

Question 1 Consider a linked chain of three nodes, such that each node contains a string. The first node contains "A", the second node contains "B", and the third node contains " C ".
a. Write \(\mathrm{C}++\) statements that create the described linked chain. Beginning with a head pointer headPtr that contains nul1ptr, create and attach a node for " C ", then create and attach a node for " B ", and finally create and attach a node for " A ".
b. Repeat part \(a\), but instead create and attach nodes in the order "A", "B", "С".

The method toVector. Recall that the method toVector retrieves the entries that are in a bag and returns them to the client within a vector. A loop within toVector adds the bag's entries to this vector. In the array-based implementation, this loop simply accesses an array of these entries. Here we must retrieve the entries from the nodes in a chain. To do that, we must move from node to node; that is, we must traverse the chain. As we visit each node, we copy its data item into the vector.

Let's write some high-level pseudocode for this loop, given the linked chain pictured in Figure 4-5.
```

Let a current pointer point to the first node in the chain
while (the current pointer is not the null pointer)
{
Assign the data portion of the current node to the next element in a vector
Set the current pointer to the next pointer of the current node
}

```

This solution requires that you keep track of the current position within the chain. Thus, you need a pointer variable-let's call it curPtr - that points to the current node.


Note: The pointer variable curPtr is analogous to the integer variable curIndex that we used in Section 3.2.4 of Chapter 3 to keep track of the current entry in an array.

Initially, curPtr must point to the first node. Because headPtr points to the first node, simply copy headPtr into curPtr by writing

Node<ItemType>* curPtr \(=\) headPtr;
Then you can use the expression curPtr->getItem() to access the data portion of the current node. After copying the data into the vector, you advance the current pointer to the next node by writing
```

curPtr = curPtr->getNext();

```

Figure 4-7 illustrates this action. If the previous assignment statement is not clear, consider
```

Node<ItemType>* temp = curPtr->getNext();
curPtr = temp;

```
and then convince yourself that the intermediate variable temp is not necessary.

FIGURE 4-7 The effect of the assignment curPtr = curPtr->getNext()


These ideas lead to the following definition of toVector:
```

temp1ate<c1ass ItemType>
vector<ItemType> LinkedBag<ItemType>::toVector() const
{
vector<ItemType> bagContents;
Node<ItemType>* curPtr = headPtr;
int counter = 0;
while ((curPtr != nu11ptr) \&\& (counter < itemCount))

```
```

    {
    bagContents.push_back(curPtr->getItem());
    curPtr = curPtr->getNext();
    counter++;
    } // end while
    return bagContents;
    } // end toVector

```

Here curPtr points to each node in a nonempty chain during the course of the loop's execution, and so the data portion of each node is accessed and assigned to the end of the vector. After the last node is accessed, curPtr becomes nul1ptr, and the loop terminates. When the bag is empty-that is, when headPtr is nu71ptr-the loop is correctly skipped. Note that the variable counter, while not necessary, provides a defense against going beyond the end of the chain.

Programming Tip: A common error in the while statement we used in toVector is to compare curPtr->getNext() instead of curPtr with nul1ptr. When curPtr points to the last node of a nonempty chain, curPtr->getNext() is nul1ptr, and so the loop would terminate before accessing the data in the last node. In addition, when the chain is empty, headPtr-and therefore-curPtr are nullptr, making the value of curPtr->getNext() undefined. Such references are incorrect and should be avoided.

The methods isEmpty and getCurrentSize. The last two methods in our core group have the following definitions, which are like the ones in the array-based implementation given in the previous chapter.
```

temp1ate<c1ass ItemType>
boo1 LinkedBag<ItemType>::isEmpty() const
{
return itemCount== 0;
} // end isEmpty
template<class ItemType>
int LinkedBag<ItemType>::getCurrentSize() const
{
return itemCount;
} // end getCurrentSize

```

\section*{Note: Testing the core methods}

Just as we did in the previous chapter, you should test each method as you define it. In fact, with little change, you can revise the program given in Listing 3-2 and use it to test the core methods that we have just defined. The bag in that program became full, but it will not here. In Section 4.4, we will examine how to change that program so that it can test multiple bag implementations.

Question 2 Why are only a few changes necessary to reuse the code in Listing 3-2? How would you implement the changes using the "find and replace" functionality of a text editor or IDE?
Question 3 Why is a LinkedBag object not concerned about becoming full?

\subsection*{4.2.3 Implementing More Methods}

Having successfully defined the core methods, we will continue to define the remaining ones in our class.

The method getFrequencyOf. To count the number of times a given object occurs in a bag, we count the number of times the object occurs in the linked chain. To do so, we need to traverse the chain and compare each of its data items with the given object. Each time we find a match, we increment a counter. When the traversal ends, we return the value of the counter.

The loop we will use is like the one we wrote in the method toVector. Here is the definition of getFrequencyOf:
```

template<class ItemType>
int LinkedBag<ItemType>::getFrequencyOf(constItemType\& anEntry) const
{
int frequency = 0;
int counter = 0;
Node<ItemType>* curPtr = headPtr;
while ((curPtr != nullptr) \&\& (counter < itemCount))
{
if (anEntry == curPtr->getItem())
{
frequency++;
} // end if
counter ++;
curPtr = curPtr->getNext();
} // end while
return frequency;
} // end getFrequencyOf

```

Question 4 Suppose that the ADT bag had an operation that displayed its contents. Write a C++ definition for such a method for the class LinkedBag.

Question 5 How many assignment operations does the method that you wrote for the previous question require?

The method contains. The discussion in Chapter 3 about the method contains applies here as well. Although the method could call the method getFrequencyOf, which we just defined, doing so usually will involve more work than is necessary. Whereas getFrequencyOf must check every entry in the bag, contains exits as soon as it finds an entry in the bag that is the same as the given one.

We observed in the previous chapter that the methods contains and remove perform the same search for a specific entry. Thus, to avoid duplicate code, we perform this search in a private method that both contains and remove can call. We declared this private method in the header file, and its definition follows.
```

// Returns either a pointer to the node containing a given entry
// or the null pointer if the entry is not in the bag.
template<class ItemType>
Node<ItemType>* LinkedBag<ItemType>::
getPointerTo(const ItemType\& target) const
{
bool found = false;
Node<ItemType>* curPtr = headPtr;

```
```

        while (!found && (curPtr != nul1ptr))
        {
        if (target == curPtr->getItem())
            found = true;
        else
            curPtr = curPtr->getNext();
        } // end while
        return curPtr;
    } // end getPointerTo
The definition of the method contains is straightforward:
template<class ItemType>
boo1 LinkedBag<ItemType>::contains(const ItemType\& anEntry) const
{
return (getPointerTo(anEntry) != nul7ptr);
} // end contains

```

\section*{? Question 6 If the pointer variable curPtr becomes nullptr in the method getPointerTo, what value does the method contains return when the bag is not empty?}

Question 7 Trace the execution of the method contains when the bag is empty.
Question 8 Revise the definition of the method getPointerTo so that the loop is controlled by a counter and the value of \(i\) temCount.
Question 9 What is a disadvantage of the definition of the method getPointerTo, as described in the previous question, when compared to its original definition?

Question 10 Why should the method getPointerTo not be made public?

The method remove. Recall that the method remove deletes one occurrence of a given entry and returns either true or false to indicate whether the removal was successful. Just as adding a new node to a linked chain is easiest at its beginning, so is removing the first node. But the entry that we need to remove is not always in the chain's first node.

Suppose that we locate the entry to delete in node \(n\). We can replace that entry with the entry in the first node, and then delete the first node. Thus, we can describe the logic for remove with the following pseudocode:
```

remove(anEntry)

```

Find the node that contains anEntry
Replace anEntry with the entry that is in the first node
Delete the first node
By using the private method getPointerTo to locate the entry that we want to delete, we can define the method remove as follows:
```

template<class ItemType>
boo1 LinkedBag<ItemType>::remove(const ItemType\& anEntry)
{
Node<ItemType>* entryNodePtr = getPointerTo(anEntry);
bool canRemoveItem = !isEmpty() \&\& (entryNodePtr != nul1ptr);
if (canRemoveItem)

```
```

    {
        // Copy data from first node to located node
        entryNodePtr->setItem(headPtr->getItem());
    // DeTete first node
    Node<ItemType>* nodeToDe1etePtr = headPtr;
    headPtr = headPtr->getNext();
    // Return node to the system
    nodeToDe1etePtr->setNext(nul1ptr);
    delete nodeToDeletePtr;
    nodeToDeletePtr = nul1ptr;
    itemCount--;
    } // end if
return canRemoveItem;
} // end remove

```

After the method remove deletes a node, the system can use this returned memory and possibly even reallocate it to your program as a result of the new operator. Suppose that this reallocation actually occurs when you ask for a new node for your linked chain. You can be sure that your new node does not still point to your linked chain, because you executed the statement nodeToDe1etePtr->setNext (nul1ptr) before you deallocated the node. Doing this and setting the variable nodeToDeletePtr to nullptr are examples of defensive programming that can avoid devastating, subtle errors later in the program. We take these steps, even though nodeToDeletePtr is a local variable that we do not use again, to clarify our intent to future programmers who revise our method.

Programming Tip: Remember that any time you allocate memory by using new, you must eventually deallocate it by using delete.

Note: For a pointer \(p\), delete \(p\) deallocates the node to which \(p\) points; it does not deallocate \(p\). The pointer \(p\) still exists, but it contains an undefined value. You should not reference \(p\) or any other pointer variable that still points to the deallocated node. To help you avoid this kind of error, you can assign nul1ptr to \(p\) after executing de1ete p. However, if variables other than \(p\) point to the deallocated node, the possibility of error still exists.

Question 11 Given the previous definition of the method remove, which entry in a bag can be deleted in the least time? Why?
Question 12 Given the previous definition of the method remove, which entry in a bag takes the most time to delete? Why?

The method clear. The method clear cannot simply set ItemCount to zero, thereby ignoring all of the entries in the linked chain. Because the nodes in the chain were allocated dynamically, clear must deallocate them. Thus, we have the following definition for this method:
```

template<class ItemType>
void LinkedBag<ItemType>::clear()

```

You must write a destructor if your class allocates memory dynamically

Situations that invoke the copy constructor

A compilergenerated copy constructor performs a shallow copy
```

{
while (headPtr != nul1ptr)
{
Node<ItemType>* nodeToDeletePtr = headPtr;
headPtr = headPtr->getNext();
// Return node to the system
nodeToDeletePtr->setNext(nul1ptr);
delete nodeToDeletePtr;
} // end while
// headPtr is nul1ptr
nodeToDe1etePtr = nul1ptr;
itemCount = 0;
} // end clear

```

The destructor. Each class has only one destructor. The destructor destroys an instance of the class, that is, an object, when the object's lifetime ends. Typically, the destructor is invoked implicitly at the end of the block in which the object was created.

Classes that use only statically allocated memory can depend on the compiler-generated destructor, as was the case for the class ArrayBag in Chapter 3. However, when a class uses dynamically allocated memory, as in the present link-based implementation, you need to write a destructor that deallocates this memory by using delete. The destructor for LinkedBag can simply call the method clear, as it uses delete to deallocate each node in the linked chain containing the bag's entries. The destructor's definition follows:
```

temp1ate<class ItemType>
LinkedBag<ItemType>::~LinkedBag()
{
clear();
} // end destructor

```

A destructor's name is a tilde \((\sim)\) followed by the class name. A destructor cannot have arguments, has no return type-not even void—and cannot use return to return a value.

Question 13 Revise the destructor in the class LinkedBag so that it does not call clear, but it instead directly deletes each node of the underlying linked chain.

The copy constructor. The second constructor in LinkedBag is the copy constructor:
LinkedBag (const LinkedBag<ItemType>\& aBag);
The copy constructor makes a copy of an object. It is invoked implicitly when you either pass an object to a function by value, return an object from a valued function, or define and initialize an object, as in

LinkedBag bag2 (bag1) ;
where bag1 exists already.
When copying an object involves only copying the values of its data members, the copy is called a shallow copy. If a shallow copy is sufficient, you can omit the copy constructor, in which case the compiler generates a copy constructor that performs a shallow copy. Such was the case in Chapter 3 for the class ArrayBag, although we did not mention it. That array-based implementation of the ADT bag used a compiler-generated copy constructor to copy both the array of bag items and the number of items.

FIGURE 4-8 (a) A linked chain and its shallow copy; (b) a linked chain and its deep copy


For our new link-based implementation, a compiler-generated copy constructor would copy only the data members itemCount and headPtr. For example, Figure \(4-8\) a pictures a linked chain and the result of this shallow copy. Both the original pointer headPtr and its copy point to the same linked chain. In other words, the chain's nodes are not copied. If you need to create a copy of the linked chain, you must write your own copy constructor. That is, a deep copy is needed, as Figure 4-8b illustrates.

Thus, the copy constructor appears as follows.
```

template<class ItemType>
LinkedBag<ItemType>::LinkedBag(const LinkedBag<ItemType>\& aBag)
{
itemCount = aBag->itemCount;
Node<ItemType>* origChainPtr = aBag->headPtr
if (origChainPtr == nul1ptr)
headPtr = nul1ptr; // Original bag is empty; so is copy
else
{
// Copy first node
headPtr = new Node<ItemType>();
headPtr->setItem(origChainPtr ->getItem());
// Copy remaining nodes
Node<ItemType>* newChainPtr = headPtr; // Last-node pointer
while (origPtr != nul1ptr)
{
origChainPtr = origChainPtr ->getNext(); // Advance pointer
// Get next item from original chain
ItemType nextItem = origChainPtr->getItem();
// Create a new node containing the next item
Node<ItemType>* newNodePtr = new Node<ItemType>(nextItem);

```
```

            // Link new node to end of new chain
            newChainPtr->setNext(newNodePtr);
            // Advance pointer to new last node
            newChainPtr = newChainPtr->getNext();
        } // end while
        newChainPtr->setNext(nul1ptr); // Flag end of new chain
    } // end if
    } // end copy constructor

```

As you can see, the copy constructor is an expensive operation. It requires traversing the original linked chain and duplicating each node visited.

\subsection*{4.3 Using Recursion in Link-Based Implementations}

It is possible, and sometimes desirable, to process linked chains recursively. This section examines how to write the iterative methods given previously as recursive ones. Such recursive methods will require the chain's head pointer as an argument. Therefore, they should not be public, because the head pointer is a private data member of the class and the client does not - and should not-have access to it. Otherwise, clients could access the linked nodes directly, thereby violating the ADT's wall.

\subsection*{4.3.1 Recursive Definitions of Methods in LinkedBag}

As an introduction to this topic, we will revise two methods in the class LinkedBag to use recursion. These methods simply traverse a chain of linked nodes without making any changes to it.

The method toVector. We begin with the method toVector because it has a straightforward recursive implementation. This operation requires us to traverse the linked chain of nodes as we copy data from the nodes into a vector. Traversal of a linked chain is an operation that occurs in many situations.

We have established that the method performing the recursion must be private and must have the head pointer as a parameter. Since the method will copy data into a vector as it traverses the linked chain, the vector must also be a parameter. We can declare this method in the private section of the class LinkedBag, as follows:
```

// Fills the vector bagContents with the data in the nodes of
// the linked chain to which curPtr points.
void fillVector(vector<ItemType>\& bagContents, Node<ItemType>* curPtr) const

```

Given a definition of this method, we could implement toVector as follows:
```

template<class ItemType>
vector<ItemType> LinkedBag<ItemType>::toVector() const
{
vector<ItemType> bagContents;
fillVector(bagContents, headPtr);
return bagContents;
} // end toVector

```

Thus, after creating a vector, toVector fills it-by calling fillVector—with the data in the chain of linked nodes whose head pointer is headPtr. Finally, toVector returns the vector to the client.

To define fillVector, we need to think recursively. If the chain is empty-that is, if curPtr is nullptr-we have nothing to do. This is the base case. In the recursive step, we first add the data curPtr->getItem() to the vector and then recursively fill the vector with the chain that begins at curPtr->getNext(). The method then has the following definition:
```

{
if (curPtr != nul1ptr)
{
bagContents.push_back(curPtr->getItem());
fil1Vector(bagContents, curPtr->getNext());
} // end if
} // end toVector

```

Even though fillvector has a reference parameter, the method is safe because it is private. A similar situation occurs in Chapter 16 and is discussed further in a Note in Section 16.3.1.

The private method getPointerTo. The method getPointerTo locates a given entry within the linked chain. To do so, it traverses the linked chain, but unlike fillVector's traversal, this traversal stops if it locates the node that contains the given entry. The iterative version of getPointerTo has one parameter, which is the given entry. We could define a private method to perform the recursionas we did for toVector-that has two parameters; the given entry and a head pointer. Then getPointerTo could call this new private method. However, since getPointerTo is itself a private method, we can revise it to have the necessary two parameters. That is the approach we will take here. Of course, we also will have to revise the calls to the method that appear in the rest of the class.

We begin by replacing the declaration of getPointerTo in the header file for the class LinkedBag with the following statements:
```

// Locates a given entry within this bag.
// Returns either a pointer to the node containing a given entry or
// the nul7 pointer if the entry is not in the bag.
Node<ItemType>* getPointerTo(const ItemType\& target,
Node<ItemType>* curPtr) const;

```

Before we forget to do so, let's revise the calls to this method that occur in the methods remove and contains. The original calls are getPointerTo(anEntry), but they now each need to be getPointerTo(anEntry, headPtr).

The recursive definition of this method has two base cases. One case occurs when the chain is empty, causing the method to return nul1ptr. The other base case occurs when we locate the desired entry at curPtr->getItem(). In this case, the method returns curPtr. The recursive step searches the chain beginning at curPtr->getNext (). Thus, the method has the following definition:
```

temp7ate<class ItemType>
Node<ItemType>* getPointerTo(const ItemType\& target,
Node<ItemType>* curPtr) const
{
Node<ItemType>* result = nul1ptr;
if (curPtr != nul1ptr)
{
if (target== curPtr->getItem())
result = curPtr;
e1se
result = getPointerTo(target, curPtr->getNext());
} // end if
return result;
} // end getPointerTo

```

Other methods that can have a recursive definition are left for you as exercises.

Question 14 Revise the method clear so that it calls a recursive method to deallocate the nodes in the chain.

\subsection*{4.4 Testing Multiple ADT Implementations}

In Chapter 3, we developed a short program, shown in Listing 3-2, that tested the core methods of our array-based bag implementation. Because we used ADT bag methods when we tested our implementation, we can use the same code-with a few changes-to test our linked implementation. You need only to change each occurrence of ArrayBag to LinkedBag and recompile the program.

But what if later you change the array-based implementation and want to retest it? You would need to change each occurrence of LinkedBag back to ArrayBag. Let's examine a way we can revise the test program to let the user decide which bag implementation needs testing. By taking advantage of polymorphism, we can do this without needing to duplicate the test code.

In Chapter 1, we described our ADT bag in the abstract class BagInterface. Since both ArrayBag and LinkedBag share public methods defined in that class, we made each a subclass of BagInterface. In our test program, if we declare the variable bagPtr as a pointer to an object that implements the methods in BagInterface, we can use an instance of either ArrayBag or LinkedBag as the object bagPtr references. We declare bagPtr as
```

BagInterface<string>* bagPtr;

```

We then ask the user which implementation to test, and if we assign either ' A ' or ' L ' to the char variable userChoice, we can instantiate an instance of the requested bag type and run the test by writing the following code:
```

if (userChoice == 'A')
{
bagPtr = new ArrayBag<string>();
cout << "Testing the Array-Based Bag:" << endl;
}
else
{
bagPtr = new LinkedBag<string>();
cout << "Testing the Link-Based Bag:" << end1;
} // end if
cout << "The initial bag is empty." << endl;
bagTester(bagPtr);
delete bagPtr;
bagPtr = nul1ptr;

```

To accommodate the data type of bagPtr, we need to change the parameter lists of the functions in our test program as follows:
```

void displayBag(BagInterface<string>* bagPtr)

```
and
```

void bagTester(BagInterface<string>* bagPtr)

```

Finally, we must change the notation used to call methods on the instances of bagPtr, since it is a pointer. For example,
bag.isEmpty ()
must be changed to
bagPtr->isEmpty()
Listing 4-4 is a complete listing of the modified test program for core methods of classes that are implementations of the ADT bag.

\section*{LISTING 4-4 A program that tests the core methods of classes that are derived from the abstract class BagInterface}
```

\#include "BagInterface.h"
\#include "ArrayBag.h"
\#include "LinkedBag.h"
\#include <iostream>
\#include <string>
\#include <cctype>

```
using namespace std;
void displayBag(BagInterface<string>* bagPtr)
\{
    cout << "The bag contains " << bagPtr->getCurrentSize()
        << " items:" << end7;
    vector<string> bagItems;
    bagPtr->toVector(bagItems);
    int numberOfEntries = bagItems.size();
    for (int \(\mathbf{i}=0\); \(\mathbf{i}\) < numberOfEntries; \(\mathbf{i + +}\) )
    \{
        cout << bagItems[i] << " ";
    \} \(/ /\) end for
    cout << endl << endl;
\} // end displayBag
void bagTester (BagInterface<string>* bagPtr)
\{
    cout << "isEmpty: returns " << bagPtr->isEmpty()
        << "; should be 1 (true)" << end7;
    string items[] = \{"one", "two", "three", "four", "five", "one"\};
    cout << "Add 6 items to the bag: " << end1;
    for (int i = 0; i < 6; i++)
    \{
        bagPtr->add(items[i]);
    \} // end for
    displayBag(bagPtr);
    cout << "isEmpty: returns " << bagPtr->isEmpty()
        << "; should be 0 (false)" << endl;
    cout << "getCurrentSize returns : " << bagPtr->getCurrentSize()
        << "; should be 6" << endl;
    cout << "Try to add another entry: add(\"extra\") returns "
        << bagPtr->add("extra") << endl;
\} // end bagTester
int main()
\{
    BagInterface<string>* bagPtr = nul1ptr;
    char userChoice;
    cout << "Enter 'A' to test the array-based implementation \(\backslash n\) ";
        << " or 'L' to test the link-based implementation: ";
    cin >> userChoice;
```

    if (toupper(userChoice) == 'A')
    {
    bagPtr = new ArrayBag<string>();
    cout << "Testing the Array-Based Bag:" << endl;
    }
    else
    {
    bagPtr = new LinkedBag<string>();
    cout << "Testing the Link-Based Bag:" << end7;
    } // end if
cout << "The initial bag is empty." << endl;
bagTester(bagPtr);
delete bagPtr;
bagPtr = nul1ptr;
cout << "Al1 done!" << endl;
return 0;
} // end main

```

\section*{Sample Output 1}

Enter 'A' to test the array-based implementation
or 'L' to test the link-based implementation: A
Testing the Array-Based Bag:
The initial bag is empty.
isEmpty: returns 1 ; should be 1 (true)
Add 6 items to the bag:
The bag contains 6 items:
one two three four five one
isEmpty: returns 0 ; should be 0 (false)
getCurrentSize returns : 6; should be 6
Try to add another entry: add("extra") returns 0
Al1 done!

\section*{Sample Output 2}

Enter 'A' to test the array-based implementation
or 'L' to test the link-based implementation: L
Testing the Link-Based Bag:
The initial bag is empty.
isEmpty: returns 1 ; should be 1 (true)
Add 6 items to the bag:
The bag contains 6 items:
one five four three two one
isEmpty: returns 0 ; should be 0 (false)
getCurrentSize returns : 6; should be 6
Try to add another entry: add("extra") returns 1
All done!

Question 15 Revise the program in Listing 4-4 so that it tests first the array-based implementation and then the link-based implementation. Ensure that the program does not have a memory leak.

\subsection*{4.5 Comparing Array-Based and Link-Based Implementations}

Typically, the various implementations that a programmer contemplates for a particular ADT have advantages and disadvantages. When you must select an implementation, you should weigh these advantages and disadvantages before you make your choice. As you will see, the decision among possible implementations of an ADT is one that you must make often. As an example of how you should proceed in general, we compare the two implementations of the ADT bag that you have seen.

The array-based implementation that you saw in Chapter 3 appears to be a reasonable approach. Arrays are easy to use, but as was already mentioned, an array has a fixed size, and it is possible for the number of items in the bag to exceed this fixed size. In practice, when choosing among implementations of an ADT, you must ask the question of whether the fixed-size restriction of an array-based implementation presents a problem in the context of a particular application. The answer to this question depends on two factors. One obvious factor is whether, for a given application, you can predict in advance the maximum number of items in the ADT at any one time. If you cannot, it is quite possible that an operation - and hence the entire program-will fail because the ADT in the context of a particular application requires more storage than the array can provide.

On the other hand, if for a given application you can predict in advance the maximum number of items in the ADT at any one time, you must explore a more subtle factor: Would you waste storage by declaring an array to be large enough to accommodate this maximum number of items? Consider a case in which the maximum number of items is large, but you suspect that this number rarely will be reached. For example, suppose that your bag could contain as many as 10,000 items, but the actual number of items in the bag rarely exceeds 50 . If you declare 10,000 array locations at compilation time, at least 9,950 of them will be wasted most of the time. In both of the previous cases, the arraybased implementation given in Chapter 3 is not desirable.

What if you use a dynamically allocated array? Because you would use the new operator to allocate storage dynamically, you will be able to provide as much storage as the bag needs (within the bounds of the particular computer, of course). Thus, you do not have to predict the maximum size of the bag. However, if you double the size of the array each time you reach the end of the array-which is a reasonable approach to enlarging the array-you still might have many unused array locations. In the example just given, you could allocate an array of 50 locations initially. If you actually have 10,000 items in your bag, array doubling will eventually give you an array of 12,800 locations, which is 2,800 more than you need. Remember also that you waste time by copying the array and then deallocating it each time you need more space.

Now suppose that your bag will never contain more than 25 items. You could allocate enough storage in the array for the bag and know that you would waste little storage when the bag contains only a few items. With respect to its size, an array-based implementation is perfectly acceptable in this case.

A link-based implementation can solve any difficulties related to the fixed size of an array-based implementation. You use the new operator to allocate storage dynamically, so you do not need to predict the maximum size of the bag. Because you allocate memory one item at a time, the bag will be allocated only as much storage as it needs. Thus, you will not waste storage.

Array-based and link-based implementations have other differences, and these differences affect both the time and memory requirements of the implementations. Any time you store a collection of data in an array or a linked chain, the data items become ordered; that is, there is a first item, a second item, and so on. This order implies that a typical item has a predecessor and a successor. In an array anArray, the location of the next item after the item in anArray [ i\(]\) is implicit-it is in anArray \([i+1]\). In a linked chain, however, you explicitly determine the location of the next item by using the pointer in the current node. This notion of an implicit versus explicit next item is one of the primary differences between an array and a linked chain. Therefore, an advantage of an array-based implementation

Arrays are easy to use, but they have a fixed size

Can you predict the maximum number of items in the ADT?

Will an array waste storage?

Increasing the size of a dynamically allocated array can waste storage and time

An array-based implementation is a good choice for a small bag

Linked chains do not have a fixed size

The item after an array item is implied; in a chain of linked nodes, an item points explicitly to the next item

An array-based implementation requires less memory than a link-based implementation

You can access array items directly with equal access time

You must traverse a linked chain to access its \(i^{\text {th }}\) node

The time to access
the \(i^{\text {th }}\) node in a linked chain depends on \(i\)
is that it does not have to store explicit information about where to find the next data item, thus requiring less memory than a link-based implementation.

Another, more important advantage of an array-based implementation is that it can provide direct access to a specified item. For example, if you use the array items to implement the ADT bag, you know that accessing either items [0] or items [49] takes the same amount of time. That is, the access time is constant for an array.

On the other hand, if you use a link-based implementation, you have no way of immediately accessing the node that contains the \(i\) th item. To get to the appropriate node, you use the next pointers to traverse the linked chain from its beginning until you reach the \(i\) th node. That is, you access the first node and get the pointer to the second node, access the second node and get the pointer to the third node, and so on until you finally access the \(i\) th node. Clearly, the time it takes you to access the first node is less than the time it takes to access the 50th node. The access time for the \(i\) th node depends on \(i\). The type of implementation chosen will affect the efficiency of the ADT's operations.

We will continue to compare various solutions to a problem throughout this book. Chapter 10 introduces a more formal way to discuss the efficiency of algorithms. Until then, our discussions will be informal.

\section*{Summary}
1. You can link objects-called nodes - to one another to form a chain of linked data. Each node contains a data item and a pointer to the next node in the chain. An external pointer variable-called the head pointer-points to the first node. The last node in the chain has nul1ptr in its pointer portion, so it points to no other node.
2. You use the new operator to dynamically allocate a new node, whereas you use the de1ete operator to deallocate a node.
3. Inserting a new node at the beginning of a linked chain or deleting the first node of a linked chain are easier to perform than insertions and deletions anywhere else in the chain. The insertion requires a change to two pointers: the pointer within the new node and the head pointer. The deletion requires a change to the head pointer and an application of the delete operator to the removed node.
4. Unlike an array, which enables you direct access to any of its elements, a linked chain requires a traversal to access a particular node. Therefore, the access time for an array is constant, whereas the access time for a linked chain depends on the location of the node within the chain.
5. When traversing a linked chain by using the pointer variable curPtr, you must be careful not to reference curPtr after it has "passed" the last node in the chain, because it will have the value nu11ptr at that point. For example, the loop
```

while (value > curPtr->getItem())
curPtr = curPtr->getNext();

```
is incorrect if value is greater than all the data values in the linked chain, because curPtr becomes nul1ptr. Instead you should write
```

while ((curPtr != nul1ptr) \&\& (value > curPtr->getItem()))
curPtr = curPtr->getNext();

```

Because C++ uses short-circuit evaluation (see Appendix A) of logical expressions, if curPtr becomes nul1ptr, the expression curPtr->getItem() will not be evaluated.
6. A class that allocates memory dynamically needs an explicit copy constructor that copies an instance of the class. The copy constructor is invoked implicitly when you pass an object to a function by value, return an object from a valued function, or define and initialize an object. If you do not define a copy constructor, the compiler will generate one for you. A compiler-generated copy constructor is sufficient only for classes that use statically allocated memory.
7. A class that allocates memory dynamically needs an explicit destructor. The destructor should use delete to deallocate the memory associated with the object. If you do not define a destructor, the compiler will generate one for you. A compiler-generated destructor is sufficient only for classes that use statically allocated memory.
8. Although you can use the new operator to allocate memory dynamically for either an array or a linked chain, you can increase the size of a linked chain one node at a time more efficiently than you can increase the size of an array. When you increase the size of a dynamically allocated array, you must copy the original array entries into the new array and then deallocate the original array.

\section*{Exercises}
1. If headPtr is a pointer variable that points to the first node of a linked chain of at least two nodes, write C++ statements that delete the second node and return it to the system.
2. Revise the public method add in the class LinkedBag so that the new node is inserted at the end of the linked chain.
3. Suppose that the class LinkedBag did not have the data member itemCount. Revise the method getCurrentSize so that it counts the number of nodes in the linked chain
a. Iteratively
b. Recursively
4. Revise the public method getFrequencyOf in the class LinkedBag so that it is recursive.
5. Add a constructor to the class LinkedBag that creates a bag from a given array of entries.
6. Specify and define a method for LinkedBag that removes a random entry from the bag.
7. Compare the number of operations required to display the data in each of the \(n\) nodes in a linked chain with the number of operations required to display each of the \(n\) items in an array.
8. Compare the number of operations required to display the data in the \(n\)th node in a linked chain with the number of operations required to display the \(n\)th item in an array.
9. Compare the array-based and link-based implementations of the ADT bag operation remove (anEntry). Consider the various locations of anEntry within the array or chain.
10. In a doubly linked chain, each node can point to the previous node as well as to the next node. Figure 4-9 shows a doubly linked chain and its head pointer. Define a class to represent a node in a doubly linked chain.
11. List the steps necessary to add a node to the beginning of the doubly linked chain shown in Figure 4-9.
12. List the steps necessary to remove the first node from the beginning of the doubly linked chain shown in Figure 4-9.

\section*{FIGURE 4-9 A doubly linked chain}


\section*{Programming Problems}
1. Add the methods union, intersection, and difference to the class ArrayBag. Exercises 6, 7, and 8 of Chapter 1 describe these operations.
2. Implement the ADT set that you specified in Programming Problem 5 of Chapter 1 by using a linked chain.
3. Implement the ADT pile that you specified in Programming Problem 6 of Chapter 1 by using a linked chain.
4. Implement the ADT polynomial that Exercise 9 in Chapter 1 describes by using a linked chain.
5. Consider a sparse implementation of the ADT polynomial that stores only the terms with nonzero coefficients. For example, you can represent the revised polynomial \(p\) in Exercise 9 of Chapter 1 with the linked chain shown in Figure 4-10.
a. Complete the sparse implementation.
b. Define a traverse operation for the ADT polynomial that will allow you to add two sparse polynomials without having to consider terms with zero coefficients explicitly.
6. Define a class DoublyLinkedBag that implements the ADT bag by using a doubly linked chain, as shown in Figure 4-9. Use the class of nodes that Exercise 10 defines.
7. Use the classes for a set or a bag, as defined in this chapter or described in the previous projects, to create a spell checker. Consult the details given in Programming Problems 13 and 14 of Chapter 3.
8. Specify and implement an ADT character string by using a linked chain of characters. Include typical operations such as finding its length, appending one string to another, finding the index of the leftmost occurrence of a character in a string, and testing whether one string is a substring of another.

FIGURE 4-10 A sparse polynomial

9. When you play a board or card game or when you use a shared computing resource, you get a turn and then wait until everyone else has had a turn. Although the number of players in a game remains relatively static, the number of users of a shared computing service fluctuates. Let's assume that this fluctuation will occur.

Design an ADT that keeps track of turns within a group of people. You should be able to add or delete people and determine whose turn occurs now.

Begin with a given group of people; assign these people an initial order. (This order can be random or specified by the user.) The first new person joining the group should get a turn after all others have had an equal number of turns. Each subsequent new person should get a turn after the person who joined the group most recently has had a turn.

Also design an ADT to represent a person. (You can be conservative with the amount of data that this ADT contains.) The data that your first ADT stores is made up of instances of the ADT person.

Implement your ADTs as \(\mathrm{C}++\) classes. Write a program that uses-and therefore tests-your ADTs completely. Your program should process several insertion and deletion operations, and demonstrate that people are given turns correctly.
10. Occasionally, a link-based structure that does not use pointers is useful. One such structure uses an array whose items are "linked" by array indexes. Figure 4-11a illustrates an array of nodes that represents the linked chain whose first node contains the string " B ", second node contains "E", and third node contains "J". Each node has two members, item and next. The next member is an integer index to the array element that contains the next node in the linked chain. Note that the next member of the last node contains -1 . The integer variable head contains the index of the first node in the chain.

The array elements that currently are not a part of the linked chain make up a free list of available nodes. These nodes form another linked chain, with the integer variable free containing the index of the first free node. To insert an item into the beginning of the original linked chain, you take a free node from the beginning of the free list and insert it into the linked chain (Figure 4-11b). When you delete an item from the linked chain, you insert the node into the beginning of the free list (Figure 4-11c). In this way, you can avoid shifting data items.

Implement the ADT bag by using this array-based linked chain.

FIGURE 4-11 (a) An array-based implementation of a linked chain; (b) after inserting " \(D\) " at the beginning of the chain; (c) after deleting " B "


\section*{Recursion as a Problem-Solving Technique}

\section*{Chapter} 5

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Prerequisites
Chapter 2 Recursion: The Mirrors

Chapter 2 presented the basic concepts of recursion, and now this chapter moves on to some extremely useful and somewhat complex applications in computer science. The recursive solutions to the problems you will see are far more elegant and concise than the best of their nonrecursive counterparts.

This chapter introduces two new concepts: formal grammars and backtracking. Formal grammars enable you to define, for example, syntactically correct algebraic

A grammar states the rules for forming the strings in a language

Symbols that grammars use
expressions, which we explore in some detail. Backtracking is a problem-solving technique that involves guesses at a solution. The chapter concludes with a discussion of the close relationship between recursion and mathematical induction; you will learn how to use mathematical induction to study properties of algorithms.

More applications of recursion appear in subsequent chapters.

\subsection*{5.1 Defining Languages}

English and C++ are two languages with which you are familiar. A language is nothing more than a set of strings of symbols from a finite alphabet. For example, if you view a C++ program as one long string of characters, you can define the set of all syntactically correct C++ programs. This set is the language
\[
C++ \text { Programs }=\{\text { string } s: s \text { is a syntactically correct } \mathrm{C}++ \text { program }\}
\]

Notice that whereas all programs are strings, not all strings are programs. A C ++ compiler is a program that, among other things, sees whether a given string is a member of the language \(C++\) Programs; that is, the compiler determines whether the string is a syntactically correct \(\mathrm{C}++\) program. Of course, this definition of \(C++\) Programs is not descriptive enough to allow the construction of a compiler. The definition specifies a characteristic of the strings in the set \(C++\) Programs: The strings are syntactically correct \(\mathrm{C}++\) programs. However, this definition does not give the rules for determining whether a string is in the set; that is, the definition does not specify what is meant by a syntactically correct C++ program.

The word "language" does not necessarily mean a programming language or a communication language. For example, the set of algebraic expressions forms a language
\[
\text { AlgebraicExpressions }=\{\text { string } s: s \text { is an algebraic expression }\}
\]

The language AlgebraicExpressions is the set of strings that meets certain rules of syntax; however, the set's definition does not give these rules.

In both examples, the rules for forming a string within the language are missing. A grammar states the rules of a language. The grammars that you will see in this chapter are recursive in nature. One of the great benefits of using such a grammar to define a language is that you can often write a straightforward recursive algorithm, based on the grammar, that determines whether a given string is in the language. Such an algorithm is called a recognition algorithm for the language.

As it is a complex task to present a grammar for the set \(C++\) Programs, we will look instead at grammars for some simpler languages, including several common languages of algebraic expressions.

\subsection*{5.1.1 The Basics of Grammars}

A grammar uses several special symbols:
- \(x \mid y\) means \(x\) or \(y\).
- \(x y\) (and sometimes \(x \cdot y\) ) means \(x\) followed by \(y\).
- <word > means any instance of word, where word is a symbol that must be defined elsewhere in the grammar.

A grammar for the language
\(C++\) Identifiers \(=\{\) string \(s: s\) is a legal \(\mathrm{C}++\) identifier \(\}\)

is simple, so we begin with it. As you know, a legal C++ identifier begins with a letter and is followed by zero or more letters and digits. In this context, the underscore ( \(\_\)) is a letter. One way to represent this definition of an identifier is with a syntax diagram, as shown in Figure 5-1.

A syntax diagram is convenient for people to use, but a grammar is a better starting point if you want to write a function that will recognize an identifier. A grammar for the language \(C++\) Identifiers is
\[
\begin{aligned}
& <\text { identifier }>=<\text { letter }>\mid<\text { identifier }><\text { letter }>\mid<\text { identifier }><\text { digit }> \\
& <\text { letter }>=\mathrm{a}|\mathrm{~b}| \ldots .\left.\mathrm{z}|\mathrm{~A}| \mathrm{B}|\ldots| \mathrm{Z}\right|_{-} \\
& <\text {digit }>=0|1| \ldots \mid 9
\end{aligned}
\]

The definition reads as follows:
An identifier is a letter, or an identifier followed by a letter, or an identifier followed by a digit.

The most striking aspect of this definition is that identifier appears in its own definition: This grammar is recursive, as are many grammars.

Given a string \(s\), you can determine whether it is in the language \(C++\) Identifiers by using the grammar to construct the following recognition algorithm: If \(s\) is of length 1 , it is in the language if the character is a letter. (This statement is the base case, so to speak.) If \(s\) is of length greater than 1 , it is in the language if the last character of \(s\) is either a letter or a digit, and \(s\) minus its last character is an identifier.

The pseudocode for a recursive valued function that determines whether a string is in the language \(C++\) Identifiers follows:
```

// Returns true if s is a legal C++ identifier;
// otherwise returns false.
isId(s: string): boolean
if (s is of length 1)
// Base case
if (sis a letter)
return true
else
return false
else if (the last character ofs is a letter or a digit)
return isId(s minus its last character) // Point X
else
return false

```

A grammar for the language of C++ identifiers

Many grammars are recursive

Figure 5-2 contains a trace of this function for the string "A2B".

The initial call is made and the function begins execution.
```

s = "A2B"

```

At point X , a recursive call is made and the new invocation of isId begins execution:


At point X , a recursive call is made and the new invocation of isId begins execution:


This is the base case, so this invocation of isId completes:


The value is returned to the calling function, which completes execution:


The value is returned to the calling function, which completes execution:


\subsection*{5.1.2 Two Simple Languages}

Now consider two more simple examples of languages, their grammars, and resulting recognition algorithms.

Palindromes. A palindrome is a string that reads the same from left to right as it does from right to left. For example, "radar" and "deed" are both palindromes. You can define the language of palindromes as follows:

Palindromes \(=\{\) string \(s: s\) reads the same left to right as right to left \(\}\)
How can you use a grammar to define the language Palindromes? You need to devise a rule that allows you to determine whether a given string \(s\) is a palindrome. In the spirit of recursive definitions, you should state this rule in terms of determining whether a smaller string is a palindrome. Your first instinct might be to choose \(s\) minus its last (or first) character for the smaller string. However, this approach does not work, because there is no relationship between the statements
\(s\) is a palindrome
and
\(s\) minus its last character is a palindrome
That is, \(s\) might be a palindrome, although \(s\) minus its last character is not, as is the case for "deed." Similarly, \(s\) minus its last character might be a palindrome, although \(s\) is not, as is the case for "deeds."

A little thought reveals that you must consider characters in pairs: There is a relationship between the statements
\(s\) is a palindrome
and
\(s\) minus its first and last characters is a palindrome
Specifically, \(s\) is a palindrome if and only if
- The first and last characters of \(s\) are the same
and
- \(s\) minus its first and last characters is a palindrome

You need a base case that you will reach after stripping away enough pairs of characters. If \(s\) has an even number of characters, you will eventually be left with two characters, and then, after you strip away another pair, you will be left with zero characters. A string of length 0 is called the empty string and is a palindrome. If \(s\) has an odd number of characters, you will eventually be left with one character, after which you cannot strip away another pair. Hence, you must have a second base case: A string of length 1 is a palindrome.

This discussion leads to the following grammar for the language of palindromes:
\[
\begin{aligned}
& <\text { pal }>=\text { empty string }|<c h>| \mathrm{a}<\text { pal }>\mathrm{a} \mid \mathrm{b}<\text { pal }>\mathrm{b}|\ldots| \mathrm{Z}<\text { pal }>\mathrm{Z} \\
& <\text { ch }>\mathrm{a}|\mathrm{~b}| \ldots|\mathrm{z}| \mathrm{A}|\mathrm{~B}| \ldots \mid \mathrm{Z}
\end{aligned}
\]

Based on this grammar, you can construct a recursive valued function for recognizing palindromes. The pseudocode for such a function follows:
```

// Returns true if the string s of letters is a palindrome; otherwise returns false.
isPalindrome(s: string): boolean
if (s is the empty string or s is of length 1)
return true
e1se if (s's first and last characters are the same letter)
return isPalindrome(s minus its first and last characters)
else
return false

```

Strings of the form \(\mathbf{A}^{n} \mathbf{B}^{n}\). The notation \(\mathrm{A}^{n} \mathrm{~B}^{n}\) represents the string that consists of \(n\) consecutive \(A\) 's followed by \(n\) consecutive \(B\) 's. Another simple language consists of such strings:
\[
A n B n=\left\{\text { string } s: s \text { is of the form } \mathrm{A}^{n} \mathrm{~B}^{n} \text { for some } n \geq 0\right\}
\]

The grammar for this language is actually very similar to the grammar for palindromes. You must strip away both the first and last characters and check to confirm that the first character is an \(A\) and the last character is a \(B\). Thus, the grammar for the language \(A n B n\) is
\[
<\text { legal_word }>=\text { empty string } \mid \mathrm{A}<\text { legal_word }>\mathrm{B}
\]

The pseudocode for a recognition function for this language follows:
```

// Returns true if s is of the form A" B}\mp@subsup{B}{}{n}\mathrm{ ; otherwise returns false.
isAnBn(s: string): boolean
if (the length of s is zero)
return true
else if (s begins with the character A and ends with the character B)
return isAnBn(s minus its first and last characters)
else
return false

```

A recursive description of a palindrome

Strings of length 0 or 1 are the base cases

A grammar for the language of palindromes

A recognition algorithm for palindromes

A grammar for the language of strings \(\mathrm{A}^{n} \mathrm{~B}^{n}\)

A recognition algorithm for strings \(\mathrm{A}^{n} \mathrm{~B}^{n}\)

内
Note: Grammars, like recursive algorithms, must have carefully chosen base cases. You must ensure that, when a string is decomposed far enough, it will always reach the form of one of the grammar's base cases.

Question 1 Consider the language of these character strings: \$, cc\$d, cccc\$dd, cccccc\$ddd, and so on. Write a recursive grammar for this language.

Processing expressions

\subsection*{5.2 Algebraic Expressions}

One of the tasks a compiler must perform is to recognize and evaluate algebraic expressions. For example, consider the \(\mathrm{C}++\) assignment statement
\[
y=x+z *(w / k+z *(7 * 6)) ;
\]

A C++ compiler must determine whether the right side is a syntactically legal algebraic expression; if so, the compiler must then indicate how to compute the expression's value.

There are several common definitions for a "syntactically legal" algebraic expression. Some definitions force an expression to be fully parenthesized - that is, to have parentheses around each pair of operands together with their operator. Thus, you would have to write \(((a * b) * c)\) rather than \(\mathrm{a} * \mathrm{~b}\) * c. In general, the stricter a definition, the easier it is to recognize a syntactically legal expression. On the other hand, conforming to overly strict rules of syntax is an inconvenience for programmers.

This section presents three different languages for algebraic expressions. The expressions in these languages are easy to recognize and evaluate but are generally inconvenient to use. However, these languages provide us with good, nontrivial applications of grammars. We will see other languages of algebraic expressions whose members are difficult to recognize and evaluate but are convenient to use. To avoid unnecessary complications, assume that you have only the binary operators ,,\(+-{ }^{*}\), and / (no unary operators or exponentiation). Also, assume that all operands in the expression are single-letter identifiers.

\subsection*{5.2.1 Kinds of Algebraic Expressions}

The algebraic expressions you learned about in school are called infix expressions. The term "infix" indicates that every binary operator appears between its operands. For example, in the expression
\[
a+b
\]
the operator + is between its operands \(a\) and \(b\). This convention necessitates associativity rules, precedence rules, and the use of parentheses to avoid ambiguity. For example, the expression
\[
a+b * c
\]
is ambiguous. What is the second operand of the + ? Is it \(b\) or is it \(\left(b^{*} c\right)\) ? Similarly, the first operand of the * could be either \(b\) or \((a+b)\). The rule that * has higher precedence than + removes the ambiguity by specifying that \(b\) is the first operand of the * and that \(\left(b^{*} c\right)\) is the second operand of the + . If you want another interpretation, you must use parentheses:
\[
(a+b) * c
\]

Even with precedence rules, an expression like
\[
a / b * c
\]
is ambiguous. Typically, / and * have equal precedence, so you could interpret the expression either as \((a / b)^{*} c\) or as \(a /(b * c)\). The common practice is to associate from left to right, thus yielding the first interpretation.

Two alternatives to the traditional infix convention are prefix and postfix expressions. Under these conventions, an operator appears either before its operands (prefix) or after its operands (postfix). Thus, the infix expression
\[
a+b
\]
is written in prefix form as
\[
+a b
\]
and in postfix form as
\[
a b+
\]

To further illustrate the conventions, consider the two interpretations of the infix expression \(a+b * c\) just considered. You write the expression
\[
a+(b * c)
\]
in prefix form as
\[
+a * b c
\]

The + appears before its operands \(a\) and \(\left({ }^{*} b c\right)\), and the * appears before its operands \(b\) and \(c\). The same expression is written in postfix form as
\[
a b c^{*}+
\]

The * appears after its operands \(b\) and \(c\), and the + appears after its operands \(a\) and ( \(\left.b c^{*}\right)\).
Similarly, you write the expression
\[
(a+b) * c
\]
in prefix form as
\[
*+a b c
\]

The * appears before its operands \((+a b)\) and \(c\), and the + appears before its operands \(a\) and \(b\). The same expression is written in postfix form as
\[
a b+c *
\]

The + appears after its operands \(a\) and \(b\), and the * appears after its operands \((a b+\) ) and \(c\).
If the infix expression is fully parenthesized, converting it to either prefix or postfix form is straightforward. Because each operator then corresponds to a pair of parentheses, you simply move the operator to the position marked by either the open parenthesis "("-if you want to convert to prefix form-or the close parenthesis ")"-if you want to convert to postfix form. This position either precedes or follows the operands of the operator. All parentheses would then be removed.

For example, consider the fully parenthesized infix expression
\[
((a+b) * c)
\]

In a prefix expression, an operator precedes its operands

In a postfix expression, an operator follows its operands

Converting to prefix form

Converting to postfix form

To convert this expression to prefix form, you first move each operator to the position marked by its corresponding open parenthesis:
\[
\begin{aligned}
& ((a b) c) \\
& \downarrow \downarrow \\
& *+
\end{aligned}
\]

Next, you remove the parentheses to get the desired prefix expression:
\[
*+a b c
\]

Similarly, to convert the infix expression to postfix form, you move each operator to the position marked by its corresponding close parenthesis:
\[
\begin{gathered}
((a b) c) \\
\downarrow \downarrow \\
+*
\end{gathered}
\]

Then you remove the parentheses:
\[
a b+c^{*}
\]

When an infix expression is not fully parenthesized, these conversions are more complex. Chapter 6 discusses the general case of converting an infix expression to postfix form.

The advantage of prefix and postfix expressions is that they never need precedence rules, association rules, or parentheses. Therefore, the grammars for prefix and postfix expressions are quite simple. In addition, the algorithms that recognize and evaluate these expressions are relatively straightforward.

\subsection*{5.2.2 Prefix Expressions}

A grammar that defines the language of all prefix expressions is
\[
\begin{aligned}
& <\text { prefix }>=<\text { identifier }>\mid<\text { operator }><\text { prefix }><\text { prefix }> \\
& <\text { operator }>=+|-|*| / \\
& <\text { identifier }>=a|b| \ldots \mid z
\end{aligned}
\]

From this grammar, you can construct a recursive algorithm that recognizes whether a string is a prefix expression. If the string is of length 1 , it is a prefix expression if and only if the string is a single lowercase letter. Strings of length 1 can be the base case. If the length of the string is greater than 1 , then for it to be a legal prefix expression, it must be of the form
\[
<\text { operator }><\text { prefix }><\text { prefix }>
\]

Thus, the algorithm must check to see whether
- The first character of the string is an operator
and
- The remainder of the string consists of two consecutive prefix expressions

The first task is trivial, but the second is a bit tricky. How can you tell whether you are looking at two consecutive prefix expressions? A key observation is that if you add any string of nonblank characters to the end of a prefix expression, you will no longer have a prefix expression. That is, if \(E\) is
a prefix expression and \(Y\) is any nonempty string of nonblank characters, then \(E Y\) cannot be a prefix expression. This is a subtle point; Exercise 17 at the end of this chapter asks you to prove it.

Given this observation, you can begin to determine whether you have two consecutive prefix expressions by identifying a first prefix expression. If you find one, the previous observation implies that only one endpoint is possible for this first expression.

If you find that the first prefix expression ends at position end1, you then attempt to find a second prefix expression beginning at position end1 +1 . If you find the second expression, you must check whether you are at the end of the string in question.

By using these ideas, you can show, for example, that +* \(a b-c d\) is a prefix expression. For \(+* a b-c d\) to be a prefix expression, it must be of the form \(+E_{1} E_{2}\), where \(E_{1}\) and \(E_{2}\) are prefix expressions. Now you can write
\[
\begin{aligned}
& E_{1}=* E_{3} E_{4}, \text { where } \\
& E_{3}=a \\
& E_{4}=b
\end{aligned}
\]

Because \(E_{3}\) and \(E_{4}\) are prefix expressions, \(E_{1}\) is a prefix expression. Similarly, you can write
\[
\begin{aligned}
& E_{2}=-E_{5} E_{6}, \text { where } \\
& E_{5}=c \\
& E_{6}=d
\end{aligned}
\]
and see that \(E_{2}\) is a prefix expression.
You can write a function to test whether an expression is a prefix expression by first constructing a recursive valued function endPre (strExp, first) to examine the expression that begins at position first of the string strExp and to locate the end of the first prefix expression it finds. If successful, the function returns the index of the end of the prefix expression. If no such prefix expression exists, endPre returns -1 . The function appears in pseudocode as follows.
```

// Finds the end of a prefix expression, if one exists.
// Precondition:The substring of strExp from the index first through the end of
// the string contains no blank characters.
// Postcondition: Returns the index of the last character in the prefix expression that
// begins at index first of strExp, or -1 if no such prefix expression exists.
endPre(strExp: string, first: integer): integer
1ast = strExp.1ength() - 1
if (first < 0 or first > 1ast)
return -1
ch = character at position first of strExp
if (ch is an identifier)
return first // Index of last character in simple prefix expression
else if (ch is an operator)
{
// Find the end of the first prefix expression
firstEnd = endPre(strExp, first + 1) // Point X
// If the end of the first prefix expression was found, find the end of the second
// prefix expression
if (firstEnd > -1)
return endPre(strExp, firstEnd + 1) // Point Y
e1se
return -1
}
else
return -1

```

If \(E\) is a prefix expression, \(E Y\) cannot be
endPre determines the end of a prefix expression

Figure 5-3 contains a trace of endPre when the initial expression is +*ab-cd.

FIGURE 5-3 Trace of endPre("+*ab-cd", 0)

The initial call endPre("+*ab-cd", 0 ) is made, and endPre begins execution:
\begin{tabular}{|ll|}
\hline first & \(=0\) \\
last & \(=6\) \\
& \\
\hline
\end{tabular}

First character of strExp is + , so at point \(X\), a recursive call is made and the new invocation of endPre begins execution:


Next character of \(\operatorname{strExp}\) is *, so at point \(X\), a recursive call is made and the new invocation of endPre begins execution:
\begin{tabular}{|c|c|c|c|c|c|}
\hline first \(=0\) & \multirow[t]{3}{*}{X} & first \(=1\) & \multirow[t]{2}{*}{} & \multirow[t]{3}{*}{first
last} & \multirow[t]{3}{*}{\(=2\)
\(=6\)} \\
\hline last \(=6\) & & last \(=6\) & & & \\
\hline \[
\begin{aligned}
& \text { firstEnd }=\text { ? } \\
& \text { X: endPre("+*ab-cd", 1) }
\end{aligned}
\] & & \[
\begin{aligned}
& \text { firstEnd }=\text { ? } \\
& \text { X: endPre("+*ab-cd", 2) }
\end{aligned}
\] & & & \\
\hline
\end{tabular}

Next character of strExp is a, which is a base case. The current invocation of endPre completes execution and returns its value:
\begin{tabular}{|c|c|c|c|c|c|c|}
\hline first \(=0\) & & first & \(=1\) & & first & \(=2\) \\
\hline last \(=6\) & X & 1ast & \(=6\) & 2 & last & \(=6\) \\
\hline firstEnd = ? & & firstEnd & \(=2\) & & return 2 & \\
\hline X: endPre("+*ab-cd", 1) & & & & & & \\
\hline
\end{tabular}

Because firstEnd \(>-1\), a recursive call is made from point \(Y\) and the new invocation of endPre begins execution:
\begin{tabular}{|ll}
\hline \begin{tabular}{l} 
first \\
last \\
firstEnd \\
X: endPre("+*ab-cd", 1)
\end{tabular} & \begin{tabular}{l}
\(=0\) \\
\(=6\)
\end{tabular} \\
\hline
\end{tabular}

Next character of \(\operatorname{str} \operatorname{Exp}\) is b , which is a base case. The current invocation of endPre completes execution and returns its value:
\begin{tabular}{|c|c|c|c|c|c|}
\hline first \(=0\) & \multirow{4}{*}{X} & first \(=1\) & & first & = 3 \\
\hline last =6 & & last \(=6\) & & last & \(=6\) \\
\hline firstEnd =? & & firstEnd =2 & 3 & 1 return 3 & \\
\hline X: endPre("+*ab-cd", 1) & & return 3 & & & \\
\hline
\end{tabular}

The current invocation of endPre completes execution and returns its value:
\begin{tabular}{|c|c|c|c|}
\hline first & \(=0\) & first & \(=1\) \\
\hline 1ast & \(=6\) & last & \(=6\) \\
\hline firstEnd & \(=3\) & firstEnd & \(=2\) \\
\hline
\end{tabular}

Because firstEnd \(>-1\), a recursive call is made from point \(Y\) and the new invocation of endPre begins execution:
\begin{tabular}{|lr|}
\hline \begin{tabular}{l} 
last \\
firstEnd \\
Y: endPre("+*ab-cd", 4) \\
return ?
\end{tabular} \\
\hline
\end{tabular}

Next character of strExp is -, so at point \(X\), a recursive call is made and the new invocation of endPre begins execution:
\begin{tabular}{|ll|}
\hline \begin{tabular}{l} 
last \\
firstEnd \\
Y: endPre("+*ab-cd", 4) \\
return?
\end{tabular} \\
\hline
\end{tabular}

Next character of \(\operatorname{strExp}\) is \(c\), which is a base case. The current invocation of endPre completes execution and returns its value:


Because firstEnd \(>-1\), a recursive call is made from point \(Y\) and the new invocation of endPre begins execution:


Next character of strExp is \(d\), which is a base case. The current invocation of endPre completes execution and returns its value:


The current invocation of endPre completes execution and returns its value:
\begin{tabular}{|c|c|c|c|c|}
\hline first & \(=0\) & & first & \(=4\) I \\
\hline last & \(=6\) & & 1 last & \(=6\) \\
\hline firstEnd & \(=3\) & 6 & firstEnd & \(=5\) \\
\hline return 6 & 4 & & return 6 & \\
\hline
\end{tabular}

The current invocation of endPre completes execution and returns its value to the original call to endPre:


Now you can use the function endPre to determine whether a string contains a prefix expression as follows:
```

// Sees whether an expression is a prefix expression.
// Precondition: strExp contains a string with no blank characters.
// Postcondition: Returns true if the expression is in prefix form; otherwise returns false.
isPrefix(strExp: string): boolean
lastChar = endPre(strExp, 0)
return (lastChar >= 0) and (lastChar == strExp.7ength() - 1)

```

Having determined that a string is a prefix expression, how can you evaluate it? Because each operator is followed by its two operands, you can look ahead in the expression for them. However, such operands can themselves be prefix expressions, which you must evaluate first. These prefix expressions are subexpressions of the original expression and must therefore be "smaller." A recursive solution to this problem seems natural.

The following function, which appears in pseudocode, evaluates a prefix expression. This algorithm is simpler than one that evaluates infix expressions.

A recognition algorithm for prefix expressions

An algorithm to evaluate a prefix expression
```

    // Returns the value of a given prefix expression.
    // Precondition: strExp is a string containing a valid prefix expression with no blanks.
    evaluatePrefix(strExp: string): float
    strLength = the length of strExp
    if (strLength == 1)
        return value of the identifier // Base case-single identifier
    else
    {
op = strExp[0] // strExp begins with an operator
// Find the end of the first prefix expression-will be the first operand
endFirst = endPre(strExp, 1)
// Recursively evaluate this first prefix expression
operand1 = evaluatePrefix(strExp[1..endFirst]);
// Recursively evaluate the second prefix expression - will be the second operand
endSecond = strLength - endFirst + 1
operand2 = evaluatePrefix(strExp[endFirst + 1, endSecond])
// Evaluate the prefix expression
return operand1 op operand2
}

```

Question 2 Write the prefix expression that represents the following infix expression: \((a / b) * c-(d+e) * f\)
Question 3 Write the infix expression that represents the following prefix expression: \(-a / b+c * \operatorname{def}\)
Question 4 Is the following string a prefix expression? +-/abc*+def * \(g h\)

\subsection*{5.2.3 Postfix Expressions}

A grammar that defines the language of all postfix expressions is
\[
\begin{aligned}
& <\text { postfix }>=<\text { identifier }>\mid<\text { postfix }><\text { postfix }><\text { operator }> \\
& <\text { operator }>=+|-|*| / \\
& <\text { identifier }>=a|b| \ldots \mid z
\end{aligned}
\]

Some calculators require that you enter two numbers before you enter the operation that you want to perform. Such calculators, in fact, require you to enter postfix expressions.

Here we shall develop an algorithm for converting a prefix expression to a postfix expression. Chapter 6 presents a nonrecursive algorithm for evaluating postfix expressions. Together, these two algorithms give you another technique for evaluating a prefix expression. To simplify the conversion algorithm, assume that, by using the prefix recognition algorithm, you have a syntactically correct prefix expression.

If you think recursively, the conversion from prefix form to postfix form is straightforward. If the prefix expression exp is a single letter, then
\[
\text { postfix }(\exp )=\exp
\]

Otherwise exp must be of the form
\[
<\text { operator }><\text { prefix } 1><\text { prefix } 2>
\]

The corresponding postfix expression is then
\[
<\text { postfix } 1><\text { postfix } 2><\text { operator }>
\]
where \(<\) prefix \(1>\) converts to \(<\) postfix \(1>\) and \(<\) prefix \(2>\) converts to \(<\) postfix \(2>\). Therefore,
```

postfix(exp) = postfix(prefix1) • postfix(prefix 2) • <operator>

```

Thus, at a high level, the conversion algorithm is
```

if (exp is a single letter)
return exp
else
return postfix(prefix1) • postfix(prefix2) • <operator>

```

The following pseudocode function convert refines this algorithm. The string pre contains the prefix expression.
```

// Converts a prefix expression to postfix form.
// Precondition: The string pre is a valid prefix expression with no blanks.
// Postcondition: Returns the equivalent postfix expression.
convert(preExp: string): string
preLength $=$ the length of preExp
ch $=$ first character in preExp
postExp = an empty string
if (ch is a lowercase letter)
// Base case-single identifier
postExp $=$ postExp • ch // Append to end of postExp
e1se // ch is an operator
\{
// pre has the form <operator> <prefix1> <prefix2>
endFirst = endPre(preExp, 1) // Find the end of prefix1
// Recursively convert prefix1 into postfix form
postExp $=$ postExp • convert(preExp[1..endFirst])
// Recursively convert prefix2 into postfix form
postExp $=$ postExp $\cdot \operatorname{convert(preExp[endFirst~+~1..preLength~-~1))~}$
postExp $=$ postExp • ch // Append the operator to the end of postExp
\}
return post

```

Question 5 Write the postfix expression that represents the following infix expression: \(\left(a^{*} b-c\right) / d+(e-f)\)

\subsection*{5.2.4 Fully Parenthesized Expressions}

Most programmers would object to using prefix or postfix notation for their algebraic expressions, so most programming languages use infix notation. However, infix notation requires precedence rules, rules for association, and parentheses to avoid ambiguity within the expressions.

An algorithm that converts a prefix expression to postfix form

A recursive algorithm that converts a prefix expression to postfix form

A grammar for the
language of fully parenthesized algebraic expressions

You can make precedence and association rules unnecessary by placing parentheses around each pair of operands together with their operator, thereby avoiding any ambiguity. A grammar for the language of all fully parenthesized infix expressions is
\[
\begin{aligned}
& <\text { infix }>=<\text { identifier }>\mid(<\text { infix }><\text { operator }><\text { infix }>) \\
& <\text { operator }>=+|-|*| / \\
& <\text { identifier }>=a|b| \ldots \mid z
\end{aligned}
\]

Although the grammar is simple, the language is rather inconvenient for programmers.
Therefore, most programming languages support a definition of algebraic expressions that includes both precedence rules for the operators and rules of association so that fully parenthesized expressions are not required. However, the grammars for defining such languages are more involved and the algorithms for recognizing and evaluating their expressions are more difficult than those you have seen in this section. Programming Problem 8 at the end of this chapter describes such a grammar without left-to-right association rules and asks you to write a recognition algorithm. Programming Problem 8 at the end of Chapter 6 presents a nonrecursive evaluation algorithm for algebraic expressions that use both precedence and left-to-right association rules.

\subsection*{5.3 Backtracking}

This section considers an organized way to make successive guesses at a solution. If a particular guess leads to a dead end, you back up to that guess and replace it with a different guess. This strategy of retracing steps in reverse order and then trying a new sequence of steps is called backtracking. You can combine recursion and backtracking to solve the following problem.

\subsection*{5.3.1 Searching for an Airline Route}

This example will introduce you to a general type of search problem. In this particular problem, you must find a path from some point of origin to some destination point. We shall solve this problem by using recursion. In the next chapter, we will solve it again without recursion.

The High Planes Airline Company (HPAir) wants a program to process customer requests to fly from some origin city to some destination city. So that we can focus on recursion, we will simplify the problem: For each customer request, just indicate whether a sequence of HPAir flights from the origin city to the destination city exists.

Imagine three input text files that specify all of the flight information for the airline as follows:
- The names of cities that HPAir serves
- Pairs of city names, each pair representing the origin and destination of one of HPAir's flights
- Pairs of city names, each pair representing a request to fly from some origin to some destination

The program should then produce output such as
```

Request is to fly from Providence to San Francisco.
HPAir flies from Providence to San Francisco.
Request is to fly from Philadelphia to Albuquerque.
Sorry. HPAir does not fly from Philadelphia to Albuquerque.
Request is to fly from Salt Lake City to Paris.
Sorry. HPAir does not serve Paris.

```


Representing the flight data. The flight map in Figure 5-4 represents the routes that HPAir flies. An arrow from city \(C_{1}\) to city \(C_{2}\) indicates a flight from \(C_{1}\) to \(C_{2}\). In this case, \(C_{2}\) is adjacent to \(C_{1}\) and the path from \(C_{1}\) to \(C_{2}\) is called a directed path. Notice that if \(C_{2}\) is adjacent to \(C_{1}\), it does not follow that \(C_{1}\) is adjacent to \(C_{2}\). For example, in Figure 5-4, there is a flight from city \(R\) to city \(X\), but not from city \(X\) to city \(R\). As you will see in Chapter 20, the map in Figure 5-4 is called a directed graph.

The nature of the search. When processing a customer's request to fly from some origin city to some destination city, you must determine from the flight map whether there is a route from the origin to the destination. For example, by examining the flight map in Figure 5-4, you can see that a customer could fly from city \(P\) to city \(Z\) by flying first to city \(W\), then to city \(Y\), and finally to city \(Z\); that is, there is a directed path from \(P\) to \(Z: P \rightarrow W, W \rightarrow Y, Y \rightarrow Z\). Thus, you must develop an algorithm that searches the flight map for a directed path from the origin city to the destination city. Such a path might involve either a single flight or a sequence of flights. The solution developed here performs an exhaustive search. That is, beginning at the origin city, the solution will try every possible sequence of flights until either it finds a sequence that gets to the destination city or it determines that no such sequence exists.

First consider how you might perform the search by hand. One approach is to start at the origin city \(C_{0}\) and select an arbitrary flight that departs from the origin city. This flight will lead you to a new city, \(C_{1}\). If city \(C_{1}\) happens to be the destination city, you are done; otherwise, you must attempt to get from \(C_{1}\) to the destination city. To do this, you select a path to travel out of \(C_{1}\). This flight will lead you to city \(C_{2}\). If \(C_{2}\) is the destination, you are done; otherwise, you must attempt to get from \(C_{2}\) to the destination city, and so on.
A recursive strategy. To fly from the origin city to the destination city by first flying from the origin city to a city \(C\) and then by flying from \(C\) to the destination has a distinct recursive flavor. We can restate this strategy as follows:

Select a city C adjacent to the origin
Fly from the origin to city C
if (C is the destination city)
Terminate- the destination is reached
else
Fly from city C to the destination
\(C_{2}\) is adjacent to \(C_{1}\) if there is a directed path from \(C_{1}\) to \(C_{2}\)

A recursive search strategy

Possible outcomes of the exhaustive search strategy

Use backtracking to recover from a wrong choice

A refinement of the recursive search algorithm

Consider the possible outcomes of applying the previous strategy:
1. You eventually reach the destination city and can conclude that it is possible to fly from the origin to the destination.
2. You reach a city \(C\) from which there are no departing flights.
3. You go around in circles. For example, from \(C_{1}\) you go to \(C_{2}\), from \(C_{2}\) you go to \(C_{3}\), and from \(C_{3}\) you go back to \(C_{1}\). You might continue this tour of the three cities forever; that is, the algorithm might not terminate.

If you always obtained the first outcome, everyone would be happy. This outcome corresponds to a base case of the recursive algorithm. If you ever reach the destination city, no additional problems of the form "fly from city \(C\) to the destination" are generated, and the algorithm terminates. However, because HPAir does not fly between all pairs of cities, you certainly cannot expect that the algorithm will always find a path from the origin city to the destination. For example, if city \(P\) in Figure 5-4 is the origin city and city \(Q\) is the destination city, the algorithm could not possibly find a path from city \(P\) to city \(Q\).

Even if there were a sequence of flights from the origin city to the destination, it would take a bit of luck for the previous strategy to discover it - the algorithm would have to select a "correct" flight at each step. For example, even though there is a way to get from city \(P\) to city \(Z\) in Figure 5-4, the algorithm might not find it and instead might reach outcome 2 or 3 . That is, suppose that from city \(P\) the algorithm chose to go to city \(R\). From city \(R\) the algorithm would have to go to city \(X\), from which there are no flights out (outcome 2). On the other hand, suppose that the algorithm chose to go to city \(W\) from city \(P\). From city \(W\), the algorithm might choose to go to city \(S\). It would then have to go to city \(T\) and then back to \(W\). From \(W\), it might once again choose to go to city \(S\) and continue to go around in circles (outcome 3).

You thus need to make the algorithm more sophisticated, so that it always finds a path from the origin to the destination, if such a path exists, and otherwise terminates with the conclusion that there is no such path. Suppose that the earlier strategy results in outcome 2: You reach a city \(C\) from which there are no departing flights. This certainly does not imply that there is no way to get from the origin to the destination; it implies only that there is no way to get from city \(C\) to the destination. In other words, it was a mistake to go to city \(C\). After discovering such a mistake, the algorithm can retrace its steps, or backtrack, to the city \(C^{\prime}\) that was visited just before city \(C\) was visited. Once back at city \(C^{\prime}\), the algorithm can select a flight to some city other than \(C\). Notice that it is possible that there are no other flights out of city \(C^{\prime}\). If this were the case, it would mean that it was a mistake to visit city \(C^{\prime}\), and thus, you would want to backtrack again - this time to the city that was visited just before city \(C^{\prime}\).

For example, you saw that, in trying to get from city \(P\) to city \(Z\) in Figure 5-4, the algorithm might first choose to go from city \(P\) to city \(R\) and then on to city \(X\). As there are no departing flights from city \(X\), the algorithm must backtrack to city \(R\), the city visited before city \(X\). Once back at city \(R\), the algorithm would attempt to go to some city other than city \(X\) but would discover that this is not possible. The algorithm would thus backtrack once more, this time to city \(P\), which was visited just before city \(R\). From city \(P\), the algorithm would choose to go to city \(W\), which is a step in the right direction!

We can resolve these problems by refining our strategy: We mark visited cities and never fly to a city that has been visited already.
```

// Discovers whether a sequence of flights from originCity to destinationCity exists.
searchR(originCity: City, destinationCity: City): boolean
Mark originCity as visited
if (originCity is destinationCity)
Terminate - the destination is reached
else
for (each unvisited city C adjacent to originCity)
searchR(C, destinationCity)

```


Now consider what happens when the algorithm reaches a city that has no unvisited city adjacent to it. For example, consider the piece of a flight map in Figure 5-5. When searchR reaches city \(M\) that is, when the argument originCity has the value \(M\)-the for loop will not be entered, because no unvisited cities are adjacent to \(M\). Hence, the searchR algorithm returns. This return has the effect of backtracking to city \(L\), from which the flight to \(M\) originated. In terms of the previous pseudocode, the return is made to the point from which the call searchR(M, destinationCity) occurred. This point is within the for loop, which iterates through the unvisited cities adjacent to \(L\); that is, the argument originCity has the value \(L\).

After backtracking from \(M\) to \(L\), the for loop will again execute. This time the loop chooses city \(K\), resulting in the recursive call searchR(K, destinationCity). From this point, the algorithm will either eventually reach the destination city-and terminate-or backtrack once again to city \(L\). If it backtracks to \(L\), the for loop will terminate, because there are no more unvisited cities adjacent to \(L\), and a return from searchR will occur. The effect is to backtrack to the city where the flight to \(L\) originated. If the algorithm ever backtracks to the origin city and no remaining unvisited cities are adjacent to it, the algorithm will terminate, and you can conclude that no sequence of flights from the origin to the destination exists. Notice that the algorithm will always terminate in one way or another, because it will either reach the destination city or run out of unvisited cities to try.

\section*{Note: Backtracking}

When searching for a sequence of flights between cities, you must take into account the possibility that the algorithm will make wrong choices. For example, the algorithm must be able to backtrack when it hits a dead end, and you must eliminate the possibility that the algorithm will cycle.

Operations on the flight map. Now consider the operations that the search algorithm must perform on the flight map. The algorithm marks cities as it visits them, determines whether a city has been visited, and determines which cities are adjacent to a given city. You can treat the flight map as an ADT that has at least these operations in addition to the search operation itself. Other desirable operations include placing data into the flight map, inserting a city adjacent to another city, displaying the flight map, displaying a list of all cities, and displaying all cities that are adjacent to

ADT flight map operations

C++ implementation of searchR
a given city. Thus, the ADT flight map could include the following operations, given here in UML notation.
```

// Reads flight information into the flight map.
+readFlightMap(cityFileName: string, flightFileName: string): void
// Displays flight information.
+displayFlightMap(): void
// Displays the names of all cities that HPAir serves.
+displayA11Cities(): void
// Displays all cities that are adjacent to a given city.
+displayAdjacentCities(aCity: City): void
// Marks a city as visited.
+markVisited(aCity: City): void
// Clears marks on all cities.
+unvisitAl1(): void
// Sees whether a city was visited.
+isVisited(aCity: City): boolean
// Inserts a city adjacent to another city in a flight map.
+insertAdjacent(aCity: City, adjCity: City): void
// Returns the next unvisited city, if any, that is adjacent to a given city.
// Returns a sentinel value if no unvisited adjacent city was found.
+getNextCity(fromCity: City): City
// Tests whether a sequence offlights exists between two cities.
+isPath(originCity: City, destinationCity: City): boolean

```

The latter operation isPath uses the searchR algorithm to make its determination. If the class Map is the implementation of the ADT flight map, its method isPath has the following C++ definition. Note that NO_CITY is a City object that Map defines as a constant for getNextCity to return in case it cannot find an unvisited adjacent city.
```

/** Tests whether a sequence of flights exists between two cities.
@pre originCity and destinationCity both exist in the flight map.
@post Cities visited during the search are marked as visited
in the flight map.
@param originCity The origin city.
@param destinationCity The destination city.
@return True if a sequence of flights exists from originCity
to destinationCity; otherwise returns false. */
bool Map::isPath(City originCity, City destinationCity)
{
bool result, done;
// Mark the current city as visited
markVisited(originCity);
// Base case: the destination is reached
if (originCity == destinationCity)
result = true;
else // Try a flight to each unvisited city
{
done = false;
City nextCity = getNextCity(originCity);

```
```

        while (!done && (nextCity != NO_CITY))
        {
            done = isPath(nextCity, destinationCity);
            if (!done)
                nextCity = getNextCity(originCity);
        } // end while
        result = done;
    } // end if
    return result;
    } // end isPath

```

Note: The subproblems that a recursive solution generates eventually must reach a base case. Failure to do so could result in an algorithm that does not terminate. Solutions that involve backtracking are particularly subject to this kind of error.

Question 6 Trace the method isPath with the map in Figure 5-6 for the following requests. Show the recursive calls and the returns from each.
- Fly from \(A\) to \(B\).
- Fly from \(A\) to \(D\).
- Fly from \(C\) to \(G\).

FIGURE 5-6 Flight map for Checkpoint Question 6


\subsection*{5.3.2 The Eight Queens Problem}

A chessboard contains 64 squares that form eight rows and eight columns. The most powerful piece in the game of chess is the queen, because it can attack any other piece within its row, within its column, or along its diagonal. The Eight Queens problem asks you to place eight queens on the chessboard so that no queen can attack any other queen.

One strategy is to guess at a solution. However, according to Section 2.6.3 of Chapter 2, there are \(g(64,8)=4,426,165,368\) ways to arrange eight queens on a chessboard of 64 squares-so many that it would be exhausting to check all of them for a solution to this problem. Nevertheless, a simple observation eliminates many arrangements from consideration: No queen can reside in a row or a column that contains another queen. Alternatively, each row and column can contain exactly one queen. Thus, attacks along rows or columns are eliminated, leaving only \(8!=40,320\) arrangements of queens to be checked for attacks along diagonals. A solution now appears more feasible.

Place eight queens on the chessboard so that no queen can attack any other queen

FIGURE 5-7 Placing one queen at a time in each column, and the placed queens' range of attack: (a) the first queen in column 1; (b) the second queen in column 2; (c) the third queen in column 3 ; (d) the fourth queen in column 4 ; (e) the five queens can attack all of column 6; (f) backtracking to column 5 to try another square for the queen; (g) backtracking to column 4 to try another square for the queen; (h) considering column 5 again


Place queens one column at a time

If you reach an impasse, backtrack to the previous column

Suppose that you provide some organization for the guessing strategy by placing one queen per column, beginning with the first square of column 1. Figure 5-7a shows this queen and its range of attack. When you consider column 2 , you eliminate its first square because row 1 contains a queen, you eliminate its second square because of a diagonal attack, and you finally place a queen in the third square of column 2, as Figure 5-7b illustrates. The black dots in the figure indicate squares that are rejected because a queen in that square is subject to attack by another queen in an earlier column. The blue dots indicate the additional squares that the new queen can attack.

We continue to place queens in this manner until we get to column 6, as Figure \(5-7 \mathrm{e}\) shows. Although the five placed queens cannot attack each other, they can attack any square in column 6 , and therefore, you cannot place a queen in column 6. You must back up to column 5 and move its queen to the next possible square in column 5, which is in the last row, as Figure 5-7f indicates. When you consider column 6 once again, there are still no choices for a queen in that column. Because you have exhausted the possibilities in column 5, you must back up to column 4. As Figure \(5-7 \mathrm{~g}\) shows, the next possible square in column 4 is in row 7 . You then consider column 5 again and place a queen in row 2 (Figure 5-7h).

How can you use recursion in the process that was just described? Consider an algorithm that places a queen in a column, given that you have placed queens correctly in the preceding columns.

First, if there are no more columns to consider, you are finished; this is the base case. Otherwise, after you successfully place a queen in the current column, you need to consider the next column. That is, you need to solve the same problem with one fewer column; this is the recursive step. Thus, you begin with eight columns, consider smaller problems that decrease in size by one column at each recursive step, and reach the base case when you have a problem with no columns.

This solution appears to satisfy the criteria for a recursive solution. However, you do not know whether you can successfully place a queen in the current column. If you can, you recursively consider the next column. If you cannot place a queen in the current column, you need to backtrack, as has already been described.

The Eight Queens problem can be solved in a variety of ways. The solution in this chapter uses two classes: a Board class to represent the chessboard and a Queen class to represent a queen on the board. A Queen object keeps track of its row and column placement and contains a static pointer to the Board. It also has operations to move to the next row and to see whether it is subject to attack. A Board object keeps track of the Queen objects currently on the board and contains operations-such as placeQueens-to perform the Eight Queens problem and display the solution.

The following pseudocode describes the algorithm for placing queens in columns, given that the previous columns contain queens that cannot attack one another:
// Places queens in eight columns.
p7aceQueens (queen: Queen) : void
if (queen's column is greater than the last column) The problem is solved
else
\{
while (unconsidered squares exist in queen's column and the problem is unsolved)
\{
Find the next square in queen's column that is not under attack by a queen in an earlier column if (such a square exists) \{

Place a queen in the square
// Try next column
placeQueens (new Queen(firstRow, queen's column + 1))
if (no queen is possible in the next column)
\{
Delete the new queen
Remove the last queen placed on the board and consider the next square in that column
\}
        \}
        \}
\}
The Eight Queens problem is initiated by the method doEightQueens, which calls placeQueens with a new queen in the upper-left corner of the board:
```

doEightQueens()
{
placeQueens(new Queen(firstRow, firstColumn))
}

```

After doEightQueens has completed, the board may display the solution, if one was found.

The solution combines recursion with backtracking

FIGURE 5-8 A solution to the Eight Queens problem


Figure 5-8 indicates the solution that the previous algorithm finds. By modifying the arguments to placeQueens, you can discover other solutions to the Eight Queens problem. The Programming Problems at the end of this chapter ask you to consider other solutions to this algorithm, as well as additional modifications.

Question 7 Consider a Four Queens problem, which has the same rules as the Eight Queens problem but uses a \(4 \times 4\) board. Find all solutions to this new problem by applying backtracking by hand.

Implementing eight queens using the STL class vector. The Board class in the solution described thus far may be represented in a number of ways. The simplest representation would be a two-dimensional array; however, such an array wastes space because only eight squares out of 64 are used. Another approach would be to use a one-dimensional array of only the squares that contain a queen. Because the algorithm uses backtracking, a dynamic array is the optimal choice. The vector container in the STL is often used in place of an array type, because it allows the number of elements to vary dynamically and provides several built-in methods. Indexing is provided with array-type subscripting or with the at method, which provides range checking as well.

As given in Listing 5-1, the class Board contains a vector of pointers to Queen objects, as well as several methods to manipulate the queens on the board. The class Queen appears in Listing 5-2.

\section*{LISTING 5-1 The header file for the class Board}
```

/** @file Board.h */
\#ifndef _BOARD
\#define _BOARD
\#include "Queen.h"
\#include <vector>
\#include <cassert>

```
```

\#include <iostream>
using namespace std;
static const int BOARD_SIZE = 8;
class Board
{
private:
vector<Queen*> queens; // Array of pointers to queens on the board
/** Sees whether a queen exists in position (inRow, inCol). */
bool isQueen(int inRow, int inCol) const;
/** Attempts to place queens on board, starting with the designated queen. */
const placeQueens(Queen* queenPtr);
/** Removes the last queen from the board, but does not deallocate it. */
void removeQueen();
/** Places a queen on the board. */
void setQueen(const Queen* queenPtr);
public:
/** Supplies the Queen class with a pointer to the board. */
Board();
/** Clears the board and removes pointer from queens. */
~Board();
/** Clears board. */
void clear();
/** Displays board. */
void display() const;
/** Initiates the Eight Queens prob7em. */
void doEightQueens();
/** @return The number of queens on the board. */
int getNumQueens() const;
/** @return A pointer to the queen at the designated index. */
const Queen* getQueen(int index) const;
}; // end Board
\#endif

```
```

LISTING 5-2 The class Queen
class Board; // Forward declaration of Board class
/** The Queen class. */
class Queen
{
public:
/** Places a queen in upper-1eft corner of board. */
Queen();

```
```

    /** Places a queen in supplied location. */
    Queen(int inRow, int inCol);
    /** @return Column number.*/
    int getCol() const;
    /** areturn Row number. */
    int getRow() const;
    /** Moves queen to next row. */
    void nextRow();
    /** Sees whether the queen is under attack by another queen.
    @return True if another queen is in the same row or the same
        diagonal; otherwise, returns false. */
    bool isUnderAttack() const;
    /** Saves a pointer to the board for al1 queens. */
    static void setBoard(const Board* bPtr);
    private:
int row; // Row (or prospective row) of queen if it is on the board
int col; // Column (or prospective column) of queen if it is on
// the board
// A17 queens share the same board
static const Board* boardPtr;
}; // end Queen

```

An implementation of p laceQueens follows:
```

bool Board::placeQueens(Queen* queenPtr)
{
// Base case: Try to place a queen in a nonexistent column.
if (queenPtr->getCo1() >= BOARD_SIZE)
{
de1ete queenPtr;
return true;
} // end if
bool isQueenP1aced = false;
while (!isQueenPlaced \&\& queenPtr->getRow() < BOARD_SIZE)
{
// If the queen can be attacked, try moving it
// to the next row in the current column
if (queenPtr->isUnderAttack())
queenPtr->nextRow();
else
{
// Put this queen on the board and try putting a
// new queen in the first row of the next column
setQueen(queenPtr);
Queen* newQueenPtr = new Queen(0, queenPtr->getCol() + 1);
isQueenPlaced = placeQueens(newQueenPtr);

```
```

            // If it wasn't possible to put the new queen in the next column,
            // backtrack by deleting the new queen and moving the last
            // queen placed down one row
            if (!isQueenP1aced)
        {
            delete newQueenPtr;
            removeQueen();
            queenPtr->nextRow();
        } // end if
        } // end if
    } // end while
    return isQueenPlaced;
    } // end placeQueens

```

\subsection*{5.4 The Relationship Between Recursion and Mathematical Induction}

A very strong relationship exists between recursion and mathematical induction. Recursion solves a problem by specifying a solution to one or more base cases and then demonstrating how to derive the solution to a problem of an arbitrary size from the solutions to smaller problems of the same type. Similarly, mathematical induction proves a property about the natural numbers by proving the property about a base case-usually 0 or 1 -and then proving that the property must be true for an arbitrary natural number \(n\) if it is true for the natural numbers smaller than \(n\).

Given the similarities between recursion and mathematical induction, it should not be surprising that induction is often employed to prove properties about recursive algorithms. What types of properties? You can, for example, prove that an algorithm actually performs the task that you intended. As an illustration, we will prove that the recursive factorial algorithm of Chapter 2 does indeed compute the factorial of its argument. Another use of mathematical induction is to prove that a recursive algorithm performs a certain amount of work. For example, we will prove that the solution to the Towers of Hanoi problem—also from Chapter 2—makes exactly \(2^{N}-1\) moves when it starts with \(N\) disks.

You can use induction to prove that a recursive algorithm either is correct or performs a certain amount of work

\subsection*{5.4.1 The Correctness of the Recursive Factorial Function}

The following pseudocode describes a recursive function that computes the factorial of a nonnegative integer \(n\) :
```

fact(n: integer): integer
if (n is 0)
return 1
else
return n * fact(n - 1)

```

You can prove that the function fact returns the values
\[
\begin{aligned}
& \operatorname{factorial(0)}=0!=1 \\
& \operatorname{factorial}(n)=n!=n \times(n-1) \times(n-2) \times \ldots \times 1 \quad \text { if } n>0
\end{aligned}
\]

The proof is by induction on \(n\).
Basis. Show that the property is true for \(n=0\). That is, you must show that fact(0) returns 1 . But this result is simply the base case of the function: fact ( 0 ) returns 1 by its definition.

You now must establish that
\[
\text { property is true for an arbitrary } k \Rightarrow \text { property is true for } k+1
\]

Inductive hypothesis. Assume that the property is true for \(n=k\). That is, assume that
\[
\text { factorial }(k)=k!=k \times(k-1) \times(k-2) \times \ldots \times 1
\]

Inductive conclusion. Show that the property is true for \(n=k+1\). That is, you must show that fact \((k+1)\) returns the value
\[
(k+1) \times k \times(k-1) \times(k-2) \times \ldots \times 2 \times 1
\]

By definition of the function fact, fact \((k+1)\) returns the value
\[
(k+1) \times \text { factorial }(k)
\]

But by the inductive hypothesis, fact \((\mathrm{k})\) returns the value
\[
k \times(k-1) \times(k-2) \times \ldots \times 2 \times 1
\]

Thus, fact \((k+1)\) returns the value
\[
(k+1) \times k \times(k-1) \times(k-2) \times \ldots \times 2 \times 1
\]
which is what you needed to show to establish that
property is true for an arbitrary \(k \Rightarrow\) property is true for \(k+1\)
The inductive proof is thus complete.

\subsection*{5.4.2 The Cost of Towers of Hanoi}

In Chapter 2, you saw the following solution to the Towers of Hanoi problem:
```

solveTowers(count, source, destination, spare)
if (count is 1)
Move a disk directly from source to destination
else
{
solveTowers(count - 1, source, spare, destination)
solveTowers(1, source, destination, spare)
solveTowers(count - 1, spare, destination, source)
}

```

We now pose the following question: If you begin with \(N\) disks, how many moves does solveTowers make to solve the problem?

Let \(\operatorname{moves}(N)\) be the number of moves made starting with \(N\) disks. When \(N=1\), the answer is easy:
\[
\operatorname{moves}(1)=1
\]

When \(N>1\), the value of \(\operatorname{moves}(N)\) is not so apparent. An inspection of the solveTowers algorithm, however, reveals three recursive calls. Therefore, if you knew how many moves solveTowers made starting with \(N-1\) disks, you could figure out how many moves it made starting with \(N\) disks; that is,
\[
\operatorname{moves}(N)=\operatorname{moves}(N-1)+\operatorname{moves}(1)+\operatorname{moves}(N-1)
\]

Thus, you have a recurrence relation for the number of moves required for \(N\) disks:
```

moves(1) = 1
moves}(N)=2\times\operatorname{moves}(N-1)+1\quad\mathrm{ if }N>

```

For example, you can determine moves(3) as follows:
\[
\begin{aligned}
\operatorname{moves}(3) & =2 \times \operatorname{moves}(2)+1 \\
& =2 \times(2 \times \operatorname{moves}(1)+1)+1 \\
& =2 \times(2 \times 1+1)+1 \\
& =7
\end{aligned}
\]

A recurrence relation for the number of moves that solveTowers requires for \(N\) disks

A closed-form formula for the number of moves that solveTowers requires for \(N\) disks mecause there are techniques for obtaining a closed-form formula from 1 . Because these tech niques are not relevant to us right now, we simply pull the formula out of the blue and use mathematical induction to prove that it is correct.

The solution to the previous recurrence relation is
\[
\operatorname{moves}(N)=2^{N}-1 \quad \text { for all } N \geq 1
\]

Notice that \(2^{3}-1\) agrees with the value 7 that was just computed for moves (3).
The proof that \(\operatorname{moves}(N)=2^{N}-1\) is by induction on \(N\).
Basis. Show that the property is true for \(N=1\). Here \(2^{1}-1=1\), which is consistent with the recurrence relation's specification that moves \((1)=1\).

You now must establish that
\[
\text { property is true for an arbitrary } k \Rightarrow \text { property is true for } k+1
\]

Inductive hypothesis. Assume that the property is true for \(N=k\). That is, assume
\[
\operatorname{moves}(k)=2^{k}-1
\]

Inductive conclusion. Show that the property is true for \(N=k+1\). That is, you must show that moves \((k+1)=2^{k+1}-1\). Now
\[
\begin{aligned}
\operatorname{moves}(k+1) & =2 \times \operatorname{moves}(k)+1 & & \text { from the recurrence relation } \\
& =2 \times\left(2^{k}-1\right)+1 & & \text { by the inductive hypothesis } \\
& =2^{k+1}-1 & &
\end{aligned}
\]
which is what you needed to show to establish that
property is true for an arbitrary \(k \Rightarrow\) property is true for \(k+1\)
The inductive proof is thus complete.
Do not get the false impression that proving properties of programs is an easy matter. These two proofs are about as easy as any will be. However, well-structured programs are far more amenable to these techniques than are poorly structured programs.

Appendix E provides more information about mathematical induction.


Note: The subtleties of some of the algorithms you encountered in this chapter indicate the need for mathematical techniques to prove their correctness. The application of these techniques during the design of the various components of a solution can help to eliminate errors in logic before they appear in the program. One such technique is mathematical induction; another is the use of loop invariants, which Appendix F discusses.

\section*{Summary}
1. A grammar is a device for defining a language, which is a set of strings of symbols. By using a grammar to define a language, you often can construct a recognition algorithm that is directly based on the grammar. Grammars are frequently recursive, thus allowing you to describe vast languages concisely.
2. To illustrate the use of grammars, we defined several different languages of algebraic expressions. These languages have their relative advantages and disadvantages. Prefix and postfix expressions, though difficult for people to use, have simple grammars and eliminate ambiguity. On the other hand, infix expressions are easier for people to use but require parentheses, precedence rules, and rules of association to eliminate ambiguity. Therefore, the grammar for infix expressions is more involved.
3. Backtracking is a solution strategy that involves both recursion and a sequence of guesses that ultimately lead to a solution. If a particular guess leads to an impasse, you retrace your steps in reverse order, replace that guess, and try to complete the solution again.
4. A close relationship between mathematical induction and recursion exists. You can use induction to prove properties about a recursive algorithm. For example, you can prove that a recursive algorithm is correct, and you can derive the amount of work it requires.

\section*{Exercises}
1. Trace the following recursive functions:
a. isPal with the string abcdeba
b. isAnBn with the string AABB
c. endPre with the expression -*/abcd
2. Consider the language that the following grammar defines:
\[
\begin{aligned}
& <S>=\$|<W>| \$<S> \\
& <W>=\mathrm{abb} \mid \mathrm{a}<W>\mathrm{bb}
\end{aligned}
\]

Write all strings that are in this language and that contain seven or fewer characters.
3. Write a recursive grammar for the language of strings of one or more letters. The first letter of each string must be uppercase, and all other letters in the string must be lowercase.
4. Consider a language of character strings that contain only dots and dashes. All strings in this language contain at least four characters and begin with either two dots or two dashes. If the first two characters are dots, the last one must be a dash; if the first two characters are dashes, the last one must be a dot. Write a recursive grammar for this language.
5. Consider a language of strings that contains only \(X\) 's, \(Y\) 's and \(Z\) 's. A string in this language must begin with an \(X\). If a \(Y\) is present in a string, it must be the final character of the string.
a. Write a recursive grammar for this language.
b. Write all the possible two-character strings of this language.
6. Consider a language of words, where each word is a string of dots and dashes. The following grammar describes this language:
\[
\begin{aligned}
& <\text { word }>=<\text { dot }>\mid<\text { dash }><\text { word }>\mid<\text { word }><\text { dot }> \\
& <\text { dot }> \\
& <\text { dash }>
\end{aligned}
\]
a. Write all three-character strings that are in this language.
b. Is the string ••••- in this language? Explain.
c. Write a seven-character string that contains more dashes than dots and is in the language. Show how you know that your answer is correct.
d. Write pseudocode for a recursive recognition function isIn(str) that returns true if the string str is in this language and returns false otherwise.
7. Consider the following grammar:
\[
\begin{array}{ll}
<\text { word }> & =R|<D>|<D><\text { word }><S> \\
<D> & =Q \mid P \\
<S> & =1
\end{array}
\]

Write pseudocode for a recursive function that returns true if a string is in this language and returns false otherwise.
8. Consider the following grammar:
\[
\begin{aligned}
& <G>=\text { empty string }|<E>|<V><E>|<E><G><V> \\
& <E>=\& \mid \# \\
& <V>=\mathrm{W} \mid \mathrm{A}
\end{aligned}
\]
a. Write pseudocode for a recursive function that returns true if a string is in this language and returns false otherwise.
b. Is the string \(\& \mathrm{~W} \# \mathrm{~W}\) in this language?
9. Let \(L\) be the language
\[
L=\left\{S: S \text { is of the form } \mathrm{A}^{n} \mathrm{~B}^{2 n}, \text { for some } n>0\right\}
\]

Thus, a string is in \(L\) if and only if it starts with a sequence of \(A\) 's and is followed by a sequence of twice as many \(B\) 's. For example, AABBBB is in \(L\), but \(\mathrm{ABBB}, \mathrm{ABBABB}\), and the empty string are not.
a. Give a grammar for the language \(L\).
b. Write a recursive function that determines whether the string strExp is in \(L\).
10. Is \(+* a-b / c++d e-f g\) a prefix expression? Explain in terms of the grammar for prefix expressions.
11. Is \(a b / c^{*} e f g^{*} h /+d-+\) a postfix expression? Explain in terms of the grammar for postfix expressions.
12. Consider the language that the following grammar defines:
\[
\begin{aligned}
& <S>=<L>\mid<D><S><S> \\
& <L>=A \mid B \\
& <D>=1 \mid 2
\end{aligned}
\]
a. Write all three-character strings that are in this language.
b. Write one string in this language that contains more than three characters.
13. Consider a language of the following character strings: The letter \(A\), the letter \(B\), the letter \(C\) followed by a string that is in the language, and the letter \(D\) followed by a string in the language. For example, these strings are in this language: \(\mathrm{A}, \mathrm{CA}, \mathrm{CCA}, \mathrm{DCA}, \mathrm{B}, \mathrm{CB}, \mathrm{CCB}, \mathrm{DB}\), and DCCB .
a. Write a grammar for this language.
b. Is CAB in this language? Explain.
c. Write a recursive recognition algorithm for this language.
14. Consider the language that the following grammar defines:
\[
<\text { word }>=\$ \mid \mathrm{a}<\text { word }>\mathrm{a} \mid \mathrm{b}<\text { word }>\mathrm{b}|\cdots| \mathrm{y}<\text { word }>\mathrm{y} \mid \mathrm{z}<\text { word }>\mathrm{z}
\]

Equivalently,
\[
L=\{s \$ \operatorname{reverse}(s): s \text { is a string of letters of length } \geq 0\}
\]

Note that this language is very similar to the language of palindromes, but there is a special middle character here.

The algorithm that this chapter gave for recognizing palindromes can be adapted easily to this language. The algorithm, which is recursive and processes the string str from both ends toward the middle, is based on the following facts:
- A string with no characters is not in the language.
- A string with exactly one character is in the language if the character is a \(\$\).
- A longer string is in the language if the ends are identical letters and the inner substring (from the second character to the next-to-last character of str) is in the language.
Describe a recursive recognition algorithm that processes the string from left to right, reading one character at a time without saving the string for future reference. Write a C++ function that implements your algorithm.

Consider the following recursive function:
```

int p(int x)
{
if (x =< 3)
return x;
else
return p(x-1) * p(x-3);
} // end p

```

Let \(m(x)\) be the number of multiplication operations that the execution of \(p(x)\) performs.
a. Write a recursive definition of \(m(x)\).
b. Prove that your answer to part \(a\) is correct by using mathematical induction.
16. Consider palindromes that consist only of lowercase letters, such as "level" and "deed," but not "RadaR," "ADA," or "101." Let \(c(n)\) be the number of palindromes of length \(n\).
a. Write a recursive definition of \(c(n)\).
b. Prove that your answer to part \(a\) is correct by using mathematical induction.
17. Prove the following for single-letter operands: If \(E\) is a prefix expression and \(Y\) is a nonempty string of nonblank characters, then \(E Y\) cannot be a legal prefix expression. (Hint: Use a proof by induction on the length of \(E\).)
18. Chapter 2 gave the following definition for \(g(n, k)\), where \(n\) and \(k\) are assumed to be nonnegative integers:
\[
g(n, k)= \begin{cases}1 & \text { if } k=0 \\ 1 & \text { if } k=n \\ 0 & \text { if } k>n \\ g(n-1, k-1)+g(n-1, k) & \text { if } 0<k<n\end{cases}
\]

Prove by induction on \(n\) that the following is a closed form for \(g(n, k)\) :
\[
g(n, k)=\frac{n!}{(n-k)!k!}
\]

\section*{Programming Problems}
1. Complete the classes Queen and Board for the Eight Queens problem.
2. Revise the program that you just wrote for the Eight Queens problem so that it answers the following questions:
a. How many backtracks occur? That is, how many times does the program remove a queen from the board?
b. How many calls to isUnderAttack are there?
c. How many recursive calls to placeQueens are there?
*3. You can begin the Eight Queens problem by placing a queen in the second square of the first column instead of in the first square. You can then call placeQueens to begin with the second column. This revision should lead you to a new solution. Write a program that finds all solutions to the Eight Queens problem. \({ }^{1}\)
4. Do you know how to find your way through a maze? After you write this program, you will never be lost again!

Assume that a maze is a rectangular array of squares, some of which are blocked to represent walls. The maze has one entrance and one exit. For example, if \(x\) 's represent the walls, a maze could appear as follows:
```

XXXXXXXXXXXXXXXXXX X
X X XXXX X
X XXXXX XXXXX XX X
X XXXXX XXXXXXX XX X
X X XX XX X
X XXXXXXXXXX XX X
XxXXXXXXXXXXOXXXXXXX

```

A creature, indicated in the previous diagram by \(o\), sits just inside the maze at the entrance (bottom row). Assume that the creature can move in only four directions: north, south, east, and west. In the diagram, north is up, south is down, east is to the right, and west is to the left. The problem is to move the creature through the maze from the entrance to the exit (top row), if possible. As the creature moves, it should mark its path. At the conclusion of the trip through the maze, you should see both the correct path and incorrect attempts. Write a program to solve this problem.

Squares in the maze have one of several states: CLEAR (the square is clear), WALL (the square is blocked and represents part of the wall), PATH (the square lies on the path to the exit), and VISITED (the square was visited, but going that way led to an impasse).

This problem uses two ADTs that must interact. The ADT creature represents the creature's current position and contains operations that move the creature. The creature should be able to move north, south, east, and west one square at a time. It should also be able to report its position and mark its trail.

The ADT maze represents the maze itself, which is a two-dimensional rectangular arrangement of squares. You could number the rows of squares from the top beginning with zero, and number the columns of squares from the left beginning with zero. You could then use a row number and a column number to uniquely identify any square within the maze. The ADT clearly needs a data structure to represent the maze. It also needs such data as the height and width of the maze given in numbers of squares; the length of a side of a square, and the row and column coordinates of both the entrance to and the exit from the maze.

The ADT maze should also contain, for example, operations that create a specific maze given descriptive data that we will detail to display a maze, determine whether a particular square is part of the wall, determine whether a particular square is part of the path, and so on.

The search algorithm and its supporting functions are outside both of the ADTs creature and maze. Thus, the maze and the creature will be arguments that you must pass to these functions. If you are at the maze's entrance, you can systematically find your way out of the maze by using the following search algorithm. This involves backtracking - that is, retracing your steps when you reach an impasse.

Step 1. First check whether you are at the exit. If you are, you're done (a very simple maze); if you are not, go to step 2.
Step 2. Try to move to the square directly to the north by calling the function goNorth (step 3).
Step 3. If goNorth was successful, you are done. If it was unsuccessful, try to move to the square directly to the west by calling the function goWest (step 4).
Step 4. If goWest was successful, you are done. If it was unsuccessful, try to move to the square directly to the south by calling the function goSouth (step 5).
Step 5. If goSouth was successful, you are done. If it was unsuccessful, try to move to the square directly to the east by calling the function goEast (step 6).
Step 6. If goEast was successful, you are done. If it was unsuccessful, you are still done, because no path exists from the entrance to the exit.

The function goNorth will examine all the paths that start at the square to the north of the present square as follows. If the square directly to the north is clear, is inside the maze, and has not been visited before, move into this square and mark it as part of the path. (Note that you are moving from the south.) Check whether you are at the exit. If you are, you're done. Otherwise, try to find a path to the exit from here by trying all paths leaving this square except the one going south (going south would put you back in the square from which you just came) as follows. Call goNorth; if it is not successful, call goWest and, if it is not successful, call goEast. If goEast is not successful, mark this square as visited, move back into the square to the south, and return.

The following pseudocode describes the goNorth algorithm:
```

goNorth(maze, creature)
if (the square to the north is clear, inside the maze, and unvisited)
{
Move to the north
Mark the square as part of the path
if (at exit)

```
```

    success = true
    else
    {
        success = goNorth(maze, creature)
        if (!success)
        {
            success = goWest(maze, creature)
            if (!success)
            {
                success = goEast(maze, creature)
                if (!success)
                    {
                        Mark square visited
                        Backtrack south
                    }
                }
        }
    }
    }
else
success = false
return success

```

The gowest function will examine all the paths that start at the square to the west of the present square as follows. If the square directly to the west is clear, is inside the maze, and has not been visited before, move into this square and mark it as part of the path. (Note that you are moving from the east.) Check whether you are at the exit. If you are, you're done. Otherwise, try to find a path to the exit from here by trying all paths leaving this square except the one going east (this would put you back in the square from which you just came) as follows. Call goNorth; if it is not successful, call goWest; and if it is not successful, call goSouth. If goSouth is not successful, mark this square as visited, move back into the square to the east, and return. The functions goEast and goSouth are analogous to goWest and goNorth.

The input data to represent a maze is simple. For example, the previously given maze is represented by the following lines of input, which can be in a text file:
\begin{tabular}{|c|c|}
\hline 207 & \(\leftarrow\) width and height of the maze in squares \\
\hline 018 & \(\leftarrow\) row and column coordinate of maze exit \\
\hline 612 & \(\leftarrow\) row and column coordinate of maze entrance \\
\hline  & \\
\hline \(x \quad \mathrm{xxxx}\) & \\
\hline x xxxxx xxxxx dx & \\
\hline  & \\
\hline x x ( xx xx & \\
\hline x xxxyxxyxx \({ }^{\text {x }}\) & \\
\hline  & \\
\hline
\end{tabular}

After the first three lines of numeric data in the file, each of the next lines corresponds to a row in the maze, and each character in a line corresponds to a column in the maze. An \(x\) indicates a blocked square (part of the wall), and a blank indicates a clear square. This notation is convenient, because you can see what the maze looks like as you design it.
5. Write a program that will recognize and evaluate prefix expressions. First design and implement a class of prefix expressions. This class should contain methods to recognize and evaluate prefix expressions. This chapter discusses the algorithms you will need to implement these methods.
6. Implement a recognition algorithm for the language in Exercise 5.
7. Implement the algorithm described in this chapter to convert a prefix expression to postfix form.
8. The following is a grammar that allows you to omit parentheses in infix algebraic expressions when the precedence rules remove ambiguity. For example, \(a+b * c\) means \(a+(b * c)\). However, the grammar requires parentheses when ambiguity would otherwise result. That is, the grammar does not permit left-to-right association when several operators have the same precedence. For example, \(a / b * c\) is illegal. Notice that the definitions introduce factors and terms.
```

$<$ expression $>=<$ term $>\mid<$ term $>+<$ term $>\mid<$ term $>-<$ term $>$
$<$ term $>=<$ factor $>\mid<$ factor $>*<$ factor $>\mid<$ factor $>\mid<$ factor $>$
$<$ factor $>=<$ letter $>\mid(<$ expression $>)$
$<$ letter $>=\mathrm{a}|\mathrm{b}| \ldots \mid \mathrm{z}$

```

The recognition algorithm is based on a recursive chain of subtasks: find an expression \(\rightarrow\) find a term \(\rightarrow\) find a factor. What makes this a recursive chain is that find an expression uses find a term, which in turn uses find a factor. Find a factor either detects a base case or uses find an expression, thus forming the recursive chain. The pseudocode for the recognition algorithm follows:
```

FIND AN EXPRESSION
// The grammar specifies that an expression is either a single term
// or a term followed by a + or a -, which then must be followed by a second term.
Find a term
if (the next symbol is + or -)
Find a term
FIND A TERM
// The grammar specifies that a term is either a single factor or
// a factor followed by a* or a /, which must then be followed by a second factor.
Find a factor
if (the next symbol is * or/)
Find a factor
FIND A FACTOR
// The grammar specifies that a factor is either a single letter (the base case) or
// an expression enclosed in parentheses.
if (the first symbol is a letter)
Done
else if (the first symbol is a '(')
{
Find an expression starting at the character after ' ('
Checkfor ')'
}
e1se
No factor exists

```

Design and implement a class of infix expressions, as described by the given grammar. Include a method to recognize a legal infix expression.

\section*{Stacks}

\section*{Chapter}

\section*{Contents}

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\section*{Prerequisites}
\(\begin{array}{ll}\text { Chapter 1 } & \text { Data Abstraction: The Walls } \\ \mathrm{C}++ \text { Interlude 1 } & \mathrm{C}++ \text { Classes }\end{array}\)

So far in this book, we have discussed data abstraction as it relates to the design of a solution, introduced \(\mathrm{C}++\) classes as a way to hide a solution's implementation, introduced resizable arrays and linked nodes as data structures used in many ADT implementations, and developed recursion as a problem-solving technique that is useful in the construction of algorithms. The primary concerns of the remainder of this book are the aspects of problem solving that involve the management of datathat is, the identification and implementation of some of the more common datamanagement operations.

Our study of data management has three goals. The first is to identify useful sets of operationsthat is, to identify abstract data types. The second goal is to examine applications that use these abstract data types. The third goal is to construct implementations for the abstract data types-that is, to develop data structures and classes. As you will discover, the nature of the operations of an abstract data type, along with the application in which you will use it, greatly influences the choice of its implementation.

The ADT bag discussed in Chapter 1 does not organize its data. However, you can organize data either by position or by value. In general, these organizations are appropriate for applications of rather different natures. For example, if an application needs to ask a question about the first person in a line, you should organize the data by position. On the other hand, if an application needs to ask a question about the employee named Smith, you should organize the data by value. Throughout the rest of this book, you will see several ADTs that use these two data organizations.

This chapter introduces a well-known ADT called a stack. You will see how the operations on a stack give it a last-in, first-out behavior. Two of the several applications of a stack that the chapter considers are evaluating algebraic expressions and searching for a path between two points. Finally, the chapter discusses the important relationship between stacks and recursion. We leave the implementations of the stack to the next chapter.

\subsection*{6.1 The Abstract Data Type Stack}

The specification of an abstract data type that you can use to solve a particular problem can emerge
be an important one: the ADT stack.

\subsection*{6.1.1 Developing an ADT During the Design of a Solution}

When you type a line of text on a keyboard, you are likely to make mistakes. If you use the Backspace key to correct these mistakes, each backspace erases the previous character entered. Consecutive backspaces are applied in sequence and so erase several characters. For instance, if you type the line
\(\mathrm{abcc} \leftarrow \mathrm{ddde} \leftarrow \leftarrow \leftarrow \mathrm{eg} \leftarrow \mathrm{fg}\)
where \(\leftarrow\) represents the backspace character, the corrected input would be
abcdefg
How can a program read the original line and get the correct input? In designing a solution to this problem, you eventually must decide how to store the input line. In accordance with the ADT approach, you should postpone this decision until you have a better idea of what operations you will need to perform on the data.

A first attempt at a solution leads to the following pseudocode:
```

// Read the line, correcting mistakes along the way
while (not end of line)
{
Read a new character ch
if (ch is nota''\leftarrow')
Add ch to the ADT
e1se
Remove from the ADT (discard) the item that was added most recently
}

```

This solution calls to attention two of the operations that the ADT will have to include:
- Add a new item to the ADT.
- Remove from the ADT the item that was added most recently.

Notice that potential trouble lurks if you type a \(\leftarrow\) when the ADT is empty-that is, when the ADT contains no characters. If this situation should occur, you have three options: (1) have the program terminate and write an error message, (2) throw an exception, or (3) have the program ignore the \(\leftarrow\) and continue. Any of these options is reasonable, so let's suppose that you decide to ignore the \(\leftarrow\) and continue. Therefore, the algorithm becomes
```

// Read the line, correcting mistakes along the way
while (not end of line)
{
Read a new character ch
if (ch is not a '\leftarrow')
Add ch to the ADT
else if (the ADT is not empty)
Remove from the ADT the item that was added most recently
else
Ignore the ' }\leftarrow
}

```

From this pseudocode, you can identify a third operation required by the ADT:
- See whether the ADT is empty.

This solution places the corrected input line in the ADT. Now suppose that you want to display the line. At first, it appears that you can accomplish this task by using the ADT operations already identified, as follows:
```

// Display the line
while (the ADT is not empty)
{
Remove from the ADT the item that was added most recently
Display....Uh-oh!
}

```

This pseudocode is incorrect for two reasons:
- When you remove an item from the ADT, the item is gone, so you cannot display it. What you should have done was to look at the item that was added to the ADT most recently. An accessor operation returns an item-enabling you to look at it-but does not change the ADT's data. Only after getting and displaying the item should you remove it from the ADT.
- The last character of the input line is the item most recently added to the ADT. You certainly do not want to write it first. The resolution of this particular difficulty is left to you as an exercise.

If we address only the first difficulty, the following pseudocode displays the input line in reverse order:

Two ADT operations that are required

The "read and correct" algorithm

ADT operation

A false start at writing the line

Reasons why the attempted solution is incorrect

The write-backward algorithm
```

// Display the line in reverse order
while (the ADT is not empty)
{
Get the item that was added to the ADT most recently and assign it to ch
Display ch
Remove from the ADT the item that was added most recently
}

```

Another required ADT operation

LIFO: The last item inserted onto a stack is the first item out

Thus, a fourth operation is required by the ADT:
- Get the item that was added to the ADT most recently.

Although you have yet to think about an implementation of the ADT, you know that you must be able to perform four specific operations. \({ }^{1}\) These operations define the required ADT, which happens to be well known: It is usually called a stack.

\subsection*{6.1.2 Specifications for the ADT Stack}

We have identified the following operations for the ADT stack:
- See whether a stack is empty.
- Add a new item to the stack.
- Remove from the stack the item that was added most recently.
- Get the item that was added to the stack most recently.

The term "stack" is intended to conjure up visions of things encountered in daily life, such as a stack of plates in the school cafeteria, a stack of books on your desk, or a stack of assignments that you need to work on. In common English usage, "stack of" and "pile of" are synonymous. To computer scientists, however, a stack is not just any old pile. A stack has the property that the last item placed on the stack will be the first item removed. This property is commonly referred to as last in, first out or simply LIFO.

A stack of plates in a cafeteria makes a very good analogy of the abstract data type stack, as Figure 6-1 illustrates. As new plates are added, the old ones drop farther into the well beneath the surface. At any particular time, only the plate last placed on the stack is above the surface and visible. This plate is at the top of the stack and is the one that must be removed next. In general, the plates are removed in exactly the opposite order from that in which they were added.

The LIFO property of stacks seems inherently unfair. Think of the poor person who finally gets the last plate on the cafeteria's stack, one that may have been placed there six years ago. Or how would you like to be the first person to arrive on the stack for a movie-as opposed to the line for a movie. You would be the last person allowed in! These examples demonstrate the reason that stacks are not especially prevalent in everyday life. The property that we usually desire in our daily lives is first in, first out, or FIFO. A queue, which you will learn about in Chapter 13, is the ADT with the FIFO property. Most people would much prefer to wait in a movie queue-as a line is called in Britain-than in a movie stack. However, while the LIFO property of stacks is not appropriate for very many everyday situations, it is precisely what is needed for a large number of problems that arise in computer science.

\section*{FIGURE 6-1 A stack of cafeteria plates}


\footnotetext{
1 The final algorithm to write the line correctly instead of in reverse order does not require additional ADT operations.
}

Notice how well the analogy holds between the ADT stack and the stack of cafeteria plates. The operations that manipulate data on the ADT stack are the only such operations, and they correspond to the only things that you can do to a stack of plates. You can determine whether the stack of plates is empty but not how many plates are on the stack; you can inspect the top plate but no other plate; you can place a plate on top of the stack but at no other position; and you can remove a plate from the top of the stack but from no other position. If you were not permitted to perform any of these operations, or if you were permitted to perform any other operations, the ADT would not be a stack.

Although the stack of cafeteria plates suggests that, as you add or remove plates, the other plates move, do not have this expectation of the ADT stack. The stack operations involve only the top item and imply only that the other items in the stack remain in sequence. Implementations of the ADT stack operations might or might not move the stack's items. The implementations given in the next chapter do not move data items.

Refining the specification of the ADT stack. The following summary of the ADT stack provides some detail about how we wish to specify its operations. The names given here for stack operations are conventional.

\section*{Abstract Data Type: Stack}

\section*{DATA}
- A finite number of objects, not necessarily distinct, having the same data type and ordered by when they were added.

\section*{Operations}
\begin{tabular}{|c|c|}
\hline Pseudocode & Description \\
\hline \multirow[t]{3}{*}{isEmpty ()} & Task: Sees whether this stack is empty. \\
\hline & Input: None. \\
\hline & Output: True if the stack is empty; otherwise false. \\
\hline \multirow[t]{3}{*}{push(newEntry)} & Task: Adds newEntry to the top of this stack. \\
\hline & Input: newEntry. \\
\hline & Output: True if the operation is successful; otherwise false. \\
\hline \multirow[t]{3}{*}{pop()} & Task: Removes the top of this stack. That is, it removes the item that was added most recently. \\
\hline & Input: None. \\
\hline & Output: True if the operation is successful; otherwise false. \\
\hline \multirow[t]{3}{*}{peek ()} & Task: Returns the top of this stack. That is, it gets the item that was added most recently. The operation does not change the stack. \\
\hline & Input: None. \\
\hline & Output: The top of the stack. \\
\hline
\end{tabular}

Figure 6-2 shows a UML diagram for a class of stacks.
Using the ADT stack in a solution. Recall that Chapter 1 urged you to focus on the specification of a module before you considered its implementation. After writing an ADT's operations in
\begin{tabular}{|l|}
\hline \multicolumn{1}{|c|}{ Stack } \\
\hline \\
\begin{tabular}{l} 
+isEmpty (): boolean \\
+push(newEntry: ItemType) : boolean \\
+pop(): boolean \\
+peek(): ItemType
\end{tabular} \\
\hline
\end{tabular}

The refined algorithms
pseudocode, you should try to use them as a check of your design. Such a test can highlight any deficiencies in your specifications or design. For example, you can use the previous stack operations to refine the algorithms developed earlier in this chapter.
```

// Reads an input line, recognizing the character ' }\leftarrow'\mathrm{ ' as a backspace that erases the
// previously typed character.
// Returns a stack of the corrected characters read.
readAndCorrect(): Stack
aStack = a new empty stack
Read newChar
while (newChar is not the end-of-line symbol)
{
if (newChar is nota ' }\leftarrow\mathrm{ ')
aStack.push(newChar)
else if (!aStack.isEmpty())
aStack.pop()
Read newChar
}
return aStack
// Displays the input line in reverse order by writing the contents of the stack aStack.
displayBackward(aStack: Stack)
while (!aStack.isEmpty())
{
newChar = aStack.peek()
aStack.pop()
Write newChar
}
Advance to new line

```

You should be able to use the stack operations without knowing their implementations or even what a stack looks like. Because working with an ADT involves building a wall around the implementation, your program can use a stack independently of the stack's implementation. As long as the program correctly uses the ADT's operations - that is, as long as it honors the operation contract-it will work regardless of how you implement the ADT.

For the program to do this, however, the operation contract must be written precisely. That is, before you implement any operations of the ADT, you should specify in detail both their requirements before use and their ultimate effect. Realize, however, that during program design, the first attempt at specification is often informal and is only later made precise by the writing
of preconditions and postconditions. For example, the previous specifications of the ADT stack leave the following questions unanswered:
- How will pop affect an empty stack?
- What will peek do when the stack is empty?

Note: Operations such as pop and peek must take reasonable action when the stack

An interface. The interface given in Listing 6-1 formalizes our specification of the ADT stack and answers the previous questions about pop and peek .
```

LISTING 6-1 A C++ interface for stacks

## LISTING 6-1 A C++ interface for stacks

```
/** @file StackInterface.h */
```

/** @file StackInterface.h */
\#ifndef _STACK_INTERFACE
\#ifndef _STACK_INTERFACE
\#define _STACK_INTERFACE
\#define _STACK_INTERFACE
temp1ate<c1ass ItemType>
temp1ate<c1ass ItemType>
class StackInterface
class StackInterface
{
{
public:
public:
/** Sees whether this stack is empty.
/** Sees whether this stack is empty.
@return True if the stack is empty, or false if not. */
@return True if the stack is empty, or false if not. */
virtual bool isEmpty() const = 0;
virtual bool isEmpty() const = 0;
/** Adds a new entry to the top of this stack.
/** Adds a new entry to the top of this stack.
@post If the operation was successful, newEntry is at the top of the stack.
@post If the operation was successful, newEntry is at the top of the stack.
@param newEntry The object to be added as a new entry.
@param newEntry The object to be added as a new entry.
@return True if the addition is successful or false if not. */
@return True if the addition is successful or false if not. */
virtual bool push(const ItemType\& newEntry) = 0;
virtual bool push(const ItemType\& newEntry) = 0;
/** Removes the top of this stack.
/** Removes the top of this stack.
@post If the operation was successful, the top of the stack
@post If the operation was successful, the top of the stack
has been removed.
has been removed.
@return True if the removal is successful or false if not. */
@return True if the removal is successful or false if not. */
virtual bool pop() = 0;
virtual bool pop() = 0;
/** Returns the top of this stack.
/** Returns the top of this stack.
@pre The stack is not empty.
@pre The stack is not empty.
@post The top of the stack has been returned, and
@post The top of the stack has been returned, and
the stack is unchanged.
the stack is unchanged.
@return The top of the stack. */
@return The top of the stack. */
virtual ItemType peek() const = 0;
virtual ItemType peek() const = 0;
}; // end StackInterface
}; // end StackInterface
\#endif

```
```

\#endif

```
```

Questions that the informal specifications of the stack leave unanswered
is empty.


An axiom is a mathematical rule

Axioms for multiplication

Axioms specify the behavior of an ADT

An example of an axiom

Question 1 If you push the letters $A, B, C$, and $D$ in order onto a stack of characters and then pop them, in what order will they be deleted from the stack?
Question 2 What do the initially empty stacks stack1 and stack2 "look like" after the following sequence of operations?

```
stack1.push(1)
stack1.push(2)
stack2.push(3)
stack2.push(4)
stack1.pop()
stackTop = stack2.peek()
stack1.push(stackTop)
stack1.push(5)
stack2.pop()
stack2.push(6)
```

Axioms (optional). Most people will understand intuitive specifications, such as those given previously for the stack operations. However, some ADTs are much more complex and less intuitive than a stack. For such ADTs, you should use a more rigorous approach to defining the behavior of their operations: You must supply a set of mathematical rules-called axioms-that precisely specify the behavior of each operation of an ADT.

An axiom is a true statement for an ADT's operation. For example, you are familiar with axioms for algebraic operations; in particular, you know the following rules for multiplication:

$$
\begin{aligned}
& (a \times b) \times c=a \times(b \times c) \\
& a \times b=b \times a \\
& a \times 1=a \\
& a \times 0=0
\end{aligned}
$$

These rules, or axioms, are true for any numeric values of $a, b$, and $c$, and describe the behavior of the multiplication operator $\times$.

In a similar fashion, you can write a set of axioms that completely describes the behavior of the operations for the ADT stack. For example,

## A newly created stack is empty

is an axiom because it is true for all newly created stacks. We can state this axiom succinctly in terms of the ADT stack operations as follows, if we represent a newly created stack by the pseudocode expression new Stack():
(new Stack()).isEmpty() = true
To formally capture the intuitive notion that the last item inserted into the stack aStack is the first item to be removed, you could write an axiom such as
(aStack.push(newItem)). pop() = aStack
That is, if you push newItem onto aStack and then pop it, you are left with the original stack aStack.
The following axioms formally define the ADT stack.

D

## Note: Axioms for the ADT stack

```
(new Stack()).isEmpty() = true
(new Stack()).pop() = false
(new Stack()).peek() = error
(aStack.push(item)).isEmpty() = false
(aStack.push(item)).peek() = item
(aStack.push(item)).pop() = true
```


### 6.2 Simple Uses of a Stack

Once you have satisfactorily specified the behavior of an ADT, you can design applications that access and manipulate the ADT's data solely in terms of its operations and without regard for its implementation. Thus, we can use the operations of the ADT stack even though we have not discussed their implementations yet. This section presents two rather simple examples for which the LIFO property of stacks is appropriate.

### 6.2.1 Checking for Balanced Braces

C++ uses curly braces, "\{" and "\}", to delimit groups of statements. For example, braces begin and end a method's body. If you treat a C++ program as a string of characters, you can use a stack to verify that a program contains balanced braces. For example, the braces in the string

```
abc{defg{ijk}{7{mn}}op}qr
```

are balanced, while the braces in the string

```
abc{def}}{ghij{k1}m
```

are not balanced. You can check whether a string contains balanced braces by traversing it from left to right. As you move from left to right, you match each successive close brace " $\}$ " with the most recently encountered unmatched open brace " $\{$ "; that is, the " $\{$ " must be to the left of the current " $\}$ ". The braces are balanced if

1. Each time you encounter a " $\}$ ", it matches an already encountered " $\{$ ".
2. When you reach the end of the string, you have matched each " $\{$ ".

The solution requires that you keep track of each unmatched " $\{$ " and discard one each time you encounter a " $\}$ ". One way to perform this task is to push each " $\{$ " encountered onto a stack and pop one off each time you encounter a " \}". Thus, a first-draft pseudocode solution is

```
for (each character in the string)
{
    if (the character is a '{')
        aStack.push('{')
    else if (the character is a'}')
        aStack.pop()
}
```

Although this solution correctly keeps track of braces, missing from it are the checks that conditions 1 and 2 are met - that is, that the braces are indeed balanced. To verify condition 1 when a " $\}$ " is encountered, you must check to see whether the stack is empty before popping from it. If it is empty,

Requirements for balanced braces

Initial draft of a solution

A detailed pseudocode solution to check a string for balanced braces
you terminate the loop and report that the string is not balanced. To verify condition 2 , you check that the stack is empty when the end of the string is reached.

Thus, the pseudocode solution to check for balanced braces in aString becomes

```
// Checks the string aString to verify that braces match.
// Returns true if aString contains matching braces, false otherwise.
checkBraces(aString: string): boolean
    aStack = a new empty stack
    balancedSoFar = true
    i = 0
    while (balancedSoFar and i < length of aString)
    {
        ch = character at position i in aString
        i++
        // Push an open brace
        if (ch is a'{')
            aStack.push('{')
        // Close brace
        else if (ch is a'}')
        {
            if (!aStack.isEmpty())
                aStack.pop() // Pop a matching open brace
            else // No matching open brace
                balancedSoFar = false
        }
        // Ignore all characters other than braces
    }
if (balancedSoFar and aStack.isEmpty())
    aString has balanced braces
e1se
    aString does not have balanced braces
```

Figure 6-3 shows the stacks that result when this algorithm is applied to several simple examples.

FIGURE 6-3 Traces of the algorithm that checks for balanced braces


Question 3 For each of the following strings, trace the execution of the balanced-braces algorithm and show the contents of the stack at each step.
a. $\mathrm{x}\{\{\mathrm{yz}\}\}\}$
b. $\{x\{y\{\{z\}\}\}$
c. $\{\{\{x\}\}\}$

Note: The push operation can fail for implementation-dependent reasons. For example, push fails if the array in an array-based implementation is full. In the spirit of failsafe programming, a function that implements this balanced-braces algorithm should check push's return value.

## Aside: A solution without stacks

It may have occurred to you that a simpler solution to this problem is possible. You need only keep a count of the current number of unmatched open braces. ${ }^{2}$ You need not actually store the open braces in a stack. However, the stack-based solution is conceptually useful as it previews more legitimate uses of stacks. For example, Exercise 7 at the end of this chapter asks you to extend the algorithm given here to check for balanced parentheses and square brackets in addition to braces.

### 6.2.2 Recognizing Strings in a Language

Consider the problem of recognizing whether a particular string is in the language
$L=\left\{s \$ s^{\prime}: s\right.$ is a possibly empty string of characters other than $\left.\$, s^{\prime}=\operatorname{reverse}(s)\right\}$
For example, the strings $\mathrm{A} \$ \mathrm{~A}, \mathrm{ABC} \$ \mathrm{CBA}$, and $\$$ are in $L$, but $\mathrm{AB} \$ \mathrm{AB}$ and ABC CB are not. (Exercise 14 in Chapter 5 introduced a similar language.) This language is like the language of palindromes that you saw in Chapter 5, but strings in this language have a special middle character.

A stack is useful in determining whether a given string is in $L$. Suppose you traverse the first half of the string and push each character onto a stack. When you reach the $\$$ you can undo the process: For each character in the second half of the string, you pop a character off the stack. However, you must match the popped character with the current character in the string to ensure that the second half of the string is the reverse of the first half. The stack must be empty when-and only when-you reach the end of the string; otherwise, one "half" of the string is longer than the other, and so the string is not in $L$.

The following algorithm uses this strategy. To avoid unnecessary complications, assume that aString contains exactly one $\$$.
// Checks the string aString to verify that it is in language L.
// Returns true if aString is in L, false otherwise.
recognizeString(aString: string) : boolean

[^5]```
```

aStack = a new empty stack

```
```

aStack = a new empty stack
// Push the characters that are before the \$ (that is, the characters in s) onto the stack
// Push the characters that are before the \$ (that is, the characters in s) onto the stack
i = 0
i = 0
ch = character at position i in aString
ch = character at position i in aString
while (ch is not a'$')
while (ch is not a'$')
{
{
aStack.push(ch)
aStack.push(ch)
i++
i++
ch = character at position i in aString
ch = character at position i in aString
}
}
// Skip the \$
// Skip the \$
i++
i++
// Match the reverse of S
// Match the reverse of S
inLanguage = true // Assume string is in language
inLanguage = true // Assume string is in language
while (inLanguage and i < length of aString)
while (inLanguage and i < length of aString)
{
{
if (!aStack.isEmpty())
if (!aStack.isEmpty())
{
{
stackTop = aStack.peek()
stackTop = aStack.peek()
aStack.pop()
aStack.pop()
ch = character at position i in aString
ch = character at position i in aString
if (stackTop equals ch)
if (stackTop equals ch)
i++ // Charactersmatch
i++ // Charactersmatch
else
else
inLanguage = false // Characters do not match (top of stack is not ch)
inLanguage = false // Characters do not match (top of stack is not ch)
}
}
else
else
inLanguage = false // Stack is empty (first half of string is shorter
inLanguage = false // Stack is empty (first half of string is shorter
// than second half)
// than second half)
}
}
if (inLanguage and aStack.isEmpty())
if (inLanguage and aStack.isEmpty())
aString is in language
aString is in language
e1se
e1se
aString is not in language

```
    aString is not in language
```

```
    ++
```

```
    ++
```

You can use an ADT's operations in an application without the distraction of implementation details

Note: In both of the preceding examples, notice how you can focus on the task at hand without the distraction of implementation details such as arrays. With less to worry about, you are less likely to make an error in your logic when you use the ADT's operations in applications such as checkBraces and recognizeString. Likewise, when you finally implement the ADT's operations in C++, you will not be distracted by these applications. In addition, because checkBraces and recognizeString do not depend on any implementation decisions that you make, they are not altered by your decisions. These remarks assume that you do not change the specifications of the ADT's operations when you implement them. However, developing software is not a linear process. You may realize during implementation that you need to refine your specifications. Clearly, changes to the specification of any module affect any alreadydesigned uses of that module.

To summarize, you can specify the behavior of an ADT independently of its implementation. Given such a specification, and without any knowledge of how the ADT will be implemented, you can design applications that use the ADT's operations to access its data.

Question 4 Trace the execution of the language-recognition algorithm described in the previous section for each of the following strings, and show the contents of the stack at each step.
a. $\mathrm{a} \$ \mathrm{a}$
b. $a b \$ a b$
c. $\mathrm{ab} \$ \mathrm{a}$
d. $a b \$ b a$

### 6.3 Using Stacks with Algebraic Expressions

This section contains two more problems that you can solve neatly by using the ADT stack. Keep in mind throughout that you are using the ADT stack to solve the problems. You can use the stack operations, but you may not assume any particular implementation. You choose a specific implementation only as a last step.

Chapter 5 presented recursive grammars that specified the syntax of algebraic expressions. Recall that prefix and postfix expressions avoid the ambiguity inherent in the evaluation of infix expressions. We will now consider stack-based solutions to the problems of evaluating infix and postfix expressions. To avoid distracting programming issues, we will allow only the binary operators *,/, + , and - , and we will disallow exponentiation and unary operators.

The strategy we shall adopt here is first to develop an algorithm for evaluating postfix expressions and then to develop an algorithm for transforming an infix expression into an equivalent postfix expression. Taken together, these two algorithms provide a way to evaluate infix expressions. This strategy eliminates the need for an algorithm that directly evaluates infix expressions, which is a somewhat more difficult problem that Programming Problem 8 considers.

### 6.3.1 Evaluating Postfix Expressions

As we mentioned in Chapter 5, some calculators require you to enter postfix expressions. For example, to compute the value of

$$
2 *(3+4)
$$

by using a postfix calculator, you would enter the sequence $2,3,4,+$, and *, which corresponds to the postfix expression

## $234+*$

Recall that an operator in a postfix expression applies to the two operands that immediately precede it. Thus, the calculator must be able to retrieve the operands entered most recently. The ADT stack provides this capability. In fact, each time you enter an operand, the calculator pushes it onto a stack. When you enter an operator, the calculator applies it to the top two operands on the stack, pops the operands from the stack, and pushes the result of the operation onto the stack. Figure 6-4 shows the action of the calculator for the previous sequence of operands and operators. The final result, 14 , is on the top of the stack.

You can formalize the action of the calculator to obtain an algorithm that evaluates a postfix expression, which is entered as a string of characters. To avoid issues that cloud the algorithm with programming details, assume that

- The string is a syntactically correct postfix expression

Simplifying

- No unary operators are present
- No exponentiation operators are present
- Operands are single lowercase letters that represent integer values

Your use of an ADT's operations should not depend on its implementation

To evaluate an infix expression, first convert it to postfix form and then evaluate the postfix expression

A pseudocode algorithm that evaluates postfix expressions

FIGURE 6-4 The effect of a postfix calculator on a stack when evaluating the expression 2 * $(3+4)$


The pseudocode algorithm is then

```
for (each character ch in the string)
{
    if (ch is an operand)
        Push the value of the operand ch onto the stack
    e1se // ch is an operator named op
    {
            // Evaluate and push the result
            operand2 = top of stack
            Pop the stack
            operand1 = top of stack
            Pop the stack
            result = operand1 op operand2
            Push result onto the stack
    }
}
```

Upon termination of the algorithm, the value of the expression will be on the top of the stack. Programming Problem 5 at the end of this chapter asks you to implement this algorithm.

Question 5 Evaluate the postfix expression $a b-c+$. Assume the following values for the identifiers: $a=7, b=3$, and $c=-2$. Show the status of the stack after each step.

### 6.3.2 Converting Infix Expressions to Equivalent Postfix Expressions

Now that you know how to evaluate a postfix expression, you will be able to evaluate an infix expression if you first can convert it into an equivalent postfix expression. The infix expressions here are the
familiar ones, such as $(a+b) * c / d-e$. They allow parentheses, operator precedence, and left-toright association.

Will you ever want to evaluate an infix expression? Certainly—you have written such expressions in programs. The compiler that translated your programs had to generate machine instructions to evaluate the expressions. To do so, the compiler first transformed each infix expression into postfix form. Knowing how to convert an expression from infix to postfix notation not only will lead to an algorithm to evaluate infix expressions, but also will give you some insight into the compilation process.

If you manually convert a few infix expressions to postfix form, you will discover three important facts:

- The operands always stay in the same order with respect to one another.
- An operator will move only "to the right" with respect to the operands; that is, if in the infix expression the operand $x$ precedes the operator $o p$, it is also true that in the postfix expression the operand $x$ precedes the operator $o p$.
- All parentheses are removed.

As a consequence of these three facts, the primary task of the conversion algorithm is determining where to place each operator.

The following pseudocode describes a first attempt at converting an infix expression to an equivalent postfix expression postfixExp:

```
Initialize postfixExp to the empty string
for (each character ch in the infix expression)
{
    switch (ch)
    {
        case ch is an operand:
            Append ch to the end of postfixExp
            break
        case ch is an operator:
            Save ch until you know where to place it
            break
        case ch is a'('ora')':
            Discard ch
            break
    }
}
```

You may have guessed that you really do not want simply to discard the parentheses, as they play an important role in determining the placement of the operators. In any infix expression, a set of matching parentheses defines an isolated subexpression that consists of an operator and its two operands. Therefore, the algorithm must evaluate the subexpression independently of the rest of the expression. Regardless of what the rest of the expression looks like, the operator within the subexpression belongs with the operands in that subexpression. The parentheses tell the rest of the expression that

You can have the value of this subexpression after it is evaluated; simply ignore everything inside.

Parentheses are thus one of the factors that determine the placement of the operators in the postfix expression. The other factors are precedence and left-to-right association.

In Chapter 5, you saw a simple way to convert a fully parenthesized infix expression to postfix form. Because each operator corresponded to a pair of parentheses, you simply moved each operator to the position marked by its close parenthesis and finally removed the parentheses.

The actual problem is more difficult, however, because the infix expression is not always fully parenthesized. Instead, the problem allows precedence and left-to-right association, and therefore

Facts about converting from infix to postfix

First draft of an algorithm to convert an infix expression to postfix form

Parentheses, operator precedence, and left-to-right association determine where to place operators in the postfix expression

Five steps in the process to convert from infix to postfix form
requires a more complex algorithm. The following is a high-level description of what you must do when you encounter each character as you read the infix string from left to right.

1. When you encounter an operand, append it to the output string postfixExp.

Justification: The order of the operands in the postfix expression is the same as the order in the infix expression, and the operands that appear to the left of an operator in the infix expression also appear to its left in the postfix expression.
2. Push each "(" onto the stack.
3. When you encounter an operator, if the stack is empty, push the operator onto the stack. However, if the stack is not empty, pop operators of greater or equal precedence from the stack and append them to postfixExp. You stop when you encounter either a "(" or an operator of lower precedence or when the stack becomes empty. You then push the current operator in the expression onto the stack. Thus, this step orders the operators by precedence and in accordance with left-to-right association. Notice that you continue popping from the stack until you encounter an operator of strictly lower precedence than the current operator in the infix expression. You do not stop on equality, because the left-to-right association rule says that in case of a tie in precedence, the leftmost operator is applied first-and this operator is the one that is already on the stack.
4. When you encounter a ")", pop operators off the stack and append them to the end of postfixExp until you encounter the matching "(".
Justification: Within a pair of parentheses, precedence and left-to-right association determine the order of the operators, and step 3 has already ordered the operators in accordance with these rules.
5. When you reach the end of the string, you append the remaining contents of the stack to postfixExp.

For example, Figure 6-5 traces the action of the algorithm on the infix expression $a-(b+c * d) / e$, assuming that the stack aStack and the string postfixExp are initially empty. At the end of the algorithm, postfixExp contains the resulting postfix expression $a b c d *+e /-$.

You can use the previous five-step description of the algorithm to develop a fairly concise pseudocode solution, which follows. The symbol • in this algorithm means concatenate (join), so

FIGURE 6-5 A trace of the algorithm that converts the infix expression $a-\left(b+c^{*} d\right) / e$ to postfix form

| ch | aStack (bottom to top) | postfixExp |  |
| :---: | :---: | :---: | :---: |
| a |  | a |  |
| - | - | a |  |
| ( | - | a |  |
| b | - | ab |  |
| + | - ${ }^{+}$ | ab |  |
| c | - + | abc |  |
| * | $-(+$ * | abc |  |
| d | $-(+$ * | abcd |  |
| ) | - ${ }^{\text {+ }}$ | abcd* | Move operators from stack to |
|  | - | abcd*+ | postfixExp until "( " |
|  | - | abcd*+ |  |
| 1 | -1 | abcd*+ | Copy operators from |
| e | -1 | abcd*+e | stack to postfixExp |
|  |  | abcd*+e/- |  |

postfixExp • x means concatenate the string currently in postfixExp and the character $x$ - that is, follow the string in postfixExp with the character $x$.

```
for (each character ch in the infix expression)
{
        switch (ch)
        {
            case operand: // Append operand to end of postfix expression-step I
            postfixExp = postfixExp • ch
            break
        case '(': // Save'('on stack-step 2
            aStack.push(ch)
            break
        case operator: // Process stack operators of greater precedence-step 3
            while (!aStack.isEmpty() and aStack.peek() is not a ' (' and
                precedence(ch) <= precedence(aStack.peek()))
            {
                Append aStack.peek() to the end of postfixExp
                aStack.pop()
            }
            aStack.push(ch) // Save the operator
            break
        case ')': // Pop stack until matching '('-step 4
            while (aStack.peek() is not a '(')
            {
                Append aStack.peek() to the end of postfixExp
                aStack.pop()
            }
            aStack.pop() // Remove the open parenthesis
            break
    }
}
// Append to postfixExp the operators remaining in the stack -step 5
while (!aStack.isEmpty())
{
    Append aStack.peek() to the end of postfixExp
    aStack.pop()
}
```

Because this algorithm assumes that the given infix expression is syntactically correct, it can ignore the return values of the stack operations. Programming Problem 7 at the end of this chapter asks you to remove this assumption.

$\square$
Note: Algorithms that evaluate an infix expression or transform one to postfix form must determine which operands apply to a given operator. Doing so allows for precedence and left-to-right association so that you can omit parentheses.

Question 6 Convert the infix expression $a / b * c$ to postfix form. Be sure to account for left-to-right association. Show the status of the stack after each step.
Question 7 Explain the significance of the precedence tests in the infix-to-postfix conversion algorithm. Why is $\mathrm{a} \geq$ test used rather than $\mathrm{a}>$ test?

A pseudocode algorithm that converts an infix expression to postfix form

Use a stack to organize an exhaustive search

Possible outcomes of the exhaustive search strategy

Use backtracking to recover from a wrong choice

### 6.4 Using a Stack to Search a Flight Map

The previous chapter described the HPAir problem, whereby we searched an airline's flight map to see whether we could fly from some origin city to some destination city. At that time, we found a recursive solution to the problem. We now will solve the problem again using a stack instead of recursion. Doing so will bring to light the close relationship between stacks and recursion.

Let's recall some of the discussion about this problem from the previous chapter. Given HPAir's flight map shown in Figure 6-6, you can fly from city $P$ to city $Z$ by flying first to city $W$, then to city $Y$, and finally to city $Z$. We need to find this path by trying every possible sequence of flights until we either find a sequence that gets to the destination city or determine that no such sequence exists. You will see that the ADT stack is useful in organizing this search.

If we were to perform the search by hand, we could start at the origin city $C_{0}$ and select an arbitrary departing flight to a new city, $C_{1}$. If city $C_{1}$ happens to be the destination city, we are done; otherwise, we must try to fly from $C_{1}$ to the destination city. By repeating this strategy, we will encounter three possible outcomes, just as you saw in the previous chapter.

1. You eventually reach the destination city.
2. You reach a city $C$ from which there are no departing flights.
3. You go around in circles forever.

Even if a sequence of flights exists from the origin city to the destination city, we might not find it and instead reach outcome 2 or 3 . Just as we did in the recursive solution, if we reached a city $C$ from which there are no departing flights-that is outcome 2 -we would backtrack to the city $C^{\prime}$ that we visited just before we visited city $C$. Once back at city $C^{\prime}$, we would select a flight to some city other than $C$. If there were no other flights out of city $C^{\prime}$, we would backtrack again-this time to the city that was visited just before city $C^{\prime}$. To avoid outcome 3, we mark the cities as we visit them, just as we did in Chapter 5, and never visit a city more than once.

For the algorithm to implement this strategy, it must maintain information about the order in which it visits the cities. First notice that when the algorithm backtracks from a city $C$, it must retreat to the city that it visited most recently before $C$. This observation suggests that you maintain the sequence of visited cities in a stack. That is, each time you decide to visit a city, you push its name onto the stack. For example, if you want to fly from city $P$ to city $Z$ in Figure 6-6, the algorithm might first choose to go from city $P$ to city $R$ and then on to city $X$. Parts $a, b$, and $c$ of Figure 6-7 illustrate the stack for these flights. You select the next city to visit from those adjacent to the city on the top of

FIGURE 6-6 A flight map


FIGURE 6-7 The stack of cities as you travel (a) from $P$; (b) to $R$; (c) to $X$; (d) back to $R$; (e) back to $P$; (f) to $W$

the stack. When you need to backtrack from the city at the top of the stack (for example, because there are no flights out of city $X$ ), you simply pop a city from the stack, as shown in Figure 6-7d. After the pop, the city on the top of the stack- $R$-is the city on the current path that you visited most recently before $X$. It is not possible to fly out of city $R$ to some city other than city $X$, so you backtrack once more, this time to city $P$, by popping the stack, as Figure 6-7e illustrates. Figure 6-7f shows the stack after the subsequent flight to city $W$.

The algorithm, as developed so far, is as follows.

```
aStack = a new empty stack
aStack.push(originCity) // Push origin city onto aStack
while (a sequence offlights from the origin to the destination has not been found)
{
    if (you need to backtrack from the city on the top of the stack)
        aStack.pop()
    else
    {
        Select a destination city C for a flight from the city on the top of the stack
        aStack.push(C)
    }
}
```

Notice that, at any point in the algorithm, the contents of the stack correspond to the sequence of flights currently under consideration. The city on the top of the stack is the city you are visiting currently, directly "below" it is the city visited previously, and so forth down to the bottom city, which is the first city visited in the sequence - that is, the origin city. In other words, within the while loop,

The stack contains a directed path from the origin city at the bottom of the stack to the city at the top of the stack. ${ }^{3}$

You can, therefore, always retrace your steps as far back through the sequence as needed.
Now consider the question of when to backtrack from the city on the top of the stack. You have already seen one case when backtracking is necessary. You must backtrack from the city on the top of the stack when there are no flights out of that city. Another time when you need to backtrack is related to the problem of going around in circles, described previously as the third possible outcome of the original strategy. You never want to visit a city that the search has already visited. As a consequence, you must backtrack from a city whenever there are no more unvisited cities to fly to. To see why, consider two cases:

- If you have visited city $C$ and it is still somewhere in the stack-that is, $i t$ is part of the sequence of cities that you are exploring currently-you do not want to visit $C$ again. Any sequence that

[^6]Backtrack when there are no more unvisited cities

Two reasons for not visiting a city more than once
goes from $C$ through $C_{1}, C_{2}, \ldots, C_{k}$, back to $C$, and then to $C^{\prime}$ might just as well skip the intermediate cities and go from $C$ directly to $C^{\prime}$.

For example, suppose that the algorithm starts at $P$ in Figure 6-6 and, in trying to find a path to $Y$, visits $W, S$, and $T$. There is now no reason for the algorithm to consider the flight from $T$ to $W$ because $W$ is already in the stack. Any city you could fly to by going from $W$ to $S$, from $S$ to $T$, and then back to $W$-such as city $Y$-you could fly to directly from $W$ without first going through $S$ and $T$. Because you do not allow the algorithm to visit $W$ a second time, it will backtrack from $S$ and $T$ to $W$ and then go from $W$ directly to $Y$. Figure 6-8 shows how the stack would appear if revisits were allowed and how it looks after backtracking when revisits are not allowed. Notice that backtracking to $W$ is very different from visiting $W$ for a second time.

- If you have visited city $C$, but it is no longer in the stack-because you backtracked from it and popped it from the stack-you do not want to visit $C$ again. This situation is subtle; consider two cases that depend on why you backtracked from the city.

If you backtracked from $C$ because there were no flights out of it, then you certainly do not ever want to try going through $C$ again. For example, if, starting at $P$ in Figure 6-6, the algorithm goes to $R$ and then to $X$, it will backtrack from $X$ to $R$. At this point, although $X$ is no longer in the stack, you certainly do not want to visit it again, because you know there are no flights out of $X$.

Now suppose that you backtracked from city $C$ because all cities adjacent to it had been visited. This situation implies that you have already tried all possible flights from $C$ and have failed to find a way to get to the destination city. There is thus no reason to go to $C$ again. For example, suppose that starting from $P$ in Figure 6-6, the algorithm executes the following sequence: Visit $R$, visit $X$, backtrack to $R$ (because there are no flights out of $X$ ), backtrack to $P$ (because there are no more unvisited cities adjacent to $R$ ), visit $W$, visit $Y$. At this point, the stack contains $P-W-Y$ with $Y$ on top, as Figure 6-8b shows. You need to choose a flight out of $Y$. You do not want to fly from $Y$ to $R$, because you have visited $R$ already and tried all possible flights out of $R$.

In both cases, visiting a city a second time does not gain you anything, and in fact, it may cause you to go around in circles.

FIGURE 6-8 The stack of cities (a) allowing revisits and (b) after backtracking when revisits are not allowed

(a)

(b)

To implement the rule of not visiting a city more than once, you simply mark a city when it has been visited. When choosing the next city to visit, you restrict consideration to unmarked cities adjacent to the city on the top of the stack. The algorithm thus becomes

```
aStack = a new empty stack
Clear marks on all cities
aStack.push(originCity) // Push origin city onto the stack
Mark the origin as visited
while (a sequence offlights from the origin to the destination has not been found)
{
    // Loop invariant: The stack contains a directed path from the origin city at
    // the bottom of the stack to the city at the top of the stack
    if (no flights exist from the city on the top of the stack to unvisited cities)
        aStack.pop() // Backtrack
    else
    {
        Select an unvisited destination city C for a flight from the city on the top of the stack
        aStack.push(C)
        Mark C as visited
    }
}
```

Finally, you need to refine the condition in the while statement. That is, you need to refine the algorithm's final determination of whether a path exists from the origin to the destination. The loop invariant, which states that the stack contains a directed path from the origin city to the city on the top of the stack, implies that the algorithm can reach an affirmative conclusion if the city at the top of the stack is the destination city. On the other hand, the algorithm can reach a negative conclusion only after it has exhausted all possibilities-that is, after the algorithm has backtracked to the origin and there remain no unvisited cities to fly to from the origin. At that point, the algorithm will pop the origin city from the stack and the stack will become empty.

With this refinement, the algorithm appears as follows:

```
// Searches for a sequence offlights from originCity to destinationCity
searchS(originCity: City, destinationCity: City): boolean
    aStack = a new empty stack
    Clear marks on all cities
    aStack.push(originCity) // Push origin onto the stack
    Mark the origin as visited
    while (!aStack.isEmpty() and destinationCity is not at the top of the stack)
    {
        // Loop invariant: The stack contains a directed path from the origin city at
        // the bottom of the stack to the city at the top of the stack
        if (noflights exist from the city on the top of the stack to unvisited cities)
        aStack.pop() // Backtrack
        e1se
        {
            Select an unvisited destination city C for a flight from the city on the top of the stack
            aStack.push(C)
            Mark C as visited
        }
    }
    if (aStack.isEmpty())
        return false // No path exists
    else
            return true // Path exists
```

Mark the visited cities

Next draft of the search algorithm

The final version of the search algorithm

FIGURE 6-9 A trace of the search algorithm, given the flight map in Figure 6-6

| Action | Reason |  |  |
| :--- | :--- | :--- | :--- |
| Push P | Initialize |  | Pontents of stack (bottom to top) |
| Push R | Next unvisited adjacent city |  | PR |
| Push X | Next unvisited adjacent city | PRX |  |
| Pop X | No unvisited adjacent city | PR |  |
| Pop R | No unvisited adjacent city | P |  |
| Push W | Next unvisited adjacent city | PW |  |
| Push S | Next unvisited adjacent city | PW S |  |
| Push T | Next unvisited adjacent city | PW S T |  |
| Pop T | No unvisited adjacent city | PW S |  |
| Pop S | No unvisited adjacent city | PW |  |
| Push Y | Next unvisited adjacent city | PW Y |  |
| Push Z | Next unvisited adjacent city | PW Y Z |  |
|  |  |  |  |

Notice that the algorithm does not specify the order of selection for the unvisited cities. It really does not matter what selection criteria the algorithm uses, because the choice will not affect the final outcome: Either a sequence of flights exists or it does not. The choice, however, will affect the specific flights that the algorithm considers. For example, suppose that the algorithm always flies to the alphabetically earliest unvisited city from the city on the top of the stack. Under this assumption, Figure 6-9 contains a trace of the algorithm's action, given the map in Figure 6-6, with $P$ as the origin city and $Z$ as the destination city. The algorithm terminates with success.

Recall from Chapter 5 the following operations that the search algorithm must perform on the flight map.

ADT flight map operations

C++ implementation of searchS

```
// Marks a city as visited.
markVisited(aCity: City): void
// Clears marks on all cities.
unvisitAl1(): void
// Returns the next unvisited city, if any, that is adjacent to a given city.
// Returns a sentinel value if no unvisited adjacent city was found.
getNextCity(fromCity: City): City
// Tests whether a sequence offlights exists between two cities.
isPath(originCity: City, destinationCity: City): boolean
```

Assuming that the class Map implements the ADT flight map and the class Stack implements the stack operations, the definition of Map's method isPath is as follows and uses the searchS algorithm. As we mentioned in Chapter 5, NO_CITY is a City object that Map defines as a constant for getNextCity to return in case it cannot find an unvisited adjacent city.

```
/** Tests whether a sequence of flights exists between two cities.
        Nonrecursive stack version.
    @pre originCity and destinationCity both exist in the flight map.
    @post Cities visited during the search are marked as visited
        in the flight map.
    @param originCity The origin city.
    @param destinationCity The destination city.
    @return True if a sequence of flights exists from originCity
    to destinationCity; otherwise returns false. */
```

```
bool Map::isPath(City originCity, City destinationCity)
{
    bool success;
    Stack aStack;
    unvisitAll(); // Clear marks on all cities
    // Push origin city onto aStack and mark it as visited
    aStack.push(originCity);
    markVisited(originCity);
    City topCity = aStack.peek();
    while (!aStack.isEmpty() && (topCity != destinationCity))
    {
        // The stack contains a directed path from the origin city
        // at the bottom of the stack to the city at the top of
        // the stack
        // Find an unvisited city adjacent to the city on the
        // top of the stack
        City nextCity = getNextCity(topCity, nextCity);
        if (nextCity == NO_CITY)
            aStack.pop(); // No city found; backtrack
        else // Visit city
        {
            aStack.push(nextCity);
            markVisited(nextCity);
        } // end if
        if (!aStack.isEmpty())
        topCity = aStack.peek();
    } // end while
    return !aStack.isEmpty();
} // end isPath
```

Programming Problem 11 at the end of this chapter provides implementation details that will enable you to complete the solution to the HPAir problem.

## Note: The STL class stack

The Standard Template Library (STL) contains the container stack, which is a template class. However, some methods of this class have slightly different names than the ones we have used here. You can use the STL stack to implement the algorithms given in this chapter until we write our own implementations in the next chapter. Alternatively, you can use the class OurStack, as described in Programming Problem 1 at the end of this chapter. You can learn more about the STL in C++ Interlude 7.

Question 8 Trace the method isPath with the map in Figure 6-10 for the following requests. Show the state of the stack after each step.
a. Fly from $A$ to $B$.
b. Fly from $A$ to $D$.
c. Fly from $C$ to $G$.

A comparison of key aspects of two search algorithms

FIGURE 6-10 Flight map for Check Point Question 8


### 6.5 The Relationship Between Stacks and Recursion

We have solved the HPAir problem twice, once in the previous chapter by using recursion and again in the previous section by using the ADT stack. We now want to look at the ways that these two approaches organize the search for a sequence of flights. You will see that the ADT stack has a hidden presence in the concept of recursion and, in fact, that stacks have an active role in most computer implementations of recursion.

Consider how the two search algorithms implement three key aspects of their common strategy.

- Visiting a new city. The recursive algorithm searchR visits a new city $C$ by calling searchR (C, destinationCity). The algorithm searchS visits city $C$ by pushing $C$ onto a stack. Notice that if you were to use the box trace to trace the execution of searchR, the call searchR (C, destinationCity) would generate a box in which the city $C$ is associated with the parameter originCity of searchR.

For example, Figure 6-11 shows both the state of the box trace for searchR and the stack for searchS at corresponding points of the search for a path from city $P$ to city $Z$ in Figure 6-6.

- Backtracking. Both search algorithms attempt to visit an unvisited city that is adjacent to the current city. Notice that this current city is the value associated with the parameter originCity in the deepest (rightmost) box of searchR's box trace. Similarly, the current city is on the top of searchs's stack. In Figure 6-11, this current city is $X$. If no unvisited cities are adjacent to the current city, the algorithms must backtrack to the previous city. The algorithm searchR backtracks by returning from the current recursive call. You represent this action in the box trace by crossing off the deepest box. The algorithm searchS backtracks by explicitly popping from its stack. For example, from the state depicted in Figure 6-11, both algorithms backtrack to city $R$ and then to city $P$, as Figure 6-12 illustrates.

FIGURE 6-11 Visiting city $P$, then $R$, then $X$ : (a) box trace versus (b) stack
(a) Box trace of recursive search:

(b) Stack-based search:


FIGURE 6-12 Backtracking from city $X$ to $R$ to $P$ : (a) box trace versus (b) stack
(a) Box trace of recursive search:

(b) Stack-based search:


- Termination.The search algorithms terminate either when they reach the destination city or when they exhaust all possibilities. All possibilities are exhausted when, after backtracking to the origin city, no unvisited adjacent cities remain. This situation occurs for searchR when all boxes have been crossed off in the box trace and a return occurs to the point of the original call to the algorithm. For searchS, no unvisited cities are adjacent to the origin when the stack becomes empty.

Thus, the two search algorithms really do perform the identical action. In fact, provided that they use the same rule to select an unvisited city-for example, traverse the current city's list of adjacent cities alphabetically-they will always visit the identical cities in the identical order. The similarities between the algorithms are far more than coincidence. In fact, it is always possible to capture the actions of a recursive function by using a stack.

An important context in which the close tie between stacks and recursion is explicitly utilized is a compiler's implementation of a recursive function. It is common for a compiler to use a stack to implement a recursive function in a manner that greatly resembles the box trace. When a recursive call to a function occurs, the implementation must remember certain information. This information consists essentially of the same local environment that you place in the boxes-values of both arguments and local variables and a reference to the point from which the recursive call was made.

During execution, the compiled program must manage these boxes of information, or activation records, just as you must manage them on paper. As the HPAir example has indicated, the operations needed to manage the activation records are those that a stack provides. When a recursive call occurs, a new activation record is created and pushed onto a stack. This action corresponds to the creation of a new box at the deepest point in the sequence. When a return is made from a recursive call, the stack is popped, bringing the activation record that contains the appropriate local environment to the top of the stack. This action corresponds to crossing off the deepest box and following the arrow back to the preceding box. Although we have greatly simplified the process, most implementations of recursion are based on stacks of activation records.

You can use a similar strategy to implement a nonrecursive version of a recursive algorithm. You might need to recast a recursive algorithm into a nonrecursive form to make it more efficient, as mentioned in Chapter 2. The previous discussion should give you a taste of the techniques for removing recursion from a program. You will encounter recursion removal as a formal topic in more advanced courses, such as compiler construction.

Typically, stacks are used to implement recursive functions

Each recursive call generates an activation record that is pushed onto a stack

You can use stacks when implementing a nonrecursive version of a recursive algorithm

## Summary

1. The ADT stack operations have a last-in, first-out (LIFO) behavior.
2. Algorithms that operate on algebraic expressions are an important application of stacks. The LIFO nature of stacks is exactly what the algorithm that evaluates postfix expressions needs to organize the operands. Similarly, the algorithm that transforms infix expressions to postfix form uses a stack to organize the operators in accordance with precedence rules and left-to-right association.
3. You can use a stack to determine whether a sequence of flights exists between two cities. The stack keeps track of the sequence of visited cities and enables the search algorithm to backtrack easily. However, displaying the sequence of cities in their normal order from origin to destination is awkward, because the origin city is at the bottom of the stack and the destination is at the top.
4. A strong relationship between recursion and stacks exists. Most implementations of recursion maintain a stack of activation records in a manner that resembles the box trace.
5. The formal mathematical study of ADTs uses systems of axioms to specify the behavior of ADT operations.

## Exercises

1. Write pseudocode statements that create a stack of the strings "Jamie", "Jane", and "Ji11" in that order with "Jamie" at the top.
2. Given the stack created in Exercise 1, in what order will three pop operations remove the strings from the stack?
3. Suppose that you have a stack aStack and an empty auxiliary stack auxStack. Show how you can do each of the following tasks by using only the ADT stack operations.
a. Display the contents of aStack in reverse order; that is, display the top last.
b. Count the number of items in aStack, leaving aStack unchanged.
c. Delete every occurrence of a specified item from aStack, leaving the order of the remaining items unchanged.
4. The diagram of a railroad switching system in Figure 6-13 is commonly used to illustrate the notion of a stack. Identify three stacks in the figure and show how they relate to one another. How can you use this system to construct any possible permutation of railroad cars?
5. Suppose that the ADT stack included a method remove $(n)$ that removes the topmost $n$ entries from a stack. Specify this method by writing comments and a header. Consider the various ways that the method could behave when the stack does not contain at least $n$ entries.
6. Section 6.1.1 described an algorithm that reads a string of characters, correcting mistakes along the way.
a. For the following input line, trace the execution of the algorithm and show the contents of the stack at each step:
```
abc}\leftarrow\textrm{de}\leftarrow\leftarrow\textrm{fg}\leftarrow\textrm{h
```


## FIGURE 6-13 Railroad switching system for Exercise 4


b. The nature of the stack-based algorithm makes it simple to display the string in reverse order, but somewhat harder to display it in its correct order. Write a pseudocode algorithm that displays the string in its correct forward order.
c. Implement the read-and-correct algorithm as a C++ function, but make the stack local to the function instead of an argument. Let the function construct a string that contains the corrected input characters in forward order, and return it.

Revise the solution to the balanced-braces problem so that the expression can contain three types of delimiters: ( ), [ ], and \{ \}.Thus, $\{\mathrm{ab}(\mathrm{c}[\mathrm{d}]) \mathrm{e}\}$ is valid, but $\{\mathrm{ab}(\mathrm{c})$ ) is not.
8. For each of the following strings, trace the execution of the language-recognition algorithm described in Section 6.2.2, and show the contents of the stack at each step.
a. $\mathrm{xy} \$ \mathrm{xy}$
b. $\mathrm{y} \$ \mathrm{yx}$
c. $\mathrm{xy} \$ \mathrm{y}$
d. $\mathrm{xy} \$ \mathrm{x}$
e. $\mathrm{xx} \$ \mathrm{xx}$
9. Write a pseudocode function that uses a stack to determine whether a string is in the language $L$, where
a. $L=\{s: s$ contains equal numbers of $A$ 's and $B$ 's $\}$
b. $L=\left\{s: s\right.$ is of the form $\mathrm{A}^{n} \mathrm{~B}^{n}$ for some $\left.n \geq 0\right\}$

Write a function that uses a stack to determine whether a string is in the language $L$, where

$$
L=\left\{s s^{\prime}: s \text { is a string of characters, } s^{\prime}=\operatorname{reverse}(s)\right\}
$$

Note: The following strings are not in the language: The empty string, a string with fewer than two characters, and a string with an odd number of characters.

Evaluate the following postfix expressions by using the algorithm given in this chapter. Show the status of the stack after each step of the algorithm. Assume the following values for the identifiers: $a=7, b=3, c=12$, $d=-5, e=1$.
a. $a b c+-$
b. $a b c-d^{*+}$
c. $a b+c-d e^{*+}$
12. Convert the following infix expressions to postfix form by using the algorithm given in this chapter. Show the status of the stack after each step of the algorithm.
a. $a-b+c$
b. $a-(b / c * d)$
c. $a /(b * c)$
d. $a / b / c-(d+e) * f$
e. $(a+b) * c$
f. $a^{*}(b / c / d)+e$
g. $a-(b+c)$
h. $a-(b+c * d) / e$
13. Execute the HPAir algorithm with the map in Figure 6-10 (see Check Point Question 8) for the following requests. Show the state of the stack after each step.
a. Fly from $A$ to $F$.
b. Fly from $A$ to $G$.
c. Fly from $F$ to $H$.
d. Fly from $D$ to $A$.
e. Fly from $I$ to $G$.
14. Section 5.1 .2 of Chapter 5 defines "palindrome." Write a pseudocode algorithm that uses a stack to test whether a given string is a palindrome.
*15. You can use the axioms for a stack to prove that the stack defined by the sequence of operations
Create an empty stack
Push a 5
Push a 7
Push a 3
Pop (the 3)
Push a 9
Push a 4
Pop (the 4)
which you can write as
(( (( (( (new Stack ()) .push(5)) .push(7)) .push(3)).pop()) .push(9)).push(4)).pop()
is exactly the same as the stack defined by the sequence
Create an empty stack
Push a 5
Push a 7
Pusha 9
which you can write as
(( (new Stack()).push(5)).push(7)).push(9)
Similarly, you can use the axioms to show that
(( (( (( (new Stack ()) .push(1)).push(2)).pop()).push(3)).pop()).pop()).isEmpty()
is true.
a. The following representation of a stack as a sequence of push operations without any pop operations is called a canonical form:
(...(new Stack()).push()).push())... ). push()

Prove that any stack is equal to a stack that is in canonical form.
b. Prove that the canonical form is unique. That is, a stack is equal to exactly one stack that is in canonical form.
c. Use the axioms to show formally that
((( (C((C( (new Stack()).push(6)).push(9)).pop()).pop()).push(2)).pop()).push(3)). push(1)).pop()).peek()
equals 3 .

## Programming Problems

1. Using the class stack in the Standard Template Library, define and test the class OurStack that is derived from StackInterface, as given in Listing 6-1. The class stack has the following methods that you can use to define the methods for OurStack.
```
stack(); // Default constructor
bool empty() const; // Tests whether the stack is empty
ItemType& top(); // Returns a reference to the top of the stack
void push(const ItemType& newEntry); // Adds newEntry to the top of the stack
void pop();
// Removes the top of the stack
```

To access stack, use the following include statement:

```
#include <stack>
```

Whenever you need a stack for any of the following problems, use the class OurStack that Programming Problem 1 asks you to write.
2. Implement the solution to the expanded balanced-braces problem in Exercise 7.
3. Write a function that uses a stack to test whether a given string is a palindrome. Exercise 14 asked you to write an algorithm for such a function.
4. Section 6.2.2 describes a recognition algorithm for the language
$L=\left\{s \$ s^{\prime}: s\right.$ is a possibly empty string of characters other than $\left.\$, s^{\prime}=\operatorname{reverse}(s)\right\}$
Implement this algorithm.
5. Design a class of postfix calculators. Use the algorithm given in this chapter to evaluate postfix expressions as entered into the calculator. Use only the operators,,$+- *$, and $/$. Assume that the postfix expressions are syntactically correct.
6. Consider simple infix expressions that consist of single-digit operands; the operators,,+- , , and $/$; and parentheses. Assume that unary operators are illegal and that the expression contains no embedded spaces. Design and implement a class of infix calculators. Use the algorithms given in this chapter to evaluate infix expressions as entered into the calculator. You must first convert the infix expression to postfix form and then evaluate the resulting postfix expression.
7. The infix-to-postfix conversion algorithm described in this chapter assumes that the given infix expression is syntactically correct. Modify Programming Problem 6 without this assumption.
8. Repeat Programming Problem 6, but use the following algorithm to evaluate an infix expression infixExp. The algorithm uses two stacks: One stack opStack contains operators, and the other stack va1Stack contains values of operands and intermediate results. Note that the algorithm treats parentheses as operators with the lowest precedence.

```
for (each character ch in infixExp)
{
    switch (ch)
    {
        case ch is an operand, that is, a digit
            va1Stack.push(ch)
            break
        case chis a '('
            opStack.push(ch)
```

```
            break
        case ch is an operator
            if (opStack.isEmpty())
                opStack.push(ch)
            else if (precedence(ch) > precedence(opStack.peek()))
                opStack.push(ch)
            e1se
            {
                while (!opStack.isEmpty() and
                    precedence(ch) <= precedence(opStack.peek())
                    Execute
            opStack.push(ch)
        }
            break
        case chis a')'
            while (opStack.peek() is not a '(')
            Execute
        opStack.pop()
        break
    }
}
while (!opStack.isEmpty())
    Execute
result = valStack.peek()
```

Note that Execute means:

```
operand2 = va1Stack.peek()
va1Stack.pop()
operand1 = va1Stack.peek()
va1Stack.pop()
op = opStack.peek()
opStack.pop()
result = operand1 op operand2
va1Stack.push(result)
```

Choose one of the following two approaches for your implementation:

- The operator stack opStack contains characters, but the operand stack va1Stack contains integers.
- The stack opStack contains integer codes that represent the operators, so both stacks contain integers.

9. The infix evaluation algorithm given in Programming Problem 8 assumes that the given infix expression is syntactically correct. Repeat Programming Problem 8 without this assumption.

Using stacks, write a nonrecursive version of the function solveTowers, as defined in Chapter 2.
11. Complete the solution to the HPAir problem. The input to the program consists of three text files, as follows:

| cityFile | Each line contains the name of a city that HPAir serves. The names are in <br> alphabetical order. |
| :--- | :--- |
| flightFile | Each line contains a pair of city names that represent the origin and destina- <br> tion of one of HPAir's flights. |
| requestFile | Each line contains a pair of city names that represent a request to fly from <br> some origin to some destination. |

You can make the following assumptions:

- Each city name contains at most 15 characters. Pairs of city names are separated by a comma.
- HPAir serves at most 20 cities.
- The input data is correct.

For example, the input files could appear as

| cityFile: | Albuquerque <br> Chicago <br> San Diego |  |
| :--- | :--- | :--- |
| flightFile: | Chicago, <br> Chicago, <br> Albuquerque, | San Diego <br> Albuquerque <br> Chicago |
| requestFile: | Albuquerque, <br> Albuquerque, | San Diego <br> Saris <br> San Diego, |

For this input, the program should produce the following output:

```
Request is to fly from Albuquerque to San Diego.
HPAir flies from Albuquerque to San Diego.
Request is to fly from Albuquerque to Paris.
Sorry. HPAir does not serve Paris.
Request is to fly from San Diego to Chicago.
Sorry. HPAir does not fly from San Diego to Chicago.
```

Begin by implementing the ADT flight map as the C++ class Map. Use the stack version of isPath. Because getNextCity is the primary operation that the search algorithm performs on the flight map, you should choose an implementation that will efficiently determine which cities are adjacent to a given city. If there are $n$ cities numbered $1,2, \ldots, n$, you can use $n$ chains of linked nodes to represent the flight map. You place a node on list $i$ for city $j$ if and only if there is a directed path from city $i$ to city $j$. Such a data structure is called an adjacency list; Figure 6-14 illustrates an adjacency list for the flight map in Figure 6-6. Chapter 20 discusses adjacency lists further when it presents ways to represent graphs. At that time, you will learn why an adjacency list is a good choice for the present program.

To simplify reading the input text files, define a class that includes the following methods:

```
// Returns a name from the next line in a text file.
+getName(): string
// Returns a pair of two names from the next line in a text file.
+getNamePair(): Pair
```

12. In the implementation of the HPAir problem (see the previous programming problem), the search for the next unvisited city adjacent to a city $i$ always starts at the beginning of the $i$ th chain in the adjacency list. This approach is actually a bit inefficient, because once the search visits a city, the city can never become unvisited. Modify the program so that the search for the next city begins where the last search left off. That is, maintain an array of try-next pointers into the adjacency list.

FIGURE 6-14 Adjacency list for the flight map in Figure 6-6 for Programming Problem 11

13. Implement an expanded version of the HPAir problem. In addition to the "from" and "to" cities, each line of input contains a flight number (an integer) and the cost of the flight (an integer). Modify the HPAir program so that it will produce a complete itinerary for each request, including the flight number of each flight, the cost of each flight, and the total cost of the trip.

For example, the input files could appear as

| cityFile: | Albuquerque |  |  |  |
| :--- | :--- | :--- | :--- | :--- |
|  | Chicago |  |  |  |
|  | San Diego |  |  |  |
| flightFile: | Chicago, | San Diego | 703 | 325 |
|  | Chicago, | Albuquerque | 111 | 250 |
|  | Albuquerque, | Chicago | 178 | 250 |
| requestFile: | Albuquerque, | San Diego |  |  |
|  | Albuquerque, | Paris |  |  |
|  | San Diego, | Chicago |  |  |

For this input, the program should produce the following output:

```
Request is to fly from Albuquerque to San Diego.
Flight #178 from Albuquerque to Chicago Cost: $250
F1ight #703 from Chicago to San Diego Cost: $325
Total Cost .............. $575
Request is to fly from Albuquerque to Paris.
Sorry. HPAir does not serve Paris.
Request is to fly from San Diego to Chicago.
Sorry. HPAir does not fly from San Diego to Chicago.
```

When the nonrecursive isPath method finds a sequence of flights from the origin city to the destination city, its stack contains the corresponding path of cities. The stumbling block to reporting this path is that the cities appear in the stack in reverse order; that is, the destination city is at the top of the stack and the origin city is at the bottom. For example, if you use the program to find a path from city $P$ to city $Z$ in Figure 6-6, the final contents of the stack will be $P-W-Y-Z$ with $Z$ on top. You want to display the origin city $P$ first, but it is at the bottom of the stack. If you restrict yourself to the stack operations, the only way that you can write the path in its correct order is first to reverse the stack by popping it onto a temporary stack and then to write the cities as you pop them off the temporary stack. Note that this approach requires that you process each city on the path twice.

Evidently, a stack is not the appropriate ADT for the problem of writing the path of cities in the correct order; the appropriate ADT is a traversable stack. In addition to the standard stack operations isEmpty, push, pop, and peek, a traversable stack includes the operation traverse. The traverse operation begins at one end of the stack and visits each item in the stack until it reaches the other end of the stack. For this project, you want traverse to begin at the bottom of the stack and move toward the top.
14. What modifications to the previous programming problem are required to find a least-cost trip for each request? How can you incorporate time considerations into the problem?

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## C++

## Exceptions

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C ++ Interlude $1 \quad \mathrm{C}++$ Classes

Exceptions provide a mechanism in C++ and other programming languages for interrupting program execution when errors, unusual circumstances, or other events occur. An exception is an object that signals the rest of the program that something unexpected has happened. Our code can react appropriately to the exception based on its type and what the exception can tell us via its methods. We handle the exception when we detect and react to it.

Some exceptions indicate mistakes in your code. By correcting those mistakes, you avoid the exceptions and no longer have to worry about them. In fact, your final code gives no indication that an exception could occur. Furthermore, if your code is entirely correct, an exception will not occur. On the other hand, you can intentionally cause an exception. In fact, the programmers who wrote the code for the $\mathrm{C}++$ Standard Library did so. At the very least, we need to know about exceptions so we can use the methods in the Standard Library.

What should we do when an exception occurs? Should we ever intentionally cause an exception in our own programs, and if so, how would we do so? These are some of the questions that this interlude will answer. This knowledge will be particularly important when we implement the ADT stack in the next chapter.

## C3.1 Background

It would be great if every time a method was called, its preconditions were met. As you have seen, that does not always happen. The client could ask a method to remove an item from a container, but the container might be empty. How do we deal with this unusual or exceptional condition? In our implementations so far, a method returns false to indicate that it is unable to perform its task. This value tells the client that some precondition was not met, or that the method has failed for some other reason. It is then the client's responsibility to check the return value to ensure that the method completed its task before continuing.

Sometimes it is not possible for a method to return a boolean value when an unusual situation prevents it from completing its task. For example, the ADT stack's method peek returns a copy of the top item on the stack. Its prototype is

```
virtual ItemType peek() const = 0;
```

What should this method do if the stack is empty? Returning a boolean value would be possible only if the stack contained boolean values, that is, if ItemType was of type boot. Regardless of the type of data in the stack, the return value will indicate a problem only if it is a special value not contained in the stack. We need a consistent mechanism to let the client know that the method could not perform its task-one that does not depend on the type of data stored in the stack. Before we get into the details of such mechanisms, let's look at another example.

## C3.1.1 A Problem to Solve

Recall from the previous C++ Interludes the video game that we are working on for our friend. Our next task is to create a function that searches for a given string in a number of boxes. The function has three parameters: an array of PlainBox<string> objects, an integer that represents the number of PlainBox objects in the array, and the string to be located in each box. The function will return a copy ${ }^{1}$ of the box that contains the string. Listing C3-1 shows our first attempt at this function.

```
LISTING C3-1 First try at the function findBox
PlainBox<string> findBox(PlainBox<string> boxes[], int size, string target)
{
    int index = 0;
    bool found = false;
    while (!found && (index < size))
    {
        if (target == boxes[index].getItem())
            found = true;
        else
            index++;
    } // end while
    return boxes[index];
} // end findBox
```

[^7]This version of findBox meets the basic requirements: It searches the array of boxes and returns the box containing the target string. We encounter a problem if a box containing the target string is not in the array. If the target is not found, the function's while loop ends when index is equal to size, which is the number of entries in the array. The function would then return boxes [size], which is undefined. Problems will occur when the client tries to use this "box."

We can repair our code by testing found to make sure we found the target before returning a box from the array. Using a simple if statement is insufficient, as we still would have the problem of what to return when the target is not found. The next section gives one way to fix our code.

## C3.2 Assertions

Chapter 3 and Appendix B define an assertion as a statement of truth about some aspect of a program's logic. You can express an assertion either as a comment or by using the $\mathrm{C}++$ function assert. By using assert, you can make assertions about variables and objects in a program. Such assertions are in the form of a boolean expression that should be true at a specific point in the program. If the assertion is false, the assert function halts execution of the program. Assertions can be used to validate method preconditions before trying to execute the body of a function or method.

To use assert in a program, you first must include its header file:

```
#include <cassert>
```

To call the assert function, you provide a boolean condition as its argument:

## assert(someBooleanCondition);

When program execution reaches this statement, the boolean condition-that is, the assertion-is tested. If the assertion is true, the program continues to execute normally. If the assertion is false, the program halts and an error message is displayed.

The assert statement is an easy way to verify a method's preconditions or postconditions. For example, when implementing the stack method peek, we could execute
assert (!isEmpty());
before trying to return the top of the stack.
Our findBox function can test whether the box was found before returning boxes[index] by calling assert:

```
assert(found);
```

If the target is not found, the assertion is false and the program halts. This prevents the program from trying to use a box that does not exist. Listing C3-2 shows the revised findBox function using assertions.

## LISTING C3-2 Revised findBox function with assertions

PlainBox<string> findBox (PlainBox<string> boxes[], int size, string target)
\{
int index $=0$;
bool found = false;

Use an assert statement to test a precondition or postcondition

```
    while (!found && (index < size))
    {
    if (target == boxes[index].getItem())
        found = true;
    else
        index++;
    } // end while
    assert(found); // Verify that there is a box to return
    return boxes[index];
} // end findBox
```

Assertions are good to use when you are testing and debugging your program. They can prevent the execution of any function or method if the preconditions have not been met. Halting your program as soon as a problematic condition becomes evident is a good debugging technique that narrows the focus of your search for errors.

As useful as they are, assertions may not be the best solution for a final shipping program. A violation of a function's preconditions may be a simple mistake that the client can fix. Such a situation does not warrant terminating program execution. We need another way to let the client know that an error or unusual condition has occurred and permit it to fix or handle the situation. We should use exceptions here.


Programming Tip: A statement of truth about some aspect of a program's logic is known as an assertion. You can express an assertion either as a comment or by using the assert function. By including assertions in your program, you facilitate the debugging process.

VideoNote
C++ exceptions

## C3.3 Throwing Exceptions

The examples we've looked at in this C++ Interlude have a common problem - how to let the client know that an error, unusual circumstance, or event occurred during execution. The two solutions presented so far-returning a boolean value or using assert to halt the program-are vastly different techniques.

Returning a boolean value requires that the client check the return value of the function to see whether the function was successful. We have no way to force the client to do so. If the client fails to check the return value, execution could continue even though the function did not complete its task. Even if the client checks the return value and finds that the function was unsuccessful, the client has no information about why the function failed. The client knows only that the function failed.

A function that uses assertions can alleviate the need for the client to check whether it completed successfully. With assertions, the program halts when there is an error, and the client cannot perform any further tasks. For an error that either is unimportant to the client's goals or is simple for the client to fix, assertions are an extreme solution.

However, an alternate way of communicating or returning information to a function's client is to throw an exception. A thrown exception bypasses normal execution, and control immediately returns to the client. The exception can contain information about the error or unusual condition that helps the client resolve the issue and possibly try the function again.

You can throw an exception by executing a throw statement with the following form:
throw ExceptionClass(stringArgument);

FIGURE C3-1 Hierarchy of C++ exception classes


Here ExceptionClass is the type of exception you want to throw, and stringArgument is an argument to the constructor of ExceptionClass that provides a more detailed description of what may have caused the exception. When a throw statement executes, a new object of type ExceptionClass is created, the exception is thrown and propagated back to the point where the function was called, and the statements in the function that follow the throw statement do not execute.

C++ has two families of exceptions that a function can throw. You can select one of these or derive a class of your own. Although the exception class in a throw statement does not need to be derived from a standard C++ exception class, it is a good programming practice to do so. Figure C3-1 shows the family of exception classes provided by C++.

Exceptions derived from runtime_error or one of its derived classes are thrown by errors that are detectable only at runtime, such as division by zero or numeric overflow. If such an exception is thrown, the program will halt, just as it will when an assertion fails. Unlike assertions, however, when a runtime_error exception is thrown, the exception can contain a message that is displayed to the user giving details about the error and the reason the program halted.

Classes derived from the logic_error exception class represent errors in the logic of the program, such as an unmet precondition or the failure to satisfy a postcondition, that you could detect before the program executes. In our findBox example, if we wanted to throw an exception to indicate the failure to find the target string, we could write:

```
if (!found)
    throw logic_error("Target not found in a box!");
```

When we throw exceptions, we first test for an error or unusual condition using an if statement. Then, if one occurred, we use the throw statement to throw an exception containing a message to our client.

To restrict the exceptions that a function or method can throw, you include a throw clause in its header. This also lets the client know that the function might throw an exception. A throw clause consists of the keyword throw followed by a list of exception types separated by commas and enclosed within parentheses. For example, here is the prototype of our findBox function with a throw clause added:

> PlainBox<string> findBox (PlainBox<string> boxes[], int size, string target) throw(logic_error);

Listing C3-3 shows the complete findBox function that throws a logic_error exception. Later in this interlude, we will derive an exception class to handle this specific error. First, however, we will look at how a client can deal with a thrown exception.

Use a throw statement to throw an exception

## LISTING C3-3 Revised findBox function that throws an exception

```
PlainBox<string> findBox(PlainBox<string> boxes[], int size,
    string target) throw(logic_error)
```

\{
int index $=0$;
bool found = false;
while (!found \&\& (index < size))
\{
if (target $==$ boxes[index].getItem())
found = true;
else
index++;
\} // end while
if (! found)
throw logic_error("Target not found in a box!");
return boxes[index];
\} // end findBox

Programming Tip: Including a throw clause in a function specification ensures that the function can throw only those exceptions. An attempt by the function to throw any other exception will result in a runtime error. Omitting a throw clause allows a function to throw any exception.

## Note: Standard exceptions

The header <stdexcept> defines several standard exceptions that can be thrown by the functions defined in the C++ Standard Library. This header contains two kinds of exceptions: runtime exceptions and logic exceptions.

The runtime exceptions are

```
overflow_error
range_error
underflow_error
An arithmetic overflow has occurred A range error has occurred An arithmetic underflow has occurred
```

The logic exceptions are

```
invalid_argument
length_error
out_of_range
```

An invalid argument was used in a function call An attempt was made to create an object that was too large The value of an argument to a function is not in range

## Note: The javadoc tag @throws

A javadoc comment that precedes a method's header should contain a separate line for each exception the method might throw. Each of these lines begins with the tag @throws, and they should be ordered alphabetically by the names of the exceptions.

## Note: When to throw an exception

- If you can resolve the unusual situation in a reasonable manner, your code should do so instead of throwing an exception.
- If several resolutions to an abnormal occurrence are possible, and you want the client to choose one, you should throw an exception.
- If you detect that a programmer has used your method incorrectly, you can throw a runtime exception. Recall that this action will halt program execution.


## C3.4 Handling Exceptions

In the previous section, we looked at how you can cause an exception when an error or unusual condition occurs. In this section, we look at exceptions from the point of view of a client and present how to react to or handle an exception.

Note: You cannot handle an exception of the standard class runtime_error or any of its descendant classes. You can handle only exceptions of the class logic_error and its derived classes.


Note: You should not throw a runtime exception simply so that the client does not have to handle it.

To handle an exception, you write code that consists of two pieces. The first piece, the try block, contains the statements that might cause or throw an exception. The second piece consists of one or more catch blocks that must immediately follow the try block. Each catch block contains code to react to or catch a particular type of exception.

Here is the general syntax for a try block followed by one catch block:

```
try
{
    < statement(s) that might throw an exception >
}
catch (ExceptionClass identifier)
{
    < statement(s) that react to an exception of type ExceptionClass >
}
```

The statements within the try block execute just as they would if the block was not there. If no exception occurs and the try block completes, execution continues with the statement after the catch block. However, if a statement within a try block causes an exception of the type specified in the catch block, the remainder of the try block is abandoned, and execution immediately transfers to the statements in the catch block. The exception now has been caught. The statements in the catch block execute, and upon their completion, execution continues with the statement immediately after the last catch block.

Note: If an exception has no applicable catch block, the function's execution ends and abnormal program termination occurs.

Note: If an exception occurs in the middle of a try block, the destructors of all objects local to that block are called. This ensures that all resources allocated in that block are released, even if the block is not completely executed. In general, you should not use the new operator inside of a try block. An exception will complicate knowing whether memory was allocated and, if it was, whether it was deallocated.

Programming Tip: An important implementation detail of a try block is that it should contain as few statements as possible-ideally, only the statement that could throw an exception. Keeping the try block small reduces the risk of either important code not being executed or allocated memory not being deallocated if an exception is thrown.

The syntax for a catch block resembles that of a function definition. It specifies both the type of exception the catch block will handle and an identifier. This identifier-called a catch block parame-ter-provides a name for the caught exception that can be used within the catch block. Although a catch block is not a function, throwing an exception is like calling a catch block as if it were a function. The catch block parameter represents the actual exception object provided by the throw statement.

The steps taken in the catch block vary from situation to situation and can be as simple as either doing nothing or displaying an error message, or as elaborate as tasks that update variables and retry the function that threw the exception. Every exception class in the C++ Standard Library defines the accessor method what, which returns a descriptive string created when the exception is thrown. The body of the catch block can display this string to provide an indication of the nature of the exception.


Note: A catch block whose parameter has the type ExceptionClass can catch exceptions of the class ExceptionClass and any of ExceptionClass's descendant classes. Specifying a catch block with a parameter of type exception catches any exception that can be thrown.

Listing C3-4 shows sample code that calls our findBox function and catches any logic_error exception thrown. The catch block displays a message to the user that something unusual has happened and creates a PlainBox object for foundBox so that the statement following the catch block executes correctly.

## LISTING C3-4 Trying the function findBox

// Create and initialize an array of boxes
PlainBox<string> myBoxes[5]; // Array of PlainBox objects
myBoxes[0] = PlainBox<string>("ring");
myBoxes[1] = PlainBox<string>("hat");
myBoxes[2] = PlainBox<string>("shirt");
myBoxes[3] = P1ainBox<string>("sock");
myBoxes[4] = PlainBox<string>("shoe");
PlainBox<string> foundBox;

```
// Try to find a box containing glasses
try
{
    foundBox = findBox(myBoxes, 5, "glasses");
}
catch(logic_error logErr)
{
    cout << logErr.what() << endl; // Display error message to user
    foundBox = PlainBox<string>("nothing"); // Fix prob7em
} // end try-catch
// Because we catch the exception and fix the problem, the following
// statement should work even if the target is not found
cout << foundBox.getItem();
```


## Output

Target not found in a box! nothing

## C3.4.1 Multiple catch Blocks

A single statement within a try block might cause more than one type of exception. Thus, a try block can have many catch blocks associated with it. When more than one catch block follows a try block, the catch block that executes is the first one-in order of appearance-whose parameter matches the thrown exception in type. Thus, the catch blocks must be ordered so that the most specific exception classes are caught before the more general exception classes. For example, the following catch blocks are in the wrong order:

```
string str = "Sarah";
try
{
    str.substr(99, 1);
    < other statements appear here . . >
}
catch (exception e) // WRONG ORDER!!
{
    cout << "Something else was caught" << endl;
}
catch (out_of_range e)
{
    cout << "out_of_range exception caught" << end1;
} // end try-catch
```

The compiler may issue a warning message similar to the following one:

```
'class std::out_of_range' : is caught by base class ('class exception')
on line n
```

This warning means that any out_of_range exception thrown in the try block will be caught by the first catch block-with the parameter of type exception-so the second catch block will never be reached. To get the code to compile without warnings, you must interchange the two catch blocks.

$!$
Programming Tip: Arrange catch blocks in order of specificity, catching the most specific one first. Since all exception classes have exception as an ancestor, try to avoid using exception in a catch block. If you must catch exceptions of type exception, do so in the last catch block.

## C3.4.2 Uncaught Exceptions

Let's see what happens when an exception is thrown but not caught. The program in Listing C3-5 encodes a string by doing a simple substitution. It replaces each letter in a string with the character that appears three positions later in the alphabet. When it reaches the end of the alphabet, it wraps around to the beginning. For example, $a$ is replaced by $d, b$ is replaced by $e$, and $w$ is replaced by $z$. At this point, it gets replacement letters by wrapping around to the beginning of the alphabet. Thus, $x$ is replaced by $a, y$ is replaced by $b$, and $z$ is replaced by $c$.

Figure C3-2 shows the flow of control when an exception occurs in this code. When the function encodeChar attempts to access the 99th character in str by the method call str. replace (99, 1, 1, newChar), an out_of_range exception is thrown. Because encodeChar does not handle the

## LISTING C3-5 A program with an uncaught exception

```
#include <iostream>
#include <string>
using namespace std;
// Encodes the character at index i of the string str.
void encodeChar(int i, string& str)
{
    int base = static_cast<int>('a');
    if (isupper(str[i]))
        base = int('A');
    char newChar = (static_cast<int>(str[i]) - base + 3) % 26 + base;
    str.replace(i, 1, 1, newChar); // Method replace can throw exception
} // end encodeChar
// Encodes numChar characters within a string.
void encodeString(int numChar, string& str)
{
    for (int j = numChar - 1; j >= 0; j-)
        encodeChar(j, str);
} // end encodeString
int main()
{
    string str1 = "Sarah";
    encodeString(99, str1);
    return 0;
} // end main
```

FIGURE C3-2 Flow of control for an uncaught exception

exception, the function terminates, and the exception propagates back to encodeString at the point where encodeChar was called. The function encodeString also does not handle the exception, so it too terminates and the exception propagates back to main. Because main is the main function of the program, it does not handle the exception, and the program execution terminates abnormally with an error message.

This code contains no indication that the function encodeChar could throw the exception out_ of_range. However, the documentation of encodeChar should indicate any exceptions it might throw. Thus, when documenting a function or method, you should also include any possible exceptions that could occur.

We can catch the exception out_of_range at any point in the sequence of function calls. For example, we could rewrite the function encodeChar as follows to catch the exception.

```
void encodeChar(int i, string& str)
{
    int base = static_cast<int>('a');
    if (isupper(str[i]))
        base = int('A');
```

```
    char newChar = (static_cast<int>(str[i]) - base + 3) % 26 + base;
    try
    {
        str.replace(i, 1, 1, newChar);
    }
    catch (out_of_range e)
    {
    cout << "No character at position " << i << endl;
} // end try-catch
} // end encodeChar
```

The call encodeString (99, str) within main causes 99 calls to encodeChar, and hence the exception is thrown 98 - str. length() times. The following output from encodeChar is the result.

```
No character at position 99
No character at position 98
No character at position 97
```

When the exception was not handled, the program terminated the first time the exception occurred. Handling the exception allows the code to continue execution.

Although the out_of_range exception is thrown in the function encodeChar, that is not necessarily the best place to handle the exception. For example, if the client had made the call encodeString (10000, str), the message printed by encodeChar would have appeared 9999 str. length() times! In this case, it makes more sense for the try and catch blocks to appear in the function encodeString and not in encodeChar, as follows:

```
void encodeString(int numChar, string& str)
{
    try
    {
        for (int i = numChar - 1; i >= 0; i--)
            encodeChar(i, str);
    }
    catch (out_of_range e)
    {
        cout << "The string does not contain " << numChar;
        cout << " characters." << endl;
        cout << e.what() << endl;
    } // end try-catch
} // end encodeString
```

Now when encodeChar throws the exception out_of_range, the exception propagates back to the function encodeString, which ends execution of the statements in the try block and executes the statements in the catch block. The expression e.what () invokes the method what for the exception e, and so it represents a string that describes the exception. For this example, the following output would be displayed.

```
The string does not contain 10 characters.
invalid string position
```

The message is printed only once, since the for loop is inside the try block, which is abandoned when the exception occurs. If the try block had been placed inside the for loop, encodeChar would be called-and therefore the exception would be thrown and handled-at each iteration of the loop. The result would be multiple copies of the message.

Programming Tip: When an exception can occur within a loop, you must decide whether to place a try block within the body of the loop-in which case an exception will be caught each time the loop cycles (as in the first example) - or place the entire loop within a try block-in which case the exception would be caught only once (as in the second example) Neither choice is better than the other, but if you make the wrong choice, the client likely will not get the necessary information.

Note: Sometimes a catch block will intentionally terminate program execution. When the block does not do so, but rather completes its execution, any statements that appear after the last catch block execute next.

## C3.5 Programmer-Defined Exception Classes

You may find that the C++ Standard Library has an exception class already defined that suits the exception needs of your program. You may also want to define your own exception class. Usually, the C++ exception class exception, or one of its derived classes, is the base class for programmerdefined exception classes. Adhering to this practice provides a standardized interface for working with exceptions. If you follow this practice, you will need to use the std namespace.

An exception class typically consists of a constructor that has a string parameter. For example, you can define the class TargetNotFoundException as

```
#include <stdexcept>
#include <string>
using namespace std;
class TargetNotFoundException: public exception
{
public :
    TargetNotFoundException(const string& message = "")
                        : exception("Target not found: " + message.c_str())
    {
    } // end constructor
}; // end TargetNotFoundException
```

The constructor provides a way for a throw statement to identify the condition that caused the exception. For example, the statement

```
throw TargetNotFoundException (target + " not found in a box!");
```

invokes the constructor of TargetNotFoundException. The message given to the constructor is returned by the method what that is inherited from the class exception. Thus, a catch block, such as the following one, can access the message:

```
catch(TargetNotFoundException except)
{
    cout << except.what() << endl;
}
```

If target has the value " $g$ lasses" when this block executes, the output is
Target not found: glasses not found in a box!

While the string message is helpful, sometimes we want to return additional information to the client to help resolve the problem. We could add a private data field to our exception class to store additional error details. A second parameter in the constructor to receive an initial value for those details, as well as a public accessor method so that the client can retrieve them, are also needed. When a method throws this exception, it would call the constructor and send it the extra data and the message as its two arguments.

## Programming Tip: Using exceptions

To throw an exception in a function:

- Determine what errors, preconditions, postconditions, and unusual conditions you must check and where they should be checked, so your function can correctly perform its task.
- Detect the error, precondition, postcondition, or unusual condition. Usually this is done with an if statement.
- If an error condition has occurred, throw an exception using the throw statement. As an argument to the exception class constructor, pass a string description of the reason for the error.
- Add a throw clause to the function header indicating the exceptions thrown by the function.
- Derive a custom exception class to better identify the conditions that caused the error. Although doing so is optional, it is desirable.

To handle an exception or use a function that could throw an exception

- Place the statement that might throw an exception in a try block. The fewer statements in the try block the better. Do not use new to create an object in a try block.
- After the try block, place catch blocks for each type of exception that can be thrown by the statements in the try block. Place the catch blocks in order from the most specific exception classes to the more general exception classes.


## Implementations of the ADT Stack

## Chapter

## 7

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| C++ Interlude 2 | Pointers, Polymorphism, and Memory Allocation |
| Chapter 4 | Link-Based Implementations |
| Chapter 6 | Stacks |
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This chapter implements the ADT stack using the techniques we described in Chapters 3 and 4 for the ADT bag. We first will use an array and then a chain of linked nodes to store the stack's entries. You should be happy to discover the simplicity and efficiency of these implementations. We then will modify these implementations to use exceptions as a way for a method to signal an unusual occurrence, such as an attempt to get an entry from an empty stack.

VideoNote
Overview of ArrayStack

### 7.1 An Array-Based Implementation

Figure 7-1a shows a sketch of an array that contains the entries in a stack. Where in the array should the top entry be? Since entries are added or removed at the top of the stack, we can avoid shifting the current entries if we anchor the bottom of the stack at index 0 and track the location of the stack's top using an index top. Then if items is the array and items [top] is the top entry, items and top can be the private data members of our class of stacks. Figure 7-1b illustrates these details. Notice that the top of the stack in this figure is at index 2 . We can either add a new entry at index[3] and then increment top or remove the top entry in index [2] by decrementing top to 1.

Our class will have a default constructor to initialize the private data members items and top. However, since we plan to store a stack's entries in statically allocated memory, the compilergenerated destructor and copy constructor will be sufficient. If, on the other hand, we were to use a dynamically allocated array, we would have to define a destructor and a copy constructor.

Listing 7-1 shows the header file for an array-based implementation of the ADT stack. Comments describing the methods are in Listing 6-1 of Chapter 6.

FIGURE 7-1 Using an array to store a stack's entries: (a) a preliminary sketch; (b) implementation details


## LISTING 7-1 The header file for an array-based stack

```
/** ADT stack: Array-based implementation.
    @file ArrayStack.h */
#ifndef _ARRAY_STACK
#define _ARRAY_STACK
#include "StackInterface.h"
const int MAX_STACK = maximum-size-of-stack;
template<class ItemType>
class ArrayStack : public StackInterface<ItemType>
{
private:
    ItemType items[MAX_STACK]; // Array of stack items
    int top; // Index to top of stack
```

```
public:
    ArrayStack(); // Default constructor
    bool isEmpty() const;
    bool push(const ItemType& newEntry);
    bool pop();
    ItemType peek() const;
}; // end ArrayStack
#include "ArrayStack.cpp"
#endif
```

The implementations of the methods that the previous header file declares are given in Listing 7-2. These definitions are straightforward, but notice that the method peek uses assert to enforce its precondition that the stack contain at least one entry. Later in this chapter, we will revise peek to throw an exception if the client calls it when the stack is empty.

## LISTING 7-2 The implementation file for an array-based stack

```
/** @file ArrayStack.cpp */
#include <cassert> // For assert
#include "ArrayStack.h" // Header file
template<class ItemType>
ArrayStack<ItemType>::ArrayStack() : top(-1)
{
} // end default constructor
// Copy constructor and destructor are supplied by the compiler
temp1ate<c1ass ItemType>
bool ArrayStack<ItemType>::isEmpty() const
{
    return top < 0;
} // end isEmpty
temp1ate<class ItemType>
boo1 ArrayStack<ItemType>::push(const ItemType& newEntry)
{
    bool result = false;
    if (top < MAX_STACK - 1) // Does stack have room for newEntry?
    {
        top++;
        items[top] = newEntry;
        result = true;
    } // end if
    return result;
} // end push
temp1ate<c1ass ItemType>
bool ArrayStack<ItemType>::pop()
```

Private data members are hidden from the client

```
{
    bool result = false;
    if (!isEmpty())
    {
        top--;
        result = true;
    } // end if
    return result;
} // end pop
template<class ItemType>
ItemType ArrayStack<ItemType>::peek() const
{
    assert (!isEmpty()); // Enforce precondition
    // Stack is not empty; return top
    return items[top];
} // end peek
// end of implementation file
```

A program that uses a stack could begin as follows:

```
#include <iostream>
#include <string>
#include "ArrayStack.h"
using namespace std;
int main()
{
    StackInterface<string>* stackPtr = new ArrayStack<string>();
    string anItem = "";
    cout << "Enter a string: ";
    cin >> anItem; // Read an item
    stackPtr->push(anItem); // Push item onto stack
    . . .
```

By implementing the stack as a class, and by declaring items and top as private, you ensure that the client cannot violate the ADT's walls. If you did not hide your implementation within a class, or if you made the array items public, the client could access the entries in items directly instead of by using the ADT stack operations. Thus, the client could access any entries in the stack, not just its top entry. You might find this capability attractive, but in fact it violates the specifications of the ADT stack. If you truly need to access all the items of your ADT randomly, do not use a stack!

Finally, note that push receives newEntry as a constant reference argument. Therefore, push uses newEntry as an alias to its actual argument, and no copy is made.

Question 1 In Chapter 6, the algorithms that appear in Section 6.2 involve strings. Under what conditions would you choose an array-based implementation for the stack in these algorithms?

Question 2 Describe the changes to the previous stack implementation that are necessary to replace the fixed-size array with a resizable array.

### 7.2 A Link-Based implementation

Many applications require a link-based implementation of a stack so that the stack can grow and shrink dynamically. Figure 7-2 illustrates such an implementation, where topPtr is a pointer to the head of the linked nodes.

As we did for the array-based implementation, we begin with a header file based upon StackInterface, as you can see in Listing 7-3. The pointer topPtr points to an instance of the class Node that we used in the implementation of the class LinkedBag in Chapter 4.

Because memory is allocated dynamically for the nodes, you must write both a copy constructor and a virtual destructor for the link-based stack implementation. As you saw in Chapter 4, if a shallow copy of the stack is sufficient, you can omit the copy constructor, in which case the compiler generates a copy constructor that performs a shallow copy. Such was the case for the array-based implementation.

If we did not write our own copy constructor for the LinkedStack class, the compiler-generated copy constructor would copy only the pointer topPtr. Thus, both topPtr and its copy would point to the same linked nodes; the stack itself would not be copied. You must write a copy constructor that explicitly makes a copy of all the nodes in a linked stack.

FIGURE 7-2 A link-based implementation of a stack


You need an explicit copy constructor and a virtual destructor

## LISTING 7-3 The header file for the class LinkedStack

```
/** ADT stack: Link-based implementation.
@file LinkedStack.h */
#ifndef _LINKED_STACK
#define _LINKED_STACK
#include "StackInterface.h"
#include "Node.h"
template<class ItemType>
class LinkedStack : public StackInterface<ItemType>
{
private:
    Node<ItemType>* topPtr; // Pointer to first node in the chain;
    // this node contains the stack's top
public:
// Constructors and destructor:
    LinkedStack(); // Defau7t constructor
    LinkedStack(const LinkedStack<ItemType>& aStack); // Copy constructor
    virtual ~LinkedStack(); // Destructor
// Stack operations:
    bool isEmpty() const;
    bool push(const ItemType& newItem);
    bool pop();
    ItemType peek() const;
}; // end LinkedStack
#include "LinkedStack.cpp"
#endif
```

Listing 7-4 gives the implementation file for LinkedStack. Note the similarity between the copy constructors of LinkedStack and LinkedBag as given in Chapter 4.

## LISTING 7-4 The implementation file for the class LinkedStack

```
/** afile LinkedStack.cpp */
#include <cassert> // For assert
#include "LinkedStack.h" // Header file
temp1ate<c1ass ItemType>
LinkedStack<ItemType>::LinkedStack() : topPtr(nu11ptr)
{
} // end default constructor
template<class ItemType>
LinkedStack<ItemType>::
LinkedStack(const LinkedStack<ItemType>& aStack)
{
```

```
    // Point to nodes in original chain
    Node<ItemType>* origChainPtr = aStack->topPtr;
    if (origChainPtr == nul1ptr)
        topPtr = nul1ptr; // Original bag is empty
    else
    {
        // Copy first node
        topPtr = new Node<ItemType> ();
        topPtr->setItem(origChainPtr->getItem());
    // Point to first node in new chain
    Node<ItemType>* newChainPtr = topPtr;
    // Copy remaining nodes
    while (origChainPtr != nul1ptr)
    {
            // Advance original-chain pointer
            origChainPtr = origChainPtr->getNext();
            // Get next item from original chain
            ItemType nextItem = origChainPtr->getItem();
            // Create a new node containing the next item
            Node<ItemType>* newNodePtr = new Node<ItemType>(nextItem);
            // Link new node to end of new chain
            newChainPtr->setNext(newNodePtr);
            // Advance pointer to new last node
            newChainPtr = newChainPtr->getNext();
        } // end while
        newChainPtr->setNext(nul1ptr); // Flag end of chain
    } // end if
} // end copy constructor
temp1ate<c1ass ItemType>
LinkedStack<ItemType>::~LinkedStack()
{
    // Pop until stack is empty
    while (!isEmpty())
        pop();
} // end destructor
template<class ItemType>
bool LinkedStack<ItemType>::isEmpty() const
{
    return topPtr == nul1ptr;
} // end isEmpty
temp1ate<c1ass ItemType>
bool LinkedStack<ItemType>::push(const ItemType& newItem)
{
    Node<ItemType>* newNodePtr = new Node<ItemType>(newItem, topPtr);
    topPtr = newNodePtr;
    newNodePtr = nul1ptr;
```

```
    return true;
} // end push
template<class ItemType>
bool LinkedStack<ItemType>::pop()
{
    bool result = false;
    if (!isEmpty())
    {
            // Stack is not empty; delete top
            Node<ItemType>* nodeToDe1etePtr = topPtr;
            topPtr = topPtr->getNext();
            // Return deTeted node to system
            nodeToDeletePtr->setNext(nul1ptr);
            delete nodeToDeletePtr;
            nodeToDeletePtr = nul1ptr;
            result = true;
    } // end if
    return result;
} // end pop
temp1ate<class ItemType>
ItemType LinkedStack<ItemType>::peek() const
{
    assert(!isEmpty()); // Enforce precondition
    // Stack is not empty; return top
    return topPtr->getItem();
} // end getTop
// end of implementation file
```

Question 3 In Chapter 6, the algorithms that appear in Section 6.2 involve strings. Under what conditions would you choose a link-based implementation?

## Note: Comparing implementations

You have seen two implementations of the ADT stack, one that uses an array and another that uses linked nodes to contain the items in a stack. The discussion in Section 4.5 of Chapter 4 applies here as well. The array-based implementation is a reasonable choice if the number of items in the stack does not exceed the fixed size of the array. For example, when we read and correct an input line, if the system allows a line length of only 80 characters, you reasonably could use a statically allocated array to represent the stack. For stacks that might be large, but often are not, the array-based implementation will waste storage. In that case, the link-based implementation is a better choice.

## Programming Tip: Initializers versus set methods for constructors

Chapter 3 suggested that you use initializers instead of assignment statements in constructors to initialize the values of a class's data members. That is still good advice, and initializers clearly show the initialization. However, if you write set methods to alter the values of the data members, you have the opportunity to ensure that the values assigned are valid. Doing so is especially important when those values are supplied by the client. In such cases, constructors can call the set methods to initialize the data members. Thus, we have the following advice when you initialize the values of data members in constructors:

- Use an initializer if the value of the data member has no restrictions.
- Use a set method if the value of the data member needs validation.

Note that even if you do not want a public set method for a particular data member, you can either define a private one for the constructor to call or have the constructor perform its own validation.

### 7.3 Implementations That Use Exceptions

Our specifications of the ADT stack appear to handle all eventualities. If you try to add an item to a full stack, the method push returns false. If you try to remove an entry from an empty stack, the method pop returns false. What happens when you try to look at the top of an empty stack? The method peek expects that you will not do this, as stated in its precondition. Our implementations, however, deal with the misuse of this method by beginning its definition with the statement

```
assert(!isEmpty());
```

If the stack is empty, assert will issue an error message and halt execution.
Instead of using assert, peek could throw an exception if its precondition is violated-that is, if it is called when the stack is empty. Let's define the class PrecondViolatedExcep that peek can use. Listings 7-5 and 7-6 show this class.

```
LISTING 7-5 The header file for the class PrecondViolatedExcep
/** @file PrecondViolatedExcep.h */
#ifndef _PRECOND_VIOLATED_EXCEP
#define _PRECOND_VIOLATED_EXCEP
#include <stdexcept>
#include <string>
using namespace std;
class PrecondViolatedExcep: public logic_error
{
public:
    PrecondViolatedExcep(const string& message = "");
}; // end PrecondViolatedExcep
#endif
```


## LISTING 7-6 The implementation file for the class PrecondViolatedExcep

```
/** afile PrecondViolatedExcep.cpp */
#include "PrecondViolatedExcep.h"
PrecondViolatedExcep::PrecondViolatedExcep(const string& message):
    logic_error("Precondition Violated Exception: " + message)
{
} // end constructor
```

We now revise the declarations of peek in both of the header files ArrayStack.h and LinkedStack. h by adding a throw clause as follows:

```
ItemType peek() const throw(PrecondViolatedExcep);
```

In the implementation files, we add the same throw clause to the header of peek and then replace the assert statement in both definitions of peek with the following if statement:

```
if (isEmpty())
    throw PrecondViolatedExcep("peek() called with empty stack");
```

Thus, the definition of peek in LinkedStack, for example, is

```
temp1ate<class ItemType>
ItemType LinkedStack<ItemType>::peek() const throw(PrecondViolatedExcep)
{
    // Enforce precondition
    if (isEmpty())
        throw PrecondViolatedExcep("peek() called with empty stack");
    // Stack is not empty; return top
    return topPtr->getItem();
} // end getTop
```

The previous revisions to peek are the same for both the array-based and link-based implementations of the stack.

Note: In the link-based implementation of the stack, both the method push and the copy constructor allocate new nodes. This allocation could fail if the system has no spare memory available. In such a case, the standard exception bad_alloc is thrown. However, in our simple examples, such an occurrence is unlikely to occur unless your computing system has a severely restricted memory size. For simplicity, we will not worry about this situation.

Question 4 Define the exception class MemoryAllocationException and then revise the definition of the method push in the class LinkedStack so that it throws this exception if it cannot allocate a new node.

## Summary

1. You can implement a stack by using an array. If the bottom of the stack is in the first element of the array, no stack entries are moved when you add or remove entries.
2. You can implement a stack by using a chain of linked nodes that has a head pointer. The first node of the chain should contain the top of the stack to provide the easiest and fastest addition and removal operations.
3. You should call assert or throw an exception to enforce the precondition for the method peek.

## Exercises

1. Discuss the advantages and disadvantages of an array-based implementation of the ADT stack as compared to a link-based implementation.
2. Consider the ADT bag, as described in Chapters 1,3 , and 4 .
a. Would you be able to implement the ADT stack by using a bag to contain its entries? Why, or why not?
b. Would you be able to implement the ADT bag by using a stack to contain its entries? Why, or why not?
3. An operation that displays the contents of a stack can be useful during program debugging. Add a display method to the ADT stack such that
a. The method uses only ADT stack operations; that is, it is independent of the stack's implementation.
b. The method assumes and uses the link-based implementation of the ADT stack.
c. The method assumes and uses the array-based implementation of the ADT stack.
4. Repeat the previous exercise, but define the method toVector instead of the method display.
5. To the ADT stack given in this chapter, add a void method remove $(\mathrm{n})$ that removes and discards the topmost n entries from a stack. Write a link-based implementation for this method.
6. Repeat the previous exercise, but write an array-based implementation instead.
7. The destructor given for the link-based implementation of the ADT stack calls pop. Although easy to write, this destructor can be inefficient due to repeated method calls. Write another implementation for the destructor that deallocates the linked nodes directly without calling pop.
8. Implement a memory-safe copy constructor for the linked stack. If a memory allocation fails, this constructor should release all memory that was allocated prior to the failure and then throw an exception.
9. Imagine an array-based implementation of the ADT stack that stores the stack's entries beginning at the end of an array. Describe how you can define the stack operations so that the push and pop operations do not require you to move existing entries in the array.
10. Imagine a link-based implementation of the ADT stack that stores the stack's top entry at the end of a chain of linked nodes. Describe how you can define the stack operations so that a traversal of the chain is not necessary.
11. Although we do not usually throw our own exception when a method or constructor in LinkedStack fails to allocate a new node, you could do so. Define the exception MemoryAllocationException, and revise the copy constructor of LinkedStack so that it throws a MemoryAllocationException when the new operator fails to allocate memory. Throwing exceptions from constructors requires great care, since you must be sure to properly deallocate any memory allocated by the constructor before throwing the exception.
Test your revised version of LinkedStack.

## Programming Problems

1. Write an implementation of the ADT stack that uses a resizable array to represent the stack items. Anytime the stack becomes full, double the size of the array. Maintain the stack's bottom entry at the beginning of the array.
2. Repeat Programming Problem 1, but maintain the stack's bottom entry at the end of the array.
3. Repeat Programming Problem 1, but maintain the stack's top entry at the beginning of the array.
4. Repeat Programming Problem 1, but maintain the stack's top entry at the end of the array.
5. Repeat any of the previous four programming problems, but after the first doubling of the array, halve the size of the array if fewer than half of the array's locations are occupied by current stack entries.
6. Suppose that instead of doubling the size of an array-based stack when it becomes full, you increase the size of the array by some positive integer $k$. Implement and demonstrate such a stack that allows the client to specify $k$ when the stack is created.
7. Repeat the previous programming problem, but each time the stack becomes full, increase the size of the array by the next value in this sequence: $3 k, 5 k, 7 k, \ldots$ for a client-specified positive integer $k$.
8. Write the implementation of the ADT stack that Exercise 10 describes.
9. The ADT stack lets you peek at its top entry without removing it. For some applications of a stack, you need to also peek at the entry beneath the top entry without removing it. Let's name such an operation peek2. If peek2 fails because the stack contains fewer than two entries, it should throw an exception. Write a link-based implementation of the ADT stack that includes both peek and peek2.
10. Repeat any of the programming problems in Chapter 6, except the first one, using an implementation of the ADT stack that this chapter describes.

## Lists

## Chapter

 8
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Programming Problems 264

## Prerequisites

Chapter 1 Data Abstraction: The Walls<br>$\mathrm{C}++$ Interlude $1 \quad \mathrm{C}++$ Classes

Ann everyday list provides a way for us to organize our thoughts, plans, or tasks. Each list has a first entry, a last entry, and often entries in between. The entries have a position within the list, and therefore they are ordered. In computer science, an ADT list provides a way to organize data. Like an everyday list, the entries in an ADT list have a position and are ordered. This order is not determined by the list itself, but rather by its client.

This chapter specifies the ADT list and gives some examples of how to use it.

VideoNote The ADT list

### 8.1 Specifying the ADT List

Consider a list that you might encounter, such as a list of chores, a list of important dates, a list of addresses, or the grocery list pictured in Figure 8-1. As you write a grocery list, where do you put new items? Assuming that you write a neat one-column list, you probably add new items to the end of the list. You could just as well add items to the beginning of the list or add them so that your list is sorted alphabetically. Regardless, the items on a list appear in a sequence. The list has one first item and one last item. Except for the first and last items, each item has a unique predecessor and a unique successor. The first item - the head or front of the list - does not have a predecessor, and the last item - the tail or end of the list-does not have a successor.

Although the six items on the list in Figure 8-1 have a sequential order, they are not necessarily sorted by name. Perhaps the items appear in the order in which they occur on the grocer's shelves, but more likely they appear in the order in which they occurred to you as you wrote the list.

Lists contain items of the same type: You can have a list of grocery items or a list of phone numbers. What can you do to the items on a list? You might count the items on the list, add an item to the list, remove an item from the list, or look at (retrieve) an item. The items on a list, together with operations that you can perform on the items, form an ADT: the ADT list.

Note: The ADT list is simply a container of items whose order you indicate and whose position you reference by number.

You reference list items by their position

You must specify the behavior of the list's operations without thinking about how you could implement them. For example, you need to decide where to add a new item and which item to retrieve or remove. The various possible decisions lead to several kinds of lists. You might decide to add, remove, and retrieve items only at the end of the list, only at the front, or at both the front and the end. The specifications of these lists, which manipulate items at one or both ends, are left as an exercise, as they are not really adequate for an actual grocery list. Instead we will discuss a more general list, one that allows you to access items anywhere on the list. That is, you might look at the item at position $i$, remove the item at position $i$, or insert an item at position $i$ on the list. Such operations are part of the ADT list. Other operations that see whether a list is empty or return the length of the list are also useful.

FIGURE 8-1
A grocery list


## Note: ADT list operations

- Test whether a list is empty.
- Get the number of entries on a list.
- Insert an entry at a given position on the list.
- Remove the entry at a given position from the list.
- Remove all entries from the list.
- Look at (get) the entry at a given position on the list.
- Replace (set) the entry at a given position on the list.

Figure 8-2 shows the UML diagram for this ADT and provides more detail for its operations.

## FIGURE 8-2 UML diagram for the ADT list

| List |
| :--- |
|  |
| +isEmpty(): boolean |
| +getLength(): integer |
| +insert(newPosition: integer, newEntry: ItemType) : boolean |
| +remove(position: integer): boolean |
| +clear(): void |
| +getEntry(position: integer): ItemType |
| +setEntry(position: integer, newEntry: ItemType): void |

To get a more precise idea of how the operations work, let's apply them to the following grocery items:

milk, eggs, butter, apples, bread, chicken

Milk will be the first item on the list, and chicken will be the last item. To begin, consider how you can construct this list by using the ADT list operations. One way is first to create an empty list aList and then use a series of insertion operations to append the items to the list one at a time, as follows:

```
aList = a new emptylist
aList.insert(1, milk)
aList.insert(2, eggs)
aList.insert(3, butter)
aList.insert(4, app1es)
aList.insert(5, bread)
aList.insert(6, chicken)
```

We assume for this simple example that the list can contain all of the items we add to it.
In the previous example, we in effect have inserted each new item at the end of the list. Nevertheless, the list's insertion operation can place new items into any position of the list, not just at its front or end. The effect of an insertion between existing items, however, is not apparent from the previous example. For instance, if you start with the previous grocery list and you perform the operation

```
aList.insert(4, nuts)
```

the list aList should become
milk, eggs, butter, nuts, apples, bread, chicken
All items that had position numbers greater than or equal to 4 before the insertion now have their position numbers increased by 1 after the insertion. In general, if a new item is inserted into position $i$, the position of each item that was at or after position $i$ is increased by 1 .

Similarly, the removal operation specifies that if an item is deleted from position $i$, the position of each item that was at a position greater than $i$ is decreased by 1 . Thus, for example, if aList is the list
milk, eggs, butter, nuts, apples, bread, chicken
and you perform the operation
aList. remove(5)
the list becomes
milk, eggs, butter, nuts, bread, chicken
All items that had position numbers greater than 5 before the removal now have their position numbers decreased by 1 after the deletion.

These examples illustrate that we can specify the effects of an ADT's operations without having to indicate how to store the data. The following summary of the ADT list provides more details for its operation contract. As usual, the specifications of these operations are the sole terms of the operation contract for the ADT list. The behavior of the operations is the only thing on which a program should depend.

## Abstract Data Type: List

Data

- A finite number of objects, not necessarily distinct, having the same data type and ordered by their positions, as determined by the client.

Operations

| Pseudocode | Description |
| :---: | :---: |
| isEmpty () | Task: Sees whether this list is empty. |
|  | Input: None. |
|  | Output: True if the list is empty; otherwise false. |
| getLength() | Task: Gets the current number of entries in this list. |
|  | Input: None. |
|  | Output: The integer number of entries currently in the list. |
| insert(newPosition, newEntry) | Task: Inserts an entry into this list at a given position. An insertion before existing entries causes the renumbering of entries that follow the new one. |
|  | Input: newPosition is an integer indicating the position of the insertion, and newEntry is the new entry. |
|  | Output: True if $1 \leq$ newPosition $\leq$ getLength () +1 and the insertion is successful; otherwise false. |


| remove(position) | Task: Removes the entry at a given position from this list. A removal before the last entry causes the renumbering of entries that follow the deleted one. <br> Input: position is the position of the entry to remove. <br> Output: True if $1 \leq$ newPosition $\leq$ getLength() and the removal is successful; otherwise false. |
| :---: | :---: |
| clear () | Task: Removes all entries from this list. Input: None. <br> Output: None. The list is empty. |
| getEntry (position) | Task: Gets the entry at the given position in this list. <br> Input: position is the position of the entry to get; $1<=$ position $<=$ getLength(). <br> Output: The desired entry. |
| setEntry (position, newEntry) | Task: Replaces the entry at the given position in this list. <br> Input: position is the position of the entry to replace; $1<=$ position $<=$ getLength(). newEntry is the replacement entry. <br> Output: None. The indicated entry is replaced. |

Note that the return values of the insertion and removal operations provide the ADT with a simple mechanism to communicate operation failure to its client. For example, if you try to remove the tenth item from a five-item list, remove will return false. Likewise, insert will return false if, for example, the list is full or position is out of range. In this way, the client can deal with error situations in an implementation-independent way.


Note: The list operations fall into the three broad categories presented earlier in this book:

- The operation insert adds data to a data collection.
- The operation remove removes data from a data collection.
- The operations isEmpty, getLength, and getEntry ask questions about the data in a data collection.

The operation setEntry replaces existing data in a data collection, so you can think of it as removing and then adding data.

Question 1 The specifications of the ADT list do not mention the case in which two or more items have the same value. Are these specifications sufficient to cover this case, or must they be revised?

Question 2 Write specifications for a list whose operations insert, remove, getEntry, and setEntry always act at the end of the list.

Axioms (optional). The previous specifications for the operations of the ADT list have been stated rather informally. For example, they rely on your knowing the meaning of "an item is at position $i$ " in

Use axioms to determine the effect of a sequence of ADT operations
a list. However, you can write a set of axioms that completely describes the behavior of the list operations. For example,

## A newly created list is empty

is an axiom because it is true for all newly created lists. You can state this axiom succinctly in terms of the ADT list operations as follows: ${ }^{1}$
(new List()).isEmpty() = true
The statement
If you insert an item $x$ into the $i^{\text {th }}$ position of a list, retrieving the $i^{\text {th }}$ item will result in $x$
is true for all lists, and so it is an axiom. You can state this axiom in terms of the ADT list operations as follows:

```
(aList.insert(i, x)).getEntry(i) = x
```

That is, getEntry returns the item x that insert has put at position i of aList.
The following axioms formally define the ADT list:


## Note: Axioms for the ADT list

1. (new List()).isEmpty() = true
2. (new List()).getLength() $=0$
3. aList.getLength() $=($ aList.insert(i, $x)$ ).getLength() -1
4. aList.getLength() $=($ aList. remove(i)). getLength() +1
5. (aList.insert(i, item)).isEmpty() = false
6. (new List()).remove(i) = false
7. (aList.insert(i, x)).remove(i) = aList
8. (new List()).getEntry(i) = error
9. (aList.insert(i, x)).getEntry(i) $=x$
10. aList.getEntry (i) $=$ (aList.insert $(i, x)$ ).getEntry $(i+1)$
11. aList.getEntry $(i+1)=(a L i s t . \operatorname{remove}(i))$.getEntry (i)
12. (new List()).setEntry (i, x) = error
13. (aList.setEntry (i, x)).getEntry (i) $=x$

A set of axioms does not make the preconditions and postconditions for an ADT's operations unnecessary. For example, the previous axioms do not describe insert's behavior when you try to insert an item into position 50 of a list of two items. One way to handle this situation is to include the restriction
$1<=$ position <= getLength() +1
in insert's precondition. Another way-which we used in our previous specifications of the ADT list-does not restrict position, but rather has the method return false if position is outside the previous range. Thus, you need both a set of axioms and a set of preconditions and postconditions to define the behavior of an ADT's operations completely.

1 The $=$ notation within these axioms denotes algebraic equality.

You can use axioms to determine the outcome of a sequence of ADT operations. For example, if aList is a list of strings and $s$ and $t$ are strings, how does the sequence of operations

```
aList.insert(1, t)
aList.insert(1, s)
```

affect aList? We will show that $s$ is the first item in this list and that $t$ is the second item by using getEntry to look at these items.

You can write the previous sequence of operations in another way as

```
(aList.insert(1, t)).insert(1, s)
```

or

```
tempList = aList.insert(1, t)
tempList.insert(1, s)
```

Now get the first and second items in the list tempList.insert(1, s), as follows:
(tempList.insert(1, s)).getEntry (1) = s
by axiom 9
and
(tempList.insert(1, s)).getEntry(2)

```
= tempList.getEntry(1) byaxiom 10
    = (aList.insert(1, t)).getEntry(1) by definition of tempList
    = t byaxiom 9
```

Thus, $s$ is the first item in the list and $t$ is the second item.

### 8.2 Using the List Operations

We now consider some simple examples of how you can use the operations of the ADT list. Recall that exploring such uses is a way to confirm your understanding of an ADT's specifications and gives you an opportunity to change your mind about your design choices prior to implementing the ADT. You can write client functions in terms of the operations that define the ADT list, even though you do not know how the list's data is stored.

Displaying the items on a list. Suppose that you want to display the items on a list. Since this task is not an ADT list operation, we write pseudocode for a client function displayList that uses the ADT operations as follows:

```
// Displays the items on the list aList.
displayList(aList)
    for (position = 1 through aList.getLength())
    {
        dataItem = aList.getEntry(position)
        Display dataItem
    }
```

Notice that as long as the ADT list is implemented correctly, the displayList function will perform its task. In this case, getEntry successfully retrieves each list item, because position's value is always valid.

VideoNote
Using the ADT list

An implementationindependent application of the ADT list

Replacing an item. Now suppose that we did not include the operation setEntry in our specification of the ADT list. To replace the entry at a given position in the list with a new one, we would need to write a client function. The following pseudocode defines such a function replace:

```
// Replaces the ith entry in the list aList with newEntry.
// Returns true if the replacement was successful; otherwise return false.
replace(aList, i, newEntry)
    success = aList.remove(i)
    if (success)
        success = aList.insert(i, newItem)
    return success
```

If remove is successful, it sets success to true. By testing success, replace will attempt the insertion only if the deletion actually occurred. Then insert sets success, which replace returns to the function that called it. If remove is unsuccessful for any reason, including an incorrect value of $\boldsymbol{i}$, it sets success to false. The replace function then ignores the insertion and returns success.

5
Note: Notice that the two previous algorithms depend only on the specifications of the list operations and not on their implementations. The algorithms do not depend on how you implement the list. They will work regardless of whether you use an array or some other data structure to store the list's data. This feature is a definite advantage of abstract data types. In addition, by thinking in terms of the available ADT operations, you will not be distracted by implementation details.

Creating a list of names in alphabetical order. Let's create a list of our friends' names as we think of them. Moreover, let's place the names on the list in alphabetical order. It is up to us to place each name into its correct position in the list. The ADT list does not choose the order of its entries.

The following pseudocode statements place the names Amy, Ellen, Bob, Drew, Aaron, and Carol in an alphabetical list. The comment at the end of each statement shows the list after the statement executes.

```
a1phaList = a new empty list
alphaList.insert(1, "Amy") // Amy
alphaList.insert(2, "Ellen") // Amy Ellen
a7phaList.insert(2, "Bob") // Amy Bob Ellen
alphaList.insert(3, "Drew") // Amy Bob Drew Ellen
alphaList.insert(1, "Aaron") // Aaron Amy Bob Drew Ellen
alphaList.insert(4, "Caro1") // Aaron Amy Bob Carol Drew Ellen
```

After initially placing Amy at the beginning of the list and Ellen at the end of the list (at position 2), we insert

- Bob between Amy and Ellen at position 2
- Drew between Bob and Ellen at position 3
- Aaron before Amy at position 1
- Carol between Bob and Drew at position 4

Later in Chapter 11, you will learn that this way of inserting each name into a collection of alphabetized names is called an insertion sort.

If we now remove the entry at position 4-Carol-by writing

```
a1phaList.remove(4)
```

Drew and Ellen will be at positions 4 and 5, respectively. Thus, a7phaList.getEntry (4) would return Drew.

Finally, suppose that we want to replace a name in this list. We cannot replace a name with just any name and expect that the list will remain in alphabetical order. Replacing Bob with Ben by writing

```
alphaList.setEntry(3, "Ben");
```

would maintain alphabetical order, but replacing Bob with Nancy would not. The list's alphabetical order resulted from our original decisions about where to place names on the list. The order did not come about automatically as a result of list operations. That is, the client, not the list, maintains the order. We could, however, design an ADT that maintains its data in alphabetical order. You will see an example of such an ADT in Chapter 12.


Question 3 Write a pseudocode function swap(aList, $\mathbf{i}, j$ ) that interchanges the items currently in positions $i$ and $j$ of a list. Define the function in terms of the ADT list operations, so that it is independent of any particular implementation of the list. Assume that the list, in fact, has items at positions $i$ and $j$. What impact does this assumption have on your solution? (See Exercise 2 at the end of this chapter.)

Question 4 What grocery list results from the following sequence of ADT list operations?
aList $=$ a new empty list
aList.insert(1, "butter")
aList.insert(1, "eggs")
aList.insert(1, "milk")
Question 5 Suppose that myList is a list that contains the five objects a b c de.
a. What does myList contain after executing myList.insert (5, w) ?
b. Starting with the original five entries, what does myList contain after executing myList.insert(6, w)?
c. Which of the operations in parts $a$ and $b$ of this question require entries in the array to shift?

### 8.3 An Interface Template for the ADT List

We now will formalize our specification of the ADT list by writing a $\mathrm{C}++$ interface, which is given in Listing 8-1.

## LISTING 8-1 A C++ interface for lists

/** Interface for the ADT 1ist
@file ListInterface.h */
\#ifndef _LIST_INTERFACE
\#define _LIST_INTERFACE
template<c1ass ItemType>
class ListInterface
/** Sees whether this list is empty.
@return True if the list is empty; otherwise returns false. */
virtual bool isEmpty() const $=0$;
/** Gets the current number of entries in this list.
areturn The integer number of entries currently in the list. */
virtual int getLength() const $=0$;
/** Inserts an entry into this list at a given position.
@pre None.
@post If $1<=$ position <= getLength() +1 and the insertion is
successful, newEntry is at the given position in the list,
other entries are renumbered according7y, and the returned
value is true.
@param newPosition The 1ist position at which to insert newEntry.
@param newEntry The entry to insert into the list.
@return True if insertion is successful, or false if not. */
virtual bool insert(int newPosition, const ItemType\& newEntry) $=0$;
/** Removes the entry at a given position from this list.
@pre None.
@post If $1<=$ position <= getLength() and the removal is successful,
the entry at the given position in the list is removed, other
items are renumbered accordingly, and the returned value is true.
@param position The list position of the entry to remove.
@return True if removal is successful, or false if not. */
virtual bool remove(int position) $=0$;
/** Removes all entries from this 1ist.
@post List contains no entries and the count of items is 0. */
virtual void clear() $=0$;
/** Gets the entry at the given position in this list.
@pre 1 <= position <= getLength().
@post The desired entry has been returned.
@param position The list position of the desired entry.
@return The entry at the given position. */
virtual ItemType getEntry(int position) const $=0$;
$/ * *$ Replaces the entry at the given position in this list.
@pre $1<=$ position <= getLength().
@post The entry at the given position is newEntry.
@param position The list position of the entry to replace.
@param newEntry The replacement entry. */
virtual void setEntry (int position, const ItemType\& newEntry) $=0$;
\}; // end ListInterface
\#endif

## Summary

1. The ADT list maintains its data by position. Each entry in a list is identified by its position, which is given by an integer, beginning with 1 . Thus, the data in a list has an order, but that order is determined by the list's client, not the list itself.
2. You can insert a new entry into a list at a position that ranges from 1 to 1 greater than the current length of the list. Thus, you can insert a new entry before the first entry, after the last entry, or between two current entries.
3. Inserting a new entry into a list renumbers any existing entries that follow the new one in the list.
4. You can remove an entry that is currently at a position that ranges from 1 to the current length of the list. Thus, you can remove the first entry, the last entry, or any interior entry.
5. Removing an entry from a list renumbers any existing entries that follow the deleted one in the list.

## Exercises

1. Consider an ADT list of integers. Write a pseudocode function that computes the sum of the integers in the list aList. The definition of your function should be independent of the list's implementation.
2. Implement the function swap, as described in Checkpoint Question 3, but remove the assumption that the $i^{\text {th }}$ and $j^{\text {th }}$ items on the list exist. Return a value that indicates whether the swap was successful.
3. Use the function swap that you wrote in Exercise 2 to write a function that reverses the order of the items in a list aList.
4. Section 8.2 describes the functions displayList and replace. Their definitions are written in terms of the ADT list operations.
a. What is an advantage and a disadvantage of the way that displayList and replace are implemented?
b. What is an advantage and a disadvantage of defining displayList and replace as operations of the ADT list?
5. Suppose that the ADT list has a method getPosition that returns the position of a given entry within the list. Write specifications for such a method.
6. Write a pseudocode function getPosition at the client level that returns the position of a given entry within a given list.
7. Suppose that the ADT list has a method contains that tests whether the list contains a given entry. Write specifications for such a method.
8. Write a pseudocode function contains at the client level that tests whether a given list contains a given entry.
9. The ADT list method remove removes from the list the entry at a given position. Suppose that the ADT list has another method remove that removes a given entry from the list. Write specifications for such a method. What does your method do if the list contains duplicate entries?
10. Write a pseudocode function remove at the client level that removes a given entry from a given list.
11. Use the axioms for the ADT list, as given in "Axioms" in Section 8.1 of this chapter, to prove that the sequence of operations

Insert A into position 2
Insert B into position 2
Insert C into position 2
has the same effect on a nonempty list of characters as the sequence
Insert C into position 2
Insert B into position 3
Insert A into position 4
12. Repeat Exercise 20 in Chapter 2, using the ADT list to implement the function $f(n)$.

## Programming Problems

1. Write pseudocode implementations of the ADT polynomial operations, as defined in Exercise 9 of Chapter 1, in terms of the ADT list operations.
2. Santa Claus allegedly keeps lists of those who are naughty and those who are nice. On the naughty list are the names of those who will get coal in their stockings. On the nice list are those who will receive gifts. Each object in this list contains a name (a string) and a list of that person's gifts (an instance of an ADT list). Design an ADT for the objects in the nice list. Specify each ADT operation by stating its purpose, describing its parameters, and writing preconditions, postconditions, and a pseudocode version of its header. Then write a template interface for the ADT that includes javadoc-style comments.

## List Implementations

## Chapter

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| Chapter 4 | Link-Based Implementations |
| C++ Interlude 3 | Exceptions |
| Chapter 8 | Lists |

Having clearly specified the operations of an ADT list in the previous chapter, we now consider data structures for implementing it. This chapter explores two implementations of the list and compares their advantages and disadvantages.

## The data members

### 9.1 An Array-Based Implementation of the ADT List

When you design an abstract data type, you concentrate on what its operations do, but you ignore how you will implement them. Now that we have clearly specified operations for the ADT list, we can implement it as a C++ class. Recall that the list operations in their UML form are

```
+isEmpty(): boolean
+getLength(): integer
+insert(newPosition: integer, newEntry: ItemType): boolean
+remove(position: integer): boolean
+clear(): void
+getEntry(position: integer): ItemType
+setEntry(position: integer, newEntry: ItemType): void
```

You need to represent the items on a list and its length. Your first thought is probably to store the list's items in an array items. In fact, you might believe that "list" is simply a fancy name for an array. This belief is not quite true, however. An array-based implementation is a natural choice because both an array and a list identify their items by number. However, the ADT list has operations such as getLength that an array does not. Later in this chapter, you will see an implementation of the ADT list that does not use an array.

With an array-based implementation, you can store a list's $k$ th entry in items [k-1]. How much of the array will the list occupy? Possibly all of it, but probably not. That is, you need to keep track of the array elements that you have assigned to the list and those that are available for use in the future. The maximum length of the array-its physical size or capacity-is a known, fixed value, such as maxItems. You can keep track of the current number of entries on the list-that is, the list's length or logical size-in a variable itemCount. An obvious benefit of this approach is that implementing the operation getLength will be easy. Thus, we could use the following statements to define the data members for an array-based implementation of the ADT list:

```
static const int DEFAULT_CAPACITY = 100;
ItemType items[DEFAULT_CAPACITY]; // Array of list items
int itemCount; // Current count of list items
int maxItems; // Maximum capacity of the list
```

Figure 9-1 illustrates these data members. To hide them from the clients of the class, we will make these data members private.

### 9.1. $\quad$ The Header File

Now that we have chosen the data members for our class ArrayList, we can write its header file, as given in Listing 9-1. We derive ArrayList from the template interface ListInterface that we developed in the previous chapter. We provide a default constructor, and since we plan to use a statically

FIGURE 9-1 An array-based implementation of the ADT list

allocated array, the compiler-generated destructor and copy constructor will be sufficient. The only other detail that ListInterface does not specify is the behavior of the methods getEntry and setEntry should their caller violate their precondition. That is, what should we do if these methods are given an out-of-bounds list position? We can throw an exception in this case, and to this end, we use the class PrecondViolatedExcep as defined in Listings 7-5 and 7-6 of Chapter 7.

Note: By defining the data member maxItems for ArrayList, we enable the class implementer to allocate the array items dynamically instead of statically. Thus, the programmer could easily define a constructor that allows the client to choose the size of the array.

## LISTING 9-1 The header file for the class ArrayList

```
/** ADT list: Array-based implementation.
    @file ArrayList.h */
#ifndef _ARRAY_LIST
#define _ARRAY_LIST
#include "ListInterface.h"
#include "PrecondViolatedExcep.h"
temp1ate<c1ass ItemType>
class ArrayList : public ListInterface<ItemType>
{
private:
    static const int DEFAULT_CAPACITY = 100;
    ItemType items[DEFAULT_CAPACITY]; // Array of list items
    int itemCount; // Current count of list items
    int maxItems; // Maximum capacity of the 7ist
public:
    ArrayList();
    // Copy constructor and destructor are supplied by compiler
    bool isEmpty() const;
    int getLength() const;
    bool insert(int newPosition, const ItemType& newEntry);
    bool remove(int position);
    void clear();
    /** @throw PrecondViolatedExcep if position < 1 or
                            position > getLength(). */
    ItemType getEntry(int position) const throw(PrecondViolatedExcep);
    /** @throw PrecondViolatedExcep if position < 1 or
                            position > getLength(). */
    void setEntry(int position, const ItemType& newEntry)
                throw(PrecondViolatedExcep);
}; // end ArrayList
#include "ArrayList.cpp"
#endif
```

Shift array entries to insert an item

### 9.1.2 The Implementation File

As usual, we will begin our implementation with the method or methods that add new entries to our con-

Overview of ArrayList tainer and any methods that help us to verify the correctness of those additions. For ArrayList, we want to define the method insert. The methods isEmpty and getLength will be easy to define and will be useful while testing insert. We also will want to call the method getEntry to check the contents of the list. And, of course, we need to define the constructor. Let's begin with these methods as our core group.

The constructor and the methods isEmpty and getLength. The default constructor initializes the data members itemCount to zero and maxItems to DEFAULT_CAPACITY:

```
template<class ItemType>
ArrayList<ItemType>::ArrayList() : itemCount(0),
                                maxItems(DEFAULT_CAPACITY)
{
} // end default constructor
```

The method isEmpty tests whether itemCount is zero, and getLength simply returns the value of itemCount:

```
template<class ItemType>
bool ArrayList<ItemType>::isEmpty() const
{
    return itemCount == 0;
} // end isEmpty
template<class ItemType>
int ArrayList<ItemType>::getLength() const
{
    return itemCount;
} // end getLength
```

The method insert. To insert a new entry at a given position in the array items, you must create room for the new entry by shifting the entries at and beyond this position toward the end of array. Figure 9-2 depicts this insertion. Part $a$ shows the list prior to adding a new entry at position 3. Starting at the end of the list, we copy the entry in items [i-1] to items [i] for values of $i$ ranging from k down to 3, as Figure 9-2b illustrates. Finally, we insert the new entry into items [newPosition - 1], or items [2] in this example. Figure 9-2c shows the list after this step.

The definition of the method insert follows:

```
template<class ItemType>
boo1 ArrayList<ItemType>::insert(int newPosition,
                                    const ItemType& newEntry)
{
    bool ableToInsert = (newPosition >= 1) &&
                                    (newPosition <= itemCount + 1) &&
                                    (itemCount < maxItems);
    if (ableToInsert)
    {
            // Make room for new entry by shifting a11 entries at
            // positions >= newPosition toward the end of the array
            // (no shift if newPosition == itemCount + 1)
            for (int pos = itemCount; pos >= newPosition; pos--)
                items[pos] = items[pos - 1];
            // Insert new entry
            items[newPosition - 1] = newEntry;
            itemCount++; // Increase count of entries
        } // end if
        return ableToInsert;
} // end insert
```

FIGURE 9-2 Shifting items for insertion: (a) the list before the insertion; (b) copy items to produce room at position 3; (c) the result


We can begin testing these methods once we write stubs for the remaining methods. To thoroughly test insert, however, we need the method getEntry, which we define next.

Question 1 Given a nonempty list that is an instance of ArrayList, at what position does an insertion of a new entry require the fewest operations? Explain.

Question 2 Describe an implementation of the method insert for ArrayList that places the entry in position 1 in the last element of the array, the entry in position 2 in the next-to-last element, and so on.

Question 3 How does the original version of insert given previously compare with the one described in Question 2 with respect to the number of operations required?

The method getEntry. This method simply needs to return the value in items [position], but we decided to enforce the method's precondition by throwing an exception if position is out of bounds. For the purpose of testing the method insert, you could write a temporary version of getEntry that does not throw an exception but just returns the desired list entry. You would then complete its definition as follows:

```
temp1ate<c1ass ItemType>
ItemType ArrayList<ItemType>::getEntry(int position) const
    throw(PrecondViolatedExcep)
```

A client of the class cannot access the class's private members directly

```
{
    // Enforce precondition
    bool ableToGet = (position >= 1) && (position <= itemCount);
    if (ableToGet)
        return items[position - 1];
    else
    {
        string message = "getEntry() called with an empty list or ";
        message = message + "invalid position.";
        throw(PrecondViolatedExcep(message));
    } // end if
} // end getEntry
```

A client. At this point, and after writing stubs for the remaining methods, you should test the core group of methods. The following main function demonstrates how you can test the methods insert and getEntry, but it assumes that you change the value of DEFAULT_CAPACITY to 5 in the class definition given in Listing 9-1:

```
int main()
{
    ListInterface<string>* 1istPtr = new ArrayList<string>();
    string data[] = {"one", "two", "three", "four", "five", "six"};
    cout << "isEmpty: returns " << listPtr->isEmpty()
        << "; should be 1 (true)" << endl;
    for (int i = 0; i < 6; i++)
    {
        if (listPtr->insert(i + 1, data[i]))
            cout << "Inserted " << listPtr->getEntry(i + 1)
                << " at position " << (i + 1) << endl;
        e1se
            cout << "Cannot insert " << data[i] << " at position " << (i + 1)
                        << endl;
    } // end for
    return 0;
} // end main
```

Note that you cannot write references such as 1istPtr->itemCount or 1istPtr->items [4], because itemCount and items are within the private portion of the class.

The method setEntry. The definition of the method setEntry is similar to the one for getEntry, so we implement it next:

```
temp1ate<c1ass ItemType>
```

temp1ate<c1ass ItemType>
void ArrayList<ItemType>::setEntry(int position, const ItemType\& newEntry)
void ArrayList<ItemType>::setEntry(int position, const ItemType\& newEntry)
throw(PrecondViolatedExcep)
throw(PrecondViolatedExcep)
{
{
// Enforce precondition
// Enforce precondition
bool ableToSet = (position >= 1) \&\& (position <= itemCount);
bool ableToSet = (position >= 1) \&\& (position <= itemCount);
if (ableToSet)
if (ableToSet)
items[position - 1] = newEntry;
items[position - 1] = newEntry;
else
else
{
{
string message = "setEntry() called with an empty list or ";
string message = "setEntry() called with an empty list or ";
message = message + "invalid position.";
message = message + "invalid position.";
throw(PrecondViolatedExcep(message));
throw(PrecondViolatedExcep(message));
} // end if
} // end if
} // end setEntry

```
} // end setEntry
```

The method remove. Now consider how to remove an entry from the list. You could blank it out, but this strategy can lead to gaps in the array, as Figure 9-3a illustrates. An array that is full of gaps has three significant problems:

- itemCount - 1 is no longer the index of the last entry in the array. You need another variable, lastPosition, to contain this index.
- Because the items are spread out, the method getEntry might have to look at every cell of the array even when only a few entries are present.
- When items[maxItems - 1] is occupied, the list could appear full, even when fewer than maxItems entries are present.

Thus, what you really need to do is shift the entries in the array so that a deletion does not leave a gap, as shown in Figure 9-3b. This data movement is like the one done by the method insert, but it occurs in the opposite direction.

The definition of remove follows:

```
temp7ate<class ItemType>
bool ArrayList<ItemType>::remove(int position)
{
    bool ableToRemove = (position >= 1) && (position <= itemCount);
    if (ableToRemove)
    {
            // Remove entry by shifting al1 entries after the one at
            // position toward the beginning of the array
            // (no shift if position == itemCount)
            for (int fromIndex = position, toIndex = fromIndex - 1;
                    fromIndex < itemCount; fromIndex++, toIndex++)
                items[toIndex] = items[fromIndex];
            itemCount--; // Decrease count of entries
    } // end if
    return ab1eToRemove;
} // end remove
```

FIGURE 9-3 $\quad$ (a) Deletion can cause a gap; (b) shift items to prevent a gap at position 3; (c) the result


The method clear. To clear all of the entries from a list, our final method simply sets itemCount to zero:

```
template<class ItemType>
void ArrayList<ItemType>::clear()
{
    itemCount = 0;
} // end clear
```

The entire implementation file for the class ArrayList is available online from the book's website.

Question 4 Although the method remove cannot remove an entry from an empty list, it does not explicitly check for one. How does this method avoid an attempted deletion from an empty list?

### 9.2 A Link-Based Implementation of the ADT List

We now consider how we can use C++ pointers instead of an array to implement the ADT list. Unlike the array-based implementation, a link-based implementation does not shift items during insertion and deletion operations. It also does not impose a fixed maximum length on the list- except, of course, as imposed by the storage limits of the system.

Once again, we need to represent the items in the list and its length. Figure 9-4 indicates one possible way to represent this data by using pointers. Here headPtr points to the linked nodes containing the items in the list, where the first node contains the entry at position 1 in the list, and so on. The integer itemCount is the current number of entries in the list. Both headPtr and itemCount will be private data members of our class.

### 9.2.1 The Header File

Let's name our class of lists LinkedList and write its header file as shown in Listing 9-2. This file has similarities to the header file for ArrayList, as given in Listing 9-1, but let's point out the differences. Here we include the header files for the classes Node and PrecondViolatedExcep, as defined in Listing 4-1 of Chapter 4 and Listing 7-5 of Chapter 7, respectively.

The private section declares the two data members-headPtr and itemCount-as well as a private method getNodeAt that we will use to locate any node within the chain of linked nodes, given its position in the chain. Unlike an array, which provides direct access to a specified element, the chain of linked nodes requires the get, insertion, and removal operations to traverse it until the specified node is reached. While you could write the necessary loop in the method getEntry, you would soon discover that the methods insert and remove will need the same loop. At that time, you should realize

FIGURE 9-4 A link-based implementation of the ADT list


## LISTING 9-2 The header file for the class LinkedList

```
/** ADT list: Link-based implementation.
    @file LinkedList.h */
#ifndef _LINKED_LIST
#define _LINKED_LIST
#include "ListInterface.h"
#include "Node.h"
#include "PrecondViolatedExcep.h"
temp1ate<c1ass ItemType>
class LinkedList : public ListInterface<ItemType>
{
private:
    Node<ItemType>* headPtr; // Pointer to first node in the chain
                            // (contains the first entry in the list)
    int itemCount; // Current count of 7ist items
    // Locates a specified node in a linked list.
    // @pre position is the number of the desired node;
    // position >= 1 and position <= itemCount.
    // @post The node is found and a pointer to it is returned.
    // @param position The number of the node to locate.
    // @return A pointer to the node at the given position.
    Node<ItemType>* getNodeAt(int position) const;
public:
    LinkedList();
    LinkedList(const LinkedList<ItemType>& aList);
    virtual ~LinkedList();
    bool isEmpty() const;
    int getLength() const;
    bool insert(int newPosition, const ItemType& newEntry);
    bool remove(int position);
    void clear();
    /** @throw PrecondViolatedExcep if position < 1 or
                                    position > getLength(). */
    ItemType getEntry(int position) const throw(PrecondViolatedExcep);
    /** @throw PrecondViolatedExcep if position < 1 or
                                    position > getLength(). */
    void setEntry(int position, const ItemType& newEntry)
                throw(PrecondViolatedExcep);
}; // end LinkedList
#include "LinkedList.cpp"
#endif
```

getNodeAt is a private method

A copy constructor and a destructor are necessary for a link-based implementation

VideoNote
Overview of LinkedList
that defining a private method to perform the traversal would be convenient. Although we declare getNodeAt now, adding the declarations of private methods to the header file after you begin the implementation file is not unusual.

The method getNodeAt is not an ADT operation, because it returns a pointer to a node; since nodes are an implementation detail, you would not want any client to call it. Such clients should be able to use the ADT without knowledge of the nodes that the implementation uses. It is perfectly reasonable for the implementation of an ADT to define variables and methods that the rest of the program should not access. You could think of the method getNodeAt as marked "for internal use only." Therefore, getNodeAt is a private method that only the implementations of the ADT's operations call.

Because we are declaring a link-based implementation, we must provide a copy constructor and a destructor for our class. The declarations of the public methods are the same as for the class ArrayList.

### 9.2.2 The Implementation File

Our core methods include the same ones that we defined for ArrayList-namely, the constructor and the methods insert, getEntry, isEmpty, and getLength. The latter two methods have the same bodies as their ArrayList versions, since they involve only the data field itemCount. However, as will become evident when we code the methods getEntry and insert, the private method getNodeAt will be necessary.

The constructor. Because the compiler-generated default constructor would not necessarily initialize headPtr and itemCount to appropriate values, you must provide your own default constructor. Initializers are sufficient to set the values of headPtr and itemCount, since we, not the client, choose these values. The constructor's definition follows:

```
template<class ItemType>
LinkedList<ItemType>::LinkedList() : headPtr(nu11ptr), itemCount(0)
{
} // end default constructor
```

This is the only constructor we will need.
The method getEntry. Like getEntry in ArrayList, we will have this method enforce its precondition by throwing an exception if position is out of bounds. Note the similarity of its definition to the array-based version given earlier in this chapter:

```
template<class ItemType>
ItemType LinkedList<ItemType>::getEntry(int position) const
                                    throw(PrecondViolatedExcep)
{
    // Enforce precondition
    bool ab1eToGet = (position >= 1) && (position <= itemCount);
    if (ableToGet)
    {
        Node<ItemType>* nodePtr = getNodeAt(position);
        return nodePtr->getItem();
    }
    else
    {
        string message = "getEntry() called with an empty list or ";
        message = message + "invalid position.";
        throw(PrecondViolatedExcep(message));
    } // end if
} // end getEntry
```

The method getNodeAt. The private method getNodeAt locates the node at a given position by traversing the chain. It then returns a pointer to the located node. The traversal begins at the first node of the chain and moves from node to node, counting as it goes, until it reaches the desired one:

```
temp1ate<c1ass ItemType>
Node<ItemType>* LinkedList<ItemType>::getNodeAt(int position) const
{
    // Debugging check of precondition
    assert( (position >= 1) && (position <= itemCount) );
    // Count from the beginning of the chain
    Node<ItemType>* curPtr = headPtr;
    for (int skip = 1; skip < position; skip++)
        curPtr = curPtr->getNext();
    return curPtr ;
} // end getNodeAt
```

Since getNodeAt is a private method and is called only by other methods within the class, those methods are responsible for honoring its precondition. Thus, we need not throw an exception if the precondition is violated. Rather, we simply use an assertion during debugging.
The method insert. As you know, you can insert a new entry into a list right before its first entry, right after its last entry, or between two adjacent entries. For the link-based implementations of the ADT bag in Chapter 4 and the ADT stack in Chapter 7, we always inserted a new node into the chain of linked nodes at the chain's beginning. The ADT list requires us to manipulate the chain in additional ways. Let's explore those ways.

Since we know how to insert a new node at the beginning of a chain of nodes, let's consider inserting a node between two existing, adjacent nodes to which the pointer variables prevPtr and curPtr point, as Figure 9-5a illustrates. If the pointer variable newNodePtr points to the new node, you make the new node point to the node that curPtr currently points to (Figure 9-5b). You also must alter the value of the pointer in the node that will precede the new node so that it points to the new node, as shown in Figure 9-5c. As the diagram suggests, you can accomplish the insertion by using the pair of assignment statements

```
newNodePtr->setNext(curPtr);
prevPtr->setNext(newNodePtr);
```

Now, how did the variables curPtr and prevPtr get appropriate values? Given newPosition, the desired position of the new entry in the list, you traverse the chain until you find the proper position for the new item. The private method getNodeAt will perform this traversal for you. Thus,

```
prevPtr = getNodeAt(newPosition - 1);
curPtr = prevPtr->getNext();
```

You then use the new operator to create a new node, to which newNodePtr points. You initialize its data portion and then insert the node into the chain, as was just described.

Figure 9-6 shows the insertion of a new node at the end of a chain. This insertion is potentially a special case because the intention of the pair of assignment statements

```
newNodePtr->setNext(curPtr);
prevPtr->setNext(newNodePtr);
```

is to insert the new node between the nodes to which curPtr and prevPtr point. If you are to insert the new node at the end of the chain, to what node should curPtr point? In this situation, it makes sense to view the value of curPtr as nullptr because, as you traverse the list, curPtr becomes nul1ptr as it moves past the end of the chain. Observe that if curPtr has the value nullptr and

Inserting a node between nodes

If curPtr is nul1ptr, inserting at the end of a chain is not a special case

Three steps to insert a new node into a chain
prevPtr points to the last node in the chain, the previous pair of assignment statements will indeed insert the new node at the end of the chain. Thus, insertion at the end of a chain is not a special case.

To summarize, the insertion process requires three high-level steps:

1. Create a new node and store the new data in it.
2. Determine the point of insertion.
3. Connect the new node to the linked chain by changing pointers.

These steps and the previous discussion are implemented in the following definition of insert:

```
template<class ItemType>
bool LinkedList<ItemType>::insert(int newPosition,
                        const ItemType& newEntry)
{
        bool ableToInsert = (newPosition >= 1) &&
                        (newPosition <= itemCount + 1);
        if (ableToInsert)
        {
            // Create a new node containing the new entry
            Node<ItemType>* newNodePtr = new Node<ItemType>(newEntry);
            // Attach new node to chain
            if (newPosition == 1)
            {
                // Insert new node at beginning of chain
                    newNodePtr->setNext(headPtr);
                    headPtr = newNodePtr;
            }
            else
            {
                    // Find node that will be before new node
                    Node<ItemType>* prevPtr = getNodeAt(newPosition - 1);
                    // Insert new node after node to which prevPtr points
                    newNodePtr->setNext(prevPtr->getNext());
                    prevPtr->setNext(newNodePtr);
            } // end if
            itemCount++; // Increase count of entries
        } // end if
        return ab1eToInsert;
} // end insert
```

If you have not already begun to test the methods in the core group, you should do so now.


Question 5 Given a nonempty list that is an instance of LinkedList, at what position does an insertion of a new entry require the fewest operations? Explain.
Question 6 In the previous method insert, the second if statement tests the value of newPosition. Should the boolean expression it tests be isEmpty() || (newPosition == 1)? Explain.

Question 7 How does the insert method enforce the precondition of getNodeAt?

The method remove. To remove any entry from a list, you must be able to delete any node from a chain of linked nodes. You already know how to delete the first node in a chain, because the

FIGURE 9-5 Inserting a new node between existing nodes of a linked chain
(a) Before the insertion of a new node

(b) After newNodePtr->setNext (curPtr) executes

(c) After prevPtr->setNext (newNodePtr) executes


FIGURE 9-6 Inserting a new node at the end of a chain of linked nodes


## FIGURE 9-7 Deleting a node from a chain

Removing an interior node

link-based implementations of the ADT bag in Chapter 4 and the ADT stack in Chapter 7 do this. Let's consider the deletion of an interior node.

As Figure 9-7 indicates, you can remove the node $N$ to which curPtr points by altering the value of the pointer in the node that precedes $N$. You need to set this pointer so that it points to the node that follows $N$, thus bypassing $N$ on the chain. (The dashed line indicates the old pointer value.) Notice that this pointer change does not directly affect node $N$. Node $N$ remains in existence, and it points to the same node that it pointed to before the removal. However, the node has effectively been removed from the chain. For example, a traversal of the chain would never reach node $N$.

To accomplish this pointer change, notice first that if you had only the pointer curPtr pointing to $N$, you would have no direct way to access the node that precedes $N$. After all, you cannot follow the links in the chain backward. Thus, you must have the pointer variable prevPtr in Figure 9-7. It points to the node that precedes $N$ and makes it possible for you to alter that node's next pointer, thereby removing node $N$ from the chain. The following assignment statement is all that you need to remove the node to which curPtr points:

```
prevPtr->setNext(curPtr->getNext());
```

Does the previous technique work for any node $N$, regardless of where in the linked chain it appears? No, it does not work if the node to be deleted is the first node in the chain, but it does work for the last node in the chain. Thus, deletion of the first node in a chain is a special case. Since we already know how to delete the first node, let's consider deleting the last node.

Figure 9-8 shows the deletion of the node to which curPtr points, which happens to be the last node. That pointer alone is insufficient, as you must have the variable prevPtr point to the next-to-last node. You set the pointer within the next-to-last node to nul1ptr. The same statement that we just used to remove an interior node will make this pointer change, since the value of curPtr->getNext() is nul1ptr:

```
prevPtr->setNext(curPtr->getNext());
```

FIGURE 9-8 Removing the last node


To remove a chain's first node, recall that you must change the value of headPtr to point to the second node. Thus, the second node becomes the chain's new first node. You make this change to headPtr by using the assignment statement

```
headPtr = headPtr->getNext();
```

As was the case for the deletion of an interior node, the pointers bypass the old first node, although it still exists. Notice also that if the node to be deleted is the only node in the chain-and thus it is both the first node and the last node-the previous assignment statement assigns the value nul1ptr to the variable headPtr. Recall that the value nu71ptr in headPtr indicates an empty chain, and so this assignment statement handles the removal of the only node in a chain correctly.

The removed node still exists, however. Assuming that curPtr points to it, changing the value of curPtr will place the removed node in a state of limbo. It will still require storage space, and the program will no longer be able to access it. You would have a memory leak. Therefore, before you change the value of curPtr, you should use the statements

```
curPtr->setNext(nul1ptr);
de1ete curPtr;
curPtr = nul1ptr;
```

to delete the node and return its memory to the system.
To summarize, the deletion process has three high-level steps:

1. Locate the node that you want to delete.
2. Disconnect this node from the linked chain by changing pointers.
3. Return the node to the system.

You will see these steps and the previous discussion implemented in the following definition of remove:

```
template<class ItemType>
bool LinkedList<ItemType>::remove(int position)
{
    bool ableToRemove = (position >= 1) && (position <= itemCount);
    if (ableToRemove)
    {
        Node<ItemType>* curPtr = nul1ptr;
        if (position == 1)
        {
            // Remove the first node in the chain
            curPtr = headPtr; // Save pointer to node
            headPtr = headPtr->getNext();
        }
        else
        {
            // Find node that is before the one to delete
            Node<ItemType>* prevPtr = getNodeAt(position - 1);
            // Point to node to delete
            curPtr = prevPtr->getNext();
            // Disconnect indicated node from chain by connecting the
            // prior node with the one after
            prevPtr->setNext(curPtr->getNext());
        } // end if
        // Return node to system
        curPtr->setNext(nul7ptr);
```

Removing the first node is a special case

Return removed nodes to the system by using delete

Three steps to delete a node from a chain

```
        delete curPtr;
        curPtr = nul1ptr;
    itemCount--; // Decrease count of entries
    } // end if
    return ableToRemove;
} // end remove
```

The method clear. To remove all entries from a list, you can call the method remove. Because the easiest entry to remove is the first one, the method clear can invoke remove (1) repeatedly until the list is empty. This action works because, after you remove the first entry in the list, the remaining entries are renumbered. Therefore, to remove all entries, you can repeatedly remove the first entry. Thus, clear has the following definition:

```
temp1ate<class ItemType>
void LinkedList<ItemType>::clear()
{
    while (!isEmpty())
        remove(1);
} // end clear
```

Question 8 The link-based implementation of the method clear contains the following loop:

```
while (!isEmpty())
    remove(1);
```

a. Can you correctly replace the loop with

```
for (int position = getLength(); position >= 1; position--)
```

remove(1) ;
b. Does your answer to part $a$ differ if you replace remove(1) with remove(position)?
c. Do your answers to parts $a$ and $b$ differ if you replace the for statement with
for (int position $=1$; position <= getLength(); position++)
The destructor. Because the method clear invokes remove repeatedly until the list is empty, and remove deallocates the nodes it removes, the destructor can simply call clear:

```
template<class ItemType>
LinkedList<ItemType>::~LinkedList()
{
    clear();
} // end destructor
```

Question 9 Revise the destructor in the class LinkedList so that it directly deletes each node of the underlying linked chain without calling either clear or remove.

The remaining methods. This link-based implementation of the ADT list cannot use a compilergenerated copy constructor, as it would copy only the data members headPtr and itemCount. Instead, it must also copy the nodes and their contents. The link-based implementations of the ADT bag in Chapter 4 and the ADT stack in Chapter 7 have copy constructors that are quite similar to the one we require for the list. Thus, we leave its definition to you as an exercise, along with the method setEntry.

### 9.2.3 Using Recursion in LinkedList Methods

Using recursion in the implementation of some of the methods in the class LinkedList can be an attractive alternative to an iterative approach. Section 4.3 of Chapter 4 developed the recursive methods fillVector and getPointerTo. Both methods considered a chain's first node and then recursively considered the rest of the chain. In general, we can state that you can process a linked chain by processing its first node and then the rest of the chain recursively. Thus, to add a new node to a chain of linked nodes at a given position, you use the following logic:

```
if (the insertion position is 1)
    Add the new node to the beginning of the chain
else
    Ignore the first node and add the new node to the rest of the chain
```

Adding to the beginning of a chain-or subchain-is the base case of this recursion. Happily, the beginning of a chain is the easiest place to make an addition.

If position is the desired position of the new node, newNodePtr points to the new node, and subChainPtr initially points to the chain and later points to the rest of the chain, we can add some detail to the previous logic, as follows:

```
if (position == 1)
{
    newNodePtr->setNext(subChainPtr)
    subChainPtr = newNodePtr
    Increment itemCount
}
else
Using recursion, add the new node at position position - 1 of the subchain pointed to by subChainPtr->getNext()
```

The private method getPointerTo in Section 4.3 returned a pointer to a given entry in a chain. Since that operation traversed the chain but did not alter it, its recursive formulation was straightforward. However, the list's method insert does alter the chain. Getting the recursive method to make these changes is the challenge.

The method insert. Let's look at the recursive implementation of the insertion operation before we describe why it works. You learned in Section 4.3 that you write a private method to perform the recursion and you write a public method-typically the one that implements the ADT's operation-to invoke this private method. Thus, we have the following method definitions:

```
// The public method insert:
temp1ate<c1ass ItemType>
bool LinkedList<ItemType>::insert(int newPosition, const ItemType& newEntry)
{
    bool ab7eToInsert = (newPosition >= 1) &&
                            (newPosition <= itemCount + 1);
    if (ableToInsert)
    {
        // Create a new node containing the new entry
        Node<ItemType>* newNodePtr = new Node<ItemType>(newEntry);
        headPtr = insertNode(newPosition, newNodePtr, headPtr);
    } // end if
    return ableToInsert;
} // end insert
```

```
// The private method insertNode:
// Adds a given node to the subchain pointed to by subChainPtr
// at a given position. Returns a pointer to the augmented subchain.
template<class ItemType>
Node<ItemType>* LinkedList<ItemType>::insertNode(int position,
                                    Node<ItemType>* newNodePtr, Node<ItemType>* subChainPtr)
{
    if (position == 1)
    {
        // Insert new node at beginning of subchain
        newNodePtr->setNext(subChainPtr);
        subChainPtr = newNodePtr;
        itemCount++; // Increase count of entries
    }
    else
    {
        Node<ItemType>* afterPtr =
            insertNode(position - 1, newNodePtr, subChainPtr->getNext());
        subChainPtr->setNext(afterPtr);
    } // end if
    return subChainPtr;
} // end insertNode
```

The private recursive method insertNode adds the new node to the subchain to which subChainPtr points at the given position. We will trace and explain its logic next.

Tracing an addition to the list's beginning. Consider the list that the chain in Figure 9-9a represents. Let's add a new entry at the beginning of the chain by using an invocation such as

```
myList->insert(1, newEntry);
```

The public method insert will create a new node containing newEntry (Figure 9-9b) and then call the private method insertNode with the invocation

```
insertNode(1, newNodePtr, headPtr);
```

The pointer in the argument headPtr is copied to the parameter subChainPtr, and so it also points to the first node in the chain, as Figure 9-9c illustrates.

The new node is linked to the beginning of the original chain, as Figure $9-9 \mathrm{~d}$ shows. Notice that subChainPtr now points to the new node at the beginning of the chain. However, headPtr is unchanged, even though it is the argument that corresponds to the parameter subChainPtr. The private method now returns the value of subChainPtr, and the public method insert assigns that value to headPtr. The chain with the completed addition appears as in Figure 9-9e.

Tracing an addition to the list's interior. What happens when the addition is not at the beginning of the original chain? Let's trace what happens when we add a new third node to the chain given in Figure $9-9$ a. The public method insert creates a new node and calls the private method insertNode with the invocation

```
insertNode(3, newNodePtr, headPtr);
```

As in the previous example, the pointer in the argument headPtr is copied to the parameter subChainPtr, and so it also points to the first node in the chain, as Figure 9-10a illustrates. Since the insertion is not at position 1 , another recursive call occurs:

```
insertNode(2, newNodePtr, subChainPtr->getNext());
```

FIGURE 9-9 Recursively adding a node at the beginning of a chain
(a) The list before any additions

(b) After the public method insert creates a new node and before it calls insertNode

(c) As insertNode (1, newNodePtr, headPtr) begins execution

(d) After the new node is linked to the beginning of the chain (the base case)

(e) After the public method insert assigns to headPtr the reference returned from insertNode


The third argument is a pointer to the chain's second node. This pointer is copied to the parameter subChainPtr, as Figure $9-10$ b depicts. We still have not reached the insertion point, so the recursive process is repeated again:

```
insertNode(1, newNodePtr, subChainPtr->getNext());
```

Now subChainPtr points to the chain's third node, as shown in Figure 9-10c, and we are at the desired position for the insertion into the chain. No other recursive call occurs, as this is the base case.

The given new node is inserted at the beginning of the subchain to which subChainPtr points, as Figure 9-10d illustrates, and then the private method insertNode returns a reference to the new node. The statement

```
Node<ItemType>* afterPtr = insertNode(1, newNodePtr, subChainPtr->getNext());
```

FIGURE 9-10 Recursively adding a node between existing nodes in a chain
(a) As insertNode(3, newNodePtr, headPtr) begins execution

(b) As the recursive call insertNode(2, newNodePtr, subChainPtr->getNext()) begins execution

(c) As the recursive call insertNode(1, newNodePtr, subChainPtr->getNext()) begins execution

(d) After a new node is linked to the beginning of the subchain (the base case)

(e) After the returned pointer is assigned to afterPtr

(f) After subChainPtr->setNext(afterPtr) executes

in the public method insert that invoked insertNode now resumes execution. The pointer to the new node returned by insertNode is assigned to afterPtr, as Figure 9-10e illustrates. At this time, subChainPtr references the second node, as it did in part $b$ of the figure. The next statement to execute in insertNode is

```
subChainPtr->setNext(afterPtr);
```

This statement changes the next pointer in the second node to reference the new node, as shown in Figure 9-10f. The private method insertNode now returns a reference to the second node. If we continue the trace, we will see that the method makes the first node point to the second node and headPtr point to the first node, even though these pointers are already in place.


Note: A recursive addition to a chain of nodes locates and remembers the nodes prior to the insertion point. After the portion of the chain that follows the insertion point is linked to the new node, the recursion links the remembered nodes back into the chain.

Question 10 Using recursion, revise the destructor in the class LinkedList so that it deletes each node of the underlying linked chain.

### 9.3 Comparing Implementations

The reasons for choosing an array-based implementation or a link-based implementation are the same as discussed in earlier chapters. Although the array-based implementation given in this chapter seems reasonable - after all, an array behaves like a list, and arrays are easy to use-it uses statically allocated memory. As such, it prevents the insert operation from adding an entry to the list if the array is full. If this restriction is not acceptable, you must either resize the array or use a link-based implementation.

The array-based and link-based implementations have other differences as well, which affect their time and memory requirements. As we discussed in Section 4.5 of Chapter 4, the time to access an array element is constant, whereas you have no way of immediately accessing the $i^{\text {th }}$ node in a linked chain. Thus, the array-based getEntry method, for example, is almost instantaneous regardless of which list item you access. A link-based getEntry, however, requires $i$ steps to access the $i^{\text {th }}$ item in the list.

You already know that the array-based implementation of the ADT list requires you to shift the data when you insert entries into or remove entries from the list. For example, if you remove the first entry of a 20-entry list, you must shift 19 entries. In general, removing the $i^{\text {th }}$ entry from a list of $n$ entries requires $n-i$ shifts. Thus, remove requires $n-1$ shifts to remove the first entry, but zero shifts to remove the last one. The method insert has similar requirements.

In contrast, you do not need to shift the data when you insert entries into or remove entries from a chain of linked nodes. Thus, if the ADT list has a link-based implementation, the methods insert and remove require essentially the same effort-regardless of the length of the list or the position of the operation within the list - once you know the point of insertion or removal. Finding this point, however, requires a list traversal, the time for which will vary depending on where in the list the operation occurs. Recall that the private method getNodeAt performs this traversal. If you examine the definition of getNodeAt, you will see that getNodeAt(i) requires $i$ assignment operations. Thus, getNodeAt's effort increases with $i$.

The time to access the $i^{\text {th }}$ node in a chain of linked nodes depends on $i$

You can access array items directly with equal access time

Insertions and removals with a link-based implementation do not require you to shift data but require a traversal

## Summary

1. Using an array results in a straightforward implementation of the ADT list, but it is somewhat more involved than the implementations of either the ADT bag or the ADT stack.
2. An array provides direct access to any of its elements, so a method such as getEntry can quickly retrieve any entry in a list.
3. Adding an entry to or removing an entry from an array-based list typically requires that other entries shift by one position within the array. This data movement degrades the time efficiency of these operations, particularly when the list is long and the position of the addition or removal is near the beginning of the list.
4. Inserting entries into or removing entries from a chain of linked nodes does not require data to shift. Thus, the methods insert and remove of a link-based implementation of the ADT list require essentially the same effort-regardless of the length of the list or the position of the operation within the list-once the point of insertion or removal is known. Finding this point, however, requires a list traversal, the time for which will vary depending on where in the list the operation occurs.
5. For a link-based implementation of the ADT list, adding to the beginning of the list and removing a list's first entry are treated as special cases.
6. Adding or removing an entry at the end of a link-based list requires a traversal of the underlying chain.
7. Adding or removing an entry anywhere within a link-based list requires a change of at most two pointers within the underlying chain.
8. The array-based getEntry method is almost instantaneous regardless of which list item you access. A linkbased getEntry, however, requires $i$ steps to access the $i^{\text {th }}$ item in the list.

## Exercises

1. Add a constructor to each of the classes ArrayList and LinkedList that creates a list containing the entries in a given array.
2. Define the method setEntry for the class LinkedList.
3. Define the copy constructor for the class LinkedList.
4. Repeat the previous exercise, but use recursion in your definition.
5. Implement the method getPosition, as described in Exercise 5 of Chapter 8, for each of the classes ArrayList and LinkedList.
6. Repeat the previous exercise, but use recursion in your definitions.
7. Implement the method contains, as described in Exercise 7 of Chapter 8, for each of the classes ArrayList and LinkedList.
8. Repeat the previous exercise, but use recursion in your definitions.
9. Implement the method remove, as described in Exercise 9 of Chapter 8, for each of the classes ArrayList and LinkedList.
10. Repeat the previous exercise, but use recursion in your definition.
11. Write a recursive definition of the private method getNodeAt for the class LinkedList.
12. A double-ended list has operations that operate at its beginning and end, in addition to the operations of the list, as given in Section 8.1 of Chapter 8. For example, you can add, remove, and get the first and last entries in a double-ended list. Write a $\mathrm{C}++$ interface for a double-ended list.

## Programming Problems

1. Write a program that thoroughly tests the classes ArrayList and LinkedList. During its execution, your program should give the user a choice of which class to test. After the results are displayed, the user should have the opportunity to choose the other class or quit.
2. Write an array-based implementation of the ADT list that expands the size of the array of list entries as needed so that the list can always accommodate a new entry.
3. Repeat the previous programming problem, but also reduce the size of the array as needed to accommodate several removals. When the size of the array is greater than 20 and the number of entries in the list is less than half the size of the array, reduce the size of the array so that it is three quarters of its current size.
4. Revise the array-based implementation of the ADT list to use recursion wherever you can.
5. Revise the link-based implementation of the ADT list to use recursion wherever you can.
6. Adding nodes to or removing nodes from a chain of linked nodes requires a special case when the operation is at the beginning of the chain. To eliminate the special case, you can add a dummy node at the beginning of the chain. The dummy node is always present but does not contain a list entry. The chain, then, is never empty, and so the head pointer never contains nullptr, even when the list is empty. Modify the class LinkedList, as presented in this chapter, by adding a dummy node to the chain.
7. Implement the ADT polynomial that Exercise 9 in Chapter 1 describes by using a list. Then write a program that adequately demonstrates your new class.
8. Implement a class of bags, as specified in Chapter 1, by using a list to contain the bag's entries. Then write a program that adequately demonstrates your new class.
9. Implement a class of stacks, as specified in Chapter 6, by using a list to contain the stack's entries. Then write a program that adequately demonstrates your new class.
10. Implement the ADT for the objects on Santa Claus's nice list, as described in Programming Problem 2 of Chapter 8. Then write a program for Santa that maintains his two lists of those who are naughty and those who are nice.
11. The popular social network Facebook was founded by Mark Zuckerberg and his classmates at Harvard University in 2004. At the time, he was a sophomore studying computer science.

Design and implement an application that maintains the data for a simple social network. Each person in the network should have a profile that contains the person's name, optional image, current status, and a list of friends. Your application should allow a user to join the network, leave the network, create a profile, modify the profile, search for other profiles, and add friends.
12. A tail pointer is a pointer variable to the last node in a chain of linked nodes. Write a link-based implementation for a class of double-ended lists, as specified in Exercise 12, that uses both a head pointer and a tail pointer.
13. The solution to the HPAir problem described in Programming Problem 11 of Chapter 6 uses an adjacency list to represent the flight map. Repeat the solution to this problem, but define the adjacency list as a list of lists instead of an array of linked chains. Use lists that have a link-based implementation.

## Algorithm Efficiency

## Chapter

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This chapter will show you how to analyze the efficiency of algorithms. The basic mathematical techniques for analyzing algorithms are central to more advanced topics in computer science and give you a way to formalize the notion that one algorithm is significantly more efficient than another. As examples, you will see analyses of some algorithms that you have studied before, including those that search data. The next chapter presents sorting algorithms, and they provide additional examples of analyzing efficiency.

### 10.1 What Is a Good Solution?

In this textbook a computer program is the final form your solutions take, so we should consider what constitutes a good computer program. Presumably, you write a program to perform some task. While performing that task, the program incurs a real and tangible cost. This cost includes such factors as the computing time and memory that the program requires, the difficulties encountered by those who use the program, and the consequences of a program that does not behave correctly.

However, the costs just mentioned do not give the whole picture. They pertain only to the life of the solution after it has been developed. In assessing whether a solution is good, you also must consider the effort required to develop the solution as well as any changes - bug fixes or extensions - to the program that are made after the program has been deployed. Each of these incurs costs, too. The total cost of a solution must take into account the value of the time of the people who analyzed, designed, coded, debugged, and tested it. A solution's cost must also include the cost of maintaining, modifying, and expanding it.

Thus, when calculating the overall cost of a solution, you must include a diverse set of factors. If you adopt such a multidimensional view of cost, it is reasonable to evaluate a solution against the following criterion:

## A solution is good if the total cost it incurs over all phases of its life is minimal.

Suppose two algorithms perform the same task, such as searching. What does it mean to compare the algorithms and conclude that one is better? The faster one is not necessarily better. Several components contribute to the cost of a computer program, including the cost of human time - the time of the people who develop, maintain, and use the program - and the cost of program execution-that is, its efficiency - measured by the amount of computer time and memory that the program requires to execute.

Today's programs are larger and more complex than ever before. Typically, many people are involved in their development. Good structure and documentation are thus of the utmost importance. In addition, the costs associated with malfunctions are high. People should not have to entrust their livelihoods-or their lives - to a program that only its authors can understand and maintain. Thus, society needs both well-structured programs and techniques for formally verifying their correctness.

At the same time, do not get the impression that a solution's execution time is no longer important. To the contrary, many situations occur for which efficiency is the prime determinant of whether a solution is even usable. However, a solution's efficiency is only one of many aspects that you must consider. If two solutions have approximately the same efficiency, other factors should dominate the comparison. However, when the efficiencies of solutions differ significantly, this difference can be the overriding concern.

You should be most concerned about efficiency when you develop the underlying algorithm. The choice of a solution's components - the objects and the design of the interactions between those objects-rather than the code you write, has the most significant impact on efficiency. This book advocates a problem-solving philosophy that views the cost of a solution as multidimensional. This philosophy is reasonable in today's world, and it likely will be reasonable in the years to come.

Note: The relative importance of the various components of a solution's cost has changed since the early days of computing. In the beginning, the cost of computer time relative to human time was extremely high. In this type of environment, one cost clearly overshadowed all others: computer resources. If two programs performed the same task, the one that required less time and memory was better.

Computing costs have dropped dramatically since the early days of computers. Thus, the value of the designers' and programmers' time is a much more significant factor than computing time in the cost of a solution. These developments have made obsolete the notion that the fastest-executing solution is always the best.

### 10.2 Measuring the Efficiency of Algorithms

Well-designed algorithms reduce the human costs of implementing the algorithm with a program, of maintaining the program, and of modifying the program. Developing good problem-solving skills and programming style has been and continues to be important. However, the efficiency of algorithms is also important. Efficiency is a criterion that you should consider when selecting an algorithm and its implementation.

The comparison of algorithms is a topic central to computer science. Measuring an algorithm's efficiency is quite important because your choice of algorithm for a given application often has a great impact. Responsive word processors, grocery checkout systems, automatic teller machines, video games, and life support systems all depend on efficient algorithms.

The analysis of algorithms is the area of computer science that provides tools for contrasting the efficiency of different algorithms. Notice the use of the term "algorithms" rather than "programs;" it is important to emphasize that the analysis concerns itself primarily with significant differences in efficiency-differences that you can usually obtain only through superior algorithms and rarely through clever tricks in coding. Reductions in computing costs due to clever coding tricks are often more than offset by reduced program readability, which increases human costs. An analysis should focus on gross differences in the efficiency of algorithms that are likely to dominate the overall cost of a solution. To do otherwise could lead you to select an algorithm that runs a small fraction of a second faster than another algorithm yet requires many more hours of your time to implement and maintain.

The efficient use of both time and memory is important. Computer scientists use similar techniques to analyze an algorithm's time and space efficiency. As none of the algorithms covered in this text has significant space requirements, our focus is primarily on time efficiency.

How do you compare the time efficiency of two algorithms that solve the same problem? One possible approach is to implement the two algorithms and run the programs. This approach has at least three fundamental problems:

- How are the algorithms coded? If algorithm $A_{1}$ runs faster than algorithm $A_{2}$, it could be the result of better programming. Thus, if you compare the running times of the programs, you are really comparing implementations of the algorithms rather than the algorithms themselves. You should not compare implementations, because they are sensitive to factors such as programming style that tend to cloud the issue of which algorithm is inherently more efficient.
- What computer should you use? The particular operations that the algorithms require can cause $A_{1}$ to run faster than $A_{2}$ on one computer, while the opposite is true on another computer. You should compare the efficiency of the algorithms independently of a particular computer.
- What data should the programs use? Perhaps the most important difficulty on this list is the selection of the data for the programs to use. There is always the danger that you could select instances of the problem for which one of the algorithms runs uncharacteristically fast. For example, when comparing a sequential search and a binary search of a sorted array, you might search for an item that happens to be the smallest item in the array. In such a case, the sequential search finds the item more quickly than the binary search because the item is first in the array and is thus the first item that the sequential search examines. Any analysis of efficiency must be independent of specific data.

To overcome these difficulties, computer scientists employ mathematical techniques that analyze algorithms independently of specific implementations, computers, or data. You begin this analysis by counting the number of significant operations in a particular solution, as the next section describes.

VideoNote
Measuring algorithmic efficiency

Consider efficiency when selecting an algorithm

A comparison of algorithms should focus on significant differences in efficiency

## Three difficulties

 with comparing programs instead of algorithmsAlgorithm analysis should be independent of specific implementations, computers, and data

Counting an algorithm's operations is a way to assess its efficiency

Displaying the data in a linked chain of $n$ nodes requires time proportional to $n$

Note: In general, you should avoid analyzing an algorithm solely by studying the running times of a specific implementation. Running times are influenced by such factors as programming style, the particular computer, and the data on which the program is run.

### 10.2.1 The Execution Time of Algorithms

Previous chapters have informally compared different solutions to a given problem by looking at the number of operations that each solution required. For example, Chapter 4 compared array-based and link-based implementations of the ADT bag. An array-based bag can access its $i^{\text {th }}$ item directly in one step, because the item is stored in items [i-1]. A link-based bag, however, must traverse a chain of linked nodes from its beginning until the $i^{\text {th }}$ node is reached and would therefore require $i$ steps.

An algorithm's execution time is related to the number of operations it requires. This is usually expressed in terms of the number, $n$, of items the algorithm must process. Counting an algorithm's operations - if possible-is a way to assess its efficiency. Let's consider a few other examples.

Traversal of linked nodes. As an example of traversing a chain of linked nodes, consider displaying the data in such a chain. If headPtr points to the first node in the chain, the following $\mathrm{C}++$ statements display its data:

```
Node<ItemType>* curPtr = headPtr; }\leftarrow1\mathrm{ assignment
while (curPtr != nullptr) }\leftarrown+1\mathrm{ comparisons
{
    cout << curPtr->getItem() < end7; }\leftarrown\mathrm{ writes
    curPtr = curPtr->getNext(); }\leftarrown\mathrm{ assignments
} // end while
```

If we have $n$ nodes, these statements require $n+1$ assignments, $n+1$ comparisons, and $n$ write operations. If each assignment, comparison, and write operation requires, respectively, $a, c$, and $w$ time units, the statements require $(n+1) \times(a+c)+n \times w$ time units. ${ }^{1}$ Thus, the time required to write $n$ nodes is proportional to $n$. This conclusion makes sense intuitively: It takes longer to display, or traverse, a linked chain of 100 nodes than it does a linked chain of 10 nodes.


Note: You must traverse a linked chain to access its $i^{\text {th }}$ node. The access time is, therefore, directly proportional to $i$. On the other hand, you can access array items directly, with equal access time for each item.

The Towers of Hanoi. Chapter 2 proved recursively that the solution to the Towers of Hanoi problem with $n$ disks requires $2^{n}-1$ moves. If each move requires the same time $m$, the solution requires $\left(2^{n}-1\right) \times m$ time units. As you will soon see, this time requirement increases rapidly as the number of disks increases.

Nested loops. Consider an algorithm that contains nested loops of the following form:

```
for ( \(i=1\) through \(n\) )
    for ( \(j=1\) through \(i\) )
        for ( \(k=1\) through 5 )
        Task T
```

[^8]If task $T$ requires $t$ time units, the innermost loop on $k$ requires $5 \times t$ time units. The loop on $j$ requires $5 \times t \times i$ time units, and the outermost loop on $i$ requires

$$
\sum_{i=1}^{n}(5 \times t \times i)=5 \times t \times(1+2+\cdots+n)=5 \times t \times n \times(n+1) / 2
$$

time units.

Question 1 How many comparisons of array items do the following loops contain?

```
for (j = 1; j <= n-1; j++)
{
    i = j + 1;
        do
        {
            if (theArray[i] < theArray[j])
                swap(theArray[i], theArray[j]);
            i++;
        } while (i <= n);
} // end for
```

Question 2 Repeat Question 1, replacing the statement $i=j+1$ with $i=j$.

### 10.2.2 Algorithm Growth Rates

As you can see, the previous examples derive an algorithm's time requirement as a function of the problem size. The way to measure a problem's size depends on the application-typical examples are the number of nodes in a linked chain, the number of disks in the Towers of Hanoi problem, the size of an array, or the number of items in a stack. Thus, we reached conclusions such as

## Algorithm A requires $n^{2} / 5$ time units to solve a problem of size $n$ <br> Algorithm $B$ requires $5 \times n$ time units to solve a problem of size $n$

The time units in the previous two statements must be the same before you can compare the efficiency of the two algorithms. If we had written

Algorithm A requires $n^{2} / 5$ seconds to solve a problem of size n
our earlier discussion indicates the difficulties with such a statement: On what computer does the algorithm require $n^{2} / 5$ seconds? What implementation of the algorithm requires $n^{2} / 5$ seconds? What data caused the algorithm to require $n^{2} / 5$ seconds?

What specifically do you want to know about the time requirement of an algorithm? The most important thing to learn is how quickly the algorithm's time requirement grows as a function of the problem size. Statements such as

## Algorithm A requires time proportional to $n^{2}$ <br> Algorithm $B$ requires time proportional to $n$

each express an algorithm's proportional time requirement, or growth rate, and enable you to compare algorithm $A$ with another algorithm $B$. Although you cannot determine the exact time requirement for either algorithm $A$ or algorithm $B$ from these statements, you can determine that for large problems, $B$ requires significantly less time than $A$. That is, $B$ 's time requirement-as a function of the problem size $n$-increases at a slower rate than $A$ 's time requirement, because $n$ increases at a slower rate than $n^{2}$. Even if $B$ actually requires $5 \times n$ seconds and $A$ actually requires $n^{2} / 5$ seconds, $B$ eventually requires significantly less time than $A$, as $n$ increases. Figure 10-1 illustrates this fact. Thus, a

Measure an algorithm's time requirement as a function of the problem size

Compare algorithm efficiencies for large problems

FIGURE 10-1 Time requirements as a function of the problem size $n$

conclusion such as " $A$ requires time proportional to $n^{22}$ " is exactly the kind of statement that characterizes the inherent efficiency of an algorithm independently of such factors as particular computers and implementations.

Figure 10-1 also shows that $A$ 's time requirement does not exceed $B$ 's until $n$ exceeds 25 . Algorithm efficiency is typically a concern for large problems only. The time requirements for small problems are generally not large enough to matter. Thus, our analyses assume large values of $n$.


Note: When comparing the efficiency of various solutions, look only at significant differences. This rule is consistent with a multidimensional view of the cost of a computer program.

### 10.2.3 Analysis and Big O Notation

Algorithm A requires time proportional to $f(n)$
Algorithm $A$ is said to be order $\boldsymbol{f}(\boldsymbol{n})$, which is denoted as $\mathbf{O}(\boldsymbol{f}(\boldsymbol{n})$ ). The function $f(n)$ is called the algorithm's growth-rate function. Because the notation uses the capital letter O to denote order, it is called the $\operatorname{Big} \mathbf{O}$ notation. If a problem of size $n$ requires time that is directly proportional to $n$, the problem is $\mathrm{O}(n)$-that is, order $n$. If the time requirement is directly proportional to $n^{2}$, the problem is $\mathrm{O}\left(n^{2}\right)$, and so on.

The following definition formalizes these ideas:

## Note: Definition of the order of an algorithm

Algorithm $A$ is order $f(n)$-denoted $\mathrm{O}(f(n))$-if constants $k$ and $n_{0}$ exist such that $A$ requires no more than $k \times f(n)$ time units to solve a problem of size $n \geq n_{0}$.

The requirement $n \geq n_{0}$ in the definition of $\mathrm{O}(f(n))$ formalizes the notion of sufficiently large problems. In general, many values of $k$ and $n$ can satisfy the definition.

The following examples illustrate the definition:

- Suppose that an algorithm requires $n^{2}-3 \times n+10$ seconds to solve a problem of size $n$. If constants $k$ and $n_{0}$ exist such that

$$
k \times n^{2}>n^{2}-3 \times n+10 \text { for all } n \geq n_{0}
$$

the algorithm is $\mathrm{O}\left(n^{2}\right)$. In fact, if $k$ is 3 and $n_{0}$ is 2 ,

$$
3 \times n^{2}>n^{2}-3 \times n+10 \text { for all } n \geq 2
$$

as Figure 10-2 illustrates. Thus, the algorithm requires no more than $k \times n^{2}$ time units for $n \geq n_{0}$ and so is $\mathrm{O}\left(n^{2}\right)$.

- Previously in Section 10.2.1, we found that displaying a linked chain's first $n$ items requires

$$
(n+1)(a+c)+n \times w
$$

time units. Because $2 \times n \geq n+1$ for $n \geq 1$,

$$
(2 \times n) \times(a+c)+n \times w \geq(n+1) \times(a+c)+n \times w \text { for } n \geq 1
$$

After factoring $n$ on the left side of the inequality, we have

$$
(2 \times a+2 \times c+w) \times n \geq(n+1) \times(a+c)+n \times w \text { for } n \geq 1
$$

Thus, this task is $\mathrm{O}(n)$. Here, $k$ is $2 \times a+2 \times c+w$ and $n_{0}$ is 1 .

- Similarly, the solution to the Towers of Hanoi problem requires $\left(2^{n}-1\right) \times m$ time units. Because

$$
m \times 2^{n}>\left(2^{n}-1\right) \times m \text { for } n \geq 1
$$

the solution is $\mathrm{O}\left(2^{n}\right)$.
The requirement $n \geq n_{0}$ in the definition of $\mathrm{O}(f(n))$ means that the time estimate is correct for sufficiently large problems. In other words, the time estimate is too small for at most a finite number of problem sizes. For example, the function $\log n$ takes on the value 0 when $n$ is 1 . Thus, the fact that $k \times \log 1$ is 0 for all constants $k$ implies an unrealistic time requirement; presumably, all algorithms require more than 0 time units, even to solve a problem of size 1 . Thus, you can discount problems of size $n=1$ if $f(n)$ is $\log n$.

FIGURE 10-2 The graphs of $3 \times n^{2}$ and $n^{2}-3 \times n+10$


Order of growth of some common functions

Intuitive interpretations of growth-rate functions

To dramatize further the significance of an algorithm's proportional growth rate, consider the table and graph in Figure 10-3. The table in part $a$ gives, for various values of $n$, the approximate values of some common growth-rate functions, which are listed in order of growth:

$$
\mathrm{O}(1)<\mathrm{O}\left(\log _{2} n\right)<\mathrm{O}(n)<\mathrm{O}\left(n \times \log _{2} n\right)<\mathrm{O}\left(n^{2}\right)<\mathrm{O}\left(n^{3}\right)<\mathrm{O}\left(2^{n}\right)
$$

The table demonstrates the relative speed at which the values of the functions grow. Figure 10-3b represents the growth-rate functions graphically. ${ }^{2}$

These growth-rate functions have the following intuitive interpretations:

```
1
```

$\log _{2} n \quad$ The time requirement for a logarithmic algorithm increases slowly as the problem size increases. If you square the problem size, you only double its time requirement. Later you will see that the recursive binary search algorithm that you studied in Chapter 2 has this behavior. Recall that a binary search halves an array and then searches one of the halves. Typical logarithmic algorithms solve a problem by solving a smaller constant fraction of the problem.

The base of the log does not affect a logarithmic growth rate, so you can omit it in a growth-rate function. Exercise 8 at the end of this chapter asks you to show why this is true.
$n$
The time requirement for an $\mathrm{O}(n)$, or linear, algorithm increases directly with the size of the problem. If you square the problem size, you also square its time requirement.
$n \times \log _{2} n \quad$ The time requirement for an $\mathrm{O}\left(n \times \log _{2} n\right)$ algorithm increases more rapidly than a linear algorithm. Such algorithms usually divide a problem into smaller problems that are each solved separately. The next chapter examines an example of this algorithm typethe merge sort.
$n^{2} \quad$ The time requirement for an $\mathrm{O}\left(n^{2}\right)$, or quadratic, algorithm increases rapidly with the size of the problem. Algorithms that use two nested loops are often quadratic. Such algorithms are practical only for small problems. The next chapter covers several quadratic sorting algorithms.
$n^{3} \quad$ The time requirement for an $\mathrm{O}\left(n^{3}\right)$, or cubic, algorithm increases more rapidly with the size of the problem than the time requirement for a quadratic algorithm. Algorithms that use three nested loops are often cubic, and are practical only for small problems.
$2^{n} \quad$ As the size of a problem increases, the time requirement for an exponential algorithm usually increases too rapidly to be practical.

If algorithm $A$ requires time that is proportional to function $f$ and algorithm $B$ requires time that is proportional to a slower-growing function $g$, it is apparent that $B$ is always significantly more efficient than $A$ for large enough problems. For large problems, the proportional growth rate dominates all other factors in determining an algorithm's efficiency.

Properties of growth-rate functions. Several mathematical properties of Big O notation help to simplify the analysis of an algorithm. As we discuss these properties, you should keep in mind that $\mathrm{O}(f(n))$ means "is of order $f(n)$ " or "has order $f(n)$." O is not a function.

[^9]FIGURE 10-3 A comparison of growth-rate functions: (a) in tabular form; (b) in graphical form
(a)

| Function | n |  |  |  |  |  |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: |
|  | 10 | 100 | 1,000 | 10,000 | 100,000 1 | 1,000,000 |
| 1 | 1 | 1 | 1 | 1 | 1 | 1 |
| $\log _{2} n$ | 3 | 6 | 9 | 13 | 16 | 19 |
| n | 10 | $10^{2}$ | $10^{3}$ | $10^{4}$ | $10^{5}$ | $10^{6}$ |
| $n \times \log _{2} n$ | 30 | 664 | 9,965 | $10^{5}$ | $10^{6}$ | $10^{7}$ |
| $\mathrm{n}^{2}$ | $10^{2}$ | $10^{4}$ | $10^{6}$ | $10^{8}$ | $10^{10}$ | $10^{12}$ |
| $\mathrm{n}^{3}$ | $10^{3}$ | $10^{6}$ | $10^{9}$ | $10^{12}$ | $10^{15}$ | $10^{18}$ |
| $2^{n}$ | $10^{3}$ | $10^{30}$ | $10^{301}$ | $10^{3,010}$ | - 1030,103 | $0310^{301,030}$ |

(b)


1. You can ignore low-order terms in an algorithm's growth-rate function. For example, if an algorithm is $\mathrm{O}\left(n^{3}+4 \times n^{2}+3 \times n\right)$, it is also $\mathrm{O}\left(n^{3}\right)$. By examining the table in Figure 10-3a, you can see that the $n^{3}$ term is significantly larger than either $4 \times n^{2}$ or $3 \times n$, particularly for large values of $n$. For large $n$, the growth rate of $n^{3}+4 \times n^{2}+3 \times n$ is the same as the growth rate of $n^{3}$. It is the growth rate of $f(n)$, not the value of $f(n)$, that is important here. Thus, even if an algorithm is $\mathrm{O}\left(n^{3}+4 \times n^{2}+3 \times n\right)$, we say that it is simply $\mathrm{O}\left(n^{3}\right)$. In general, you can usually conclude that an algorithm is $\mathrm{O}(f(n))$, where $f$ is a function similar to the ones listed in Figure 10-3.

Some properties of growth-rate functions

An algorithm can require different times to solve different problems of the same size

An array-based getEntry is $\mathrm{O}(1)$

A link-based getEntry is $\mathrm{O}(n)$
2. You can ignore a multiplicative constant in the high-order term of an algorithm's growth-rate function. For example, if an algorithm is $\mathrm{O}\left(5 \times n^{3}\right)$, it is also $\mathrm{O}\left(n^{3}\right)$. This observation follows from the definition of $\mathrm{O}(f(n))$, if you let $k=5$.
3. $\mathbf{O}(f(n))+\mathbf{O}(g(n))=\mathbf{O}(f(n)+g(n))$. You can combine growth-rate functions. For example, if an algorithm is $\mathrm{O}\left(n^{2}\right)+\mathrm{O}(n)$, it is also $\mathrm{O}\left(n^{2}+n\right)$, which you write simply as $\mathrm{O}\left(n^{2}\right)$ by applying property 1 . Analogous rules hold for multiplication.
These properties imply that you need only an estimate of the time requirement to obtain an algorithm's growth rate; you do not need an exact statement of an algorithm's time requirement, which is fortunate because deriving the exact time requirement is often difficult and sometimes impossible.

Worst-case and average-case analyses. A particular algorithm might require different times to solve different problems of the same size. For example, the time that an algorithm requires to search $n$ items might depend on the nature of the items. Usually you consider the maximum amount of time that an algorithm can require to solve a problem of size $n$-that is, the worst case. A worst-case analysis concludes that algorithm $A$ is $\mathrm{O}(f(n))$ if, in the worst case, $A$ requires no more than $k \times f(n)$ time units to solve a problem of size $n$ for all but a finite number of values of $n$. Although a worst-case analysis can produce a pessimistic time estimate, such an estimate does not mean that your algorithm is always slow. Instead, you have shown that the algorithm is never slower than your estimate. Realize, however, that an algorithm's worst case might happen rarely, if at all, in practice.

An average-case analysis attempts to determine the average amount of time that an algorithm requires to solve problems of size $n$. In an average-case analysis, $A$ is $\mathrm{O}(f(n))$ if the average amount of time that $A$ requires to solve a problem of size $n$ is no more than $k \times f(n)$ time units for all but a finite number of values of $n$. Average-case analysis is, in general, far more difficult to perform than worst-case analysis. One difficulty is determining the relative probabilities of encountering various problems of a given size; another is determining the distributions of various data values. Worst-case analysis is easier to calculate and is thus more common.

Note: While manipulating the Big O notation, remember that $\mathrm{O}(f(n))$ represents an inequality. It is not a function but simply a notation that means "is of order $f(n)$ " or "has order $f(n)$."
$?$
Question 3 What order is an algorithm that has as a growth-rate function
a. $8 \times n^{3}-9 \times n$
b. $7 \times \log _{2} n+20$
c. $7 \times \log _{2} n+n$

### 10.2.4 Keeping Your Perspective

Before continuing with additional analyses of specific algorithms, a few words about perspective are appropriate. You know that you can access the $n^{\text {th }}$ item in an array directly. This access is independent of $n$ : accessing the $100^{\text {th }}$ item takes the same time as it does to access the first item in the array. For an ADT list of $n$ items, the array-based implementation of the retrieval operation getEntry is $\mathrm{O}(1)$. However, the link-based implementation of getEntry requires $n$ steps to traverse a chain until it reaches the $n^{\text {th }}$ item, and so is $\mathrm{O}(n)$.

Throughout the course of an analysis, you should always keep in mind that you are interested only in significant differences in efficiency. Is the difference in efficiency for the two implementations of getEntry significant? As the size of the list grows, the link-based implementation might require more time to retrieve the desired entry, because its node can be farther away from the beginning of the chain. In contrast, regardless of how large the list is, the array-based implementation always requires the same constant amount of time to retrieve any particular item. Thus, no matter what time difference is significant to you, it is reached if the list is large enough. In this example, observe that the difference in efficiency for the two implementations is worth considering only when the problem is large enough. If the list never has more than 25 items, for example, the difference in the implementations is not significant at all.

Now consider an application-such as a word processor's spelling checker-that frequently retrieves items from a list but rarely inserts or removes an item. Because getEntry for an array-based list is faster than a link-based list's getEntry, you should choose an array-based implementation of the list for the application. On the other hand, if an application requires frequent insertions and removals but rarely retrieves an item, you should choose a link-based implementation of the list. The most appropriate implementation of an ADT for a given application strongly depends on how frequently the application performs these operations. More examples of this point are presented in the next chapter.

The response time of some ADT operations, however, can be crucial, even if you seldom use them. For example, an air traffic control system could include an emergency operation to resolve the impending collision of two airplanes. Clearly, this operation must occur quickly, even if it is rarely used. Thus, before you choose an implementation for an ADT, you should know what operations a particular application requires, approximately how often the application performs each operation, and the response times that the application requires of each operation.

In the next section we compare a searching algorithm that is $\mathrm{O}(n)$ with one that is $\mathrm{O}\left(\log _{2} n\right)$. While it is true that an $\mathrm{O}\left(\log _{2} n\right)$ searching algorithm requires significantly less time on large arrays than an $\mathrm{O}(n)$ algorithm requires, on small arrays-say $n<25$-the time requirements might not be significantly different at all. In fact, it is entirely possible that, because of factors such as the size of the constant $k$ in the definition of Big O , the $\mathrm{O}(n)$ algorithm runs faster on small problems. It is only on large problems that the slower growth rate of an algorithm necessarily gives it a significant advantage. Figure 10-1 illustrated this phenomenon.

Thus, in general, if the maximum size of a given problem is small, the time requirements of any two solutions for that problem likely will not differ significantly. If you know that your problem size will always be small, do not overanalyze; simply choose the algorithm that is easiest to understand, verify, and code.

Frequently, when evaluating an algorithm's efficiency, you have to weigh carefully the trade-offs between a solution's execution time requirements and its memory requirements. You are rarely able to make a statement as strong as "This approach is the best one for performing the task." A solution that requires a relatively small amount of computer time often also requires a relatively large amount of memory. It may not even be possible to say that one solution requires less time than another. Solution $A$ may perform some components of the task faster than solution $B$, while solution $B$ performs other components of the task faster than solution $A$. Often you must analyze the solutions in light of a particular application.

In summary, it is important to examine an algorithm for both style and efficiency. The analysis should focus only on gross differences in efficiency and not reward coding tricks that save milliseconds. Any finer differences in efficiency are likely to interact with coding issues, which you should not allow to interfere with the development of your programming style. If you find an algorithm that is significantly more efficient than others, you should select it, unless you know that the maximum problem size is quite small. If you are solving only small problems, it is possible that a less efficient algorithm would be more appropriate. That is, other factors, such as the simplicity of the algorithm,

When choosing an implementation of an ADT, consider how frequently particular ADT operations occur in a given application

Some seldom-used but critical operations must be efficient

If the problem size is always small, you can probably ignore an algorithm's efficiency

Weigh the trade-offs between an algorithm's time requirements and its memory requirements

Compare algorithms for both style and efficiency

## Analysis of an

 algorithm's time efficiency focuses on large problemsSequential search. Worst case: O(n); average case: $\mathrm{O}(n)$; best case: $\mathrm{O}(1)$
could become more significant than minor differences in efficiency. In fact, analyzing the time efficiency of an algorithm implicitly assumes that an algorithm is used to solve large problems. This assumption allows you to focus on growth rates, because regardless of other factors, an algorithm with a slow growth rate requires less time than an algorithm with a fast growth rate, provided that the problems to be solved are sufficiently large.


Note: If a problem is small, do not overanalyze it. In such a situation, the primary concern should be simplicity. For example, if you are searching an array that contains only a small number of items-say, fewer than 25 -a simple $\mathrm{O}(n)$ algorithm such as a sequential search is appropriate.

### 10.2.5 The Efficiency of Searching Algorithms

As another example of analyzing the time efficiency of an algorithm, consider the efficiency of two search algorithms: the sequential search and the binary search of an array.

Sequential search. In a sequential search of an array of $n$ items, you look at each item in turn, beginning with the first one, until either you find the desired item or you reach the end of the data collection. In the best case, the desired item is the first one that you examine, so only one comparison is necessary. Thus, in the best case, a sequential search is $O(1)$. In the worst case, the desired item is the last one you examine, so $n$ comparisons are necessary. Thus, in the worst case, the algorithm is $\mathrm{O}(n)$. In the average case, you find the desired item in the middle of the collection, making $n / 2$ comparisons. Thus, the algorithm is $\mathrm{O}(n)$ in the average case.

What is the algorithm's order when you do not find the desired item? Does the algorithm's order depend on whether the initial data is sorted? These questions are left for you in Checkpoint Question 4 at the end of this section.

Binary search. Is a binary search of an array more efficient than a sequential search? The binary search algorithm, which Chapter 2 presented, searches a sorted array for a particular item by repeatedly dividing the array in half. The algorithm determines which half the item must be in-if it is indeed present-and discards the other half. Thus, the binary search algorithm searches successively smaller arrays: The size of a given array is approximately one-half the size of the array previously searched.

At each division, the algorithm makes a comparison. How many comparisons does the algorithm make when it searches an array of $n$ items? The exact answer depends, of course, on where the soughtfor item resides in the array. However, you can compute the maximum number of comparisons that a binary search requires - that is, the worst case. The number of comparisons is equal to the number of times that the algorithm divides the array in half. Suppose that $n=2^{k}$ for some $k$. The search requires the following steps:

1. Inspect the middle item of an array of size $n$.
2. Inspect the middle item of an array of size $n / 2$.
3. Inspect the middle item of an array of size $n / 2^{2}$, and so on.

To inspect the middle item of an array, you must first divide the array in half. If you halve an array of $n$ items, then divide one of those halves in half, and continue dividing halves until only one item remains, you have performed $k$ divisions. This is true because $n / 2^{k}=1$. (Remember, we assumed that $n=2^{k}$.) In the worst case, the algorithm performs $k$ divisions and, therefore, $k$ comparisons. Because $n=2^{k}$,

$$
k=\log _{2} n
$$

Thus, the algorithm is $\mathrm{O}\left(\log _{2} n\right)$ in the worst case when $n=2^{k}$.
What if $n$ is not a power of 2 ? You can easily find the smallest $k$ such that

$$
2^{k-1}<n<2^{k}
$$

(For example, if $n$ is 30 , then $k=5$, because $2^{4}=16<30<32=2^{5}$.) The algorithm still requires at most $k$ divisions to obtain a subarray with one item. Now it follows that

$$
\begin{aligned}
& k-1<\log _{2} n<k \\
& k<1+\log _{2} n<k+1 \\
& k=1+\log _{2} n \text { rounded down }
\end{aligned}
$$

Thus, the algorithm is still $\mathrm{O}\left(\log _{2} n\right)$ in the worst case when $n \neq 2^{k}$. In general, the algorithm is $\mathrm{O}\left(\log _{2} n\right)$ in the worst case for any $n$.

Is a binary search faster than a sequential search? Much faster! For example $\log _{2} 1,000,000 \approx 19$, so a binary search of 1 million sorted items requires at most 20 comparisons, but a sequential search of the same items can require 1 million comparisons! For large arrays, the binary search has an enormous advantage over a sequential search.

Realize, however, that maintaining the array in sorted order requires an overhead cost, which can be substantial. The next chapter examines the cost of sorting an array.

Note: If you are searching a very large array, an $\mathrm{O}(n)$ algorithm is probably too inefficient to use.


Question 4 Consider a sequential search of $n$ data items.
a. If the data items are sorted into ascending order, how can you determine that your desired item is not in the data collection without always making $n$ comparisons?
b. What is the order of the sequential search algorithm when the desired item is not in the data collection? Do this for both sorted and unsorted data, and consider the best, average, and worst cases.
c. Show that if the sequential search algorithm finds the desired item in the data collection, the algorithm's order does not depend upon whether or not the data items are sorted.

## SUMMARY

1. Using Big O notation, you measure an algorithm's time requirement as a function of the problem size by using a growth-rate function. This approach enables you to analyze the efficiency of an algorithm without regard for such factors as computer speed and programming skill that are beyond your control.
2. When you compare the inherent efficiency of algorithms, you examine their growth-rate functions when the problems are large. Only significant differences in growth-rate functions are meaningful.
3. Worst-case analysis considers the maximum amount of work an algorithm requires on a problem of a given size, while average-case analysis considers the expected amount of work that it requires.
4. Analyzing an algorithm's time requirement will help you to choose an implementation for an abstract data type. If your application frequently uses particular ADT operations, your implementation should be efficient for at least those operations.

## Exercises

1. Using Big O notation, indicate the time requirement of each of the following tasks in the worst case. Describe any assumptions that you make.
a. After arriving at a party, you shake hands with each person there.
b. Each person in a room shakes hands with everyone else in the room.
c. You climb a flight of stairs.
d. You slide down the banister.
e. After entering an elevator, you press a button to choose a floor.
f. You ride the elevator from the ground floor up to the $n$th floor.
g. You read a book twice.

Describe a way to climb from the bottom of a flight of stairs to the top in time that is no better than $\mathrm{O}\left(n^{2}\right)$.
3. Using Big O notation, indicate the time requirement of each of the following tasks in the worst case.
a. Computing the sum of the first $n$ even integers by using a for loop
b. Displaying all $n$ integers in an array
c. Displaying all $n$ integers in a sorted linked chain
d. Displaying all $n$ names in an array of linked chains
e. Displaying one array element
f. Displaying the last integer in a linked chain
g. Searching an array of $n$ items for a particular value by using a sequential search
h. Searching an array of $n$ items for a particular value by using a binary search
i. Adding an item to a stack of $n$ items
j. Adding an item to a bag of $n$ items
4. Suppose that your implementation of a particular algorithm appears in $\mathrm{C}++$ as

```
for (int pass = 1; pass <= n; pass++)
{
    for (int index = 0; index < n; index++)
    {
        for (int count = 1; count < 10; count++)
        {
        } // end for
    } // end for
} // end for
```

The previous code shows only the repetition in the algorithm, not the computations that occur within the loops. These computations, however, are independent of $n$. What is the Big O of the algorithm? Justify your answer.
5. Consider the following $\mathrm{C}++$ function f , which calls the function swap. Assume that swap exists and simply swaps the contents of its two arguments. Do not be concerned with f's purpose.

```
void f(int theArray[], int n)
{
    for (int j = 0; j < n; ++j)
    {
        int i = 0;
        while (i <= j)
        {
            if (theArray[i] < theArray[j])
```

```
                swap(theArray[i], theArray[j]);
            i++;
        } // end while
    } // end for
} // end f
```

How many comparisons does f perform?
6. For large arrays, and in the worst case, is a sequential search faster than a binary search? Explain.
7. Show that any polynomial $f(x)=c_{n} x^{n}+c_{n-1} x^{n-1}+\cdots+c_{1} x+c_{0}$ is $\mathrm{O}\left(x^{n}\right)$.
8. Show that for all constants $a, b>1, f(n)$ is $\mathrm{O}\left(\log _{a} n\right)$ if and only if $f(n)$ is $\mathrm{O}\left(\log _{b} n\right)$. Thus, you can omit the base when you write $\mathrm{O}(\log n)$. Hint: Use the identity $\log _{a} n=\log _{b} n / \log _{b} a$ for all constants $a, b>1$.
9. Show that $7 n^{2}+5 n$ is $\operatorname{not} \mathrm{O}(n)$.
10. Consider an array of length $n$ containing positive and negative integers in random order. Write $\mathrm{C}++$ code that rearranges the integers so that the negative integers appear before the positive integers. Your solution should use
a. $\mathrm{O}\left(n^{2}\right)$ operations
b. $\mathrm{O}(n)$ operations

## Programming Problems

For the following programming problems, you need to time a section of code in $C++$. For example, the following statements time the execution of the function doSomething:

```
#include <ctime>
clock_t start = clock();
doSomething();
clock_t finish = clock();
doub7e overa11Time = static_cast<doub7e>(finish - start) / CLOCKS_PER_SEC;
```

1. Consider the following two loops:
```
// Loop A
for (i = 1; i <= n; i++)
    for (j = 1; j <= 10000; j++)
        sum = sum + j;
// Loop B
for (i = 1; i <= n; i++)
    for (j = 1; j <= n; j++)
        sum = sum + j;
```

What is the Big O of each loop? Design and implement an experiment to find a value of $n$ for which Loop B is faster than Loop A.
2. Repeat the previous project, but use the following for Loop B:

```
// Loop B
for (i = 1; i <= n; i++)
    for (j = 1; j <= n; j++)
        for (k = 1; k <= j; k++)
            sum = sum + k;
```

[^10]3. Write a C++ program that implements the following three algorithms and times them for various values of $n$. The program should display a table of the run times of each algorithm for various values of $n$.

```
// Algorithm A
sum = 0
for i = 1 to n
    sum = sum + i
```

```
// Algorithm B
```

// Algorithm B
sum = 0
sum = 0
for i = 1 to n
for i = 1 to n
{
{
for j = 1 to i
for j = 1 to i
sum = sum + 1
sum = sum + 1
}

```
}
```

For each of the following problems, you will need an implementation of the ADT bag whose method remove removes a random entry instead of a specific one.
4. Suppose that you have several numbered billiard balls on a pool table. At each step you remove a billiard ball from the table. If the ball removed is numbered $n$, you replace it with $n$ balls whose number is $n / 2$, where the division is truncated to an integer. For example, if you remove the 5 ball, you replace it with five 2 balls. Using Big O notation, predict the time requirement for this algorithm when initially the pool table contains only the $n$ ball.

Write a program that simulates this process. Use a bag of positive integers to represent the balls on the pool table. Time the actual execution of the program for various values of $n$ and plot its performance as a function of $n$. Compare your results with your predicted time requirements.
5. Repeat the previous project, but instead replace the $n$ ball with $n$ balls randomly numbered less than $n$.
6. In mythology, the Hydra was a monster with many heads. Every time the hero chopped off a head, two smaller heads would grow in its place. Fortunately for the hero, if the head was small enough, he could chop it off without two more growing in its place. To kill the Hydra, all our hero needed to do was to chop off all the heads.

Write a program that simulates the Hydra. Instead of heads, we will use strings. A bag of strings, then, represents the Hydra. Every time you remove a string from the bag, delete the first letter of the string and put two copies of the remaining string back into the bag. For example, if you remove HYDRA, you add two copies of YDRA to the bag. If you remove a one-letter word, you add nothing to the bag. To begin, read one word from the keyboard and place it into an empty bag. The Hydra dies when the bag becomes empty.

Using Big O notation, predict the time requirement for this algorithm in terms of the number $n$ of characters in the initial string. Then time the actual execution of the program for various values of $n$ and plot its performance as a function of $n$.

## Sorting Algorithms and Their Efficiency

## Chapter

 11
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Chapter 10 Algorithm Efficiency

This chapter examines the important topic of sorting data. You will study some simple algorithms, which you may have seen before, and some more sophisticated recursive algorithms. Sorting algorithms provide varied and relatively easy examples of the analysis of efficiency.

## The sorts in this

 chapter are internal sorts
### 11.1 Basic Sorting Algorithms

Sorting is a process that organizes a collection of data into either ascending ${ }^{1}$ or descending order. The need for sorting arises in many situations. You may simply want to sort a collection of data before including it in a report. Often, however, you must perform a sort as an initialization step for certain algorithms. For example, searching for data is one of the most common tasks performed by computers. When the collection of data to be searched is large, an efficient technique for searching-such as the binary search algorithm-is desirable. However, the binary search algorithm requires that the data be sorted. Thus, sorting the data is a step that must precede a binary search on a collection of data that is not already sorted. Good sorting algorithms, therefore, are quite valuable.

You can organize sorting algorithms into two categories. An internal sort requires that the collection of data fit entirely in the computer's main memory. The algorithms in this chapter are internal sorting algorithms. You use an external sort when the collection of data will not fit in the computer's main memory all at once but must reside in secondary storage, such as on a disk. Chapter 21 examines external sorts.

The data items to be sorted might be integers, character strings, or even objects. It is easy to imagine the results of sorting a collection of integers or character strings, but consider a collection of objects. If each object contains only one data member, sorting the objects is really no different from sorting a collection of integers. However, when each object contains several data members, you must know which data member determines the order of the entire object within the collection of data. This data member is called the sort key. For example, if the objects represent people, you might want to sort on their names, their ages, or their zip codes. Regardless of your choice of sort key, the sorting algorithm orders entire objects based on only one data member, the sort key.

For simplicity, the examples in this chapter sort quantities such as numbers or strings. All algorithms in this chapter sort the data into ascending order. Modifying these algorithms to sort data into descending order is simple. Finally, each example assumes that the data resides in an array.

### 11.1.1 The Selection Sort

Imagine some data that you can examine all at once. To sort it, you could select the largest item and put it in its place, select the next largest and put it in its place, and so on. For a card player, this process is analogous to looking at an entire hand of cards and ordering it by selecting cards one at a time in their proper order. The selection sort formalizes these intuitive notions. To sort an array into ascending order, you first search it for the largest item. Because you want the largest item to be in the last position of the array, you swap the last item with the largest item, even if these items happen to be identical. Now, ignoring the last-and largest-item of the array, you search the rest of the array for its largest item and swap it with its last item, which is the next-to-last item in the original array. You continue until you have selected and swapped $n-1$ of the $n$ items in the array. The remaining item, which is now in the first position of the array, is in its proper order, so it is not considered further.

Figure 11-1 provides an example of a selection sort. Beginning with five integers, you select the largest-37-and swap it with the last integer-13. (As the items in this figure are ordered, they are shaded in blue. This convention will be used throughout this chapter.) Next you select the largest integer - 29 - from among the first four integers in the array and swap it with the next-to-last integer in the array-13. Notice that the next selection-14 -is already in its proper position, but the algorithm ignores this fact and performs a swap of 14 with itself. It is more efficient in general to occasionally perform an unnecessary swap than it is to continually ask whether the swap is necessary. Finally, you select the 13 and swap it with the item in the second position of the array-10. The array is now sorted into ascending order.

[^11]FIGURE 11-1 A selection sort of an array of five integers

Gray elements are selected;
blue elements comprise the sorted portion of the array.

| Initial array: | 29 | 10 | 14 | 37 | 13 |
| :---: | :---: | :---: | :---: | :---: | :---: |
| After 1st swap: | 29 | 10 | 14 | 13 | 37 |
| After 2nd swap: | 13 | 10 | 14 | 29 | 37 |
| After 3rd swap: | 13 | 10 | 14 | 29 | 37 |
| After 4th swap: | 10 | 13 | 14 | 29 | 37 |

Question 1 Trace the selection sort as it sorts the following array into ascending order: 208040256030

Question 2 Repeat the previous question, but instead sort the array into descending order.

Listing 11-1 gives a C++ function, and the two functions it calls, that performs a selection sort on an array theArray of $n$ items, assuming that ItemType is the data type of the array's entries.

## LISTING 11-1 An implementation of the selection sort

```
/** Finds the largest item in an array.
    @pre The size of the array is >= 1.
    @post The arguments are unchanged.
    @param theArray The given array.
    @param size The number of elements in theArray.
    @return The index of the largest entry in the array. */
int findIndexofLargest(const ItemType theArray[], int size);
/** Sorts the items in an array into ascending order.
    @pre None.
    @post The array is sorted into ascending order; the size of the array
        is unchanged.
    @param theArray The array to sort.
    @param n The size of theArray. */
void selectionSort(ItemType theArray[], int n)
{
    // 1ast = index of the last item in the subarray of items yet
    // to be sorted;
    // largest = index of the largest item found
```

```
    for (int last = n - 1; last >= 1; last--)
    {
        // At this point, theArray[last+1..n-1] is sorted, and its
        // entries are greater than those in theArray[0..1ast].
        // Select the largest entry in theArray[0..7ast]
        int largest = findIndexofLargest(theArray, last+1);
        // Swap the largest entry, theArray[largest], with
        // theArray[last]
        std::swap(theArray[largest], theArray[last]);
    } // end for
} // end selectionSort
int findIndexofLargest(const ItemType theArray[], int size)
{
    int indexSoFar = 0; // Index of largest entry found so far
    for (int currentIndex = 1; currentIndex < size; currentIndex++)
    {
        // At this point, theArray[indexSoFar] >= al1 entries in
        // theArray[0..currentIndex - 1]
        if (theArray[currentIndex] > theArray[indexSoFar])
            indexSoFar = currentIndex;
    } // end for
    return indexSoFar; // Index of largest entry
} // end findIndexofLargest
```

Analysis. As you can see from the previous algorithm, sorting in general compares, exchanges, or moves items. As a first step in analyzing such algorithms, you should count these operations. Generally, such operations are more expensive than ones that control loops or manipulate array indexes, particularly when the data to be sorted is more complex than integers or characters. Thus, our approach ignores these incidental operations. You should convince yourself that by ignoring such operations we do not affect our final result. (See Exercise 1.)

Clearly, the for loop in the function selectionSort executes $n-1$ times. Thus, selectionSort calls each of the functions findIndexofLargest and swap $n-1$ times. Each call to findIndexofLargest causes its loop to execute last times (that is, size - 1 times when size is last +1 ). Thus, the $n-1$ calls to findIndexofLargest, for values of 1ast that range from $n-1$ down to 1 , cause the loop in findIndexofLargest to execute a total of

$$
(n-1)+(n-2)+\cdots+1=n \times(n-1) / 2
$$

times. Because each execution of findIndexofLargest's loop performs one comparison, the calls to findIndexofLargest require

$$
n \times(n-1) / 2
$$

comparisons.
The $n-1$ calls to swap result in $n-1$ exchanges. Each exchange requires three assignments, or data moves. Thus, the calls to swap require

$$
3 \times(n-1)
$$

moves.

Together, a selection sort of $n$ items requires

$$
n \times(n-1) / 2+3 \times(n-1)=n^{2} / 2+5 \times n / 2-3
$$

major operations. By applying the properties of the growth-rate functions given in Chapter 10, you can ignore low-order terms to get $\mathrm{O}\left(n^{2} / 2\right)$ and then ignore the multiplier $1 / 2$ to get $\mathrm{O}\left(n^{2}\right)$. Thus, the selection sort is $\mathrm{O}\left(n^{2}\right)$.

Although a selection sort does not depend on the initial arrangement of the data, which is an advantage of this algorithm, it is appropriate only for small $n$ because $\mathrm{O}\left(n^{2}\right)$ grows rapidly. While the algorithm requires $\mathrm{O}\left(n^{2}\right)$ comparisons, it requires only $\mathrm{O}(n)$ data moves. A selection sort could be a good choice over other approaches when data moves are costly but comparisons are not. Such might be the case if each data item is lengthy but the sort key is short.


Note: If you are sorting a very large array, an $\mathrm{O}\left(n^{2}\right)$ algorithm is probably too inefficient to use.

### 11.1.2 The Bubble Sort

The next sorting algorithm is one that you may have seen already. The bubble sort compares adjacent items and exchanges them if they are out of order. This sort usually requires several passes over the data. During the first pass, you compare the first two items in the array. If they are out of order, you exchange them. You then compare the items in the next pair-that is, in positions 2 and 3 of the array. If they are out of order, you exchange them. You proceed in the same manner, comparing and exchanging items two at a time, until you reach the end of the array.

Figure 11-2a illustrates the first pass of a bubble sort of an array of five integers. You compare the items in the first pair-29 and 10-and exchange them because they are out of order. Next you consider the second pair-29 and 14-and exchange these items because they are out of order. The items in the third pair- 29 and 37 -are in order, and so you do not exchange them. Finally, you exchange the items in the last pair- 37 and 13 .

Although the array is not sorted after the first pass, the largest item has "bubbled" to its proper position at the end of the array. During the second pass of the bubble sort, you return to the beginning of the array and consider pairs of items in exactly the same manner as the first pass. You do

Selection sort is $\mathrm{O}\left(n^{2}\right)$

When you order successive pairs of items, the largest item bubbles to the top (end) of the array

FIGURE 11-2
The first two passes of a bubble sort of an array of five integers


The bubble sort usually requires several passes through the array
not, however, include the last - and largest-item of the array. That is, the second pass considers the first $n-1$ items of the array. After the second pass, the second-largest item in the array will be in its proper place in the next-to-last position of the array, as Figure 11-2b illustrates. Now, ignoring the last two items, which are in order, you continue with subsequent passes until the array is sorted.

Although a bubble sort requires at most $n-1$ passes to sort the array, fewer passes might be possible to sort a particular array. Thus, you could terminate the process if no exchanges occur during any pass. The following C++ function bubbleSort uses a boolean variable to signal when an exchange occurs during a particular pass. The function uses the previous swap function.

## LISTING 11-2 An implementation of the bubble sort

```
/** Sorts the items in an array into ascending order.
    @pre None.
    @post theArray is sorted into ascending order; n is unchanged.
    @param theArray The given array.
    @param n The size of theArray. */
```

void bubbleSort(ItemType theArray[], int n)
\{
bool sorted = false; // False when swaps occur
int pass = 1 ;
while (!sorted \&\& (pass < $n$ ))
\{
// At this point, theArray[n+1-pass..n-1] is sorted
// and all of its entries are > the entries in theArray[0..n-pass]
sorted = true; // Assume sorted
for (int index $=0$; index < n - pass; index++)
\{
// At this point, al1 entries in theArray[0..index-1]
// are <= theArray[index]
int nextIndex = index + 1;
if (theArray[index] > theArray[nextIndex])
\{
// Exchange entries
std::swap(theArray[index], theArray[nextIndex]);
sorted $=$ false; // Signal exchange
\} // end if
\} // end for
// Assertion: theArray[0..n-pass-1] < theArray[n-pass]
pass++;
\} // end while
\} // end bubbleSort

Question 3 Trace the bubble sort as it sorts the following array into ascending order: 253020804060.

Question 4 Repeat the previous question, but instead sort the array into descending order.

Analysis. As was noted earlier, the bubble sort requires at most $n-1$ passes through the array. Pass 1 requires $n-1$ comparisons and at most $n-1$ exchanges; pass 2 requires $n-2$ comparisons and at most $n-2$ exchanges. In general, pass $i$ requires $n-i$ comparisons and at most $n-i$ exchanges. Therefore, in the worst case, a bubble sort will require a total of

$$
(n-1)+(n-2)+\cdots+1=n \times(n-1) / 2
$$

comparisons and the same number of exchanges. Recall that each exchange requires three data moves. Thus, altogether there are

$$
2 \times n \times(n-1)=2 \times n^{2}-2 \times n
$$

major operations in the worst case. Therefore, the bubble sort algorithm is $\mathrm{O}\left(n^{2}\right)$ in the worst case.
The best case occurs when the original data is already sorted: bubbleSort uses one pass, during which $n-1$ comparisons and no exchanges occur. Thus, the bubble sort is $\mathrm{O}(n)$ in the best case.

### 11.1.3 The Insertion Sort

Imagine once again arranging a hand of cards, but now you pick up one card at a time and insert it into its proper position; in this case, you are performing an insertion sort. Chapter 8 introduced the insertion sort algorithm in the context of a list of alphabetical names. In that example, we repeatedly called the method insert to add a string into its proper sorted order in a list.

You can use the insertion sort strategy to sort items that reside in an array. This version of the insertion sort partitions the array into two regions: sorted and unsorted, as Figure 11-3 depicts. Initially, the entire array is the unsorted region, just as the cards dealt to you sit in an unsorted pile on the table. At each step, the insertion sort takes the first item of the unsorted region and places it into its correct position in the sorted region. This step is analogous to taking a card from the table and inserting it into its proper position in your hand. The first step, however, is trivial: Moving theArray [0] from the unsorted region to the sorted region really does not require moving data. Therefore, you can omit this first step by considering the initial sorted region to be theArray[0] and the initial unsorted region to be theArray[1..n-1]. The fact that the items in the sorted region are sorted among themselves is an assertion ${ }^{2}$ of the algorithm. Because at each step the size of the sorted region grows by 1 and the size of the unsorted region shrinks by 1 , the entire array will be sorted when the algorithm terminates.

Figure 11-4 illustrates an insertion sort of an array of five integers. Initially, the sorted region is theArray [0], which is 29 , and the unsorted region is the rest of the array. You take the first item in the unsorted region-the 10-and insert it into its proper position in the sorted region. This insertion

2 As Appendix F discusses, this particular assertion is an invariant.

FIGURE 11-3 An insertion sort partitions the array into two regions


Bubble sort: Worst case: O( $n^{2}$ ) Best case: O(n)

Take each item from the unsorted region and insert it into its correct order in the sorted region

FIGURE 11-4 An insertion sort of an array of five integers

| Initial array: | 29 | 10 | 14 | 37 | 13 | Copy 10 |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: |
|  | 4 |  |  |  |  |  |
|  | 29 | 29 | 14 | 37 | 13 | Shift 29 |
|  | 10 | 29 | 14 | 37 | 13 | Insert 10; copy 14 |
|  | 1 |  |  |  |  |  |
|  | 10 | 29 | 29 | 37 | 13 | Shift 29 |
|  | 10 | 14 | 29 | 37 | 13 | Insert 14; copy 37, insert 37 on top of itself |
|  | 10 | 14 | 29 | 37 | 13 | Copy 13 |
|  |  |  |  | , | - |  |
|  | 10 | 14 | 14 | 29 | 37 | Shift 37, 29, 14 |
| Sorted array: | 10 | 13 | 14 | 29 | 37 | Insert 13 |

requires you to shift array entries to make room for the inserted item. You then take the first item in the new unsorted region-the 14-and insert it into its proper position in the sorted region, and so on.

Question 5 Trace the insertion sort as it sorts the array in Check point Question 3 into ascending order.

Question 6 Repeat the previous question, but instead sort the array into descending order.

Listing 11-3 contains a C++ function that performs an insertion sort on an array of $n$ items.

## LISTING 11-3 An implementation of the insertion sort

```
/** Sorts the items in an array into ascending order.
    @pre None.
    @post theArray is sorted into ascending order; n is unchanged.
    @param theArray The given array.
    @param n The size of theArray. */
void insertionSort(ItemType theArray[], int n)
{
    // unsorted = first index of the unsorted region,
    // loc = index of insertion in the sorted region,
    // nextItem = next item in the unsorted region.
    // Initially, sorted region is theArray[0],
    // unsorted region is theArray[1..n-1].
    // In general, sorted region is theArray[0..unsorted-1],
    // unsorted region theArray[unsorted..n-1]
```

```
    for (int unsorted = 1; unsorted < n; unsorted++)
    {
        // At this point, theArray[0..unsorted-1] is sorted.
        // Find the right position (loc) in theArray[0..unsorted]
        // for theArray[unsorted], which is the first entry in the
        // unsorted region; shift, if necessary, to make room
        ItemType nextItem = theArray[unsorted];
        int loc = unsorted;
        while ((loc > 0) && (theArray[loc - 1] > nextItem))
        {
            // Shift theArray[loc - 1] to the right
            theArray[loc] = theArray[loc - 1];
        } // end for
        // At this point, theArray[loc] is where nextItem belongs
        theArray[loc] = nextItem; // Insert nextItem into sorted region
        loc--;
    } // end while
} // end insertionSort
```

Analysis. The outer while loop in the function insertionSort executes $n-1$ times. This loop contains an inner for loop that executes at most unsorted times for values of unsorted that range from 1 to $n-1$. Thus, in the worst case, the algorithm's comparison occurs

$$
1+2+\cdots+(n-1)=n \times(n-1) / 2
$$

times. In addition, the inner loop moves data items at most the same number of times.
The outer loop moves data items twice per iteration, or $2 \times(n-1)$ times. Together, there are

$$
n \times(n-1)+2 \times(n-1)=n^{2}+n-2
$$

major operations in the worst case.
Therefore, the insertion sort algorithm is $\mathrm{O}\left(n^{2}\right)$ in the worst case. For small arrays-say, fewer than 25 items-the simplicity of the insertion sort makes it an appropriate choice. For large arrays, however, an insertion sort can be prohibitively inefficient-unless the array is already sorted. In that case, the inner for loop exits immediately, making the sort $\mathrm{O}(n)$ in its best case.

### 11.2 Faster Sorting Algorithms

The previous sorting algorithms are all you really need if you have to sort small arrays or if you need to sort a large array once. For extremely large arrays-particularly ones that must be updated and sorted again-you need faster algorithms. We now examine a few of these more sophisticated techniques.

### 11.2.1 The Merge Sort

Two important divide-and-conquer sorting algorithms, merge sort and quick sort, have elegant

The insertion sort is $\mathrm{O}\left(n^{2}\right)$ in the worst case

The insertion sort is $\mathrm{O}(n)$ in the best case

## $>$

## VideoNote

Faster sorting algorithms recursive formulations and are highly efficient. The presentations here are in the context of sorting

Halve the array, recursively sort its halves, and then merge the halves

The merge sort requires a second array as large as the original array

FIGURE 11-5 A merge sort with an auxiliary temporary array

arrays, but-as you will see in Chapter 21-the merge sort generalizes to external files. It will be convenient to express the algorithms in terms of the array theArray [first. .last].

The merge sort is a recursive sorting algorithm that always gives the same performance, regardless of the initial order of the array items. Suppose that you divide the array into halves, sort each half, and then merge the sorted halves into one sorted array, as Figure 11-5 illustrates. In the figure, the halves $\langle 1,4,8\rangle$ and $<2,3>$ are merged to form the array $\langle 1,2,3,4,8\rangle$. This merge step compares an item in one half of the array with an item in the other half and moves the smaller item to a temporary array. This process continues until there are no more items to consider in one half. At that time, you simply move the remaining items to the temporary array. Finally, you copy the temporary array back into the original array.

Although the merge step of the merge sort produces a sorted array, how do you sort the array halves prior to the merge step? The merge sort sorts the array halves by using a merge sort-that is, by calling itself recursively. Thus, the pseudocode for the merge sort is

```
// Sorts theArray[first..7ast] by
// 1. Sorting the first half of the array
// 2. Sorting the second half of the array
// 3. Merging the two sorted halves
mergeSort(theArray: ItemArray, first: integer, last: integer)
    if (first < last)
{
    mid = (first + last) / 2 // Getmidpoint
    // Sort theArray[first..mid]
    mergeSort(theArray, first, mid)
    // Sort theArray[mid+1..7ast]
    mergeSort(theArray, mid + 1, last)
```

```
    // Merge sorted halves theArray[first..mid] and theArray[mid+1..1ast]
    merge(theArray, first, mid, last)
}
// If first >= 1ast, there is nothing to do
```

Clearly, most of the effort in the merge sort algorithm is in the merge step, but does this algorithm actually sort? The recursive calls continue dividing the array into pieces until each piece contains only one item; obviously an array of one item is sorted. The algorithm then merges these small pieces into larger sorted pieces until one sorted array results. Figure 11-6 illustrates both the recursive calls and the merge steps in a merge sort of an array of six integers.

The C++ functions in Listing 11-4 implement the merge sort algorithm. To sort an array theArray of $n$ items, you would invoke the function mergeSort by writing

```
mergeSort(theArray, 0, n - 1)
```

FIGURE 11-6 A merge sort of an array of six integers


## LISTING 11-4 An implementation of the merge sort

const int MAX_SIZE = maximum-number-of-items-in-array;
/** Merges two sorted array segments theArray[first..mid] and
theArray[mid+1..7ast] into one sorted array.
@pre first <= mid <= 1ast. The subarrays theArray[first..mid] and
theArray [mid+1..1ast] are each sorted in increasing order.
@post theArray[first.. 1ast] is sorted.
@param theArray The given array.
@param first The index of the beginning of the first segment in theArray.
@param mid The index of the end of the first segment in theArray;
mid +1 marks the beginning of the second segment.
@param last The index of the last element in the second segment in theArray.

```
@note This function merges the two subarrays into a temporary
    array and copies the result into the original array theArray. */
void merge(ItemType theArray[], int first, int mid, int last)
{
    ItemType tempArray[MAX_SIZE]; // Temporary array
    // Initialize the local indices to indicate the subarrays
    int first1 = first; // Beginning of first subarray
    int last1 = mid; // End of first subarray
    int first2 = mid + 1; // Beginning of second subarray
    int last2 = 1ast; // End of second subarray
    // While both subarrays are not empty, copy the
    // smaller item into the temporary array
    int index = first1; // Next available location in tempArray
    while ((first1 <= 1ast1) && (first2 <= 1ast2))
    {
        // At this point, tempArray[first..index-1] is in order
        if (theArray[first1] <= theArray[first2])
        {
            tempArray[index] = theArray[first1];
            first1++;
        }
        else
            {
                tempArray[index] = theArray[first2];
                first2++;
            } // end if
            index++;
    } // end while
    // Finish off the first subarray, if necessary
    while (first1 <= 1ast1)
    {
        // At this point, tempArray[first..index-1] is in order
        tempArray[index] = theArray[first1];
        first1++;
        index++;
    } // end while
    // Finish off the second subarray, if necessary
    while (first2 <= last2)
    {
        // At this point, tempArray[first..index-1] is in order
            tempArray[index] = theArray[first2];
            first2++;
            index++;
    } // end for
    // Copy the result back into the original array
    for (index = first; index <= last; index++)
            theArray[index] = tempArray[index];
} // end merge
```

```
/** Sorts the items in an array into ascending order.
    @pre theArray[first..last] is an array.
    @post theArray[first..last] is sorted in ascending order.
    @param theArray The given array.
    @param first The index of the first element to consider in theArray.
    @param last The index of the last element to consider in theArray.*/
void mergeSort(ItemType theArray[], int first, int last)
{
    if (first < last)
    {
        // Sort each half
        int mid = first + (last - first) / 2; // Index of midpoint
        // Sort left half theArray[first..mid]
        mergeSort(theArray, first, mid);
        // Sort right half theArray[mid+1..1ast]
        mergeSort(theArray, mid + 1, 1ast);
        // Merge the two halves
        merge(theArray, first, mid, last);
    } // end if
} // end mergeSort
```

Question 7 By drawing a diagram like the one shown in Figure 11-6, trace the merge sort as it sorts the following array into ascending order: 253020804060 .
Question 8 Show that the merge sort algorithm satisfies the four criteria of recursion that Chapter 2 describes.

Analysis. Because the merge step of the algorithm requires the most effort, let's begin the analysis there. Each merge step merges theArray[first..mid] and theArray [mid+1..7ast]. Figure 11-7 provides an example of a merge step that requires the maximum number of comparisons. If the total number of items in the two array segments to be merged is $n$, then merging the segments requires at most $n-1$ comparisons. (For example, in Figure 11-7 six items are in the segments and five comparisons

FIGURE 11-7 A worst-case instance of the merge step in a merge sort

| first mid |  |  |  | 1 ast |  |  |  | Merge the halves: |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| theArray: | 1 | 2 | 8 |  | 4 | 5 | 6 |  |
|  |  |  |  |  |  |  |  | a. $1<4$, so move 1 from theArray[first. .mid] to tempArray b. $2<4$, so move 2 from theArray[first. .mid] to tempArray c. $8>4$, so move 4 from theArray [mid+1. . 1 ast] to tempArray d. $8>5$, so move 5 from theArray [mid+1. . 1ast] to tempArray e. $8>6$, so move 6 from theArray [mid+1. . 1 ast] to tempArray |
| tempArray: | 1 | 2 | 4 | 5 | 6 | 8 |  | f. theArray [mid+1. 7 last] is finished, so move 8 to tempArray |

FIGURE 11-8 Levels of recursive calls to mergeSort, given an array of eight items


Level 0: mergesort 8 items

Level 1: 2 calls to mergesort with 4 items each

Level 2: 4 calls to mergesort with 2 items each

Level 3: 8 calls to mergesort with 1 item each
are required.) In addition, there are $n$ moves from the original array to the temporary array, and $n$ moves from the temporary array back to the original array. Thus, each merge step requires $3 \times n-1$ major operations.

Each call to mergeSort recursively calls itself twice. As Figure 11-8 illustrates, if the original call to mergeSort is at level 0 , two calls to mergeSort occur at level 1 of the recursion. Each of these calls then calls mergeSort twice, so four calls to mergeSort occur at level 2 of the recursion, and so on. How many levels of recursion are there? We can count them, as follows.

Each call to mergeSort halves the array. Halving the array the first time produces two pieces. The next recursive calls to mergeSort halve each of these two pieces to produce four pieces of the original array; the next recursive calls halve each of these four pieces to produce eight pieces, and so on. The recursive calls continue until the array pieces each contain one item - that is, until there are $n$ pieces, where $n$ is the number of items in the original array. If $n$ is a power of $2\left(n=2^{k}\right)$, the recursion goes $k=\log _{2} n$ levels deep. For example, in Figure 11-8, there are three levels of recursive calls to mergeSort because the original array contains eight items and $8=2^{3}$. If $n$ is not a power of 2 , there are $1+\log _{2} n$ (rounded down) levels of recursive calls to mergeSort.

The original call to mergeSort (at level 0 ) calls merge once. Then merge merges all $n$ items and requires $3 \times n-1$ operations, as was shown earlier. At level 1 of the recursion, two calls to mergeSort, and hence to merge, occur. Each of these two calls to merge merges $n / 2$ items and requires $3 \times(n / 2)-1$ operations. Together these two calls to merge require $2 \times(3 \times(n / 2)-1)$ or $3 \times n-2$ operations. At level $m$ of the recursion, $2^{m}$ calls to merge occur; each of these calls merges $n / 2^{m}$ items and so requires $3 \times$ $\left(n / 2^{m}\right)-1$ operations. Together the $2^{m}$ calls to merge require $3 \times n-2^{m}$ operations. Thus, each level of the recursion requires $\mathrm{O}(n)$ operations. Because there are either $\log _{2} n$ or $1+\log _{2} n$ levels, the merge sort is $\mathrm{O}(n \times \log n)$ in both the worst and average cases. You should look at Figure 10-3 in the previous chapter to convince yourself that $\mathrm{O}(n \times \log n)$ is significantly faster than $\mathrm{O}\left(n^{2}\right)$.

The merge sort is $\mathrm{O}(n \times \log n)$

Although the merge sort is an extremely efficient algorithm with respect to time, it does have one drawback: The merge step requires an auxiliary array. This extra storage and the necessary copying of entries are disadvantages.

### 11.2.2 The Quick Sort

Consider the first two steps of the pseudocode function kSma 11 that solves the problem of finding the $k^{\text {th }}$ smallest item of the array theArray[first. . 1ast] and was discussed in Section 2.4.4 of Chapter 2:

FIGURE 11-9 A partition about a pivot


Choose a pivot value p from anArray [first.. 1 ast]
Partition the values of anArray[first..7ast] aboutp
Recall that this partition, which is pictured again in Figure 11-9, has the property that all items in $S_{1}=$ theArray[first..pivotIndex -1] are less than or equal to the pivot $p$, and all items in $S_{2}=$ theArray[pivotIndex +1..7ast] are greater than or equal to $p$. Though this property does not imply that the array is sorted, it does imply an extremely useful fact: The items within $S_{1}$ remain within $S_{1}$ when the array is properly sorted, although their positions relative to one another may change. Similarly, the items within $S_{2}$ will remain within $S_{2}$ when the array is sorted, although their relative positions may change. Finally, the pivot item remains in its position in the final, sorted array.

The partition induces relationships among the array items that are the ingredients of a recursive solution. Arranging the array items around the pivot $p$ generates two smaller sorting problems-sort the left section of the array $\left(S_{1}\right)$, and sort the right section of the array $\left(S_{2}\right)$. The relationships between the pivot and the array items imply that once you solve the left and right sorting problems, you will have solved the original sorting problem. That is, partitioning the array before making the recursive calls places the pivot in its correct position and ensures that when the smaller array segments are sorted their items will be in the proper relation to the rest of the array. Also, the quick sort algorithm will eventually terminate: The left and right sorting problems are indeed smaller problems and are each closer than the original sorting problem to the base case-which is an array containing one item - because the pivot is not part of either $S_{1}$ or $S_{2}$.

A first draft of pseudocode for the quick sort algorithm follows:

```
// Sorts theArray[first..7ast].
quickSort(theArray: ItemArray, first: integer, last: integer): void
    if (first < last)
    {
        Choose a pivot item p from theArray[first..7ast]
    Partition the items of theArray[first..last] aboutp
    // The partition is theArray[first..pivotIndex..last]
    quickSort(theArray, first, pivotIndex - 1) // Sort S S
    quickSort(theArray, pivotIndex + 1, last) // Sort S2
    }
    // If first >= last, there is nothing to do
```

Partitioning the array. Now consider the partition function that both kSma11 and quickSort must call. Partitioning an array section about a pivot item is actually the most difficult part of these two problems.

The quick sort partitions an array into items that are less than or equal to the pivot and those that are greater than or equal to the pivot

Partitioning places the pivot in its correct position within the array

Place your chosen pivot in theArray[1ast] before partitioning

The partition function will receive an array segment theArray[first. . ]ast] as an argument. The function must arrange the items of the array segment into two regions: $S_{1}$ contains the items less than or equal to the pivot and $S_{2}$ contains the items greater than or equal to the pivot. Thus, as you saw in Figure 11-9, $S_{1}$ is the segment theArray[first.. pivotIndex - 1], and $S_{2}$ is the segment theArray[pivotIndex + 1..7ast].

What pivot should you use? Various strategies exist for making this choice, as you will see later, so let's assume for now that we have chosen a pivot and go on to develop the partition. Regardless of which pivot you choose, let's swap it with the last entry theArray [last] to get it out of the way while we partition the array. Figure 11-10a shows an array after this step. Starting at the beginning of the array and moving toward the end (left to right in the figure), look for the first entry that is greater than or equal to the pivot. In Figure 11-10b, that entry is 5 and occurs at the index indexFromLeft. In a similar fashion, starting at the next-to-last entry and moving toward the beginning of the array (right to left in the figure), look for the first entry that is less than or equal to the pivot. In Figure 11-10b, that entry is 2 and occurs at the index indexFromRight. Now, if indexFromLeft is less than indexFromRight, swap the two entries at those indices. Figure 11-10c shows the result of this step. The 2, which is less than

FIGURE 11-10 A partitioning of an array during a quick sort

(c)

(d)


5 indexFromRight
(e)

(f)

(g)

(h)

the pivot, has moved toward the beginning of the array, while the 5 , which is greater than the pivot, has moved in the opposite direction.

Continue the searches from the left and from the right. Figure 11-10d shows that the search from the left stops at 4 and the search from the right stops at 1 . Since indexFromLeft is less than indexFromRight, swap 4 and 1. The array now appears as in Figure 11-10e. Entries equal to the pivot are allowed in either piece of the partition.

Continue the searches again. Figure 11-10f shows that the search from the left stops at 6, while the search from the right goes beyond the 6 to stop at 1 . Since indexFromLeft is not less than indexFromRight, no swap is necessary and the searches end. The only remaining step is to place the pivot between the subarrays $S_{1}$ and $S_{2}$ by swapping a[indexFromLeft] and a [last], as Figure 11-10g shows. The completed partition appears in Figure 11-10h.

Note that the previous searches must not go beyond the ends of the array. Soon you will see a convenient way to implement this requirement.

Entries equal to the pivot. Notice that both of the subarrays $S_{1}$ and $S_{2}$ can contain entries equal to the pivot. This might seem a bit strange to you. Why not always place any entries that equal the pivot into the same subarray? Such a strategy would tend to make one subarray larger than the other. However, to enhance the quick sort's performance, we want the subarrays to be as nearly equal in size as possible.

Notice that both the search from the left and the search from the right stop when they encounter an entry that equals the pivot. This means that rather than leaving such entries in place, they are swapped. It also means that such an entry has a chance of landing in each of the subarrays.

Selecting a pivot. Ideally, the pivot should be the median value in the array, so that the subarrays $S_{1}$ and $S_{2}$ each have the same - or nearly the same - number of entries. One way to find the median value is to sort the array and then get the value in the middle. But sorting the array is the original problem, so this circular logic is doomed. So instead of getting the best pivot by finding the median of all values in the array, we will at least try to avoid a bad pivot.

We will take as our pivot the median of three entries in the array: the first entry, the middle entry, and the last entry. One way to accomplish this task is to sort only those three entries and use the middle entry of the three as the pivot. Figure 11-11 shows an array both before and after its first, middle, and last entries are sorted. The pivot is the 5 . This pivot selection strategy is called median-of-three pivot selection.

This pivot selection scheme assumes that the array has at least three entries. If you have only three entries, the pivot selection sorts them, so there is no need for the partition method or for a quick sort. Thus, we now assume that the array contains at least four entries.

The following pseudocode describes how to sort the first, middle, and last entries in an array of at least four entries. For example, it will transform the array in Figure 11-11a to the one in Figure 11-11b.

FIGURE 11-11 Median-of-three pivot selection: (a) The original array; (b) the array with its first, middle, and last entries sorted
(a)

| $\mathbf{5}$ | 8 | 6 | 4 | $\mathbf{9}$ | 3 | 7 | 1 | $\mathbf{2}$ |
| :--- | :--- | :--- | :--- | :--- | :--- | :--- | :--- | :--- |

(b)

| $\mathbf{2}$ | 8 | 6 | 4 | $\mathbf{5}$ | 3 | 7 | 1 | $\mathbf{9}$ |
| :--- | :--- | :--- | :--- | :--- | :--- | :--- | :--- | :--- |

FIGURE 11-12 (a) The array with its first, middle, and last entries sorted; (b) the array after positioning the pivot and just before partitioning
(a)

(b)

// Arranges the first, middle, and last entries in an array into ascending order.
sortFirstMiddleLast(theArray: ItemArray, first: integer, mid: integer, last: integer): void
if (theArray[first] > theArray[mid])
Interchange theArray[first] and theArray[mid]
if (theArray[mid] > theArray[last])
Interchange theArray[mid] and theArray[1ast]
if (theArray[first] > theArray[mid])
Interchange theArray[first] and theArray[mid]
Adjusting the partition algorithm. Median-of-three pivot selection suggests some minor adjustments to our partitioning scheme. Previously, we swapped the pivot with the last entry in the array prior to partitioning. But here the first, middle, and last entries in the array are sorted, so we know that the last entry is at least as large as the pivot. Thus, the last entry belongs in the subarray $S_{2}$. We can simply leave the last entry in place. To get the pivot out of the way, we can swap it with the next-to-last entry, a[last - 1], as Figure 11-12 shows. Therefore, the partition algorithm can begin its search from the right at index last - 2 .

Also notice that the first entry is at least as small as the pivot, and so it belongs in the subarray $S_{1}$. Thus, we can leave the first entry in place and have the partition algorithm begin its search from the left at index first +1 . Figure 11-12b shows the status of the array at this point, just prior to partitioning.

This scheme provides a side benefit that simplifies the loops for the two searches. The search from the left looks for an entry that is greater than or equal to the pivot. That search will terminate because, at worst, it will stop at the pivot. The search from the right looks for an entry that is less than or equal to the pivot. That search will terminate because, at worst, it will stop at the first entry. Thus, the loops need not do anything special to prevent the searches from going beyond the ends of the array.

After the search loops end, we need to position the pivot between the subarrays $S_{1}$ and $S_{2}$. We do this by swapping the entries $a[$ indexFromLeft] and a[1ast - 1].


Note: The quick sort rearranges the entries in an array during the partitioning process. Each partition places one entry-the pivot-in its correct sorted position. The entries in each of the two subarrays that are before and after the pivot will remain in their respective subarrays.

```
The following pseudocode describes the partitioning algorithm for an array of at least four entries:

The partition algorithm
```

// Partitions theArray[first..7ast].

```
// Partitions theArray[first..7ast].
partition(theArray: ItemArray, first: integer, last: integer): integer
partition(theArray: ItemArray, first: integer, last: integer): integer
    // Choose pivot and reposition it
    // Choose pivot and reposition it
    mid = first + (last - first) / 2
    mid = first + (last - first) / 2
    sortFirstMiddleLast(theArray, first, mid, last)
    sortFirstMiddleLast(theArray, first, mid, last)
    Interchange theArray[mid] and theArray[1ast - 1]
    Interchange theArray[mid] and theArray[1ast - 1]
    pivotIndex = 1ast - 1
    pivotIndex = 1ast - 1
    pivot = theArray[pivotIndex]
    pivot = theArray[pivotIndex]
    // Determine the regions S}\mp@subsup{S}{1}{}\mathrm{ and }\mp@subsup{S}{2}{
    // Determine the regions S}\mp@subsup{S}{1}{}\mathrm{ and }\mp@subsup{S}{2}{
    indexFromLeft = first + 1
    indexFromLeft = first + 1
    indexFromRight = 1ast - 2
    indexFromRight = 1ast - 2
    done = false
    done = false
    while (not done)
    while (not done)
    {
    {
        // Locate first entry on left that is \geq pivot
        // Locate first entry on left that is \geq pivot
        while (theArray[indexFromLeft] < pivot)
        while (theArray[indexFromLeft] < pivot)
            indexFromLeft = indexFromLeft + 1
            indexFromLeft = indexFromLeft + 1
        // Locate first entry on right that is \leq pivot
        // Locate first entry on right that is \leq pivot
        while (theArray[indexFromRight] > pivot)
        while (theArray[indexFromRight] > pivot)
            indexFromRight = indexFromRight - 1
            indexFromRight = indexFromRight - 1
        if (indexFromLeft < indexFromRight)
        if (indexFromLeft < indexFromRight)
        {
        {
            Move theArray[firstUnknown] into S
            Move theArray[firstUnknown] into S
            Interchange theArray[indexFromLeft] and theArray[indexFromRight]
            Interchange theArray[indexFromLeft] and theArray[indexFromRight]
            indexFromLeft = indexFromLeft + 1
            indexFromLeft = indexFromLeft + 1
            indexFromRight = indexFromRight - 1
            indexFromRight = indexFromRight - 1
        }
        }
        e1se
        e1se
            done = true
            done = true
    }
    }
    // Place pivot in proper position between S}\mp@subsup{S}{1}{}\mathrm{ and }\mp@subsup{S}{2}{}\mathrm{ , and mark its new location
    // Place pivot in proper position between S}\mp@subsup{S}{1}{}\mathrm{ and }\mp@subsup{S}{2}{}\mathrm{ , and mark its new location
    Interchange theArray[pivotIndex] and theArray[indexFromLeft]
    Interchange theArray[pivotIndex] and theArray[indexFromLeft]
    pivotIndex = indexFromLeft
    pivotIndex = indexFromLeft
    return pivotIndex
```

```
    return pivotIndex
```

```

Question 9 Trace the quick sort's partitioning algorithm as it partitions the following array: 381640391227

A function for the quick sort. Before completing the C++ code for the quick sort, we need to think about small arrays. You have seen that the array should contain at least four entries before you call the partition method. But simply agreeing to use the quick sort only on large arrays is not enough. The pseudocode just given for the quick sort shows that partitioning even a large array will eventually lead to a recursive call that involves an array as small as two entries. The code for the quick sort needs to screen out these small arrays and use another way to sort them. An insertion sort is a good choice for small arrays. In fact, using it instead of the quick sort on arrays of as many as ten entries is reasonable. The function in Listing 11-5 implements the quick sort with these
        array. 381640391227
observations in mind. It assumes a constant MIN_SIZE that specifies the size of the smallest array on which we will use a quick sort.

\section*{LISTING 11-5 A function that performs a quick sort}
```

/** Sorts an array into ascending order. Uses the quick sort with
median-of-three pivot selection for arrays of at least MIN_SIZE
entries, and uses the insertion sort for other arrays.
@pre theArray[first..last] is an array.
@post theArray[first..last] is sorted.
@param theArray The given array.
@param first The first element to consider in theArray.
@param last The last element to consider in theArray. */
void quickSort(ItemType theArray[], int first, int last)
{
if (last - first + 1 < MIN_SIZE)
{
insertionSort(theArray, first, 1ast);
}
else
{
// Create the partition: S1 | Pivot | S2
int pivotIndex = partition(theArray, first, last);
// Sort subarrays S1 and S2
quickSort(theArray, first, pivotIndex - 1);
quickSort(theArray, pivotIndex + 1, last);
} // end if
} // end quickSort

```

Analysis. The major effort in the quickSort function occurs during the partitioning step. Partitioning will require no more than \(n\) comparisons, and so, like merging, it will be an \(\mathrm{O}(n)\) task. The ideal situation occurs when the pivot moves to the center of the array, so the two subarrays that the partition forms are the same size. If every recursive call to quickSort forms a partition with equal-sized subarrays, the quick sort will be like a merge sort in that the recursive calls halve the array. Thus, fewer recursive calls to quickSort occur. As in the previous analysis of the merge sort, you can conclude that there are either \(\log _{2} n\) or \(1+\log _{2} n\) levels of recursive calls to quickSort. Each call to quickSort involves \(m\) comparisons and at most \(m\) exchanges, where \(m\) is the number of items in the subarray to be sorted. Clearly \(m \leq n-1\). Thus, a quick sort would be \(\mathrm{O}(n \log n)\), and this would be its best case.

This ideal situation might not always occur, however. It is possible that each partition has one empty subarray. Although one recursive call will have nothing to do, the other call must sort \(n-1\) entries instead of \(n / 2\). This occurrence is the worst case because the nonempty subarray decreases in size by only 1 at each recursive call to quickSort, and so the maximum number of recursive calls to quickSort will occur. The result is \(n\) levels of recursive calls instead of \(\log n\). Thus, in the worst case, quick sort is \(\mathrm{O}\left(n^{2}\right)\).

A formal analysis of the average-case behavior of a quick sort would show that it is \(\mathrm{O}(n \log n)\). While the merge sort is always \(\mathrm{O}(n \log n)\), the quick sort can be faster in practice and does not require
the additional memory that merge sort needs for merging. Moreover, on large arrays, you can expect a quick sort to run significantly faster than an insertion sort. However, in its worst case, a quick sort will require roughly the same amount of time as an insertion sort.

The fact that the quick sort's average-case behavior is far better than its worst-case behavior distinguishes it from the other sorting algorithms considered in this chapter. If the original arrangement of data in the array is "random," the quick sort performs at least as well as any known sorting algorithm that involves comparisons.

The efficiency of a merge sort is somewhere between the possibilities for a quick sort: Sometimes the quick sort is faster, and sometimes the merge sort is faster. While the worst-case behavior of a merge sort is of the same order of magnitude as a quick sort's average-case behavior, in most situations a quick sort will run somewhat faster than a merge sort. However, in its worst case, the quick sort will be significantly slower than the merge sort.

Note: The choice of pivots affects the quick sort's efficiency. Some pivot-selection schemes can lead to worst-case behavior if the array is already sorted or nearly sorted. In practice, nearly sorted arrays can occur more frequently than you might imagine. Fortunately, median-of-three pivot selection avoids worst-case behavior for sorted arrays.

Note: The quick sort is often used to sort large arrays, as it is usually extremely fast in practice, despite its unimpressive theoretical worst-case behavior. Although a worst-case situation is not typical, even if the worst case occurs, the quick sort's performance is acceptable for moderately large arrays.

Note: The quick sort is appropriate when you are confident that the data in the array to be sorted is arranged randomly. Although the quick sort's worst-case behavior is \(\mathrm{O}\left(n^{2}\right)\), the worst case rarely occurs in practice.

\section*{ASIDE: Quick sort versus merge sort and ksmal1}

The quickSort and mergeSort algorithms are similar in spirit, but whereas quickSort does its work before its recursive calls, mergeSort does its work after its recursive calls. That is, while quickSort has the form
```

quickSort(theArray, first, 1ast)
if (first < last)
{
Prepare array for recursive calls quickSort ( $S_{1}$ region of theArray)

```
```

    quickSort(S region of theArray)
    }

```
mergeSort has the general form
```

mergeSort(theArray, first, last)
if (first < last)
{
mergeSort(Left half of theArray)
mergeSort(Right half of theArray)
Tidy up array after the recursive calls
}

```

The preparation in quickSort is to partition the array into regions \(S_{1}\) and \(S_{2}\). The algorithm then sorts \(S_{1}\) and \(S_{2}\) independently, because every item in \(S_{1}\) belongs to the left of every item in \(S_{2}\). In mergeSort, on the other hand, no work is done before the recursive calls: The algorithm sorts each half of the array with respect to itself. However, the algorithm must still deal with the interaction between the items in the two halves. That is, the algorithm must merge the two halves of the array after the recursive calls.

Now recall the pseudocode function kSma11, described in Chapter 2, that returns the \(k^{\text {th }}\) smallest value in an array:
```

kSma11(k, theArray, first, 1ast): ItemType
Choose a pivot value p from anArray[first. . 1ast]
Partition the values of anArray[first.. 1ast] about p
if (k < pivotIndex - first + 1)
return kSmall(k, anArray, first, pivotIndex - 1)
else if ( $k==$ pivotIndex - first +1 )
return $p$
else
return kSmal1 (k - (pivotIndex - first + 1) , anArray,
pivotIndex + 1, 1ast)

```

The recursive calls to kSmall act only on the section of the array that contains the desired item, and it is not called at all if the desired item is the pivot. On the other hand, quickSort is called recursively on both unsorted sections of the array. Figure 11-13 illustrates this difference.

Question 10 Suppose that you sort a large array of integers by using a merge sort. Next you use a binary search to determine whether a given integer occurs in the array. Finally, you display all of the integers in the sorted array.
a. Which algorithm is faster, in general: the merge sort or the binary search? Explain in terms of Big O notation.
b. Which algorithm is faster, in general: the binary search or displaying the integers? Explain in terms of Big O notation.

FIGURE 11-13 kSma11 versus quickSort


\subsection*{11.2.3 The Radix Sort}

The radix sort is included here because it is quite different from the other sorts we've described, as it does not compare the array's entries.

Imagine again that you are sorting a hand of cards. This time you pick up the cards one at a time and arrange them by rank into 13 possible groups in this order: \(2,3, \ldots, 10, \mathrm{~J}, \mathrm{Q}, \mathrm{K}, \mathrm{A}\). Combine these groups and place the cards face down on the table so that the 2 s are on top and the aces are on the bottom. Now pick up the cards one at a time and arrange them by suit into four possible groups in this order: clubs, diamonds, hearts, and spades. When taken together, the groups result in a sorted hand of cards.

A radix sort uses this idea of forming groups and then combining them to sort a collection of data. The sort treats each data item as a character string. As a first simple example of a radix sort, consider this collection of three-letter strings:
ABC, XYZ, BWZ, AAC, RLT, JBX, RDT, KLT, AEO, TLJ

The sort begins by organizing the data according to the rightmost (least significant) letters. Although none of the strings ends in \(A\) or \(B\), two strings end in \(C\). Place those two strings into a group. Continuing through the alphabet, you form the following groups:
\[
(\mathrm{ABC}, \mathrm{AAC})(\mathrm{TLJ})(\mathrm{AEO})(\mathrm{RLT}, \mathrm{RDT}, \mathrm{KLT})(\mathrm{JBX})(\mathrm{XYZ}, \mathrm{BWZ})
\]

The strings in each group end with the same letter, and the groups are ordered by that letter. In addition, the strings within each group retain their relative order from the original list of strings.

Now combine the groups into one as follows. Take the items in the first group in their present order, follow them with the items in the second group in their present order, and so on. The following group results:
ABC, AAC, TLJ, AEO, RLT, RDT, KLT, JBX, XYZ, BWZ

\section*{Group strings by their rightmost letter}

Combine the groups

Next, form new groups as you did before, but this time use the middle letter of each string instead of the last letter:

Group the strings by their middle letter

Combine the groups

Group the strings by their first letter

Sorted strings
(AAC) (ABC, JBX) (RDT) (AEO) (TLJ, RLT, KLT) (BWZ) (XYZ)
Now the strings in each group have the same middle letter, and the groups are ordered by that letter. As before, the strings within each group retain their relative order from the previous group of all strings.

Combine these groups into one group, again preserving the relative order of the items within each group:

\author{
AAC, ABC, JBX, RDT, AEO, TLJ, RLT, KLT, BWZ, XYZ
}

Now form new groups according to the first letter of each string:
(AAC, ABC, AEO) (BWZ) (JBX) (KLT) (RDT, RLT) (TLJ) (XYZ)

Finally, combine the groups, again maintaining the relative order within each group:
AAC, ABC, AEO, BWZ, JBX, KLT, RDT, RLT, TLJ, XYZ

The strings are now in sorted order.
In the previous example, all character strings had the same length. If the character strings have varying lengths, you can treat them as if they were the same length by padding them on the right with blanks as necessary.

To sort numeric data, the radix sort treats a number as a character string. You can treat numbers as if they were padded on the left with zeros, making them all appear to be the same length. You then form groups according to the right-most digits, combine the groups, form groups according to the next-to-last digits, combine them, and so on, just as you did in the previous example. Figure 11-14 shows a radix sort of eight integers.

FIGURE 11-14 A radix sort of eight integers
```

0123,2154,0222,0004, 0283, 1560, 1061, 2150
(1560,2150) (1061) (0222) (0123,0283) (2154,0004)
1560, 2150, 1061, 0222, 0123, 0283, 2154,0004
(0004) (0222,0123) (2150, 2154) (1560, 1061) (0283)
0004, 0222, 0123, 2150, 2154, 1560, 1061,0283
(0004, 1061) (0123, 2150, 2154) (0222,0283) (1560)
0004, 1061, 0123, 2150, 2154, 0222, 0283, 1560
(0004, 0123, 0222,0283) (1061, 1560) (2150, 2154)
0004, 0123, 0222, 0283, 1061, 1560, 2150, 2154
Original integers
Grouped by fourth digit Combined
Grouped by third digit Combined
Grouped by second digit Combined Grouped by first digit Combined (sorted)

```

The following pseudocode describes the algorithm for a radix sort of \(n\) decimal integers of \(d\) digits each:
```

// Sorts nd-digit integers in the array theArray.
radixSort(theArray: ItemArray, n: integer, d: integer): void
for (j = d down to 1)
{
Initialize 10 groups to empty
Initialize a counter for each group to 0
for (i = 0 through n - 1)

```
```

    {
        k = jth digit of theArray[i]
        Place theArray [i] at the end of group k
        Increase kth counter by l
    }
    Replace the items in theArray with all the items in group 0,
        followed by all the items in group 1, and so on.
    }

```

Question 11 Trace the radix sort as it sorts the following array into ascending order: 38121600401239341234272433335432

Analysis. From the pseudocode for the radix sort, you can see that this algorithm requires \(n\) moves each time it forms groups and \(n\) moves to combine them again into one group. The algorithm performs these \(2 \times n\) moves \(d\) times. Therefore, the radix sort requires \(2 \times n \times d\) moves to sort \(n\) strings of \(d\) characters each. However, notice that no comparisons are necessary. Thus, radix sort is \(\mathrm{O}(n)\).

Despite its efficiency, the radix sort has some difficulties that make it inappropriate as a generalpurpose sorting algorithm. For example, to perform a radix sort of strings of uppercase letters, you need to accommodate 27 groups - one group for blanks and one for each letter. If the original data collection contains \(n\) strings, each group must be able to hold \(n\) strings. For large \(n\), this requirement demands substantial memory if you use arrays for both the original data and the resulting groups. However, you can save memory by using a chain of linked nodes for each of the 27 groups. Thus, a radix sort is more appropriate for a chain than for an array.

\subsection*{11.3 A Comparison of Sorting Algorithms}

Figure 11-15 summarizes the time required in the worst case and average case for the sorting algorithms that appear in this chapter. For reference purposes, two other algorithms - the tree sort and heap sort-are included here, even though you will not study them until Chapters 15 and 17, respectively.

FIGURE 11-15 Approximate growth rates of time required for eight sorting algorithms
\begin{tabular}{|c|c|c|}
\hline & Worst case & Average case \\
\hline Selection sort & \(\mathrm{n}^{2}\) & \(\mathrm{n}^{2}\) \\
\hline Bubble sort & \(\mathrm{n}^{2}\) & \(\mathrm{n}^{2}\) \\
\hline Insertion sort & \(\mathrm{n}^{2}\) & \(\mathrm{n}^{2}\) \\
\hline Merge sort & \(n \times \log n\) & \(n \times \log n\) \\
\hline Quick sort & \(\mathrm{n}^{2}\) & \(n \times \log n\) \\
\hline Radix sort & n & n \\
\hline Tree sort & \(\mathrm{n}^{2}\) & \(n \times \log n\) \\
\hline Heap sort & \(n \times \log n\) & \(n \times \log n\) \\
\hline
\end{tabular}


Note: The Standard Template Library (STL) provides several sort functions in the library header <algorithm>. Consult C++ Interlude 7 for more information about the STL and these functions.

Even though the radix sort is \(\mathrm{O}(n)\), it is not appropriate as a general-purpose sorting algorithm

Selection sort Insertion sort Merge sort Quick sort Tree sort Heap sort

\section*{Summary}
1. The selection sort, bubble sort, and insertion sort are all \(\mathrm{O}\left(n^{2}\right)\) algorithms. Although in a particular case one might be faster than another, for large problems they all are slow. For small arrays, however, the insertion sort is a good choice.
2. The quick sort and merge sort are two very efficient recursive sorting algorithms. In the average case, the quick sort is among the fastest known sorting algorithms. However, the quick sort's worst-case behavior is significantly slower than the merge sort's. Fortunately, the quick sort's worst case rarely occurs in practice. The merge sort is not quite as fast as the quick sort in the average case, but its performance is consistently good in all cases. The merge sort has the disadvantage of requiring extra storage equal to the size of the array to be sorted.
3. The radix sort is unusual in that it does not sort the array entries by comparing them. Thus, it is not always applicable, making it inappropriate as a general-purpose sorting algorithm. However, when the radix sort is applicable, it is an \(\mathrm{O}(n)\) algorithm.

\section*{Exercises}
1. This chapter's analysis of the selection sort ignored operations that control loops or manipulate array indices. Revise this analysis by counting all operations, and show that the algorithm is still \(\mathrm{O}\left(n^{2}\right)\).
2. Trace the insertion sort as it sorts the following array into ascending order:

208040256040
3. Trace the selection sort as it sorts the following array into ascending order:

7122441932
4. Trace the bubble sort as it sorts the following array into descending order:

122351034
5. Apply the selection sort, bubble sort, and insertion sort to
a. An array sorted into descending order: 8642
b. An array sorted into ascending order: 2468
6. How many comparisons would be needed to sort an array containing 25 entries using the bubble sort in
a. The worst case?
b. The best case?
7. Find an array that makes the bubble sort exhibit its worst behavior.
8. Revise the function selectionSort so that it sorts an array of instances of a class according to one int data member, which is the sort key. Assume that the class contains a member method getSortKey that returns the integer sort key.
9. Write recursive versions of selectionSort, bubbleSort, and insertionSort.
10. Trace the merge sort algorithm as it sorts the following array into ascending order. List the calls to mergeSort and to merge in the order in which they occur.
\[
20804025 \quad 60 \quad 30
\]
11. When sorting an array by using a merge sort,
a. Do the recursive calls to mergeSort depend on the values in the array, the number of items in the array, or both? Explain.
b. In what step of mergeSort are the items in the array actually swapped (that is, sorted)? Explain.
12. Trace the quick sort algorithm as it sorts the following array into ascending order. List the calls to quicksort and to partition in the order in which they occur.

20804025601015
13. Suppose that you remove the call to merge from the merge sort algorithm to obtain
```

// Mystery algorithm for theArray [0..n-1].
mystery(theArray: ItemArray, n: integer): void
if ( }\textrm{n}>1
{
mystery(lefthalf(theArray))
mystery(righthalf(theArray))
}

```

What does this new algorithm do?
14. How many recursive calls are necessary when quickSort sorts an array of size \(n\) if you use median-of-three pivot selection?
15. Describe an iterative version of mergeSort.
16. One criterion used to evaluate sorting algorithms is stability. A sorting algorithm is stable if it does not exchange items that have the same sort key. Thus, items with the same sort key (possibly differing in other ways) will maintain their positions relative to one another. For example, you might want to take an array of students sorted by name and re-sort it by year of graduation. Using a stable sorting algorithm to sort the array by year will ensure that within each year the students will remain sorted by name. Some applications mandate a stable sorting algorithm. Others do not. Which of the sorting algorithms described in this chapter are stable?
17. When we discussed the radix sort, we sorted a hand of cards by first ordering the cards by rank and then by suit. To implement a radix sort for this example, you could use two characters to represent a card, if you used the letter \(T\) to represent a 10. For example, S2 is the 2 of spades and HT is the 10 of hearts.
a. Trace the radix sort for this example.
b. Suppose that you did not use \(T\) to represent a 10 - that is, suppose that H 10 is the 10 of hearts-and that you padded the two-character strings on the right with a blank to form three-character strings. How would a radix sort order the entire deck of cards in this case?

\section*{Programming Problems}
1. Add a counter to the functions insertionSort and mergeSort that counts the number of comparisons that are made. Run the two functions with arrays of various sizes. At what size does the difference in the number of comparisons become significant? How does this size compare with the size that the orders of these algorithms predict?
2. Revise the function quickSort so that it always chooses the first item in the array as the pivot. Add a counter to the function partition that counts the number of comparisons that are made. Compare the behavior of the revised function with the original one, using arrays of various sizes. At what size array does the difference in the number of comparisons become significant? For which pivot selection strategy does the difference in the number of comparisons become significant?
3. Various partitioning strategies are possible for quick sort. What other strategies can you think of? How do they compare to the two studied in the previous programming problem?
4. Consider the ADT list, as described in Chapters 8 and 9 . Add a method sort to the list.
5. Repeat Programming problem 4 , but assume a link-based implementation of the ADT list. Use a merge sort in the definition of the method sort. Implement any other sorting algorithms that are appropriate for this implementation of a list.
6. Repeat Programming Problem 4, but assume a link-based implementation of a list of strings. Use a radix sort in the definition of the method sort.
7. You can sort a large array of integers that are in the range 1 to 100 by using an array count of 100 items to count the number of occurrences of each integer in the array. Fill in the details of this sorting algorithm, which is called a bucket sort, and write a C++ function that implements it. What is the order of the bucket sort? Why is the bucket sort not useful as a general sorting algorithm?
8. The Shell sort (named for its inventor, Donald Shell) is an improved insertion sort. Rather than always exchanging adjacent items - as in an insertion sort - the Shell sort can exchange items that are far apart in the array. The Shell sort arranges the array so that every \(h^{\text {th }}\) item forms a sorted subarray in a decreasing sequence of values. For example, if \(h\) is 5 , every fifth item forms a sorted subarray. Ultimately, if \(h\) is 1 , the entire array will be sorted.

One possible sequence of \(h\) 's begins at \(n / 2\) and halves \(n\) until it becomes 1 . By using this sequence, and by replacing 1 with \(h\) and 0 with \(h-1\) in insertionSort, we get the following function for the Shell sort:
```

void she11Sort(ItemType theArray[], int n)
{
for (int h = n / 2; h > 0; h = h / 2)
{
for (int unsorted = h; unsorted < n; unsorted++)
{
DataItem nextItem = theArray[unsorted];
int loc = unsorted;
while ( (loc >= h) \&\& (theArray[loc - h] > nextItem) )
{
theArray[loc] = theArray[loc - h];
loc = 1oc - h;
} // end while
theArray[loc] = nextItem;
} // end for
} // end for
} // end shel7Sort

```

Add a counter to the functions insertionSort and she11Sort that counts the number of comparisons that are made. Run the two functions with arrays of various sizes. At what size does the difference in the number of comparisons become significant?
9. Write a program to display the running time of the sorts described in this chapter. Test the sorts on arrays of various sizes. Arrays of the same size should contain identical entries. Use the function clock from <ctime> to time each sort. See the beginning of the programming problems in Chapter 10 for an example of how to time code.

\section*{C++ Interlude}

\section*{Class Relationships and Reuse}

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Prerequisites
Chapter 1 Data Abstraction: The Walls
C++ Interlude 1 C++ Classes
C++ Interlude 2 Pointers, Polymorphism, and Memory Allocation
Chapter 6 ..... Stacks
Chapter 8 ..... Lists
\(C_{++}\)classes provide a way to enforce the walls of data abstraction by encapsulating an abstract data type's data and operations. An object-oriented approach, however, goes well beyond encapsulation. Inheritance, polymorphism, and containment allow you to define new classes from existing classes. This interlude describes techniques that make collections of reusable software components possible. Realize that much more can and should be said about these techniques. Consider this interlude to be a refinement and expansion of the introduction to this material begun in C++ Interlude 1.

\section*{C4.1 Inheritance Revisited}

When you think of inheritance, you might imagine a bequest of one million dollars from some long-lost wealthy relative. In the object-oriented world, however, inheritance describes the ability of a class to derive properties from a previously defined class. These properties are like the genetic characteristics you received from your parents: Some traits are the same, some are similar but different, and some are new.

FIGURE C4-1 Inheritance: Relationships among timepieces


A class can derive the behavior and structure of another

A digital alarm clock is a digital clock

Inheritance enables the reuse of existing classes

\section*{A derived class} inherits the members of its base

Inheritance, in fact, is a relationship among classes. One class can derive the behavior and structure of another class. For example, Figure C4-1 illustrates some relationships among various timepieces. Digital clocks, for example, include the clock in the dashboard of your car, the clock on the sign of the downtown bank, and the clock on your microwave oven. All digital clocks have the same underlying structure and perform operations such as

\section*{Set the time}

Advance the time
Display the time
A digital alarm clock is a digital clock that also has alarm methods, such as

\section*{Set the alarm}

Enable the alarm
Sound the alarm
Silence the alarm
That is, a digital alarm clock has the structure and operations of a digital clock and, in addition, has an alarm and operations to manipulate the alarm.

You can think of the group of digital clocks and the group of digital alarm clocks as classes. The class of digital alarm clocks is a derived class, or subclass, of the class of digital clocks. The class of digital clocks is a base class, or superclass, of the class of digital alarm clocks. Inheritance enables you to reuse software components when you define a new class. For example, you can reuse your design and implementation of an analog clock when you design a cuckoo clock.

In C++, a derived class inherits all of the members of its base class, except the constructors and destructor. That is, a derived class has the data members and methods of the base class in addition to the members it defines. A derived class can also revise any inherited method. For example, according to Figure C4-1, a cuckoo clock is a descendant of an analog clock, like the one on a classroom wall.


The cuckoo clock inherits the structure and behavior of the analog clock, but revises the way it reports the time each hour by adding a cuckoo.

Sometimes a derived class has more than one base class. For example, as Figure C4-2 illustrates, you can derive a class of clock radios from the class of digital clocks and the class of radios. This type of relationship is known as multiple inheritance. Multiple inheritance must be used with care, since it is possible for the base classes to have similarly named methods. For example, suppose that both digital clocks and radios implemented a method turnOn. When the client calls the turnOn method in the digital clock radio, you have no way to know which turnOn method is called. In this textbook, if multiple inheritance is used, the derived class inherits code from only one base class; any other bases classes should be abstract base classes.

In C++ Interlude 1, we used plain boxes, toy boxes, and magic boxes as objects. While designing the class of toy boxes, ToyBox, we decided that a toy box is simply a plain box with a color. This realization is significant in that PlainBox-the class of plain boxes-already existed. Thus, we could let PlainBox be a base class of ToyBox, and we implemented ToyBox without reinventing the basic box. Toward that end, Listing C4-1 recalls the definition of the class PlainBox from Listing C1-3.

\section*{LISTING C4-1 The class PlainBox, originally given in Listing C1-3}
```

template<class ItemType>;
class PlainBox
{
private:
ItemType item;
public:
PlainBox();
PlainBox(const ItemType\& theItem);
void setItem(const ItemType\& theItem);
ItemType getItem() const;
}; // end P7ainBox

```

Our derived class ToyBox inherited all the members of the class PlainBox-except the constructors and destructor-and made some changes. We added both a data member that held the box's color and methods to access and set the color.

You can add as many new members to a derived class as you like. Although you cannot revise an ancestor's private data members and should not reuse their names, you can redefine other inherited methods. A method in a derived class redefines a nonvirtual method in the base class if the two methods have the same name and parameter declarations.

When the derived class MagicBox was implemented in C++ Interlude 1, we redefined the method getItem so that it always returns the first item stored in the box. Figure C4-3 illustrates the relationship between PlainBox and MagicBox.

Inheritance reduces the effort necessary to add features to an existing object

A derived class can add new members to those it inherits

A derived class can redefine an inherited method of its base class

FIGURE C4-3 The derived class MagicBox inherits members of the base class PlainBox and redefines and adds members


Listing C4-2 shows the declaration of the class MagicBox, which was originally given in Listing C1-7 of C++ Interlude 1.

\section*{LISTING C4-2 The class MagicBox, originally given in Listing C1-7}
template<c1ass ItemType>;
class MagicBox : public PlainBox<ItemType>
\{
private:
bool firstItemStored;
pub7ic:
MagicBox() ;
MagicBox(const ItemType\& theItem);
void setItem(const ItemType\& theItem);
\}; // end MagicBox

An instance of a derived class has all the behaviors of its base class

A derived class inherits private members from the base class, but cannot access them directly

A derived class's methods can call the base class's public methods

Adding a colon and public PlainBox after class MagicBox indicates that PlainBox is a base class of MagicBox or, equivalently, that MagicBox is a derived class of PlainBox.

An instance of the class MagicBox has two data members-item, which is inherited, and firstItemStored, which is new. Because an instance of a derived class can invoke any public method in the base class, an instance of MagicBox has all of the methods that PlainBox defines; new constructors; a new, compiler-generated destructor; and a redefined method setItem. Although an instance of a derived class contains copies of inherited data members, the code for inherited methods is not copied.

A derived class cannot access the private members of the base class directly by name, even though they are inherited. Inheritance does not imply access. After all, you can inherit a locked vault but be unable to open it. In the current example, the data member item of Plain Box is private, so you can reference it only within the definition of the class PlainBox and not within the definition of ToyBox or MagicBox. However, the class ToyBox or MagicBox can use PlainBox's public methods setItem and getItem to set or obtain the value of item indirectly.

Within the implementation of MagicBox, you can use the public methods that MagicBox inherits from PlainBox. For example, the redefined method setItem calls the inherited version of the method setItem, which you indicate by writing PlainBox<ItemType> : : setItem. This notation is necessary to differentiate between the two versions of the method. Thus, you can access a base member, even though it has been redefined, by using the scope resolution operator :: . The implementation of MagicBox's setItem method is then
template<c1ass ItemType>
void MagicBox<ItemType>: : setItem(const ItemType\& theItem)
```

{
if (!firstItemStored)
{
PlainBox<ItemType>::setItem(theItem);
firstItemStored = true; // Box now has magic
} // end if
} // end setItem

```

Clients of a derived class also can invoke the public members of the base class. For example, if you write
```

MagicBox<string> myMagicBox("Volleyba11");
string magicBoxItem = myMagicBox.getItem();

```
myMagicBox's item, "Volleyba11", is returned by the method getItem that MagicBox inherits from PlainBox. If a new method has the same name as an ancestor method-setItem, for exampleinstances of the new class will use the new method, while instances of the ancestor class will use the original method. Therefore, if myPlainBox is a PlainBox object, the call myPlainBox.setItem(...) will invoke PlainBox's setItem, whereas myMagicBox.setItem(...) will invoke MagicBox's setItem, as Figure C4-4 illustrates. Because the compiler can determine which form of setItem to use at compilation time-as opposed to at execution time - this situation is called early binding, as introduced in \(\mathrm{C}++\) Interlude 2. Another term for early binding is static binding.

\section*{Note: Inheritance and the execution order of constructors and destructors}

A derived class's constructor executes after the base class's constructor. For example, if you define an instance of MagicBox, its constructor executes after PlainBox's constructor. The destructor of a derived class executes before the destructor of the base class. For example, MagicBox's destructor executes before PlainBox's destructor. This is true for constructors and destructors that you write as well as those generated by the compiler.

Early binding can lead to problems. For example, the statements
```

PlainBox<string>* cardBoxPtr = new PlainBox<string>("Queen of Hearts");
cardBoxPtr->setItem("Jack of Spades");

```
invoke P1ainBox's version of setItem. Unfortunately, if cardBoxPtr points to an instance of Magic-Box-for example, if you wrote
```

PlainBox<string>* cardBoxPtr = new MagicBox<string>("Queen of Hearts");

```

FIGURE C4-4 Early, or static, binding: The compiler determines which version of a method to invoke

Clients of a derived class can invoke the base class's public methods

Early binding can cause problems

the statement
```

cardBoxPtr->setItem("Jack of Spades");

```
still invokes PlainBox's version of setItem, instead of MagicBox's version. In this case, the compiler determined the version of the method to invoke from the type of the pointer cardBoxPtr instead of from the type of object to which cardBoxPtr points.

Simply redefining setItem in MagicBox-as we did earlier-is insufficient. As was briefly discussed in Section C1.5 of C++ Interlude 1 and Section C2.4 of C++ Interlude 2, you also need to tell the compiler that derived classes of PlainBox might revise setItem so that the compiler can make accommodations for such an occurrence. You do this by making setItem a virtual method in P1ainBox. Recall that to make a particular method virtual, you simply write the keyword virtual before the method's declaration within the definition of the base class. For example, the class PlainBox appears as
```

class PlainBox
{
public:
// Everything as before, except setItem
virtual ItemType setItem() const;
}; // end P7ainBox

```

The implementation of setItem is the same as given earlier.
Now, when cardBoxPtr points to an instance of MagicBox, the statement
```

cardBoxPtr->setItem("Jack of Spades");

```
invokes MagicBox's version of setItem. Thus, the appropriate version of a method is decided at execution time, instead of at compilation time, based on the type of object to which cardBoxPtr points. Recall from C++ Interlude 2 that this situation is called late binding. Another term for late binding is dynamic binding. A method such as setItem is called polymorphic. That is, the outcome of a particular operation depends upon the objects on which the operation acts. We also say that MagicBox's version of setItem overrides PlainBox's version.

A virtual method in a derived class can override a virtual method in the base class if they have the same declarations. Overriding a method is similar to redefining a method. However, you can override only virtual methods. Note that you can omit virtual in the derived class. Any method in a derived class that has the same declaration as an ancestor's virtual method-such as setItem-is also virtual without explicit identification, although tagging it with virtual is desirable stylistically.

Every class has a virtual method table (VMT), which remains invisible to the programmer. For each method in the class, the VMT contains a pointer to the actual instructions that implement the method. For a method that is not virtual, the compiler can establish this pointer, since it knows the method's definition. For a virtual method, the compiler can not complete the VMT. Instead, a call to a constructor during program execution sets the pointer, as Figure C4-5 illustrates. That is, the constructor establishes within the VMT pointers to the versions of the virtual methods that are appropriate for the object. Thus, the VMT is the mechanism that enables late binding.

\section*{C4.1.1 Public, Private, and Protected Sections of a Class}

In addition to its public and private sections, a class can have a protected section. By creating a protected section, a class designer can hide members from a class's clients but make them available to a derived class. That is, a derived class can reference the protected members of its base class directly, but clients of the base class or derived class cannot.

FIGURE C4-5 Virtual method tables when cardBoxPtr points to (a) an instance of PlainBox; (b) an instance of MagicBox
(a) PlainBox<string>* cardBoxPtr = new PlainBox<string>("Queen of Hearts");

AP1ainBox object

(b) PlainBox<string>* cardBoxPtr = new MagicBox<string>("Queen of Hearts");


For example, PlainBox has a private member item, which the derived class MagicBox cannot reference directly. If instead you declared item as protected, MagicBox would be able to access item directly by name. Clients of MagicBox or PlainBox, however, would not have direct access to item. We could then revise our implementation of MagicBox's setItem method as follows:
```

template<class ItemType>
void MagicBox<ItemType>::setItem(const ItemType\& theItem)
{
if (!firstItemStored)
{
item = theItem; // item has protected access
firstItemStored = true; // Box now has magic
} // end if
} // end setItem

```

The following summary distinguishes among the public, private, and protected sections of a class, and Figure C4-6 illustrates their access.

FIGURE C4-6 Access to public, private, and protected sections of a class by a client and a derived class


In general, a class's data members should be private

\section*{Note: Access categories of a class}
- Public members can be used by anyone.
- Private members can be used only by methods of the class.
- Protected members can be used only by methods of the class or any derived class.

\section*{Note: Stylistic guideline}

As a general stylistic guideline to maintain information hiding and to keep derived classes from coupling directly to base class data members, you should make all data members of a class private and, if required, provide indirect access to them by defining accessor or mutator methods that are either public or protected. Although a class's public members are available to anyone, its protected members are available exclusively to either its own methods or the methods of a derived class.

\section*{C4.1.2 Public, Private, and Protected Inheritance}

Several kinds of inheritance are possible. Regardless of the kind of inheritance, a derived class can access all of the base class's public and protected members, but not its private members. You can control how a class's inherited members are passed to subsequent derived classes by specifying one of three kinds of inheritance. You begin the definition of the derived class as
class DerivedClass : kindOfInheritance BaseClass
where kindOfInheritance is one of public, private, or protected. The inheritance that you saw earlier in this interlude is public inheritance. The following summary describes the three kinds of inheritance.

\section*{Note: Kinds of inheritance}
- Public inheritance: Public and protected members of the base class remain, respectively, public and protected members of the derived class.
- Protected inheritance: Public and protected members of the base class are protected members of the derived class.
- Private inheritance: Public and protected members of the base class are private members of the derived class.

In all cases, private members of a base class remain private to the base class and cannot be accessed by a derived class.

Of the three types of inheritance, public inheritance is the most important and the one that we will use most often in subsequent chapters. You use public inheritance to extend the definition of a class. You use private inheritance to implement one class in terms of another class. Protected inheritance is not often used, so we will not cover it.

The following section describes when it is appropriate to use public and private inheritance.

\section*{C4.1.3 Is-a and As-a Relationships}

As you just saw, inheritance provides for ancestor/descendant relationships among classes. Other relationships are also possible. When designing new classes from existing ones, it is important to

identify their relationship so that you can determine whether to use inheritance and, if so, the kind of inheritance that best reflects the relationship. Three basic kinds of relationships are possible. In this section we discuss is- \(a\) and \(a s-a\) relationships, which involve inheritance. We discuss an important third relationship, has-a, in Section C4.2.

Is-a relationships. Earlier in this interlude, we used public inheritance to derive the class MagicBox from PlainBox. You should use public inheritance only when an is-a relationship exists between two classes of objects. In this example, a magic box is a plain box, as Figure C4-7 illustrates. That is, whatever is true of the base class PlainBox is also true of the derived class MagicBox. Wherever you can use an object of type PlainBox, you can also use an object of type MagicBox. This feature is called object type compatibility. In general, a derived class is type-compatible with all of its ancestor classes. Thus, you can use an instance of a derived class instead of an instance of its base class, but not the other way around.

In particular, the object type of an argument in a call to a method can be a descendant of the object type of the corresponding parameter. For example, suppose your program uses the classes PlainBox and MagicBox and contains the following ordinary function, which is not a member of any class:
```

void displayBoxItem(PlainBox<string> thing)
{
cout << "The item stored in the box is "
<< thing.getItem() << ".\n";
} // end disp7ayBoxItem

```

\section*{If you define myPlainBox and myMagicBox as}

P1ainBox<string> myP1ainBox("Basketba11");
MagicBox<string> myMagicBox("Volleyba11");
the following calls to displayBoxItem are legal:
```

displayBoxItem(myPlainBox); // myPTainBox's item is displayed
displayBoxItem(myMagicBox); // myMagicBox's item is displayed

```

The first call is unremarkable because both the argument myPlainBox and the parameter thing have the same data type. The second call is more interesting: The data type of the argument myMagicBox is a descendant of the data type of the parameter thing. Because a magic box is a plain box, it can behave like a plain box. That is, myMagicBox can perform plain box behaviors, so you can use myMagicBox anywhere you can use myPlainBox. Note that object type compatibility applies to both value and reference arguments.
\(\boldsymbol{A s} \boldsymbol{a} \boldsymbol{a}\) relationships. Consider the relationship between the class Stack, which implements the ADT stack, and the class List, which implements the ADT list. Because a Stack object should not be able to have List behaviors, public inheritance is inappropriate in this case. A stack is not a list. When public inheritance is inappropriate, if your class needs access to the protected members of another

Public inheritance implements an is-a relationship

You can use an instance of a derived class anywhere you can use an instance of the base class

\section*{Because a magic} box is a plain box, you can use it anywhere you can use a plain box

If the relationship between two classes is not \(i s-a\), you should not use public inheritance

Has-a, or containment, means a class has an object as a data member
classes; that is, you can use private inheritance.
To derive Stack from List using private inheritance, you begin the definition of the class Stack with the header
```

class Stack : private List

```

\section*{class Stack : private List}

List's public members-and protected members, if List has any-would be private members of Stack. Thus, within the implementation of Stack, you would be able to use a list to store the stack's
entries and manipulate them by using List's methods. We say that Stack implements the ADT stack Stack. Thus, within the implementation of Stack, you would be able to use a list to store the stack's
entries and manipulate them by using List's methods. We say that Stack implements the ADT stack as a list. Both the descendants and clients of Stack are unable to access any members of List. Thus, the underlying list is hidden from the clients of the stack.

If public inheritance is inappropriate and your class does not need access to the protected mem-
In of another class, or if you do not need to redefine methods in that class, a has-a relationship is
If public inheritance is inappropriate and your class does not need access to the protected mem-
bers of another class, or if you do not need to redefine methods in that class, a has-a relationship is preferable and simpler to use. We consider that next.

\section*{C4.2 Containment: Has-a Relationships}

A ball-point pen has \(a\) ball as its point, as Figure C4-8 illustrates. Although you would want to use Ba11 in your definition of a class Pen, you should not use public inheritance, because a pen is not a Ball in your definition of a class Pen, you should not use public inheritance, because a pen is not a
ball. In fact, you do not use inheritance at all to implement a has-a relationship. Instead, you can define a data member point-whose type is Ball-within the class Pen as
class or if you need to redefine methods in that class, you can form an as-a relationship between your
```

c7ass Pen

```
c7ass Pen
{
{
private:
private:
    Ba11 point;
    Ba11 point;
}; // end Pen
```

}; // end Pen

```

Recall that a data member can be an instance of any class other than the one currently being definedPen, in this example. An instance of Pen has, or contains, an instance of Ba11. Thus, another name for the has- \(a\) relationship is containment.

As you continue to examine the ball's relationship to the pen, you may realize that if the pen is destroyed, the ball should also be destroyed, since it no longer has a purpose. When an instance of an object cannot exist independently of the containing class, this type of containment is called

\section*{FIGURE C4-8 A pen has \(a\) or contains a ball}

composition. We have seen this in our link-based implementations of an ADT. For example, when an instance of LinkedBag, LinkedStack, or LinkedList is destroyed, the nodes used to store the container's objects are no longer needed and are destroyed also. In the array-based implementations of our ADTs, the array containing the items is destroyed.

Another form of containment is aggregation. In an aggregate relationship the contained item can exist independently of the containing class. Our PlainBox class has an instance of ItemType named item. If our plain box is destroyed, the program can continue to use item. Another example of aggregation can be found in the implementations of our ADT bag, ADT stack, and ADT list: The items stored in a bag, stack, or list continue to exist even after we are finished using the container and it has been destroyed.

Programming Tip: Favor containment over inheritance. Use public inheritance when an is- \(a\) relationship exists, that is, when an instance of a derived class can substitute for an instance of the base class. Use private inheritance when you want to reuse a part of another class's implementation.

\section*{Note: Containment and the execution order of constructors and destructors}

The constructors for a class's member objects execute in the order listed in the class definition before the class's constructor. For example, Ball's constructor executes before Pen's constructor. The destructors of a class's member objects execute after the class's destructor. For example, Ba17's destructor executes after Pen's.

\section*{Note: Execution order of constructors and destructors}

Considering both inheritance and containment, a class's constructors execute in the following order:
1. Its base class constructor executes.
2. Constructors of its member objects execute in the declaration order.
3. The body of its constructor executes.

A class's destructors execute in the opposite order:
1. The body of its destructor executes.
2. Destructors of its member objects execute in the declaration order.
3. Its base class destructor executes.

\section*{C4.3 Abstract Base Classes Revisited}

Imagine a CD player (CDP) and a DVD player (DVDP). Both devices share several characteristics. Each involves a compact disc. You can insert, remove, play, and stop such discs. You can also skip forward or skip backward from the current position on the disc. Some of these operations are essentially the same for both devices, while others - in particular, the play operation-are different but similar.

If you were specifying both devices, you might begin by describing the common operations.
// Inserts a disc into the player.
Disc transport operations

A pure virtual method has an undefined body
```

// Removes a disc from the player.
+remove()
// Plays the disc.
+p7ay()
// Stops playing the disc.
+stop()
// Skips ahead to another section of the disc.
+skipForward()
// Skips back to an earlier section of the disc.
+skipBackward()

```

These operations could constitute a generic disc player (GDP).
If GDP, CDP, and DVDP were classes, GDP could be the base class of CDP and DVDP, as Figure C4-9 illustrates. While GDP could implement operations such as insert and remove that would be suitable for both a CDP and a DVDP, it could only indicate that these devices have a p 7 ay operation. So CDP, for example, inherits the operations provided by GDP but overrides the play operation to suit CDs, as Figure C4-10 illustrates. If necessary, CDP could override any of GDP's operations or define additional ones. We can make similar comments about DVDP. Thus,
- A CDP is a GDP that plays sound.
- A DVDP is a GDP that plays sound and video.

Because GDP cannot implement its play operation, we would not want instances of it. So GDP is simply a class without instances that forms the basis of other classes. If a class never has instances, its methods need not be implemented. Such methods, however, must be virtual so that derived classes can supply their own implementations. A virtual method with an undefined body is called a pure virtual method and is written as
virtual prototype \(=0\);
within a class definition. The implementation of a pure virtual method is deferred to a derived class.

FIGURE C4-9 CDP and DVDP have an abstract base class



Chapter 1 defined an abstract base class as a class that contains at least one virtual method that has no implementation. We can now restate that definition as a class that contains at least one pure virtual method. An abstract base class has no instances and is used only as the basis of other classes. Thus, the GDP class is an abstract base class, because it does not implement all of its methods. Any derived class that fails to implement all of the pure virtual methods is also an abstract base class and cannot have instances.

An abstract base class can provide a constructor, which cannot be pure because constructors cannot be virtual. A destructor in an abstract base class should not be pure, because derived classes will call it, and thus, it should be implemented even if it simply has an empty body. Generally, virtual destructors are recommended.

As you have seen, an abstract base class is a great place to specify an ADT. By defining the methods that make up the ADT's public interface as pure virtual methods, we force any derived class to implement these methods, ensuring a common interface across implementations. For example, Listing C4-3 contains the abstract base class for boxes from C++ Interlude 1 with the addition of a virtual destructor.

\section*{LISTING C4-3 An abstract class that is an interface for the ADT box}

A class that contains at least one pure virtual method is an abstract base class

An abstract base class has descendants but no instances

We can indicate that our class PlainBox is derived from BoxInterface by changing its class header to
class PlainBox : public BoxInterface<ItemType>
We can also define a new class, Jewe 1 ryBox, that provides an alternate implementation of the methods in BoxInterface. Its class header would be
class Jewe1ryBox : pub7ic BoxInterface<ItemType>
Figure C4-11 shows the family of box classes.
```

template <class ItemType>

```
template <class ItemType>
class BoxInterface
class BoxInterface
{
{
public:
public:
    virtual ~BoxInterface() { } // Empty implementation
    virtual ~BoxInterface() { } // Empty implementation
    virtual void setItem(const ItemType& theItem) = 0;
    virtual void setItem(const ItemType& theItem) = 0;
    virtual ItemType getItem() const = 0;
    virtual ItemType getItem() const = 0;
}; // end BoxInterface
```

}; // end BoxInterface

```

FIGURE C4-11 UML class diagram of the family of box classes


If we define several pointer variables of type BoxInterface, we can have them point to instances of any of the derived classes:
```

BoxInterface<string>* someBoxPtr = new PlainBox<string>("Ace");
BoxInterface<string>* anotherBoxPtr = new Jewe1ryBox<string>("Emerald");
BoxInterface<string>* funBoxPtr = new MagicBox<string>("Stone");

```

For each of the objects created, we can call the setItem or getItem method, and the implementation specific to the instantiated class will execute, even though the pointers are of type BoxInterface:
```

cout << someBoxPtr->getItem() << endl; // Uses PTainBox getItem
cout << anotherBoxPtr->getItem() << endl; // Uses JewelryBox getItem
cout << funBoxPtr->getItem() << end1; // Uses MagicBox getItem, which
// is PlainBox getItem

```

The key points about abstract base classes are summarized next.

\section*{Note: Abstract base classes}

An abstract base class:
- By definition is a class that contains at least one pure virtual method.
- Is used only as the basis for derived classes and thus defines a minimum interface for its descendants.
- Has no instances.
- Should, in general, omit implementations except for the destructor and methods that provide access to private data members. That is, virtual methods in an abstract base class usually should be pure.
- Must implement any virtual method that is not pure, thereby providing a default implementation if a derived class chooses not to supply its own.

\section*{Sorted Lists and Their Implementations}

\section*{Chapter} 12

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0ne of the most frequently performed computing tasks is the maintenance of a collection of data in some specified order. Many examples immediately come to mind, such as students placed in order by their names, baseball players listed in order by their batting averages, and corporations listed in order by their assets. The items in these examples are in sorted order. In contrast, the items on a grocery list might occur in the order in which they appear on the grocer's shelves, but they are probably not sorted by name. These items are ordered simply by their positions within the list.

The ADT sorted list maintains items in sorted order

The ADT sorted list maintains its entries in sorted order. The problem of maintaining sorted data involves more than simply sorting the data. Often you need to insert some new data item into its proper, sorted place. Similarly, you often need to remove some data item. For example, suppose your university maintains an alphabetical list of the students who are currently enrolled. The registrar must add names to and remove names from this list because students constantly enroll in and leave school. These operations should preserve the sorted order of the data.

This chapter specifies the ADT sorted list and then implements it in several ways. A major goal of this chapter is to show you how we can develop a class of sorted lists from an existing class of lists.

\subsection*{12.1 Specifying the ADT Sorted List}

The ADT list leaves it up to the client to arrange the entries in a given collection. The client can maintain the entries in any order that meets its needs. Recall that to use the insert operation of the ADT list, you must specify both the new entry and its desired position within the list. The ADT sorted list differs from the ADT list in that a sorted list inserts and removes items by their values and not by their positions. For example, the sorted list insertion operation determines the proper position for the new entry according to its value. Since a sorted list is responsible for organizing its entries, it has no operation that inserts a new entry by position. If you were allowed to specify a new entry's position, you might destroy the order of the sorted list's entries.


Note: The ADT sorted list is a container of items that determines and maintains the order of its entries by their values. For simplicity, we will allow the sorted list to contain duplicate items. Insisting that the sorted list contain only unique items is somewhat more complicated, and we will leave this variation as an exercise.

What other operations do we want for this ADT? We've already mentioned that you can add an entry to the sorted list. Since the sorted list determines the position of a new entry, you could ask the ADT for this position. That is, you could ask for the position of an existing entry or for the position in which a proposed entry would occur if you added it to the list. You could also ask the ADT for the entry at a particular position. And clearly you should be able to remove an entry. We could remove a particular entry given its value, but we could also remove an entry given its position in the sorted list. We will include both of these removal operations.

Let's itemize the sorted list's operations-including common ones like getting the current number of entries - and specify them more carefully. Note that none of the operations on a sorted list that involve the position of an entry destroys the list's sorted order.


Note: ADT sorted list operations
- Test whether a sorted list is empty.
- Get the number of entries in a sorted list.
- Insert an entry into a sorted list.
- Remove a given entry from a sorted list.
- Remove the entry at a given position from a sorted list.
- Remove all entries from the sorted list.
- Look at (get) the entry at a given position in the list.
- Get the position in a sorted list of a given entry.

\section*{FIGURE 12-1 UML diagram for the ADT sorted list}
\begin{tabular}{l} 
SortedList \\
\hline \\
\hline +isEmpty(): boolean \\
+getLength(): integer \\
+insertSorted(newEntry: ItemType) : void \\
+removeSorted(newEntry: ItemType): boolean \\
+remove(position: integer): boolean \\
+clear(): void \\
+getEntry(position: integer): ItemType \\
+getPosition(newEntry: ItemType): integer \\
\hline
\end{tabular}

Figure 12-1 shows the UML diagram for this ADT, and the following summary of its operations provides even more details for its operation contract:

\section*{Abstract Data Type: Sorted List}

\section*{Data}
- A finite number of objects, not necessarily distinct, having the same data type and ordered by their values.

\section*{Operations}

\section*{Pseudocode}
insertSorted(newEntry)
removeSorted(anEntry)
getPosition(anEntry)

\section*{Description}

Task: Inserts an entry into this sorted list in its proper order so that the list remains sorted.
Input: newEntry is the new entry.
Output: None.
Task: Removes the first or only occurrence of anEntry from this sorted list.
Input: anEntry is the entry to remove.
Output: Returns true if anEntry was located and removed, or false if not. In the latter case, the list remains unchanged.

Task: Gets the position of the first or only occurrence of anEntry in this sorted list.
Input: anEntry is the entry to be located.
Output: Returns the position of anEntry if it occurs in the sorted list. Otherwise, returns the position where anEntry would occur in the list, but as a negative integer.
The following operations behave as they do for the ADT list and are described in Chapter 8:
```

isEmpty()
getLength()
remove(position)
clear()
getEntry(position)

```

The first two methods are straightforward, but let's look more closely at getPosition. Given an entry in the sorted list, the method getPosition returns the entry's position number within the list, as you would expect. We number the entries beginning with 1 , just as we do for the ADT list. But what if the given entry is not in the sorted list? In this case, getPosition returns the position number where the entry belongs in the list. The returned number is negative, however, to signal that the entry is not in the list. For example, if missingObject is not in the sorted list but would belong at position 3, getPosition(missingObject) will return -3 .

The sorted list also has some, but not all, of the operations of an ADT list. We have already mentioned that adding an entry at a given position is not possible, because otherwise the client could destroy the order of the sorted list. For the same reason, the list's setEntry method is not available to a sorted list. The other operations of the ADT list, however, are useful for a sorted list as well, including the ones that retrieve or remove the entry at a given position. The methods getEntry and remove each have a position number as a parameter, but they will not alter the relative order of the entries in the sorted list.

Question 1 The specifications of the ADT sorted list do not mention the case in which two or more items have the same value. Are these specifications sufficient to cover this case, or must they be revised?

Question 2 Write specifications for the operation insertSorted when the sorted list must not contain duplicate entries.

\subsection*{12.1.1 An Interface Template for the ADT Sorted List}

The \(\mathrm{C}++\) interface given in Listing 12-1 formalizes our initial specifications of the ADT sorted list.

\section*{LISTING 12-1 A C++ interface for sorted lists}
```

/** Interface for the ADT sorted list
@file SortedListInterface.h */
\#ifndef _SORTED_LIST_INTERFACE
\#define SORTED_LIST_INTERFACE
template<class ItemType>
class SortedListInterface
{
public:
/** Inserts an entry into this sorted list in its proper order
so that the list remains sorted.
@pre None.
@post newEntry is in the list, and the list is sorted.
@param newEntry The entry to insert into the sorted list. */
virtual void insertSorted(const ItemType\& newEntry) = 0;
/** Removes the first or on7y occurrence of the given entry from this
sorted list.
@pre None.

```
```

        @post If the removal is successful, the first occurrence of the
            given entry is no longer in the sorted list, and the returned
            value is true. Otherwise, the sorted list is unchanged and the
            returned value is false.
        @param anEntry The entry to remove.
        @return True if removal is successful, or false if not. */
        virtua1 bool removeSorted(const ItemType& anEntry) = 0;
    /** Gets the position of the first or only occurrence of the given
        entry in this sorted list. In case the entry is not in the list,
        determines where it should be if it were added to the list.
    @pre None.
    @post The position where the given entry is or belongs is returned.
        The sorted list is unchanged.
    @param anEntry The entry to locate.
    @return Either the position of the given entry, if it occurs in the
        sorted list, or the position where the entry would occur, but as a
        negative integer. */
    virtual int getPosition(const ItemType& anEntry) = 0;
    // The following methods are the same as those given in ListInterface
// in Listing 8-1 of Chapter 8 and are completely specified there.
/** Sees whether this list is empty. */
virtual bool isEmpty() const = 0;
/** Gets the current number of entries in this list. */
virtual int getLength() const = 0;
/** Removes the entry at a given position from this list. */
virtual bool remove(int position) = 0;
/** Removes all entries from this 1ist. */
virtual void clear() = 0;
/** Gets the entry at the given position in this list. */
virtual ItemType getEntry(int position) const = 0;
}; // end SortedListInterface
\#endif

```

\subsection*{12.1.2 Using the Sorted List Operations}

Example. To demonstrate the operations of the ADT sorted list that the previous section specifies, we first create a sorted list of strings. We begin by declaring and allocating a sorted list, where we assume that SortedList is an implementation of the operations specified by the interface SortedListInterface:

SortedListInterface<string>* nameListPtr = new SortedList<string>();
Next, we add names in an arbitrary order, realizing that the ADT will organize them alphabetically:
```

nameListPtr->insertSorted("Jamie");

```
nameListPtr->insertSorted("Brenda");
```

nameListPtr->insertSorted("Sarah");
nameListPtr->insertSorted("Tom");
nameListPtr->insertSorted("Carlos");

```

The sorted list now contains the following entries:
Brenda
Carlos
Jamie
Sarah
Tom
Assuming the list just given, here are some examples of the operations on the sorted list:
nameListPtr->getPosition("Jamie") returns 3, the position of Jamie in the list nameListPtr->getPosition("Jil1") returns -4, because Jill belongs at position 4 in the list nameListPtr->getEntry (2) returns Carlos, because he is at position 2 in the list

Now remove Tom and the first name in the list by writing
```

nameList.remove("Tom");
nameList.remove(1);

```

The list now contains
Carlos
Jamie
Sarah
Removing the last entry, Tom, did not change the positions of the other entries in the list, but removing the first entry did. Carlos is now at position 1 , instead of 2.

Note: The ADT sorted list can add, remove, or locate an entry, given the entry as an argument. The sorted list has several operations that are the same as ADT list operations, namely getEntry (by position), removeSorted (by position), clear, getLength, and isEmpty. However, a sorted list will not let you add or replace an entry by position.

Question 3 Suppose that wordListPtr points to an unsorted list of words. Using the operations of the ADT list and the ADT sorted list, create a sorted list of these words.
Question 4 Assuming that the sorted list you created in the previous question is not empty, write C++ statements that
a. Display the last entry in the sorted list.
b. Add the sorted list's first entry to the sorted list again.

\subsection*{12.2 A Link-Based Implementation}

As with all ADTs, you have a choice of several ways in which to implement the sorted list. You could store a sorted list's entries in, for example, an array, a chain of linked nodes, an instance of a vector, or an instance of an ADT list. We first will consider a chain of linked nodes and then an instance of an ADT list. Finally, we will use inheritance to develop a completely different implementation.

\subsection*{12.2.1 The Header File}

An implementation that uses a chain of linked nodes to store the entries in a sorted list has several details in common with the link-based implementation of the ADT list that you studied in Chapter 9. In particular, it has the same data fields, similar constructors, several methods that are the same, and the same definition of the class Node. Let's begin with the header file, which is given in Listing 12-2. The class LinkedSortedList declares three new public methods for the sorted list-insertSorted, removeSorted, and getPosition-each of which takes a list entry as its argument. As you will see, to implement these methods, you will need the private method getNodeBefore. It accepts an entry as its argument and returns a pointer to the node just prior to the one that contains-or should contain-the entry. Since the definitions of the remaining methods are the same as those in LinkedList, as given in Chapter 9, and two of them involve the position of an entry, you will also need the private method getNodeAt as defined in Chapter 9. As you will see, we will use the private method copyChain to provide a recursive implementation for the copy constructor.

\section*{LISTING 12-2 The header file for the class LinkedSortedList.}
```

/** ADT sorted list: Link-based implementation.
@file LinkedSortedList.h */
\#ifndef _LINKED_SORTED_LIST
\#define _LINKED_SORTED_LIST
\#include "SortedListInterface.h"
\#include "Node.h"
\#include "PrecondViolatedExcep.h"
temp1ate<c1ass ItemType>
class LinkedSortedList : public SortedListInterface<ItemType>
{
private:
Node<ItemType>* headPtr; // Pointer to first node in the chain
int itemCount; // Current count of list items
// Locates the node that is before the node that should or does
// contain the given entry.
// @param anEntry The entry to find.
// @return Either a pointer to the node before the node that contains
// or should contain the given entry, or nullptr if no prior node exists.
Node<ItemType>* getNodeBefore(const ItemType\& anEntry) const;
// Locates the node at a given position within the chain.
Node<ItemType>* getNodeAt(int position) const;
// Returns a pointer to a copy of the chain to which origChainPtr points.
Node<ItemType>* copyChain(const Node<ItemType>* origChainPtr);
public:
LinkedSortedList();
LinkedSortedList(const LinkedSortedList<ItemType>\& aList);
virtual ~LinkedSortedList();

```
```

    void insertSorted(const ItemType& newEntry);
    bool removeSorted(const ItemType& anEntry);
    int getPosition(const ItemType& newEntry) const;
    // The following methods are the same as given in ListInterface:
    bool isEmpty() const;
    int getLength() const;
    bool remove(int position);
    void clear();
    ItemType getEntry(int position) const throw(PrecondViolatedExcep);
    }; // end LinkedSortedList
\#include "LinkedSortedList.cpp"
\#endif

```

\subsection*{12.2.2 The Implementation File}

The default constructor and destructor for our class LinkedSortedList have practically the same definitions as they do in the class LinkedList. We will, therefore, leave them to you to implement. You could also model the copy constructor after the one in LinkedList, but we will use recursion in our definition here. We begin by having the copy constructor call the private method copyChain, which will have a recursive definition:
```

template<class ItemType>
LinkedSortedList<ItemType>::
LinkedSortedList(const LinkedSortedList<ItemType>\& aList)
{
headPtr = copyChain(aList.headPtr);
} // end copy constructor

```

The private method copyChain begins by testing its pointer argument. If it contains nul1ptr, the pointer to the copy of the chain is set to nullptr and itemCount is set to zero. Otherwise, the method creates a new node containing the data from the first node of the given chain. The method then recursively inserts the new node into the copy of the chain. After each insertion, the method increments itemCount and returns a pointer to the new chain. The definition of copyChain is
```

temp1ate<c1ass ItemType>
Node<ItemType>* LinkedSortedList<ItemType>::
copyChain(const Node<ItemType>* origChainPtr)
{
Node<ItemType>* copiedChainPtr;
if (origChainPtr == nul1ptr)
{
copiedChainPtr = nul1ptr;
itemCount = 0;
}
else
{
// Build new chain from given one
Node<ItemType>* copiedChainPtr =
new Node<ItemType>(origChainPtr->getItem());
copiedChainPtr->setNext(copyChain(origChainPtr->getNext()));
itemCount++;
} // end if
return copiedChainPtr;
} // end copyChain

```

FIGURE 12-2 Places to insert strings into a sorted chain of linked nodes


The method insertSorted. Adding an entry to a sorted list requires that you find where in the list the new entry belongs. Since the entries are sorted, you compare the new entry with the entries in the sorted list until you reach an entry that is not smaller than the new entry. Figure 12-2 depicts a chain of linked nodes, each containing a string that is sorted alphabetically. The figure shows where the additional strings "Ally", "Cathy", "Luke", "Sue", and "Tom" would be inserted into the chain and the comparisons that would have to occur to arrive at those locations.

You can see from the figure that, in a string comparison, Ally is less than Bob, and so it would be inserted at the beginning of the chain. To see where to insert Luke, you would find that Luke is greater than both Bob and Jill but less than Mike. Thus, Luke belongs before Mike in the chain. Sue, on the other hand, is already in one of the nodes. You would discover that Sue is greater than Bob, Jill, and Mike but not greater than Sue. So you would insert the new entry Sue just before the existing entry Sue. Finally, Tom is greater than all the current names in the list, so you would add it to the end of the chain.


Note: Given a sorted list with entries in ascending order, you insert a new entry just before the first entry that is not smaller than the new entry.

Recall from earlier chapters that you add a new node to the beginning of a chain differently than at other points in the chain. Adding to the beginning is easy, since headPtr references the first node in the chain. To add anywhere else, you need a pointer to the node that will ultimately occur before the new node. Thus, while you traverse the chain of linked nodes to discover where the new entry belongs, you must retain a reference to the node prior to the one under consideration.

The following high-level algorithm describes our strategy:
```

// Adds a new entry to the sorted list.
add(newEntry)

```

Allocate a new node containing newEntry
Search the chain until either you find a node containing newEntry or you pass the point where it should be
Let prevPtr point to the node before the insertion point
if (the chain is empty or the new node belongs at the beginning of the chain) Add the new node to the beginning of the chain
e1se
Insert the new node after the node referenced by prevPtr
Increment the length of the sorted list

Assuming the private method getNodeBefore that we specified in the header file, we can use the previous algorithm to define the insertSorted method, as follows:
```

template<class ItemType>
void LinkedSortedList<ItemType>::insertSorted(const ItemType\& newEntry)
{
Node<ItemType>* newNodePtr = new Node<ItemType>(newEntry);
Node<ItemType>* prevPtr = getNodeBefore(newEntry);
if (isEmpty() || (prevPtr == nul1ptr)) // Add at beginning
{
newNodePtr->setNext(headPtr);
headPtr = newNodePtr;
}
else // Add after node before
{
Node<ItemType>* aftPtr = prevPtr->getNext();
newNodePtr->setNext(aftPtr);
prevPtr->setNext(newNodePtr);
} // end if
itemCount++;
} // end insertSorted

```

The private method getNodeBefore. We still need to implement the private method getNodeBefore to locate the node that is before the node that should or does contain the given entry. We will need two pointers as we traverse the list. Clearly, we need a pointer to the current node so we can compare its entry to the desired entry. But we also must retain a pointer to the previous node, because it is this pointer that the method returns. In the following implementation, these pointers are curPtr and prevPtr.
```

temp1ate<class ItemType>
Node<ItemType>* LinkedSortedList<ItemType>::
getNodeBefore(const ItemType\& anEntry) const
{
Node<ItemType>* curPtr = headPtr;
Node<ItemType>* prevPtr = nul1ptr;
while ( (curPtr != nul1ptr) \&\& (anEntry > curPtr->getItem()) )
{
prevPtr = curPtr;
curPtr = curPtr->getNext();
} // end while
return prevPtr;
} // end getNodeBefore

```

The programming problems at the end of this chapter ask you to complete this implementation of the sorted list.

Note: Since the ADTs sorted list and list share many of the same operations, portions of their implementations are identical.

Question 5 In the while statement of the method getNodeBefore, how important is the order of the two boolean expressions that the operator \&\& joins? Explain.
Question 6 What does getNodeBefore return if the sorted list is empty? How can you use this fact to simplify the implementation of the method insertSorted given previously?

Question 7 Suppose that you use the previous method insertSorted to add an entry to a sorted list. If the entry is already in the list, where in the list will the method insert it? Before the first occurrence of the entry, after the first occurrence of the entry, after the last occurrence of the entry, or somewhere else?
Question 8 What would be the answer to the previous question if you changed \(>\) to \(>=\) in the while statement of the method getNodeBefore?

\subsection*{12.2.3 The Efficiency of the Link-Based Implementation}

The performance of insertSorted depends on the efficiency of the method getNodeBefore. The latter method locates the insertion point by traversing the chain of nodes. This traversal is \(\mathrm{O}(n)\), making the addition to a sorted list an \(\mathrm{O}(n)\) operation. With the exception of isEmpty and getLength, which are \(\mathrm{O}(1)\) operations, the sorted-list methods are \(\mathrm{O}(n)\). We leave these observations for you to show as Exercise 11.

\subsection*{12.3 Implementations That Use the ADT List}

As we noted earlier in this chapter, the link-based implementation of the ADT sorted list repeats much of the corresponding implementation of the ADT list. Can we avoid this duplication of effort and reuse portions of the list's implementation? The answer to this question is yes, as you will soon see.

You can certainly use the ADT list to create and maintain an alphabetical list of strings. It is natural, then, to consider using the ADT list when implementing the ADT sorted list. Basically, you can do this using one of three techniques:
- Containment
- Public inheritance
- Private inheritance

In most cases, one or more of these three approaches will not be applicable to a particular problem; usually, one of the approaches will be best. However, we will use the sorted list to demonstrate all three techniques.

\subsection*{12.3.1 Containment}

A sorted list can maintain its entries within a list, as Figure 12-3 illustrates. You use a list as a data field within the class that implements the sorted list. As noted in C++ Interlude 4, this approach uses a

FIGURE 12-3 An instance of a sorted list that contains a list of its entries


FIGURE 12-4 SortedListHasA is composed of an instance of the class LinkedList

type of containment called composition and illustrates the has-a relationship between the class of sorted lists and the class of lists. If we name our new class SortedListHasA and use an instance of LinkedList, as given in Listing 9-2 of Chapter 9, Figure 12-4 illustrates the relationship between these two classes using UML notation.

The header file. Listing 12-3 declares the class SortedListHasA so that it has a private data member that is an instance of LinkedList containing the items in the sorted list. Our class will derive from the interface SortedListInterface, as given earlier in Listing 12-1.

\section*{LISTING 12-3 The header file for the class SortedListHasA}
```

/** ADT sorted list using the ADT list.
@file SortedListHasA.h */
\#ifndef _SORTED_LIST_HAS_A
\#define _SORTED_LIST_HAS_A
\#include "SortedListInterface.h"
\#include "ListInterface.h"
\#include "Node.h"
\#include "PrecondViolatedExcep.h"
template<class ItemType>
class SortedListHasA : public SortedListInterface<ItemType>
{
private:
ListInterface<ItemType>* 1istPtr;
public:
SortedListHasA();
SortedListHasA(const SortedListHasA<ItemType>\& sList);
virtua1 ~SortedListHasA();
void insertSorted(const ItemType\& newEntry);
bool removeSorted(const ItemType\& anEntry);
int getPosition(const ItemType\& newEntry) const;
// The following methods have the same specifications
// as given in ListInterface in Chapter 8:

```
```

    bool isEmpty() const;
    int getLength() const;
    bool remove(int position);
    void clear();
    ItemType getEntry(int position) const throw(PrecondViolatedExcep);
    }; // end SortedListHasA
\#include "SortedListHasA.cpp"
\#endif

```

The constructors and destructor. The constructor creates an instance of LinkedList and assigns a reference to it to the pointer variable 1istPtr:
```

template<class ItemType>
SortedListHasA<ItemType>::SortedListHasA()
{
1istPtr = new LinkedList<ItemType>();
} // end default constructor

```

The copy constructor creates a new list and then copies the entries in the given sorted list to the new list:
```

template<class ItemType>
SortedListHasA<ItemType>::
SortedListHasA(const SortedListHasA<ItemType>\& sList)
{
listPtr = new LinkedList<ItemType>();
for (int position = 1; position <= sList.getLength(); position++)
listPtr->insert(position, sList.getEntry(position));
} // end copy constructor

```

Note that we call LinkedList's method insert. While this particular definition is not very efficient, a recursive \(\mathrm{O}(n)\) implementation is possible. Exercise 7 at the end of this chapter asks you to investigate this possibility.

The destructor simply calls the method clear, which will deallocate the list:
```

template<class ItemType>
SortedListHasA<ItemType>::~SortedListHasA()
{
clear();
} // end destructor

```

The method insertSorted. The implementations of the remaining methods of the ADT sorted list are brief, as the list does most of the work. To add a new entry to the sorted list, we first use the method getPosition, which is an operation of the sorted list. We assume that it is already implemented, even though we have not written it yet. Recall that getPosition finds the position of an existing entry within a sorted list, or the position at which we should insert a new entry that does not occur in the sorted list. The method sets the sign of the integer it returns to indicate whether the entry exists in the list already. Since we allow duplicate entries in our sorted list, we can ignore this sign. Notice that the following implementation uses the standard method fabs to discard this sign. It also uses the insert method of the ADT list.
```

template<class ItemType>
void SortedListHasA<ItemType>::insertSorted(const ItemType\& newEntry)

```
```

{
int newPosition = fabs(getPosition(newEntry));
listPtr->insert(newPosition, newEntry);
} // end insertSorted

```

The method removeSorted. The method removeSorted can call getPosition to determine where in the sorted list the given entry exists or belongs. This time, however, we do need to know whether the entry exists in the list. If it does not exist, we cannot remove it, so removeSorted will return false when this is the case. Otherwise, the returned position will be positive, and removeSorted can call the list's method remove to delete the entry. We leave the implementation of removeSorted to you in Checkpoint Question 11.

The method getPosition. Implementing getPosition is somewhat harder than implementing the previous two methods. To decide where in a sorted list an entry is or belongs, we need to compare it to the entries already in the list, beginning with the first one. If the given entry is in the list, we obviously compare entries until we find a match. However, if it is not in the list, we want to stop the search at the point where it belongs in the sorted list. We take advantage of the sorted order of the objects by using logic similar to that described in Section 12.2.2

Checkpoint Question 12 asks you to define getPosition. You can do so by using the methods getLength and getEntry.

Question 9 Repeat Checkpoint Question 7 using the method insertSorted of the class SortedListHasA.
Question 10 Can a client of SortedListHasA invoke the method insert of the ADT list? Explain.

Question 11 Define the method removeSorted for the class SortedListHasA.
Question 12 Define the method getPosition for the class SortedListHasA.

The remaining methods. Each of the methods isEmpty, getLength, remove, clear, and getEntry of the ADT sorted list has the same specifications as in the ADT list. Each can simply invoke the corresponding list method. For example, the method remove has the following implementation in SortedListHasA:
```

temp1ate<class ItemType>
bool SortedListHasA<ItemType>::remove(int position)
{
return listPtr->remove(position);
} // end remove

```

Efficiency issues. Except perhaps for some subtle logic in getPosition, you can write the previous implementation quickly and with few, if any, errors. Saving human time is an attractive feature of using an existing class to build another. But does the implementation use computer time efficiently? In this particular implementation, several methods invoke getPosition, so their efficiency depends on getPosition's efficiency.

The method getPosition calls the list method getLength, which is an \(\mathrm{O}(1)\) operation. Therefore, we need not be concerned with it. On the other hand, a loop examines the entries in the list one at a time by invoking getEntry until the desired entry is located. Thus, the efficiency of getPosition depends in part on the efficiency of getEntry. However, the efficiency of getEntry depends upon which implementation of the ADT list you use. For the link-based implementation, getEntry is \(\mathrm{O}(n)\). Since getPosition invokes getEntry within a loop, getPosition is \(\mathrm{O}\left(n^{2}\right)\) in the worst case. Note that each time getEntry retrieves the next entry in the list, it starts its search at the beginning of the chain. This is the cause of getPosition's inefficiency.

Question 13 Suppose that instead of using LinkedList in the implementation of SortedListHasA, you used ArrayList. What Big O would describe the performance of the method getPosition?

Question 14 Give an advantage and a disadvantage of using containment in the implementation of the class SortedListHasA.

Figures 12-5 and 12-6 summarize the efficiencies of the operations for array-based and link-based implementations of the ADTs list and sorted list. Confirmation of these results is left as an exercise. As you can see, the implementation of the sorted list using containment is easy to write but is not very efficient if the underlying list uses a chain of linked nodes.

FIGURE 12-5 The worst-case efficiencies of ADT list operations for array-based and linkbased implementations
\begin{tabular}{|l|c|c|}
\hline \multicolumn{1}{|c}{ ADT List Operation } & \multicolumn{1}{c}{ Array-based } & Link-based \\
\hline insert(newPosition, newEntry) & \(\mathrm{O}(n)\) & \(\mathrm{O}(n)\) \\
remove(position) & \(\mathrm{O}(n)\) & \(\mathrm{O}(n)\) \\
getEntry(position) & \(\mathrm{O}(1)\) & \(\mathrm{O}(n)\) \\
setEntry(position, newEntry) & \(\mathrm{O}(1)\) & \(\mathrm{O}(n)\) \\
clear() & \(\mathrm{O}(1)\) & \(\mathrm{O}(n)\) \\
getLength(), isEmpty() & \(\mathrm{O}(1)\) & \(\mathrm{O}(1)\) \\
\hline
\end{tabular}

FIGURE 12-6 The worst-case efficiencies of the ADT sorted list operations when implemented using an instance of the ADT list
\begin{tabular}{ll|l|}
\hline \multicolumn{1}{c|}{ ADT Sorted List Operation } & \multicolumn{2}{c}{} \\
\hline & \multicolumn{2}{c}{\begin{tabular}{c} 
List Implementation \\
Array-based \\
Link-based
\end{tabular}} \\
\hline insertSorted(newEntry) & \(\mathrm{O}(n)\) & \(\mathrm{O}\left(n^{2}\right)\) \\
removeSorted(anEntry) & \(\mathrm{O}(n)\) & \(\mathrm{O}\left(n^{2}\right)\) \\
getPosition(anEntry) & \(\mathrm{O}(n)\) & \(\mathrm{O}\left(n^{2}\right)\) \\
getEntry(position) & \(\mathrm{O}(1)\) & \(\mathrm{O}(n)\) \\
remove(givenPosition) & \(\mathrm{O}(n)\) & \(\mathrm{O}(n)\) \\
clear() & \(\mathrm{O}(1)\) & \(\mathrm{O}(n)\) \\
getLength(), isEmpty() & \(\mathrm{O}(1)\) & \(\mathrm{O}(1)\) \\
\hline
\end{tabular}

\section*{Note: Using containment to implement the ADT sorted list}

When you use an instance of the ADT list to contain the entries in the ADT sorted list, you must use the list's operations to access the sorted list's entries, instead of accessing them directly. Such an implementation of the sorted list is easy to write but is inefficient when the underlying list uses a chain of linked nodes to store its entries.

An is-a relationship implies public inheritance

\subsection*{12.3.2 Public Inheritance}

A list is a container of items that you reference by position number. If you maintained those items in sorted order, would you have a sorted list? Ignoring name differences, most operations for the ADT list are almost the same as the corresponding operations for the ADT sorted list. The insertion and removal operations differ, however, and the ADT sorted list has an additional operation, getPosition.

You can insert an item into a sorted list by first using getPosition to determine the position in the sorted list where the new item belongs. You then use the list's insert operation to insert the item into that position in the list. You use a similar approach to remove an item from a sorted list.

Thus it appears that a sorted list is a list, so let's see whether we can use public inheritance. Figure 12-7 shows the class SortedListIsA as a descendant of the class LinkedList. This diagram represents an implementation in C++ using public inheritance. SortedListIsA inherits LinkedList's members, adds the method getPosition, and revises the insertion and removal operations.

The class SortedListIsA now has useful operations of the ADT list, such as getEntry, remove, clear, isEmpty, and getLength-which it inherits from the class List—in addition to the methods insertSorted, removeSorted, and getPosition. However, it also inherits insert and setEntry from LinkedList. By using either of these two position-oriented methods, a client could destroy the order of a sorted list. To prevent this from occurring, SortedListIsA must override them.

FIGURE 12-7 SortedListIsA as a descendant of LinkedList


The header file. Listing \(12-4\) shows a header file for the class SortedListIsA. Note again that we are using public inheritance. You might wonder why we are not using SortedListInterface. Doing so would require us to use multiple inheritance, which, while possible, is beyond our present scope. Since LinkedList is derived from ListInterface, you can think of SortedListIsA as a grandchild of ListInterface.

\section*{LISTING 12-4 A header file for the class SortedListIsA}
```

/** ADT sorted list using ADT list.
@file SortedListIsA.h */
\#ifndef _SORTED_LIST_IS_A
\#define _SORTED_LIST_IS_A
\#include "LinkedList.h"
\#include "Node.h"
\#include "PrecondViolatedExcep.h"
temp1ate<c1ass ItemType>
class SortedListIsA : pub1ic LinkedList<ItemType>

```
```

{
public:
SortedListIsA();
SortedListIsA(const SortedListIsA<ItemType>\& sList);
virtual ~SortedListIsA();
void insertSorted(const ItemType\& newEntry);
bool removeSorted(const ItemType\& anEntry);
int getPosition(const ItemType\& anEntry) const;
// The inherited methods remove, clear, getEntry, isEmpty, and
// getLength have the same specifications as given in ListInterface.
// The following methods must be overridden to disable their
// effect on a sorted list:
bool insert(int newPosition, const ItemType\& newEntry);
void setEntry(int position, const ItemType\& newEntry)
throw(PrecondViolatedExcep);
}; // end SortedListIsA
\#include "SortedListIsA.cpp"
\#endif

```

Method definitions. Let's examine the implementations of the sorted list methods. The default constructor and destructor are straightforward; the copy constructor invokes LinkedList's copy constructor by using an initializer:
```

template<class ItemType>
SortedListIsA<ItemType>: :SortedListIsA()
{
} // end default constructor
template<class ItemType>
SortedListIsA<ItemType>::
SortedListIsA(const SortedListIsA<ItemType>\& sList):
LinkedList<ItemType>(sList)
{
} // end copy constructor
template<class ItemType>
SortedListIsA<ItemType>:: SortedListIsA()
{
} // end destructor

```

The method insertSorted first calls getPosition to get the intended position of the new entry, ignores the sign of this position, and uses the result and LinkedList's insert method to complete its task. Note that since we will override insert in SortedListIsA, we must be careful not to call that implementation, but to call the base class implementation instead.
```

temp1ate<c1ass ItemType>
void SortedListIsA<ItemType>::insertSorted(const ItemType\& newEntry)
{
int newPosition = fabs(getPosition(newEntry));

```

Call LinkedList's version of insert
// We need to call the LinkedList version of insert, since the
// SortedListIsA version does nothing but return false
LinkedList<ItemType>::insert(newPosition, newEntry);
\} // end insertSorted
The method removeSorted uses an approach similar to that of insertSorted, but it does not ignore the sign of the position returned by getPosition. Insertion happens regardless of whether the new entry is a duplicate of another one already in the sorted list. However, you cannot remove an entry that is not in the list!
```

template<class ItemType>
bool SortedListIsA<ItemType>: :removeSorted(const ItemType\& anEntry)
{
boo1 ableToRemove = fa1se;
if (!LinkedList<ItemType>::isEmpty())
{
int position = getPosition(anEntry);
ab1eToRemove = position > 0;
if (ableToRemove)
ableToRemove = LinkedList<ItemType>::remove(position);
} // end if
return ableToRemove;
} // end removeSorted

```

The method getPosition uses the list's method getEntry to access each entry in the list sequentially until it either finds a match or reaches the point where the entry being sought would belong if it were in the list.
```

temp1ate<class ItemType>
int SortedListIsA<ItemType>::getPosition(const ItemType\& anEntry) const
{
int position = 1;
int length = LinkedList<ItemType>::getLength();
while ( (position <= length) \&\&
(anEntry > LinkedList<ItemType>::getEntry(position)) )
{
position++;
} // end while
if ( (position > length) ||
(anEntry != LinkedList<ItemType>::getEntry(position)) )
{
position = -position;
} // end if
return position;
} // end getPosition

```

Programming Tip: Notice that we have preceded the calls to methods inherited from LinkedList with the LinkedList<Item>: : namespace reference. When we called insert from insertSorted, we had to ensure that we called LinkedList's version of insert and not the overriding version in SortedListIsA. After all, the latter method does nothing but return false. Notice, however, that we do not override the methods remove, clear, getEntry, isEmpty, and getLength. Because these methods do not appear
explicitly in the header file for SortedListIsA, if you were to call them without the qualifier LinkedList<Item>: :, the C++ compiler would assume that they were client functions instead of class methods. You would get a syntax error.

An alternate way of calling these methods is to precede their calls by this-> instead of LinkedList<Item>: :. For example, you could write
if (!this->isEmpty())
instead of
if (!LinkedList<ItemType>::isEmpty())

Overridden methods. Recall that we have inherited two methods-insert and setEntry-from LinkedList that we do not want as public methods of the ADT sorted list. Although we called insert in our definition of insertSorted, we do not want clients to be able to destroy the sorted order of our entries by inserting an item at any specific position within a sorted list. To this end, we override insert so that it always returns false:
```

template<class ItemType>
bool SortedListIsA<ItemType>::
insert(int newPosition, const ItemType\& newEntry)
{
return false;
} // end insert

```

This approach ensures that a client of SortedListIsA can never use insert to add items by position to an object of SortedListIsA. We override setEntry in a similar manner.


Note: Recall that public inheritance enables you to use a derived-class object anywhere that you can use a base-class object. But you cannot use a sorted list in the same way that you can use a list and expect to maintain the values stored in the sorted list in sorted order. Moreover, the value-oriented nature of a sorted list does not support all of a list's operations. Therefore, a sorted list is not really a list, and hence public inheritance is not appropriate.

Programming Tip: You should use public inheritance only when two classes have an is-a relationship.


Question 15 What would have happened if you preceded the call to insert in the method insertSorted by this-> instead of LinkedList<ItemType>: :? Explain.

Efficiency issues. If you compare the implementation of SortedListIsA with the one for SortedListHasA, you will notice slight differences in the syntax, but the basic logic of the methods is the same. Both classes use methods of the ADT list to define their operations. The performance of corresponding methods is virtually the same.

Overriding insert prevents insertions into a sorted list by position

A sorted list is not a list; do not use public inheritance

\subsection*{12.3.3 Private Inheritance}

If you do not have an is- \(a\) relationship between your new class and an existing class, you should not use public inheritance. Instead, if you want to inherit members from the existing class, you can use private inheritance. Private inheritance enables you to use the methods of a base class without giving a client access to them.

The header file in Listing 12-5 declares the class SortedListAsA. Note that it is derived from the class LinkedList using private inheritance. Except for this addition and the name of the class, this header file is identical to the one in Listing 12-3 for SortedListHasA.

\section*{LISTING 12-5 The header file for the class SortedListAsA}
```

/** ADT sorted list using ADT 1ist.
@file SortedListAsA.h */
\#ifndef _SORTED_LIST_AS_A
\#define _SORTED_LIST_AS_A
\#include "SortedListInterface.h"
\#include "ListInterface.h"
\#include "Node.h"
\#include "PrecondViolatedExcep.h"
template<class ItemType>
class SortedListAsA : public SortedListInterface<ItemType>,
private LinkedList<ItemType>
{
public:
SortedListAsA();
SortedListAsA(const SortedListAsA<ItemType>\& sList);
virtual ~SortedListAsA();
< The rest of the public section is the same as in SortedListHasA in Listing 12-3. >
. . .
}; // end SortedListAsA
\#include "SortedListAsA.cpp"
\#endif

```

Because clients of SortedListAsA do not have access to the methods of the base class LinkedList, you need to provide a complete set of sorted list operations. This was not necessary for public inheritance. However, this lack of access also means that we do not have to override inherited methods such as setEntry that we do not want clients to have.

The implementation of SortedListAsA can use the public \({ }^{1}\) members of LinkedList. In fact, the methods insertSorted, removeSorted, and getPosition have the same implementations as they did with public inheritance. The remaining methods simply call the corresponding methods in LinkedList. For example, the method getEntry has the following definition:
```

temp1ate<c1ass ItemType>
ItemType SortedListAsA<ItemType>::getEntry(int position) const
throw(PrecondViolatedExcep)

```
\{
```

    return LinkedList<ItemType>::getEntry(position);
    ```
\} // end getEntry

With both private inheritance ( \(a s-a\) ) and containment (has-a), LinkedList is hidden from the clients of the sorted list. That is, you can use either private inheritance or containment to hide the underlying list in the implementation of the sorted list. Realize, however, that unlike public inheritance, private inheritance does not allow you to use an instance of SortedListAsA wherever you can use an instance of LinkedList; that is, SortedListAsA and LinkedList are not object-type compatible. The UML diagram in Figure 12-8 shows the inheritance relationship between the two classes. The notation <<implementation>> is a stereotype and is used to indicate that the SortedListAsA class is implemented in terms of the LinkedList class. You can use a stereotype to identify a unique characteristic of an element in any UML diagram.

FIGURE 12-8 The SortedListAsA class implemented in terms of the LinkedList class


\section*{Summary}
1. The ADT sorted list maintains its entries in sorted order. It, not the client, determines where to place an entry.
2. The ADT sorted list can add, remove, or locate an entry, given the entry as an argument.
3. The ADT sorted list has several operations that are the same as the corresponding operations of the ADT list. However, a sorted list will not let you add or replace an entry by position.
4. A chain of linked nodes provides a reasonably efficient implementation of the sorted list.
5. A class of sorted lists that has a list as a data field is said to use containment. Although such a class is easy to write, its efficiency can suffer if the implementation of the ADT list is inefficient.
6. Although it seems like a sorted list is a list, deriving a class of sorted lists from a class of lists using public inheritance is not appropriate. Doing so requires you to override some methods of the ADT list that a client of the sorted list class should not be able to use. Additionally, public inheritance would make the two classes object-type compatible, but you cannot use a sorted list anywhere that a list is used.
7. Private inheritance provides a reasonable alternative to containment as an approach to using the ADT list in the implementation of the ADT sorted list. Usually, however, containment is preferable.

\section*{Exercises}
1. Consider the classes Sphere and Bal1, as outlined here:
```

class Sphere
{
public:
doub7e getArea() const; // Surface area
void displayStatistics() const;
}; // end Sphere
class Ba11 : public Sphere
{
public:
doub7e getArea() const; // Cross-sectional area
void displayStatistics() const;
. . .
}; // end Ba77

```

Suppose that the implementation of each version of displayStatistics invokes the method getArea.
a. If mySphere is an instance of Sphere and myBa11 is an instance of Ba11, which version of getArea does each of the following calls to displayStatistics invoke? Explain your answer.
mySphere.displayStatistics();
myBa11.displayStatistics();
b. If the statements
```

Sphere* spherePtr;
Ba17* bal1Ptr;

```
declare spherePtr and ba11Ptr, which version of getArea does each of the following calls to displayStatistics invoke? Explain your answer.
```

spherePtr->displayStatistics();

```
spherePtr = bal1Ptr;
spherePtr->displayStatistics();
bal1Ptr->displayStatistics();
2. Define and implement a class Pen that has an instance of Ball as one of its members. Provide several members for the class Pen, such as the data member color and methods isEmpty and write.
3. Consider the following classes:
- LandVehicle represents a vehicle that travels on land. Its public methods include whee 1Count and speed.
- MotorizedLandVehicle represents a land vehicle that has a motor. Its public methods include engineCapacity and fue1Type.
a. Which of the methods mentioned previously can the implementation of speed invoke?
b. Which of the methods mentioned previously can the implementation of engineCapacity invoke?
4. Assume the classes described in Exercise 3 and consider a main method that contains the following statements:
```

LandVehicle 1andVeh;
MotorizedLandVehicle motorVeh;

```
a. Which of these objects can invoke the method wheelCount?
b. Which of these objects can invoke the method fuelType?
5. Consider the following classes:
```

class Expr
{
private:
char Array[MAX_STRING+1];
public:
int getLength() const;
virtual void display() const;
. .
}; // end Expr
class AlgExpr : public Expr
{
public:
bool isExpression() const;
bool isBlank(int first, int last) const;
}; // end AlgExpr
class InfixExpr : public AlgExpr
{
public:
bool isExpression() const;
int valueOf() const;
void display() const;
protected:
int endFactor(int first, int last) const;
int endTerm(int first, int last) const;
int endExpression(int first, int last) const;
private:
Stack<int> values;
Stack<char> operators;
}; // end InfixExpr

```

The class A7gExpr represents algebraic expressions, including prefix, postfix, and infix expressions. Its method isExpression simply examines the expression for valid characters but does not consider the order of the characters.

The class InfixExpr represents infix expressions. Its isExpression calls isB7ank, and its display calls valueOf.
a. Should isBTank be public, protected, or private? Explain.
b. If inExp is an instance of InfixExpr in the main method, can inExp invoke endExpression? Explain.
c. What small change(s) would you make to the classes to ensure that the correct version of isExpression is called?
6. Assume the classes described in Exercise 5 and consider a main method that contains the statements
```

Expr exp;
AlgExpr aExp;
InfixExpr inExp;

```
a. Which of these objects can correctly invoke the method getLength?
b. Which of these objects can correctly invoke the method isExpression?
c. Which of these objects can correctly invoke the method valueOf?
d. Give an example of object-type compatibility by writing a method declaration and a call to it that could appear in this main method.
7. Recall from Section 12.3.1 the copy constructor for the class SortedListHasA. Its definition uses a loop. Replace that definition with a recursive one, and show that it is \(\mathrm{O}(n)\).
8. Consider an ADT front list, which restricts insertions, removals, and retrievals to the first item in the list. Define and implement a class for the ADT stack that is a descendant of FrontList.
9. Define an abstract base class Person that describes a typical person. Next, define a derived class Student that describes a typical student. Finally, derive from Student a class GradStudent for a typical graduate student.
10. Design and implement the following classes:
a. An abstract base class Employee that represents a generic employee. Include methods to retrieve information about an employee.
b. A subclass of Employee called Hour 7yEmployee that describes an employee who gets paid by the hour. Include a public method called getPay that returns the pay of the employee for that month and any other relevant methods.
c. A subclass of Employee called NonHour 7yEmployee that describes an employee who gets paid a fixed salary every month. Include a public method called getPay that returns the pay of the employee for that month. Include any other relevant methods.
11. Confirm the results in Figures 12-5 and 12-6.
12. Imagine an unknown implementation of an ADT sorted list of integers. This ADT organizes its items into ascending order. Suppose that you have just read \(n\) integers into a one-dimensional array of integers called data. Write some \(\mathrm{C}++\) statements that use the ADT sorted list operations to sort the array into ascending order.
13. Write pseudocode that merges two sorted lists into a new third sorted list by using only ADT sorted list operations.
14. Define a set of axioms for the ADT sorted list and use them to prove that the sorted list of characters, which is defined by the sequence of operations
```

sList = an empty sorted list
sList.insertSorted('S')
sList.insertSorted('T')
sList.insertSorted('R')
sList.removeSorted('T')

```
is exactly the same as the sorted list defined by the sequence
```

sList = an empty sorted list
sList.insertSorted('T')
sList.insertSorted('R')
sList.removeSorted('T')
sList.insertSorted('S')

```

\section*{Programming Problems}
1. Complete the implementation of the class LinkedSortedList.
2. Define a class for an array-based implementation of the ADT sorted list. Consider a recursive implementation for getPosition. Should insertSorted and removeSorted call getPosition?
3. Complete the implementation of the class SortedListHasA.
4. Complete the implementation of the class SortedListAsA.
5. Consider the class FrontList that Exercise 8 describes. Implement FrontList in each of the following ways:
a. Store the list's entries in an instance of LinkedList.
b. Derive FrontList from LinkedList using public inheritance.
c. Derive FrontList from LinkedList using private inheritance.
6. The class LinkedList, as described in Sections 9.2.1 and 9.2.2 of Chapter 9, does not contain a method getPosition that returns the position number of a given entry. Define a descendant of LinkedList that has a method getPosition as well as methods that insert and remove items by their values instead of their positions. This new insertion method should always insert the new entry at the beginning of the list. Although the items in this list are not sorted, the new ADT is analogous to the ADT sorted list, which contains the method getPosition.
7. Consider an ADT circular list, which is like the ADT list but treats its first entry as if it were immediately after its last entry. For example, if a circular list contains six items, retrieval or removal of the eighth item actually involves the list's second item. Let insertion into a circular list, however, behave exactly like insertion into a list. Define and implement the ADT circular list as a derived class of LinkedList.
8. Programming Problem 13 in Chapter 6 describes the ADT traversable stack. In addition to the standard stack operations-isEmpty, push, pop, and peek-a traversable stack includes the operation traverse. This operation begins at the bottom of the stack and displays each item in the stack until it reaches the top of the stack.

Define and implement the ADT traversable stack as a derived class of ArrayStack, as given in Section 7.1 of Chapter 7.
9. Define a template interface for the ADT sorted list that is derived from ListInterface. Then define the class SortedListHasA that is derived from your new interface.
10. Because algebraic expressions are character strings, you can derive a class of algebraic expressions from a class of strings. Define such a class. Include an isExpression operation that uses the recognition algorithm given in Programming Problem 8 of Chapter 5 and a valueOf operation that uses the evaluation algorithm given in Programming Problem 8 of Chapter 6 .

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\section*{Queues and Priority Queues}

\section*{Chapter}

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\section*{Prerequisites}
\begin{tabular}{ll} 
Chapter 6 & Stacks \\
Chapter 8 & Lists (incidental) \\
Chapter 12 & Sorted Lists and Their Implementations (incidental)
\end{tabular}

Whereas a stack's behavior is characterized as last in, first out, a queue's behavior is characterized as first in, first out. This chapter defines the queue's operations and discusses several uses of them. As you will see, queues are common in everyday life. Their first-in, first-out behavior makes them appropriate ADTs for situations that involve waiting. Queues are also important in simulation, a technique for analyzing the behavior of complex systems. This chapter uses a queue to model the behavior of people in a line.

Sometimes the importance of an object depends on criteria other than when it is placed into a container. In such cases, you can assign each object a priority and organize them according to their priorities instead of chronologically. The priority queue will do this for you.

This chapter explores the two ADTs queue and priority queue. We will consider their implementations in the next chapter.

The ADT queue
FIFO: The first item inserted into a queue is the first item out

\subsection*{13.1 The ADT Queue}

A queue is like a line of people. The first person to join a line is the first person served, that is, the first to leave the line. New items enter a queue at its back, or rear, and items leave a queue from its front. Operations on a queue occur only at its two ends. This characteristic gives a queue its first-in, first-out (FIFO) behavior. In contrast, you can think of a stack as having only one end, because all operations are performed at the top of the stack. This characteristic gives a stack its last-in, first-out behavior.

As an abstract data type, the queue has the following operations:

\section*{Note: ADT queue operations}
- Test whether a queue is empty.
- Add a new entry to the back of the queue.
- Remove the entry at the front of the queue (the entry that was added earliest).
- Get the entry that was added earliest to the queue.

Queues occur in everyday life

Queues have applications in computer science

Queues are appropriate for many real-world situations. You wait in a queue-that is, a line-to buy a movie ticket, to check out at the bookstore, or to use an automatic teller machine. The person at the front of the queue is served, while new people join the queue at its back. Even when you call an airline to check your flight's status, your call actually enters a queue while you wait for the next available agent.

Queues also have applications in computer science. When you print an essay, the computer sends lines faster than the printer can print them. The lines are held in a queue for the printer, which removes them in FIFO order. If you share the printer with other computers, your request to print enters a queue to wait its turn.

Since all of these applications involve waiting, people study them to see how to reduce the wait. Such studies are called simulations, and they typically use queues. Later, this chapter examines a simulation of a line of customers at a bank.

The following operation contract specifies the ADT queue in more detail, and Figure 13-1 shows a UML diagram for the class Queue. Note that the conventional names of the operations that add or remove an entry are, respectively, enqueue (pronounced " \(\mathrm{N}-\mathrm{Q}\) ") and dequeue (pronounced "D-Q").

\section*{FIGURE 13-1 UML diagram for the class Queue}
\begin{tabular}{|l|}
\hline \multicolumn{1}{|c|}{ Queue } \\
\hline \\
\hline +isEmpty(): boolean \\
+enqueue(newEntry: ItemType) : boolean \\
+dequeue(): boolean \\
+peekFront() : ItemType
\end{tabular}

\section*{Abstract Data Type: Queue}

DATA
- A finite number of objects, not necessarily distinct, having the same data type and ordered by when they were added.

\section*{Operations}

Pseudocode
isEmpty ()

enqueue(newEntry)
dequeue()
peekFront()

\section*{Description}

Task: Sees whether this queue is empty. Input: None. Output: True if the queue is empty; otherwise false.
Task: Adds newEntry at the back of this queue.
Input: newEntry.
Output: True if the operation is successful; otherwise false.
Task: Removes the front of this queue. That is, removes the item that was added earliest.
Input: None.
Output: True if the operation is successful; otherwise false.
Task: Returns the front of this queue. That is, gets the item that was added earliest. The operation does not change the queue.
Input: None.
Output: The front of the queue.

Figure 13-2 illustrates the effect of these operations on a queue of integers. Notice that enqueue adds an item at the back of the queue and that peekFront looks at the item at the front of the queue, whereas dequeue removes the item at the front of the queue.

FIGURE 13-2 Some queue operations
\begin{tabular}{|c|c|c|}
\hline Operation & Front & Queue after operation \\
\hline & Front & \\
\hline aQueue = an empty queue & & \(\nabla\) \\
\hline aQueue.enqueue (5) & & 5 \\
\hline aQueue.enqueue(2) & & 52 \\
\hline aQueue.enqueue (7) & & 527 \\
\hline aQueue. peekFront() & & 527 (Returns 5) \\
\hline aQueue.dequeue() & & 27 \\
\hline aQueue.dequeue() & & 7 \\
\hline
\end{tabular}

Question 1 If you add the letters \(A, B, C\), and \(D\) in sequence to a queue of characters and then remove them, in what order will they leave the queue?

Question 2 What do the initially empty queues queue1 and queue2 "look like" after the following sequence of operations?
```

queue1.enqueue(1)
queue1.enqueue(2)
queue2.enqueue(3)
queue2.enqueue(4)
queue1.dequeue()
queueFront = queue2.peekFront()
queue1.enqueue(queueFront)
queue1.enqueue(5)
queue2.dequeue()
queue2.enqueue(6)

```

Compare these results with Checkpoint Question 2 in Chapter 6.

An interface. Listing 13-1 contains a C++ template that completes our specification of the ADT queue.

\section*{LISTING 13-1 A C++ interface for queues}
```

/** @file QueueInterface.h */
\#ifndef _QUEUE_INTERFACE
\#define _QUEUE_INTERFACE
template<class ItemType>
class QueueInterface
{
public:
/** Sees whether this queue is empty.
@return True if the queue is empty, or false if not. */
virtual bool isEmpty() const = 0;
/*** Adds a new entry to the back of this queue.
@post If the operation was successful, newEntry is at the
back of the queue.
@param newEntry The object to be added as a new entry.
@return True if the addition is successful or false if not. */
virtual bool enqueue(const ItemType\& newEntry) = 0;
/** Removes the front of this queue.
@post If the operation was successful, the front of the queue
has been removed.
@return True if the removal is successful or false if not. */
virtual bool dequeue() = 0;
/** Returns the front of this queue.
@pre The queue is not empty.
@post The front of the queue has been returned, and the
queue is unchanged.
@return The front of the queue. */
virtual ItemType peekFront() const = 0;
}; // end QueueInterface
\#endif

```

\subsection*{13.2 Simple Applications of the ADT Queue}

This section presents two simple applications of the ADT queue. The applications use the ADT queue operations independently of their implementations.

\subsection*{13.2.1 Reading a String of Characters}

When you enter characters at a keyboard, the system must retain them in the order in which you typed them. It could use a queue for this purpose, as the following pseudocode indicates:
```

// Read a string of characters from a single line of input into a queue
aQueue = a new empty queue
while (not end of line)
{
Read a new character into ch
aQueue.enqueue(ch)
}

```

Once the characters are in a queue, the system can process them as necessary. For example, if you had typed the integer 247-without any mistakes, but possibly preceded or followed by blanks-the queue would contain digits and possibly blanks. The system could convert the digits 2, 4, and 7 into the decimal value 247 by computing \(10 \times(10 \times 2+4)+7\).

The following pseudocode function performs this conversion in general:
```

// Converts digits in a queue aQueue into a decimal integer
getInteger(aQueue: Queue): integer
// Get first digit, ignoring any leading blanks
do
{
aQueue.dequeue(ch)
} while (ch is blank)
// Assertion: ch contains first digit
// Compute the integer n from digits in queue
n = 0
done = false
do
{
n=10*n+(integer that ch represents)
done = aQueue.isEmpty()
if (!done)
aQueue.dequeue(ch)
} while (!done and ch is a digit)
return n

```

\subsection*{13.2.2 Recognizing Palindromes}

Recall from Section 5.1.2 in Chapter 5 that a palindrome is a string of characters that reads the same from left to right as it does from right to left. In Chapter 6, you learned that you could use a stack to reverse the order of occurrences. You should realize by now that you can use a queue to preserve the order of occurrences. Thus, you can use both a queue and a stack to see whether a string is a palindrome.

As you traverse a string from left to right, you can add each character to both a queue and a stack. Figure 13-3 illustrates the result of this action for the string "abcbd", which is not a

You can use a queue in conjunction with a stack to recognize palindromes

FIGURE 13-3 The results of inserting the characters \(a, b, c, b, d\) into both a queue and a stack
\begin{tabular}{|c|c|c|c|}
\hline String: & Queue: & Stack & \\
\hline abcbd & \(\mathrm{a} b \mathrm{~b}\) b d & d
b & \multirow[t]{3}{*}{\(\leftarrow\) Top} \\
\hline & \(\uparrow\) ¢ & c & \\
\hline & Front Back & b
a & \\
\hline
\end{tabular}
palindrome. You can see that the first character in the string is at the front of the queue and the last character in the string is at the top of the stack. Thus, characters removed from the queue will occur in the order in which they appear in the string, and characters removed from the stack will occur in the opposite order.

Knowing this, you can compare the characters at the front of the queue and the top of the stack. If the characters are the same, you can remove them. You repeat this process until the stack and the queue become empty, in which case the original string is a palindrome, or the two characters are not the same, in which case the string is not a palindrome.

The following is a pseudocode version of a nonrecursive recognition algorithm for the language of palindromes:
```

// Tests whether a given string is a palindrome.
isPalindrome(someString: string): boolean
// Create an empty queue and an empty stack
aQueue = a new empty queue
aStack = a new empty stack
// Add each character of the string to both the queue and the stack
7ength = length of someString
for (i = 1 through 1ength)
{
nextChar = ith character of someString
aQueue.enqueue(nextChar)
aStack.push(nextChar)
}
// Compare the queue characters with the stack characters
charactersAreEqua1 = true
while (aQueue is not empty and charactersAreEqua1)
{
queueFront = aQueue.peekFront()
stackTop = aStack.peek()
if (queueFront equals stackTop)
{
aQueue.dequeue()
aStack.pop()
}
e1se
charactersAreEqua1 = false
}
return charactersAreEqual

```

Question 3 Trace the palindrome-recognition algorithm described in this section for each of the following strings of characters:
a. abcda
b. radar

Question 4 Improve the palindrome-recognition algorithm described in this section by adding the first length / 2 characters to the queue and then pushing the remaining characters onto the stack.

\subsection*{13.3 The ADT Priority Queue}

Imagine a person who visits a hospital's emergency room (ER). When any patient enters the hospital, the staff creates a record about that person in a database for later retrieval by nurses, doctors, and the billing department. In addition, the staff must keep track of the ER patients and decide when each person will receive care.

What ADT should the ER staff use for their patients? The ADT sorted list would facilitate the treatment of ER patients in alphabetical order by name or in numerical order by ID number. A queue would enable treatment of patients in the order of arrival. In either case, Ms. Zither, who was just rushed to the ER with acute appendicitis, would have to wait for Mr. Able to have a splinter removed. Clearly, the ER staff should assign some measure of urgency, or priority, to the patients waiting for treatment. The next available doctor should treat the patient with the highest priority. The ADT that the ER staff needs should produce this patient on request.

Another example of the use of priorities is your list of daily or weekly tasks. Suppose that your "to do" list for this week contains the following items:

\section*{Send a birthday card to Aunt Mabel.}

Start the research paper for world history.
Finish reading Chapter 13 of Walls and Mirrors.
Make plans for Saturday night.
When you consult your list, you most likely will attend to the task that, for you, has the highest priority.

A priority value indicates, for example, a patient's priority for treatment or a task's priority for completion. What quantity should you use for this priority value? Many reasonable possibilities exist, including a simple ranking from 1 to 10 . Let's arbitrarily decide that the largest priority value indicates the highest priority. The priority value becomes a part of the item that you insert into an ADT. You then ask the ADT for the item that has the highest priority.

Such an ADT is known as a priority queue. More formally, a priority queue is an ADT that provides the following operations:

You can organize data by priorities

You usually prioritize your list of tasks

\section*{Note: ADT priority queue operations}
- Test whether a priority queue is empty.
- Add a new entry to the priority queue in its sorted position based on priority value.
- Remove from the priority queue the entry with the highest priority value.
- Get the entry in the priority queue with the highest priority value.

A priority queue orders by priority values

The following operation contract specifies the ADT priority queue in more detail, and Figure 13-4 shows a UML diagram for the class PriorityQueue.

\section*{Abstract Data Type: Priority Queue}

\section*{Data}
- A finite number of objects, not necessarily distinct, having the same data type and ordered by priority.

\section*{Operations}

Pseudocode
isEmpty()
add(newEntry)
remove()
peek()

\section*{Description}

Task: Sees whether this priority queue is empty.
Input: None.
Output: True if the priority queue is empty; otherwise false.
Task: Adds newEntry to this priority queue.
Input: newEntry.
Output: True if the operation is successful; otherwise false.
Task: Removes the entry with the highest priority from this priority queue. Input: None.
Output: True if the operation is successful; otherwise false.
Task: Returns the entry in this priority queue with the highest priority. The operation does not change the priority queue.
Input: None.
Output: The entry with the highest priority.

FIGURE 13-4 UML diagram for the class PriorityQueue
\begin{tabular}{|l|}
\hline \multicolumn{1}{|c|}{ PriorityQueue } \\
\hline \\
+isEmpty(): boolean \\
+add(newEntry: ItemType) : boolean \\
+remove(): boolean \\
+peek(): ItemType
\end{tabular}

\subsection*{13.3.1 Tracking Your Assignments}

Professors and bosses like to assign tasks for us to do by certain dates. Using a priority queue, we can organize these assignments in the order in which we should complete them. Suppose that we order the assignments by their due dates. A task with the earliest due date will have the highest priority. We can define a class Assignment of tasks that includes a data field date representing a task's due date. Figure 13-5 shows a diagram of such a class.

\section*{FIGURE 13-5 UML diagram for the class Assignment}
\begin{tabular}{|l|}
\hline \multicolumn{1}{|c|}{ Assignment } \\
\hline \begin{tabular}{l} 
course-the course code \\
task-a description of the assignment \\
date-the due date
\end{tabular} \\
\hline \begin{tabular}{l} 
+getCourseCode(): string \\
+getTask(): string \\
+getDueDate(): string
\end{tabular} \\
\hline
\end{tabular}

The following pseudocode shows how you could use a priority queue to organize your assignments and other responsibilities so that you know which one to complete first:
```

assignmentLog = a new priority queue using due date as the priority value
project = a new instance of Assignment
essay = a new instance of Assignment
task = a new instance of Assignment
errand = a new instance of Assignment
assignmentLog.add(project)
assignmentLog.add(essay)
assignmentLog.add(task)
assignmentLog.add(errand)
cout << "I should do the following first: "
cout << assignmentLog.peek()

```

\subsection*{13.4 Application: Simulation}

Simulation-a major application area for computers-is a technique for modeling the behavior of both natural and human-made systems. Generally, the goal of a simulation is to generate statistics that summarize the performance of an existing system or to predict the performance of a proposed system. In this section we will consider a simple example that illustrates one important type of simulation.

A problem to solve. Ms. Simpson, president of the First City Bank of Springfield, has heard her customers complain about how long they have to wait for service at the branch located in a downtown grocery store. Because she fears losing those customers to another bank, she is considering whether to hire a second teller for that branch.

Before Ms. Simpson hires another teller, she would like an approximation of the average time a customer has to wait for service from that branch's only teller. Ms. Simpson heard you were great at solving problems and has come to you for help. How can you obtain this information for Ms. Simpson?

Considerations. You could stand with a stopwatch in the bank's lobby all day, but that task is not particularly exciting. Besides, you should use an approach that also allows Ms. Simpson to predict how much improvement she could expect if the bank hired a given number of additional tellers. She certainly does not want to hire the tellers on a trial basis and then monitor the bank's performance before making her final decision.

Simulation models the behavior of systems

Simulated time

Sample arrival and transaction times

You conclude that the best way to obtain the information needed is to use a computer model to simulate the behavior of the bank. The first step in simulating a system such as a bank is to construct a mathematical model that captures the relevant information about the system. For example, how many tellers does the bank employ? How often do customers arrive? How long do the customers' transactions take?

If the model accurately describes the real-world system, a simulation can derive accurate predictions about the system's overall performance. For example, a simulation could predict the average time a customer has to wait before receiving service. A simulation can also evaluate proposed changes to the real-world system, such as predicting the effect of hiring more tellers at the bank. A large decrease in the time predicted for the average wait of a customer might justify the cost of hiring additional tellers.

After discussing the problem with Ms. Simpson, you decide that you want the simulation to determine
- The average time a customer waits to begin service from the current single teller
- The decrease in customer wait time with each new teller added

Simulation time and events. Central to a simulation is the concept of simulated time. Envision a stopwatch that measures time elapsed during a simulation. For example, suppose that the model of the bank specifies only one teller. At time 0 , which is the start of the banking day, the simulated system would be in its initial state with no customers. As the simulation runs, the stopwatch ticks away units of time-perhaps minutes-and certain events occur. At time 20, the bank's first customer arrives. Because there is no line, the customer goes directly to the teller and begins her transaction, which will take about 6 minutes to complete. At time 22, a second customer arrives. Because the first customer has not yet completed her transaction, the second customer must wait in line. At time 26, the first customer completes her transaction and the second customer can begin his. Figure 13-6 illustrates these four times in the simulation.

To gather the information you need, you run this simulation for a specified period of simulated time. During the course of the run, you need to keep track of certain statistics, such as the average time a customer has to wait for service. Notice that in the small example of Figure 13-6, the first customer had to wait 0 minutes to begin a transaction and the second customer had to wait 4 minutes to begin a transaction-an average wait of 2 minutes.

One point not addressed in the previous discussion is how to determine when certain events occur. For example, why did we say that the first customer arrived at time 20 and the second at time 22? After studying real-world systems like our bank, mathematicians learned to model events such as the arrival of people by using techniques from probability theory. This statistical information is incorporated into the mathematical model of the system and is used to generate events in a way that reflects the real world. The simulation uses these events and is thus called an eventdriven simulation. Note that the goal is to reflect the long-term average behavior of the system rather than to predict occurrences of specific events. This goal is sufficient for the needs of our simulation.

Although the techniques for generating events to reflect the real world are interesting and important, they require a good deal of mathematical sophistication. Therefore, we simply assume that we already have a list of events available for our use. In particular, for the bank problem, we assume that a file contains the time of each customer's arrival-an arrival event-and the duration of that customer's transaction once the customer reaches the teller. For example, the data
\begin{tabular}{cc} 
Arrival time & Transaction length \\
\cline { 1 - 1 } 20 & 6 \\
22 & 4 \\
23 & 2 \\
30 & 3
\end{tabular}

FIGURE 13-6 A bank line at time (a) 0; (b) 20; (c) 22; (d) 26
(a)

(b)

indicates that the first customer arrives 20 minutes into the simulation and her transactiononce begun-requires 6 minutes; the second customer arrives 22 minutes into the simulation, and his transaction requires 4 minutes; and so on. Assume that the input file is ordered by arrival time.

The use of a data file with predetermined event information is common in simulations. It allows us to try many different scenarios or bank teller configurations with the same set of events to ensure a fair comparison.

Notice that the file does not contain departure events; the data does not specify when a customer will complete the transaction and leave. In fact, the departure time of a customer cannot be determined until the simulation is run, so the simulation must determine when departures occur. By using the arrival time and the transaction length, the simulation can easily determine the time at which a

The results of a simulation
customer departs. To compute the departure time, we add the length of the transaction to the time when the customer begins the transaction.

For example, if we run the simulation by hand with the previous data, we would compute the departure times as follows:

A first attempt at a simulation algorithm

\section*{Time}

20
22
\(35 \quad\) Customer 4 departs

> 26 Customer 1 departs; customer 2 begins transaction Determine customer 2 departure event is at time 30
> \(30 \quad\) Customer 2 departs; customer 3 begins transaction Determine customer 3 departure event is at time 32 Customer 4 enters bank and stands at end of line Customer 3 departs; customer 4 begins transaction Determine customer 4 departure event is at time 35
> Event
> Customer 1 enters bank and begins transaction Determine customer 1 departure event is at time 26 Customer 2 enters bank and stands at end of line Customer 3 enters bank and stands at end of line

30
32

A customer's wait time is the elapsed time between arrival in the bank and the start of the transaction, that is, the amount of time the customer spends in line. The average of this wait time over all the customers is the statistic that you want to obtain.

To summarize, this simulation is concerned with two kinds of events:

\section*{Note: Kinds of events in an event-driven simulation}
- Arrival events indicate the arrival at the bank of a new customer. The input file specifies the times at which the arrival events occur. As such, they are externally generated events. When a customer arrives at the bank, one of two things happens. If the teller is idle when the customer arrives, the customer goes to the teller and begins the transaction immediately. If the teller is busy, the new customer must stand at the end of the line and wait for service.
- Departure events indicate the departure from the bank of a customer who has completed a transaction. The simulation determines the times at which the departure events occur. Thus, they are internally generated events. When a customer completes the transaction, he or she departs and the next person in line-if there is one-begins a transaction. tion, hat is, he amount ofine the custorner spends in line. The average of this wait time over all the

Event loop. The main tasks of an algorithm that performs a simulation are to repeatedly determine the times at which events occur and to process the events when they do occur. In simulation and gaming applications, this process is referred to as the event loop. The algorithm is stated at a high level as follows:
```

// Initialize
currentTime = 0
Initialize the line to "no customers"
while (currentTime <= time of the final event)
{

```
```

    if (an arrival event occurs at time currentTime)
        Process the arrival event
    if (a departure event occurs at time currentTime)
        Process the departure event
    // When an arrival event and departure event occur at the same time,
    // arbitrarily process the arrival event first
    currentTime++
    }

```

But do you really want to increment currentTime by 1 ? You would for a time-driven simulation, where you would determine arrival and departure times at random and compare those times to currentTime. Video games use this approach, since events can occur or need to be processed in almost every unit of time, which is typically a frame. In such a case, you would increment currentTime by 1 to simulate the ticking of a clock.

Recall, however, that this simulation is event driven, so you have a file of predetermined arrival times and transaction times. Because you are interested only in those times at which arrival and departure events occur, and because no action is required between events, you can advance currentTime from the time of one event directly to the time of the next.

Thus, you can revise the pseudocode solution as follows:
```

Initialize the line to "no customers"
while (events remain to be processed)
{
currentTime = time of next event
if (event is an arrival event)
Process the arrival event
else
Process the departure event

```
        // When an arrival event and a departure event occur at the same time,
        // arbitrarily process the arrival event first
\}

You must determine the time of the next arrival or departure so that you can implement the statement
```

currentTime $=$ time of next event

```

To make this determination, you must maintain an event list. An event list contains all arrival and departure events that will occur but have not occurred yet. The times of the events in the event list are in ascending order, and thus the next event to be processed is always at the beginning of the list. The algorithm simply gets the event from the beginning of the list, advances to the time specified, and processes the event. The difficulty, then, lies in successfully managing the event list.

Managing and processing customers and events. As customers arrive, they go to the back of the line. The current customer, who was at the front of the line, is being served, and it is this customer that you remove from the system next. It is thus natural to use a queue, bankQueue, to represent the line of customers in the bank. For this problem, the only information that you must store in the queue about each customer is the time of arrival and the length of the transaction.

Arrival events and departure events are ordered by time, and we always want to remove and process the next event that should occur-the highest-priority event. The ADT priority queue is used in this way. Our events can be stored in the priority queue eventListPQueue. We can initialize eventListPQueue with the arrival events in the simulation data file and later add the departure events as they are generated.

A time-driven simulation simulates the ticking of a clock

An event-driven simulation considers only the times of certain events, in this case, arrivals and departures

First revision of the simulation algorithm

An event list contains all future arrival events and departure events

Two tasks are required to process each event

The algorithm for arrival events

A new customer always enters the queue and is served while at the queue's front

The algorithm for departure events

But how can you determine the times for the departure events? Observe that the next departure event always corresponds to the customer that the teller is currently serving. As soon as a customer begins service, the time of his or her departure is simply
\[
\text { time of departure }=\text { time service begins }+ \text { length of transaction }
\]

Recall that the length of the customer's transaction is in the event list, along with the arrival time. Thus, as soon as a customer begins service, you place a departure event corresponding to this customer in the event list. Figure 13-7 illustrates a typical instance of an arrival event and a departure event used in this simulation.

FIGURE 13-7 A typical instance of (a) an arrival event; (b) a departure event


Now consider how you can process an event when it is time for the event to occur. You must perform two general types of actions:
- Update the bank line: Add or remove customers.
- Update the event list: Add or remove events.

To summarize, you process an arrival event as follows:
```

// TO PROCESSAN ARRIVAL EVENT
// Update the event list
Remove the arrival event for customer C from the event list
// Update the bank line
if (bank line is empty and teller is available)
{
Departure time of customer C is current time + transaction length
Add a departure event for customer C to the event list
Mark the teller as unavailable
}
else
Add customer C to the bank line

```

When customer \(C\) arrives at the bank, if the line is empty and the teller is not serving another customer, customer \(C\) can go directly to the teller. The wait time is 0 and you insert a departure event into the event list. If other customers are in line, or if the teller is assisting another customer, customer \(C\) must go to the end of the line.

You process a departure event as follows:
// TO PROCESSA DEPARTURE EVENT
// Update the event list
Remove the departure event from the event list
// Update the bank line
if (bank line is not empty)
```

{
Remove customer C from the front of the bank line
Customer C begins transaction
Departure time of customer C is current time + transaction length
Add a departure event for customer C to the event list
}
else
Mark the teller as available.

```

When a customer finishes a transaction and leaves the bank, if the bank line is not empty, the next customer \(C\) leaves the line and goes to the teller. You insert a departure event for customer \(C\) into the event list.

You can now combine and refine the pieces of the solution into an algorithm that performs the simulation by using the ADTs queue and priority queue:
```

// Performs the simulation.
simulate(): void
Create an empty queue bankQueue to represent the bank line
Create an empty priority queue eventListPQueue for the event list
tellerAvailable = true
// Create and add arrival events to event list
while (data file is not empty)
{
Get next arrival time a and transaction time t from file
newArrivalEvent = a new arrival event containing a and t
eventListPQueue.add(newArriva1Event)
}
// Event loop
while (eventListPQueue is not empty)
{
newEvent = eventListPQueue.peek()
// Get current time
currentTime = time of newEvent
if (newEvent is an arrival event)
processArriva1(newEvent, eventListPQueue, bankQueue)
else
processDeparture(newEvent, eventListPQueue, bankQueue)
}
// Processes an arrival event.
processArrival(arriva1Event: Event, eventListPQueue: PriorityQueue,
bankQueue: Queue)
// Remove this event from the event list
eventListPQueue.remove()
customer = customer referenced in arrivalEvent
if (bankQueue.isEmpty() \&\& te11erAvailable)
{
departureTime = currentTime + transaction time in arrivalEvent
newDepartureEvent = a new departure event with departureTime
eventListPQueue.add(newDepartureEvent)
tellerAvailable = false
}
else
bankQueue.enqueue(customer)

```

The final pseudocode for the event-driven simulation
```

// Processes a departure event.
+processDeparture(departureEvent: Event, eventListPQueue: PriorityQueue,
bankQueue: Queue)
// Remove this event from the event list
eventListPQueue.remove()
if (!bankQueue.isEmpty())
{
// Customer at front of line begins transaction
customer = bankQueue.peek()
bankQueue.dequeue()
departureTime = currentTime + transaction time in customer
newDepartureEvent = a new departure event with departureTime
eventListPQueue.add(newDepartureEvent)
}
else
tel1erAvailable = true

```

Figure 13-8 begins a trace of this algorithm for the data given earlier and shows the changes to the queue and priority queue. Checkpoint Question 6 asks you to complete the trace. There are several more implementation details that must be decided, such as how to represent customers and events. Programming Problem 6 at the end of this chapter asks you to complete the implementation of this simulation.

FIGURE 13-8 A trace of the bank simulation algorithm for the data
206
224
232
303


Question 5 In the bank simulation problem, why is it impractical to read the entire input file and create a list of all the arrival and departure events before the simulation begins?

Question 6 Complete the hand trace of the bank-line simulation that Figure 13-8 began. Show the state of the queue and the event list at each step.

\subsection*{13.5 Position-Oriented and Value-Oriented ADTs}

Of the abstract data types that we have seen so far, the stack, the list, and the queue have operations defined in terms of the positions of their data items. We call them position-oriented ADTs. Stacks and queues greatly restrict the positions that their operations can affect; only their end positions can be accessed. The list removes this restriction.

The ADT sorted list is an example of a value-oriented ADT. When a new item is added to a sorted list, the sorted list determines the item's placement according to its value. Unlike a list, a sorted list cannot be told where to place its items. The same is true of a priority queue, since it orders its entries according to their priorities, which are values that are a part of each entry.

Although stacks and queues order their entries, and you do not specify where a stack or a queue should place its items, the entries' positions are not determined according to their values. Thus, we do not classify stacks and queues as value oriented.

Note: The ADTs stack, queue, list, sorted list, and priority queue all order their entries. The stack, queue, and list are position oriented; the sorted list and priority queue are value oriented. The ADT bag does not order its entries; it is neither position oriented nor value oriented.

Stacks are really quite similar to queues. This similarity becomes apparent if you pair off their operations, as follows:
- Stack isEmpty and queue isEmpty see whether any items exist in the ADT.
- push and enqueue insert a new item into one end (the top and back, respectively) of the ADT.
- pop and dequeue: The pop operation removes the most recent item, which is at the top of the stack, and dequeue removes the first item, which is at the front of the queue.
- Stack peek and queue peekFront: The peek operation retrieves the most recent item, which is at the top of the stack, and peekFront retrieves the first item at the front of the queue.

The ADT list allows you to insert into, remove from, and inspect an item at any position of the list. Thus, it has the most flexible operations of the three position-oriented ADTs. You can view the list operations as general versions of the stack and queue operations as follows:
- getLength: If you ignore the restriction that the stack and queue versions of isEmpty can tell only when an item is present, you obtain an operation that can count the number of items that are present.
- insert: If you ignore the restriction that push and enqueue can insert new items into only one position, you obtain an operation that can insert a new item into any position of the list.
- remove: If you ignore the restriction that pop and dequeue can remove items from only one position, you obtain an operation that can remove an item from any position of the list.
- getEntry: If you ignore the restriction that peek and peekFront can retrieve items from only one position, you obtain an operation that can retrieve the item from any position of the list.

Operations for the ADTs list, stack, and queue involve the position of items

A comparison of stack and queue operations

ADT list operations generalize stack and queue operations

Question 7 For each of the following situations, which of these ADTs (1 through 6) would be most appropriate? (1) a queue; (2) a stack; (3) a list; (4) a sorted list; (5) a priority queue; (6) none of these
a. The customers at a6deli counter who take numbers to mark their turn
b. An alphabetic list of names
c. Integers that need to be sorted
d. The boxes in a box trace of a recursive function
e. A grocery list ordered by the occurrence of the items in the store
f. The items on a cash register tape
g. A word processor that allows you to correct typing errors by using the Backspace key
h. A program that uses backtracking
i. A list of ideas in chronological order
j. Airplanes that stack above a busy airport, waiting to land
k. People who are put on hold when they call for customer service
1. An employer who fires the most recently hired person

\section*{Summary}
1. The definition of the queue operations gives the ADT queue first-in, first-out (FIFO) behavior.
2. Models of real-world systems often use queues. The event-driven simulation in this chapter used a queue to model a line of customers in a bank.
3. Central to a simulation is the notion of simulated time. In a time-driven simulation, simulated time is advanced by a single time unit, whereas in an event-driven simulation, simulated time is advanced to the time of the next event. To implement an event-driven simulation, you maintain an event list that contains events that have not yet occurred. The event list is ordered by the time of the events, so that the next event to occur is always at the head of the list.
4. A priority queue has operations to retrieve or remove the item with the highest priority.
5. ADTs are classified as either position oriented or value oriented. A position-oriented ADT organizes its items according to their positions, whereas a value-oriented ADT organizes items by their values.

\section*{Exercises}
1. Consider the palindrome-recognition algorithm described in Section 13.2.2. Is it necessary for the algorithm to look at the entire queue and stack? That is, can you reduce the number of times that the loop must execute?
2. Consider the language
\[
L=\left\{s \$ s^{\prime}: s \text { is a possibly empty string of characters other than } \$, s^{\prime}=\operatorname{reverse}(s)\right\}
\]
as defined in Chapter 6. Write a recognition algorithm for this language that uses both a queue and a stack. Thus, as you traverse the input string, you insert each character of \(s\) into a queue and each character of \(s^{\prime}\) into a stack. Assume that each input string contains exactly one \(\$\).
3. What is the output of the following pseudocode, where num 1 , num 2 , and num 3 are integer variables?
```

num1 = 5
num2 = 1
num3 = 4
aQueue.enqueue(num2)
aQueue.enqueue(num3)
aQueue.dequeue()
aQueue.enqueue(num1 - num2)
num1 = aQueue.peek()
aQueue.dequeue()
num2 = aQueue.peek()
aQueue.dequeue()
cout << num2 << " " << num1 << " " << num3 << endl

```
4. Revise the infix-to-postfix conversion algorithm of Chapter 6 so that it uses a queue to represent the postfix expression.
5. Write a client function that returns the back of a queue while leaving the queue unchanged. This function can call any of the methods of the ADT queue. It can also declare new Queue objects. The return type is ItemType, and it accepts a Queue object as a parameter.
6. Consider a queue implementation that uses the ADT list to represent the items in the queue. Discuss the efficiency of the queue's insertion and removal operations when the ADT list's implementation is
a. Array based
b. Link based
7. An operation that displays the contents of a queue can be useful during program debugging. Add a display operation to the ADT queue such that display uses only ADT queue operations, so it is independent of the queue's implementation.
8. Write a \(\mathrm{C}++\) template interface PriorityQueueInterface for the ADT priority queue
9. Consider a slight variation of the ADT queue. In this variation, new items can be added to and removed from either end. This ADT is commonly called a double-ended queue, or deque. Specify each method of the deque by stating the method's purpose; by describing its parameters; and by writing preconditions, postconditions, and a pseudocode version of its header. Then write a C++ template interface for these methods that includes javadoc-style comments.
10. Use a deque, as described in the previous exercise, to solve the read-and-correct problem given in Section 6.1 .1 of Chapter 6. In that problem, you enter text at a keyboard and correct typing mistakes by using the Backspace key. Each backspace erases the most recently entered character. Your pseudocode solution should provide a corrected string of characters in the order in which they were entered at the keyboard.
11. With the following data, hand-trace the execution of the bank-line simulation that this chapter describes. Each line of data contains an arrival time and a transaction time. Show the state of the queue and the event list at each step.
\begin{tabular}{ll}
5 & 9 \\
7 & 5 \\
14 & 5 \\
30 & 5 \\
32 & 5 \\
34 & 5
\end{tabular}

Note that at time 14, there is a tie between the execution of an arrival event and a departure event.
12. In the solution to the bank simulation problem, can the event list be a queue? Can the event list be a list or sorted list?
13. Consider the stack-based search of the flight map in the HPAir problem of Chapter 6. You can replace the stack that searchS uses with a queue. That is, you can replace every call to push with a call to enqueue, every call to pop with a call to dequeue, and every call to peek with a call to peekFront. Trace the resulting algorithm when you fly from \(P\) to \(Z\) in the flight map in Figure 6-6. Indicate the contents of the queue after every operation on it.
14. As Chapter 6 pointed out, you can define ADT operations in a mathematically formal way by using axioms. Consider the following axioms for the ADT queue, where aQueue is an arbitrary queue and item is an arbitrary queue item:
```

(new Queue()).isEmpty() = true
(new Queue()).dequeue() = false
(new Queue()).peekFront() = error
((new Queue()).enqueue(item)).dequeue() = new Queue()
((new Queue()).enqueue(item)).peekFront() = item
(aQueue.enqueue(item)).isEmpty() = false
(aQueue.enqueue(item)).dequeue() = true

```

If aQueue is not empty,
(aQueue.enqueue(item)).dequeue() = (aQueue.dequeue()). enqueue(item)
and
(aQueue.enqueue(item)).peekFront() = aQueue.peekFront()
a. Note the recursive nature of the definition of peekFront. What is the base case? What is the recursive step? What is the significance of the isEmpty test? Why is the queue operation peekFront recursive in nature while the stack operation peek for the ADT stack is not?
b. The representation of a stack as a sequence of push operations without any pop operations is called a canonical form. (See Exercise 15 in Chapter 6.) Is there a canonical form for the ADT queue? That is, can you represent a queue as a sequence of enqueue operations without any dequeue operations? Prove your answer.

\section*{Programming Problems}
1. Using the class queue in the Standard Template Library, define and test the class OurQueue that is derived from QueueInterface, as given in Listing 13-1. The class queue has the following methods that you can use to define the methods for OurQueue:
```

queue(); // Defau7t constructor
bool empty() const; // Tests whether the queue is empty
void push(const ItemType\& newEntry); // Adds newEntry to the back of the queue
void pop(); // Removes the front of the queue
ItemType\& front(); // Returns a reference to the front of the queue

```

To access queue, use the following include statement:
```

\#include <queue>;

```
2. Using the class priority_queue in the Standard Template Library, define and test the class OurPriorityQueue that is derived from PriorityQueueInterface, as developed in Exercise 8. The class priority_queue has the following methods that you can use to define the methods for OurPriorityQueue:
```

priority_queue();
bool empty() const;
void push(const ItemType\& newEntry);
void pop();
ItemType\& top();
// Defau7t constructor
// Tests whether the priority queue is empty
// Adds newEntry to the priority queue
// Removes the entry having the highest priority
// Returns a reference to the entry having the
// highest priority

```

To access priority_queue, use the following include statement:
\#include <priority_queue>;
Whenever you need a queue or a priority queue for any of the following problems, use the classes OurQueue and OurPriorityQueue that Programming Problems 1 and 2 ask you to write.
3. Implement the palindrome-recognition algorithm described in Section 13.2.2.
4. Implement the recognition algorithm that you wrote to solve Exercise 2 using the classes OurQueue, as described in Programming Problem 1, and OurStack, as described in Programming Problem 1 of Chapter 6.
5. Implement the radix sort of an array by using a queue for each group. The radix sort is discussed in Section 11.2.3 of Chapter 11.
6. Implement the event-driven simulation of a bank that this chapter described. A queue of arrival events will represent the line of customers in the bank. Maintain the arrival events and departure events in a priority queue, sorted by the time of the event. Use a link-based implementation for the event list.

The input is a text file of arrival and transaction times. Each line of the file contains the arrival time and required transaction time for a customer. The arrival times are ordered by increasing time.

Your program must count customers and keep track of their cumulative waiting time. These statistics are sufficient to compute the average waiting time after the last event has been processed. Display a trace of the events executed and a summary of the computed statistics (the total number of arrivals and average time spent waiting in line). For example, the input file shown in the left columns of the following table should produce the output shown in the right column.

Input file
\begin{tabular}{ll}
1 & 5 \\
2 & 5 \\
4 & 5 \\
20 & 5 \\
22 & 5 \\
24 & 5 \\
26 & 5 \\
28 & 5 \\
30 & 5 \\
88 & 3
\end{tabular}

Output from processing file on left
\begin{tabular}{ll} 
Simulation Begins & \\
Processing an arrival event at time: & 1 \\
Processing an arrival event at time: & 2 \\
Processing an arrival event at time: & 4 \\
Processing a departure event at time: & 6 \\
Processing a departure event at time: & 11 \\
Processing a departure event at time: & 16 \\
Processing an arrival event at time: & 20 \\
Processing an arrival event at time: & 22 \\
Processing an arrival event at time: & 24 \\
Processing a departure event at time: & 25 \\
Processing an arrival event at time: & 26 \\
Processing an arrival event at time: & 28 \\
Processing an arrival event at time: & 30
\end{tabular}
Processing a departure event at time: ..... 30
Processing a departure event at time: ..... 35
Processing a departure event at time ..... 40
Processing a departure event at time ..... 45
Processing a departure event at time: ..... 50
Processing an arrival event at time: ..... 88
Processing a departure event at time: 91Simulation Ends
Final Statistics:
Total number of people processed: 10
Average amount of time spent waiting: 5.6
7. Modify and expand the event-driven simulation program that you wrote in Programming Problem 6.
a. Add an operation that displays the event list, and use it to check your hand trace in Exercise 11.
b. Add some statistics to the simulation. For example, compute the maximum wait in line, the average length of the line, and the maximum length of the line.
c. Modify the simulation so that it accounts for three tellers, each with a distinct line. You should keep in mind that there should be
- Three queues, one for each teller
- A rule that chooses a line when processing an arrival event (for example, enter the shortest line)
- Three distinct departure events, one for each line
- Rules for breaking ties in the event list

Run both this simulation and the original simulation on several sets of input data. How do the statistics compare?
d. The bank is considering the following change: Instead of having three distinct lines (one for each teller), there will be a single line for the three tellers. The person at the front of the line will go to the first available teller. Modify the simulation of part \(c\) to account for this variation. Run both simulations on several sets of input data. How do the various statistics compare (averages and maximums)? What can you conclude about having a single line as opposed to having distinct lines?
8. The people who run the Motor Vehicle Department (MVD) have a problem. They are concerned that people do not spend enough time waiting in lines to appreciate the privilege of owning and driving an automobile. The current arrangement is as follows:
- When people walk in the door, they must wait in a line to sign in.
- Once they have signed in, they are told either to stand in line for registration renewal or to wait until they are called for license renewal.
- Once they have completed their desired transaction, they must go and wait in line for the cashier.
- When they finally get to the front of the cashier's line, if they expect to pay by check, they are told that all checks must get approved. To do this, it is necessary to go to the check-approver's table and then reenter the cashier's line at the end.

Write an event-driven simulation to help the MVD gather statistics. Each line of input will contain
- A desired transaction code ( \(L\) for license renewal, \(R\) for registration renewal)
- A method-of-payment code ( \(\$\) for cash, \(C\) for check)
- An arrival time (integer)
- A name

Write out the specifics of each event (when, who, what, and so on). Then display these final statistics:
- The total number of license renewals and the average time spent in MVD (arrival until completion of payment) to renew a license
- The total number of registration renewals and the average time spent in MVD (arrival until completion of payment) to renew a registration

Incorporate the following details into your program:
- Define the following events: arrive, sign in, renew license, renew registration, and cope with the cashier (make a payment or find out about check approval).
- In the case of a tie, let the order of events be determined by the list of events just given-that is, arrivals have the highest priority.
- Assume that the various transactions take the following amounts of time:
\begin{tabular}{ll} 
Sign in & 10 seconds \\
Renew license & 90 seconds \\
Register automobile & 60 seconds \\
See cashier (payment) & 30 seconds \\
See cashier (check not approved) & 10 seconds
\end{tabular}
- As ridiculous as it may seem, the people waiting for license renewal are called in alphabetical order. Note, however, that people are not pushed back once their transactions have started.
- For the sake of this simulation, you can assume that checks are approved instantly. Therefore, the rule for arriving at the front of the cashier's line with a check that has not been approved is to go to the back of the cashier's line with a check that has been approved.

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\section*{Queue and Priority Queue Implementations}

\section*{Chapter}

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\section*{Prerequisites}

Chapter 13 Queues and Priority Queues

The previous chapter discussed the ADT queue at length and introduced the ADT priority queue. This chapter will present several implementations of the queue and compare their advantages and disadvantages.

We will consider an implementation of the priority queue, but we can write a much more efficient implementation later, after we explore the ADT heap in Chapter 17.

\subsection*{14.1 Implementations of the ADT Queue}

Like stacks, queues can have an array-based or a link-based implementation. However, we also can use an implementation of the ADT list to define a class of queues. Such a class is easy to write, as most of the work is done by the class of lists. If you needed a class of queues right away, you could use this approach. The result would not be as time efficient as possible, however.

VideoNote
Overview of LinkedQueue

We will begin by using an instance of the class LinkedList, as given in Listing 9-2 of Chapter 9, to store the queue's entries. We will then write a link-based implementation, as it is a bit more straightforward than the array-based one, which we will investigate last.

\subsection*{14.1.1 An Implementation That Uses the ADT List}

You can use the ADT list to contain the items in a queue. This approach is especially useful if you need to define a class of queues quickly and already have a class of lists. Figure 14-1 illustrates such a queue. If the item in position 1 of a list represents the front of the queue, you can implement the method dequeue as the list operation remove (1) and the method peekFront as the list operation getEntry (1). Similarly, if you let the item at the end of the list represent the back of the queue, you can implement the operation enqueue (newEntry) as the list operation insert (getLength() +1 , newEntry).

FIGURE 14-1 An implementation of the ADT queue that stores its entries in a list


The header file containing such a class definition for the ADT queue appears in Listing 14-1, and the implementation file is in Listing 14-2.

\section*{LISTING 14-1 The header file for the class ListQueue}
```

/*** ADT queue: ADT list implementation.
afile ListQueue.h */
\#ifndef _LIST_QUEUE
\#define _LIST_QUEUE
\#include "QueueInterface.h"
\#include "LinkedList.h"
\#include "PrecondViolatedExcep.h"
template<class ItemType>
class ListQueue : public QueueInterface<ItemType>
{
private:
LinkedList<ItemType>* listPtr; // Pointer to list of queue items
public:
ListQueue();
ListQueue(const ListQueue\& aQueue);
~ListQueue();

```
```

    bool isEmpty() const;
    boo1 enqueue(const ItemType& newEntry);
    bool dequeue();
    /** @throw PrecondViolatedExcep if queue is empty. */
    ItemType peekFront() const throw(PrecondViolatedExcep);
    }; // end ListQueue
\#include "ListQueue.cpp"
\#endif

```

\section*{LISTING 14-2 The implementation file for the class ListQueue}
```

/** ADT queue: ADT 1ist implementation.
@file ListQueue.cpp */
\#include "ListQueue.h" // Header file
temp1ate<c1ass ItemType>
ListQueue<ItemType>::ListQueue()
{
1istPtr = new LinkedList<ItemType>();
} // end default constructor
temp1ate<c1ass ItemType>
ListQueue<ItemType>::ListQueue(const ListQueue\& aQueue) :
1istPtr(aQueue.1istPtr)
{
} // end copy constructor
template<class ItemType>
ListQueue<ItemType>::~ListQueue()
{
} // end destructor
template<class ItemType>
boo1 ListQueue<ItemType>::isEmpty() const
{
return listPtr->isEmpty();
} // end isEmpty
temp1ate<c1ass ItemType>
boo1 ListQueue<ItemType>::enqueue(const ItemType\& newEntry)
{
return listPtr->insert(listPtr->getLength() + 1, newEntry);
} // end enqueue
template<class ItemType>
boo1 ListQueue<ItemType>::dequeue()
{
return listPtr->remove(1);
} // end dequeue
temp1ate<c1ass ItemType>
ItemType ListQueue<ItemType>::peekFront() const
throw(PrecondViolatedExcep)

```
```

    if (isEmpty())
        throw PrecondViolatedExcep("peekFront() called with empty queue.");
    // Queue is not empty; return front
    return listPtr->getEntry(1);
    } // end peekFront
// end of implementation file

```
\{

Of particular note in the implementation file are the definitions of the constructor, copy constructor, and destructor. The constructor creates a new instance of a list. Omitting this step is an easy oversight to make, leading to incorrect execution instead of a syntax error. The copy constructor uses an initializer, 1istPtr(aQueue. 1istPtr), to invoke the list's copy constructor. Finally, even though the destructor has an empty body, the list's destructor will be invoked.

Exercise 8 at the end of this chapter asks you to consider the efficiency of this implementation.

\subsection*{14.1.2 A Link-Based Implementation}

A link-based implementation of a queue uses a chain of linked nodes, much like the other link-based implementations that you have seen. However, the queue presents a challenge, since we must be able to not only remove entries from its front but also add them to its back. Removing a node from the beginning of a linked chain is easy, but to add a new node to the chain's end, we need a pointer to the chain's last node. One way to accomplish this is to begin at the first node and traverse the chain until we reach the last one. A much more efficient approach uses a tail pointer to reference the end of the chain-just as the head pointer references the beginning of the chain. Figure 14-2 illustrates a chain of linked nodes that has both head and tail pointers. Like the head pointer frontPtr, backPtr is external to the chain.

A linear linked chain or a circular linked chain can represent a queue

FIGURE 14-2 A chain of linked nodes with head and tail pointers


Figure 14-3 shows that you can actually get by with one external pointer-to the back-if you make the last node point to the first one. This data structure is a circular chain of linked nodes. Notice that the nodes in a circular chain have next pointers that never contain nul1ptr. We will refer to a chain like the one in Figure 14-2 as a linear chain, regardless of how many external pointers it has. Such a chain does have a node whose next pointer is nullptr.

Programming Problem 1 at the end of the chapter asks you to consider the details of the circular chain implementation. Here we will develop an implementation of the ADT queue using a chain that has both head and tail pointers, as illustrated in Figure 14-2.

FIGURE 14-3 A circular chain of linked nodes with one external pointer


Note: If you use a linear chain with only a head pointer to implement a queue, the enqueue operation will be inefficient. Each addition to the queue requires a traversal to the end of the chain. As the queue increases in length, the traversal time - and hence enqueue's time requirement-will increase.

Listing 14-3 shows the header file for our class definition.

\section*{LISTING 14-3 The header file for the class LinkedQueue}
```

/** ADT queue: Link-based implementation.
@file LinkedQueue.h */
\#ifndef _LINKED_QUEUE
\#define _LINKED_QUEUE
\#include "QueueInterface.h"
\#include "Node.h"
\#include "PrecondViolatedExcep.h"
template<class ItemType>
class LinkedQueue : public QueueInterface<ItemType>
{
private:
// The queue is implemented as a chain of linked nodes that has
// two external pointers, a head pointer for the front of the queue
// and a tail pointer for the back of the queue.
Node<ItemType>* backPtr;
Node<ItemType>* frontPtr;
public:
LinkedQueue();
LinkedQueue(const LinkedQueue\& aQueue);
~LinkedQueue();
bool isEmpty() const;
boo1 enqueue(const ItemType\& newEntry);
bool dequeue();

```
/** @throw PrecondViolatedExcep if the queue is empty */
ItemType peekFront() const throw(PrecondViolatedExcep);
\}; // end LinkedQueue
\#include "LinkedQueue.cpp"
\#endif

The method enqueue. Inserting a new node, to which newNodePtr points, at the back of the chain that represents the queue requires three pointer changes: the next pointer in the new node, the next pointer in the current back node, and the external pointer backPtr. Figure 14-4 illustrates these changes during the addition of an item to a nonempty queue and indicates the order in which they can occur. The statements to perform this addition are
```

newNodePtr->setNext(nu11ptr);
backPtr->setNext(newNodePtr);
backPtr = newNodePtr;

```

FIGURE 14-4 Adding an item to a nonempty queue


The addition of an item to an empty queue is a special case, as Figure \(14-5\) illustrates. If newNodePtr points to the new node, the following statements add the node to the empty chain:
```

frontPtr = newNodePtr;
backPtr = newNodePtr

```

These statements easily follow from the realization that the chain has only one node, which is both the first and last node in the chain.

FIGURE 14-5 Adding an item to an empty queue: (a) before enqueue; (b) after enqueue


Thus, we have the following definition for the method enqueue:
```

template<class ItemType>
bool LinkedQueue<ItemType>::enqueue(const ItemType\& newEntry)
{
Node<ItemType>* newNodePtr = new Node<ItemType>(newEntry);
// Insert the new node
if (isEmpty())
frontPtr = newNodePtr; // The queue was empty
else
backPtr->setNext(newNodePtr); // The queue was not empty
backPtr = newNodePtr; // New node is at back
return true;
} // end enqueue

```

The method dequeue. Removing the front of the queue involves deleting the first node of the chain. This is an easier operation than removing the last node, which, fortunately, we do not have to do. Figure 14-6 illustrates the removal of the front item from a queue that contains more than one item. Notice that you need to change only the external pointer frontPtr. Removal from a queue of one item is a special case that sets the external pointers backPtr and frontPtr to nul1ptr.

The following definition of the method dequeue shows how to implement these two cases:
```

template<class ItemType>
boo1 LinkedQueue<ItemType>::dequeue()
{
bool result = false;
if (!isEmpty())
{
// Queue is not empty; remove front
Node<ItemType>* nodeToDeletePtr = frontPtr;
if (frontPtr == backPtr)
{ // Special case: one node in queue
frontPtr = nul1ptr;
backPtr = nul1ptr;
}
else
frontPtr = frontPtr->getNext();
// Return deleted node to system
nodeToDe1etePtr->setNext(nul1ptr);
delete nodeToDeletePtr;
nodeToDeletePtr = nul1ptr;
result = true;
} // end if
return result;
} // end dequeue

```

FIGURE 14-6 Removing an item from a queue of more than one item


Retrieval. The method peekFront simply returns the value frontPtr->getItem() after checking that the queue is not empty. We leave the rest of this implementation to you as an exercise. (See Exercises 1 through 3 at the end of this chapter.)

Question 1 Why is a tail pointer desirable when you use a chain of linked nodes to implement a queue?

Question 2 If you use a circular chain that has only a tail pointer, as Figure 14-3 illustrates, how do you access the data in the first node?

Question 3 If the ADT queue had a method clear that removed all entries from a queue, what would its definition be in the previous link-based implementation?

\subsection*{14.1.3 An Array-Based Implementation}

For applications in which a fixed-sized queue does not present a problem, you can use an array to represent a queue. A naive array-based implementation of a queue might include the following definitions, as Figure 14-7a illustrates:
```

const int MAX_QUEUE = maximum size of queue;
ItemType items[MAX_QUEUE]; // Array of queue items
int front; // Index to front of queue
int back; // Index to back of queue

```

FIGURE 14-7 \(\quad\) (a) A naive array-based implementation of a queue; (b) rightward drift can cause the queue to appear full
items
(a)

\(\longleftarrow\) Array indices

\section*{items}
(b)


Here front and back are the indices of the front and back entries, respectively, in the queue. Initially, front is 0 and back is -1 . To add a new item to the queue, you increment back and place the item in items [back]. To remove an item, you simply increment front. The queue is empty whenever back is less than front. The queue is full when back equals MAX_QUEUE -1 .

The problem with this strategy is rightward drift - that is, after a sequence of additions and removals, the items in the queue will drift toward the end of the array, making it appear full. In other words, back could equal MAX_QUEUE - 1 even when the queue contains only a few items. Figure 14-7b illustrates this situation.

One possible solution to this problem is to shift array entries to the left, either after each removal from the queue or whenever back equals MAX_QUEUE -1 . This solution guarantees that the queue can always contain up to MAX_QUEUE items. Shifting is not really satisfactory, however, as it would dominate the cost of the implementation.

Question 4 Suppose that we change the naive array-based implementation of a queue pictured in Figure 14-7 so that the back of the queue is in items [0]. Although repeated removals from the front would no longer cause rightward drift, what other problem would this implementation cause?

A much more elegant solution is possible by viewing the array as circular, as Figure 14-8 illustrates. To remove an item, you increment the queue index front, and to insert an item, you increment

A naive array-based implementation of a queue

VideoNote
Overview of ArrayQueue

Rightward drift can cause a queue-full condition even though the queue contains few entries

Shifting entries to compensate for rightward drift is expensive

\section*{A circular array} eliminates rightward drift

FIGURE 14-8 A circular array as an implementation of a queue

back. Figure 14-9 illustrates the effect of a sequence of three queue operations on front, back, and the array. Notice that front and back "advance" clockwise around the array.

When either front or back advances past MAX_QUEUE -1 , it should wrap around to 0 . This wraparound eliminates the problem of rightward drift, which occurred in the previous naive implementation, because here the circular array has no end. You obtain the wraparound effect of a circular queue by using modulo arithmetic (that is, the \(\mathrm{C}++\%\) operator) when incrementing front and back. For example, you can add newEntry to the queue by using the statements

Adding to a queue

Deleting from a queue
```

back = (back + 1) % MAX_QUEUE;
items[back] = newEntry;

```

Notice that if back equaled MAX_QUEUE - 1 before the addition of newItem, the first statement, back \(=(\) back +1\() \%\) MAX_QUEUE, would have the effect of wrapping back around to index 0 . Similarly, you can remove the entry at the front of the queue by using the statement
```

front = (front + 1) % MAX_QUEUE;

```

To initialize the queue, you set front to 0 and back to MAX_QUEUE - 1 .

FIGURE 14-9 The effect of three consecutive operations on the queue in Figure 14-8
```

dequeue(); dequeue(); enqueue(9);

```


FIGURE 14-10 (a) front passes back when the queue becomes empty; (b) back catches up to front when the queue becomes full


The only difficulty with this scheme is detecting when the queue is empty or full. It seems reasonable to select as the queue-empty condition
front is one slot ahead of back
as this appears to indicate that front "passes" back when the queue becomes empty, asFigure 14-10a depicts. However, it is also possible that this condition signals a full queue: Because the queue is circular, back might in fact "catch up" with front as the queue becomes full. Figure 14-10b illustrates this situation.

Obviously, you need a way to distinguish between the two situations. One way is to keep a count of the number of items in the queue. Before adding an item to the queue, you check whether the count is equal to MAX_QUEUE; if it is, the queue is full. Before removing an item from the queue, you check whether the count is equal to zero; if it is, the queue is empty.
The header file. Listing 14-4 contains the header file for an array-based implementation of the ADT queue that uses a circular array as just described. Because the data is stored in statically allocated memory, the compiler-generated destructor and copy constructor are sufficient. \({ }^{1}\)
front and back cannot be used to distinguish between queue-full and queue-empty conditions

By counting queue items, you can detect queue-full and queue-empty conditions

\section*{LISTING 14-4 The header file for the class ArrayQueue}
```

/** ADT queue: Circular array-based implementation.
@file ArrayQueue.h */
\#ifndef _ARRAY_QUEUE
\#define _ARRAY_QUEUE
\#include "QueueInterface.h"
\#include "PrecondViolatedExcep.h"
const int MAX_QUEUE = 50;
template<class ItemType>
class ArrayQueue : public QueueInterface<ItemType>
{
private:
ItemType items[MAX_QUEUE]; // Array of queue items
int front; // Index to front of queue
int back; // Index to back of queue
int count; // Number of items currently in the queue
public:
ArrayQueue();
// Copy constructor and destructor supplied by compiler
bool isEmpty() const;
bool enqueue(const ItemType\& newEntry);
bool dequeue();
/** @throw PrecondViolatedExcep if queue is empty. */
ItemType peekFront() const throw(PrecondViolatedExcep);
}; // end ArrayQueue
\#include "ArrayQueue.cpp"
\#endif

```

The implementation file. Listing 14-5 contains the definitions of ArrayQueue's methods as they appear in the implementation file.

\section*{LISTING 14-5 The implementation file for the class ArrayQueue}
```

/** ADT queue: Circular array-based implementation.
@file ArrayQueue.cpp */
\#include "ArrayQueue.h" // Header file
template<class ItemType>
ArrayQueue<ItemType>::ArrayQueue() : front(0), back(MAX_QUEUE - 1), count(0)
{
} // end default constructor
temp1ate<c1ass ItemType>
bool ArrayQueue<ItemType>::isEmpty() const
{
return count == 0;
} // end isEmpty

```
```

template<class ItemType>
boo1 ArrayQueue<ItemType>::enqueue(const ItemType\& newEntry)
{
bool result = false;
if (count < MAX_QUEUE)
{
// Queue has room for another item
back = (back + 1) % MAX_QUEUE;
items[back] = newEntry;
count++;
result = true;
} // end if
return result;
} // end enqueue
template<class ItemType>
bool ArrayQueue<ItemType>::dequeue()
{
bool result = false;
if (!isEmpty())
{
front = (front + 1) % MAX_QUEUE;
count--;
result = true;
} // end if
return result;
} // end dequeue
temp1ate<c1ass ItemType>
ItemType ArrayQueue<ItemType>::peekFront() const
throw(PrecondViolatedExcep)
{
// Enforce precondition
if (isEmpty())
throw PrecondViolatedExcep("peekFront() called with empty queue");
// Queue is not empty; return front
return items[front];
} // end peekFront

```

Question 5 If the ADT queue had a method clear that removed all entries from a queue, what would its definition be in the previous array-based implementation?

Variations. Several commonly used variations of the previous circular-array approach do not require a count of the number of entries in the queue. One approach uses a boolean variable isFull to distinguish between the full and empty conditions. The expense of maintaining this variable is about the same as that of maintaining a counter, however. A faster implementation declares MAX_QUEUE + 1 locations for the array items, but uses only MAX_QUEUE of them for queue items. You sacrifice one

An isFull flag can replace the counter

Using an extra array location is more time efficient

Fixed size versus dynamic size

Reuse of an already implemented class saves you time

FIGURE 14-11 A more time-efficient circular implementation: (a) a full queue; (b) an empty queue
(a)

(b)

array location and let front be the index of the location before the front of the queue. As Figure 14-11 illustrates, the queue is full if
```

front equals (back + 1) % (MAX_QUEUE + 1)

```
but the queue is empty if

\section*{front equals back}

This approach does not have the overhead of maintaining a counter or boolean variable, and so is more efficient of time. Programming Problems 3 and 4 discuss these two alternate implementations further.

\subsection*{14.1.4 Comparing Implementations}

We have suggested implementations of the ADT queue that store the queue's entries in either an instance of the ADT list, a chain of linked nodes that has both a head pointer and a tail pointer, a circular chain that has only one external pointer, an array, or a circular array. You have seen the details of three of these implementations. All of our implementations of the ADT queue are ultimately either array based or link based.

The reasons for making the choice between array-based and link-based implementations are the same as those discussed in earlier chapters. The discussion here is similar to the one in Section 4.5 of Chapter 4. We repeat the highlights here in the context of queues.

An implementation based on a statically allocated array prevents the enqueue operation from adding an item to the queue if the array is full. Such a queue is appropriate for many data structures, such as buffers, within an operating system. If this restriction is not acceptable, you must use either a dynamically allocated array or a link-based implementation.

Suppose that you decide to use a link-based implementation. Should you choose the implementation that uses a linked chain or the one that uses a link-based implementation of the ADT list? Because a linked chain actually represents the items in the ADT list, using the ADT list to represent a queue is not as efficient as using a linked chain directly. However, the ADT list approach is much simpler to write.

If you decide to use a linked chain instead of the ADT list to represent the queue, should you use a linear chain or a circular chain? We leave this question for you to answer in Programming Problem 1.

\subsection*{14.2 An Implementation of the ADT Priority Queue}

While we could place the entries of a priority queue into either an array or a chain of linked nodes, doing so would be like implementing the ADT sorted list, which we did in Chapter 12. Rather than repeating those implementations, we can use a sorted list to contain the entries of a priority queue. The header file in Listing 14-6 defines a class of priority queues that has as a private data member an instance of the class LinkedSortedList, as given in Listing 12-2 of Chapter 12.
```

LISTING 14-6 A header file for the class SL_PriorityQueue.
/*** ADT priority queue: ADT sorted list implementation.
@file SL_PriorityQueue.h */
\#ifndef _PRIORITY_QUEUE
\#define _PRIORITY_QUEUE
\#include "PriorityQueueInterface.h"
\#include "LinkedSortedList.h"
\#include "PrecondViolatedExcep.h"
temp7ate<c7ass ItemType>
class SL_PriorityQueue : public PriorityQueueInterface<ItemType>
{
private:
LinkedSortedList<ItemType>* slistPtr; // Pointer to sorted 7ist of
// items in the priority queue
public:
SL_PriorityQueue();
SL_PriorityQueue(const SL_PriorityQueue\& pq);
~SL_PriorityQueue();
bool isEmpty() const;
bool add(const ItemType\& newEntry);
bool remove();
/** @throw PrecondViolatedExcep if priority queue is empty. */
ItemType peek() const throw(PrecondViolatedExcep);
}; // end SL_PriorityQueue
\#include "SL_PriorityQueue.cpp"
\#endif

```

Before we can implement the class SL_PriorityQueue, we need to consider how the sorted list behaves. LinkedSortedList maintains the entries in a sorted list in sorted order. It does so by comparing the entries with each other. For a priority queue, we want these comparisons to be based on the priorities of the entries. Let's assume that each entry is an object that has its priority as a data member. You will learn in C++ Interlude 5 how to make a comparison operator, such as \(>\), base its comparison on specific aspects of its two operands. For now, let's assume that the comparisons are made according to priority. Note that the highest-priority value can be either the largest value or the smallest value, according to the application at hand.

To give you an idea of how to implement this class, let's look at the definitions of the add and remove operations:

\footnotetext{
template<class ItemType>
boo1 SL_PriorityQueue<ItemType>::add(const ItemType\& newEntry)
}
```

{
slistPtr->insertSorted(newEntry);
return true;
} // end add
temp7ate<c7ass ItemType>
bool SL_PriorityQueue<ItemType>::remove()
{
// The highest-priority item is at the end of the sorted list
return slistPtr->remove(slistPtr->getLength());
} // end remove

```

We will leave the remainder of this implementation to you as an exercise. Note, however, that we will discuss a more time-efficient implementation in Chapter 17, when we introduce the ADT heap.

Question 6 Define the method peek for the sorted list implementation of the ADT priority queue.

\section*{Summary}
1. The enqueue and dequeue operations for a queue require efficient access to both ends of the queue. Therefore, \(a\) link-based implementation of a queue uses a linear linked chain that has both a head pointer and a tail pointer. Alternatively, you can use a circular linked chain.
2. An array-based implementation of a queue is prone to rightward drift. This phenomenon can make a queue look full when it really is not. Shifting the items in the array is one way to compensate for rightward drift. A more efficient solution uses a circular array.
3. If you use a circular array to implement a queue, you must be able to distinguish between the queue-full and queue-empty conditions. You can make this distinction by counting the number of items in the queue, using a boolean variable isFul1, or leaving one array location empty.
4. You can use an array, a chain of linked nodes, or a sorted list to contain the entries in a priority queue. A more time-efficient implementation is possible by using the ADT heap, which we will consider later in this book.

\section*{Exercises}
1. Implement the copy constructor for the class LinkedQueue that is declared in Listing 14-3. Hint: Look at the copy constructor for the ADT stack in Listing 7-4 of Chapter 7.
2. Repeat the previous exercise, but implement a memory-safe copy constructor instead. If a memory allocation fails, this constructor should release all memory that was allocated prior to the failure and then throw an exception.
3. The destructor for the class LinkedQueue that is declared in Listing 14-3 could repeatedly call dequeue. Although easy to write, this destructor can be inefficient due to repeated method calls. Write another implementation for the destructor that deallocates the linked chain directly without calling dequeue.
4. An operation that displays the contents of a queue can be useful during program debugging. Add a display operation to the ADT queue such that display assumes and uses the link-based implementation of the ADT queue.
5. Consider a method called getNumberOfElements that returns the number of elements in a queue without changing the queue. The return type is int, and it does not accept any parameters.
a. Add getNumberOfElements to the array-based ADT queue given in this chapter.
b. Add getNumberOfElements to the link-based ADT queue given in this chapter.
6. Show that the implementation of a queue that uses an extra array element as described in Section 14.1.3 is more efficient than the counter variation, even though it appears to perform two extra additions in the test:
```

front equals (back + 1) % (MAX_QUEUE + 1)

```
7. Consider the class ListQueue that implements the ADT queue by using the ADT list to represent the items in the queue. Discuss the efficiency of the queue's enqueue and dequeue operations when the ADT list's implementation is
a. Array based
b. Link based
8. The class ListQueue, as given in Listing 14-1, maintains the queue's front at the end of a list of the queue's entries and has the back of the queue at the beginning of that list. Note that the list is an object of the class LinkedList. What is the impact on the efficiency of the operations enqueue and dequeue if we were to maintain the queue's front at the beginning of the list and the queue's back at the list's end?
9. Complete the implementation of the class SL_PriorityQueue, as described in Section 14.2.
10. Implement the class SL_PriorityQueue, assuming that low values indicate a higher priority. Such a priority queue is useful within operating systems and other similar applications. When the underlying sorted list has a link-based implementation, using the lowest value for the highest priority makes the priority queue operations more efficient. Describe why this benefit is true.

\section*{Programming Problems}
1. Write a link-based implementation of a queue that uses a circular linked chain to represent the items in the queue. You will need a single tail pointer. When you are done, compare your implementation to the one given in this chapter that uses a linear linked chain with two external pointers. Which implementation is easier to write? Which is easier to understand? Which is more efficient?
2. Write an array-based implementation of a queue that uses a resizable, circular array to represent the items in the queue.
3. Revise the array-based implementation of a queue given in this chapter. Instead of counting the number of items in the queue, use a boolean variable isFul1 to distinguish between the full and empty conditions.
4. This chapter described another array-based implementation of a queue that uses no special data member-such as count or isFull (see the previous programming problem)-to distinguish between the full and empty conditions. In this implementation, you declare MAX_QUEUE + 1 locations for the array items, but use only MAX_QUEUE of them for queue items. You sacrifice one array location by making front the index of the location before the front of the queue. The queue is full if front equals (back +1 ) \% (MAX_QUEUE +1 ), but the queue is empty if front equals back. Implement this array-based approach.
5. Exercise 9 in the previous chapter defined the double-ended queue, or deque. Implement the ADT deque by using a circular array to contain the items in the deque.
6. Repeat the previous programming problem, but maintain the deque's entries in a linked chain.
7. Implement the ADT deque, as described in Exercise 9 of the previous chapter, as a derived class of ArrayQueue, as given in Listings 14-4 and 14-5.

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\section*{Overloaded \\ Operators and Friend Access}

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\(C_{++ \text {provides many operators, such as }+,-,=,==,>,<\text {, and } \ll \text {, that are defined for }}\) primitive data types. For instance, when working with integers, the symbol + represents the addition operation, \(=\) is used for the assignment operation, and \(==\) is used for equality comparisons. When we define new C++ data types using classes, these symbols do not have a definition that addresses how to add or compare two objects of the class. This C++ Interlude describes the tools that C++ provides you for extending the current operator definitions so that they can be applied to classes you have created.

\section*{C5.1 Overloaded Operators}

The standard arithmetic operators in C++ actually have multiple meanings. Although the addition operators in the expressions \(2+3\) and \(2.0+3.0\) appear to be the same, they in fact are not. Because integers such as 2 and 3 have internal representations that differ from floating-point numbers such as 2.0 and 3.0 , the algorithm to add two integers must differ from the algorithm to add two floating-point numbers. C++ could use two different symbols to designate integer addition and floating-point addition, but instead it uses only one symbol, + . The actual meaning of the + operator- that is, the type of addition it designates - is implied by the data type of its operands. An operator with more than one meaning is overloaded and is an example of a simple form of polymorphism.

An overloaded operator has more than one meaning

Two lists are equal if they have identical lengths and items

You saw earlier that you define new data types within \(\mathrm{C}++\) by using classes. Clients of such data types should be able to use them as naturally as the standard data types of the language. In particular, a client should be able to combine instances of a class with \(\mathrm{C}++\) operators in meaningful ways. To enable a particular operator to operate correctly on instances of a class, you typically must define a new meaning for the operator; that is, you must overload it.

Suppose that myList and yourList are two instances of LinkedList, as given in Listing 9-2 of Chapter 9, and you write
if (myList == yourList)
cout << "The lists are equa1. \n";
You must provide LinkedList with a definition of \(==\), since the compiler will not provide a default interpretation. To this end, let us state that the list myList is equal to the list yourList if
- myList and yourList have the same size, and
- Every item on myList is the same as the corresponding item on yourList

To overload an operator, you define an operator method whose name has the form
operatorsymbol
where symbol is the operator that you want to overload. For the == operator, you name the method operator \(==\) and declare one argument: the object that will appear on the right-hand side of the operator. The current object represents the object on the left-hand side of the operator. For LinkedList, you would add the declaration
bool operator \(==(\) const LinkedList<ItemType>\& rightHandSide) const;
to the class definition in the header file.
To understand this notation, realize that we may also have declared a client function isEqua1To that compares a parameter of type LinkedList with the current object to determine equality. The declaration for such a function would look like this:
bool isEqualTo(const LinkedList<ItemType>\& rightHandSide) const;
To use isEqualTo to compare the lists myList and yourList, you would write
```

if (myList.isEqualTo(yourList))

```
    cout << "The lists are equal. \n";

You can treat operator== exactly as you do isEqua1To; that is, you can write
if (myList.operator==(yourList))
cout << "The lists are equal. \(\backslash n\) ";
because operator== is simply a method name. However, you can also use the more natural shorthand notation myList == yourList, and the compiler will understand that you mean the operator== method.

The LinkedList implementation of this method is
```

template <class ItemType>

```
bool LinkedList<ItemType>::operator==(const
LinkedList<ItemType>\& rightHandSide) const
\{
bool isEqual = true; // Assume equal
// First check whether the number of items is the same
if (itemCount != rightHandSide.getLength())
isEqual = false;
```

    else
    { // Then compare items
        Node<ItemType>* leftSidePtr = headPtr;
        Node<ItemType>* rightSidePtr = rightHandSide.headPtr;
        while ((1eftSidePtr != nul1ptr) && (rightSidePtr != nul1ptr) && isEqual)
        {
            ItemType 1eftItem = 1eftSidePtr->getItem();
            ItemType rightItem = rightSidePtr->getItem();
            isEqual = (leftItem == rightItem);
            1eftSidePtr = 1eftSidePtr->getNext();
            rightSidePtr = rightSidePtr->getNext();
        } // end while
    } // end if
    return isEqual;
    } // end operator==

```

Note that this method depends on the \(==\) operator for items in the list. If these items are themselves instances of a class, that class must overload \(==\). You can overload the relational operators ( \(<,<=,>\), >=) in a similar manner.

Now suppose that you want all implementations of ListInterface, as given in Listing 8-1 of Chapter 8, to provide an implementation of the operator \(==\). You would need to add the declaration
virtual bool operator= \(=(\) const ListInterface<ItemType \(>\&\) rightHandSide ) const \(=0\);
to the interface ListInterface. Each implementation of ListInterface would then provide a definition of \(==\), allowing you to compare any two ListInterface objects. The challenge of making this comparison is that only ListInterface methods can be used for the list on the right-hand side, thereby affecting the efficiency of the equality check when comparing lists having two different implementations. For example, if we had an array-based list that we wished to compare to a linkbased list, we could not take advantage of the node structure of the underlying linked chain when we accessed the entries. Instead, we would need to use getEntry to access entries, but each call to getEntry would iterate through the list from its beginning until it reaches the desired entry.

The implementation of other relational operators is completed in a similar way.

\section*{C5.1.1 Overloading = for Assignment}

Overloading the assignment operator (=) is similar to overloading the equality operator, but it presents additional concerns. Once again, suppose that myList and yourList are two instances of LinkedList. If you place several items in the list yourList and then write
```

myList = yourList;

```
you would expect myList to be an exact copy of yourList. Without an overloaded assignment operator, however, you would get a shallow copy of yourList instead of a deep copy, as Figure 4-8 in Chapter 4 illustrates. Although a shallow copy might be sufficient for a statically allocated data structure, a deep copy is necessary for a dynamically allocated one such as the chain of linked nodes in the link-based implementation LinkedList. A shallow copy of a LinkedList object-yourListwould copy only the data members itemCount (the length of the list yourList) and headPtr (the pointer to yourList's first item). The items in the list would not be copied.

To provide an assignment operator for the class LinkedList, you would add the declaration

\footnotetext{
LinkedList<ItemType>\& operator=(const LinkedList<ItemType>\& rightHandSide);
}

Without an overloaded assignment operator, you get a shallow copy

The assignment operator must first deallocate myList, the object on the left-hand side

Make the assignment operator fail-safe by checking for a special case
to the class definition in LinkedList's header file. The argument rightHandSide represents the object to be copied-that is, the object that will appear on the right-hand side of the assignment operator. The method is not void, but instead returns a value to accommodate assignments such as
```

myList = yourList = theirList

```

As you will see, the method returns a reference to the invoking object.
In implementing this method, you must deal with a few subtleties. Suppose that myList and yourList each contain several items. If you write
```

myList = yourList;

```
what happens to the items that were in myList? You might not care, as long as myList ultimately contains a copy of the items in yourList. You should care, however, if you have a link-based list. Before you can copy the items in yourList to myList, you need to deallocate the nodes in myList. Failure to do so results in a memory leak-that is, memory that was allocated to myList is not returned to the system and is inaccessible. Thus, the assignment of yourList to myList must take these steps:
```

Deallocate memory assigned to myList
for (each item in yourList)
{
Allocate a new node for myList
Set the node's data portion to a copy of the item in yourList
}

```

The tasks here are like those that you implemented when you wrote the destructor and copy constructor for LinkedList. To deallocate memory currently used by LinkedList nodes, you can call clear. When copying the underlying linked chain of nodes, you can avoid redundancy with the copy constructor by defining a private or protected method, copyListNodes, that both the copy constructor and the assignment operator can call.

Now suppose you write
```

myList = myList;

```

Notice what our previous pseudocode tells you to do here: Deallocate myList and then make a copy of myList. After you have deallocated myList, there is nothing left to copy! Your implementation should test for this special case by asking whether the invoking object on the left side of the = operator is the same as the object on the right. If it is, your assignment operator should do nothing - it should be fail-safe.

You can make this test by comparing the addresses of the two list objects rather than by comparing the items on the lists. In \(\mathrm{C}++\), this is a pointer to the invoking object. Thus, you write
```

if (this != \&rightHandSide)

```
to compare the addresses of the objects on the left and right side of the \(=\). Notice the placement of the \& symbol; when used in this manner, it means "the address of." We would then read the above statement as "if the address of this object is not equal to the address of rightHandSide, then ...". In this textbook, this is the only time when it is necessary to use the \& symbol as the "address of" operator.

The link-based implementation of the overloaded assignment operator for LinkedList follows:
```

template <class ItemType>
LinkedList<ItemType>\& LinkedList<ItemType>: :operator=(const
LinkedList<ItemType>\& rightHandSide)
{
// Check for assignment to self
if (this != \&rightHandSide)

```
```

    {
        this->clear(); // Deal7ocate 7eft-hand side
        copyListNodes(rightHandSide); // Copy 7ist nodes
        itemCount = rightHandSide.itemCount; // Copy size of 7ist
    } // end if
    return *this;
    } // end operator=

```

The assignment operator = requires that we return an object of the same type as rightHandSide, so we must return a LinkedList object. The returned object is placed into the left-hand side of the assignment statement. Since our implementation of the operator = places a copy of rightHandSide into the current, or invoking, object on the left-hand side, we should return the current object, which is represented by *this.

Note: In \(\mathrm{C}++\), this is a pointer to the invoking object. Therefore, *this is the invoking object.

If you declare and initialize an object in the same statement, the compiler will invoke the copy constructor and not the = operator. For example, the declaration of myList in

LinkedList<string> myList = yourList;
constructs the object myList. This statement is equivalent to:
```

LinkedList<string> myList(yourList);

```

\section*{C5.1.2 Overloading + for Concatenation}

As we discussed in the previous section, the + operator is used for both integer and floating-point addition. In \(\mathrm{C}++\), the + operator is also used as a concatenation operator to append strings. This ability illustrates one of the concerns some programmers have about overloading operators-a symbol can represent several different operations. These operations not need be closely related to each other, so reading code that uses overloaded operators can be difficult.

Suppose you were asked to maintain code someone else wrote and found the following statement in a method:
```

theThing = myThing + yourThing;

```

Can you quickly determine whether the + performs scalar addition-adding two integers or floatingpoint numbers-or concatenates two strings? Without knowing the data type of the variables myThing and yourThing, you cannot determine what operation + performs. In a well-written program, most variables are declared close to their use, thus helping to clarify the statement, but this is not always the case.

When you choose to overload an operator, you should not change the meaning of the symbol. Your definition should extend the symbol's current operation only for use with a new data type-your class. When working with lists, the concatenation definition of the + symbol applies better than scalar addition, so let's see what is involved in appending one list onto another. Let's assume that theThing, myThing, and yourThing in the previous statement are list objects.

Link-based lists. If we choose to overload the + operator for the list-based implementation of the ADT list, we would add the following declaration to the class definition:

LinkedList<ItemType>\&
operator+(const LinkedList<ItemType>\& rightHandSide) const;

The implementation of this concatenation operator has the following logic:
```

concatList = a new, empty instance of LinkedList
concatList.itemCount = itemCount + rightHandSide.itemCount
1 eftChain = a copy of the chain of nodes in this list
rightChain = a copy of the chain of nodes in the list rightHandSide
Set the last node of 7eftChain to point to the first node of rightChain
concatList.headPtr = 1eftChain.headPtr
return concatList

```

Now, the statement
```

theThing = myThing + yourThing;

```
assigns the returned list concatList to theThing using the assignment operator defined in the previous section.

After implementing the + operator, we should also implement the \(+=\) operator, since it provides a shorthand for self-assignment using the + operator. This can be a more efficient method, as we do not need to make a copy of our list before appending the copy of rightHandSide.

Array-based lists. An implementation of the + operator for array-based lists has different challenges. For example, which of the following should the capacity of the resulting list be?
- The sum of the capacities of the two lists that are operands
- The capacity of the first (left-hand) list
- The number of actual items in the combined lists
- Twice the number of actual items in the combined lists to allow for growth

After the new ArrayList object has been created, we can copy the entries from the first list into the first array elements and then place the entries in the second list into the elements that follow.

\section*{Note: Some guidelines for overloading operators}
- Overloading an operator is the same as writing a method whose name is

\section*{operatorsymbol}
where symbol is the operator that you want to overload.
- Overloaded operators are typically placed in the public section of the class declaration.
- When you decide to overload an operator, be sure that the new operation closely matches how the operator is already used. If that is not possible, consider creating a method whose name is not operatorsymbol to perform the operation.
- Commonly overloaded operators are the assignment \((=)\), equality ( \(==\) and \(!=\) ), and relational ( \(<,<=,>,>=\) ) operators.
- You can overload any C++ operator except
.*:: ?: sizeof
- You cannot define new operators by overloading symbols that are not already C++ operators.
- You cannot change the standard precedence of a \(\mathrm{C}++\) operator.
- You cannot change the number of arguments for an overloaded operator, since they represent the operator's operands.
- At least one operand of an overloaded operator must be an instance of a class.

\section*{C5.2 Friend Access and Overloading <<}

In \(\mathrm{C}++\), a class can provide additional access to its private and protected parts by declaring other functions and classes as friends. Declaring a nonmember function as a friend to a class allows that function to access all of the private and protected members of the class. Let's look at an example in which we need to declare a function as a friend of our class.

The output stream operator \(\ll\) is used to output a stream of data to either the display or a file. We commonly stream integers, floating-point numbers, and strings to the display or a file, using \(\ll\). Occasionally, you may even have streamed a pointer to the display to help with troubleshooting or just to see what happened. \(\mathrm{C}++\) has defined the stream operator's action for each of those data types.

Trying to stream a list to the display using statements such as
```

LinkedList<string> myList;
cout << myList;

```
results in a syntax error because there is no << operation defined for the LinkedList class. Notice that the syntax of the previous cout statement differs from the syntax of previous binary operators we have overloaded. The expression cout \(\ll\) myList is shorthand for
```

cout.operator<<(myList)

```

Thus, when we define \(\ll\) for our LinkedList class, we are actually overloading the operator<< method in the ostream class of which cout is an instance. Therefore, we need to define an operation for \(\ll\) in a different manner than the way we overloaded previous operators.

Before discussing how to overload the output stream operator, let's consider what our list should look like if it is displayed or written to a file. If our list contained the four items Ace, Jack, Queen, and King, we could display them on a single line

Ace Jack Queen King
or on multiple lines with a position label:

> 1 Ace
> 2 Jack
> 3 Queen
> 4 King

For our example, let's choose the second form of output. We could use public methods to display the list in this form, but we would need to call getEntry-which iterates through the list from the beginning-to display each item. It is more efficient to access the data directly. Thus, in defining \(\ll\) for our LinkedList class, we should grant the operator<< method in the ostream class access to the private and protected data of LinkedList.

We grant this access by making the operator \(\ll\) method a friend of LinkedList by placing the following two statements in its declaration in the header file:
```

template<class friendItemType>
friend ostream\& operator<<(ostream\& outputStream,
const LinkedList< friendItemType >\& outputList);

```

The keyword friend indicates that this method has access to private and protected data members and methods of our class as well as to those that are public. Notice that the method returns an ostream object and that the first parameter is also an ostream object. This is because we are overloading an operator method from the ostream class. Also, since the operator is not part of the LinkedList class, we cannot use the same template type ItemType in the declaration; we must declare another template type-friendItemType in our example.

Functions and classes can be friends of a class

A friend method

Friend classes

The class
LinkedList is a friend of ListNode

A friend of a base class is not a friend of a derived class

We again must use the different template type when we implement the method:
```

template<class friendItemType>
ostream\& operator<< (ostream\& outStream,
const LinkedList<friendItemType>\& outputList)
{
int position = 1;
Node<friendItemType>* curPtr = outputList.headPtr;
while (curPtr != nul1ptr)
{
outStream << position << "\t" << curPtr->getItem() << endl;
curPtr = curPtr->getNext();
position++;
} // end while
return outStream;
} // end operator<<

```

Because the operator \(\ll\) method is already part of the ostream class, we don't need to include it in our namespace by writing LinkedList<ItemType>: : before the method name.

Note that the method accesses the private data member headPtr of the class LinkedList. Although this access violates the principle of information hiding, it is done in a controlled fashion: The class LinkedList explicitly grants this access by declaring the method as a friend.

We can extend the concept of friend methods to create friend classes. Granting a class access as a friend of another class allows all of the methods contained in the friend class to have access to the private and protected parts of the granting class. This feature is particularly useful when one class is used in the implementation of another class. For example, we have seen link-based implementations of ADTs that have a node declared as a separate class. For example, the LinkedList class uses the class Node. An alternative is to implement the node as a class whose data members and methods are private and to declare the LinkedList class as a friend class. For example, you could define a node for the ADT list as follows:
```

temp1ate<class ItemType>
class ListNode // A node on the list
{
private:
ItemType item; // A data item on the list
Node<ItemType> *next; // Pointer to next node
Node();
Node(const ItemType\& nodeItem, Node<ItemType>* nextNode);
// Friend class - can access private parts
friend class LinkedList<ItemType>;
}; // end ListNode

```

The class LinkedList has the same access privileges to the node's data members item and next as ListNode does. Friends of a base class do not have access to the private and protected parts added by a derived class-only to the private and protected members that appear in the base class.

\section*{Note: Some key points about friends}
- Friend methods can access the private and protected parts of the class.
- Friend methods are not members of the class.
- When a class is declared as a friend of a class \(C\), all of its methods have access to the private and protected parts of the class \(C\).
- Friendship is not inherited. The private and protected members declared in a derived class are not accessible by friends of the base class.

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\section*{Chapter}

\section*{Trees}

\section*{Contents}

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The previous chapters discussed ADTs whose operations fit into at least one of these general categories:
- Operations that insert data into a data collection
- Operations that remove data from a data collection
- Operations that ask questions about the data in a data collection

The ADTs list, stack, and queue are all position oriented, and their operations have the form
- Insert a data item into the \(i^{\text {th }}\) position of a data collection.
- Remove a data item from the \(i^{\text {th }}\) position of a data collection.
- Ask a question about the data item in the \(i^{\text {th }}\) position of a data collection.

As you have seen, the ADT list allows the value of \(i\) to range from 1 to the number of items in the list, while the ADTs stack and queue are more restrictive. For example, the operations of the ADT stack are limited to inserting into, removing from, and asking a question about one end - the top-of the stack. Thus, although they differ with respect to the flexibility of their operations, lists, stacks, and queues manage an association between data items and positions.

As Section 13.5 of Chapter 13 discussed, the ADT sorted list is value oriented. Its operations are of the form
- Insert a data item containing the value \(x\).
- Remove a data item containing the value \(x\).
- Ask a question about a data item containing the value \(x\).

Although these operations, like position-oriented operations, fit into the three general categories of operations listed earlier - they insert data, remove data, and ask questions about data-they are based on the values of data items instead of their positions.

Lists, stacks, and queues are linear in their organization of data, in that items are one after another. In this chapter, we organize data in a nonlinear, hierarchical form, whereby an item can have more than one immediate successor. We present two major ADTs: the binary tree and the binary search tree. As you will see, the binary tree is a position-oriented ADT, but it is not linear. Thus, you will not reference items in a binary tree by using a position number. Our discussion of the ADT binary tree provides an important background for the more useful binary search tree, which is a value-oriented ADT. Although a binary search tree is also not linear, it has operations similar to those of a sorted list, which is linear.

\subsection*{15.1 Terminology}

Trees are hierarchical

You use trees to represent relationships. Previous chapters informally used tree diagrams to represent the relationships between the calls of a recursive algorithm. For example, the diagram of the rabbit algorithm's recursive calls in Figure 2-19 of Chapter 2 is actually a tree. Each call to rabbit is represented by a box, or node, or vertex, in the tree. The lines between the nodes (boxes) are called edges. For this tree, the edges indicate recursive calls. For example, the edges from \(\operatorname{rabbit}(7)\) to \(\operatorname{rabbit}(6)\) and \(\operatorname{rabbit(5)}\) indicate that subproblem \(\operatorname{rabbit(7)}\) makes calls to \(\operatorname{rabbit}(6)\) and \(\operatorname{rabbit(5).}\)

All trees are hierarchical in nature. Intuitively, "hierarchical" means that a "parent-child" relationship exists between the nodes in the tree. If an edge is between node \(n\) and node \(m\), and node \(n\) is above node \(m\) in the tree, then \(n\) is the parent of \(m\), and \(m\) is a child of \(n\). In the tree in Figure 15-1a, nodes \(B\) and \(C\) are children of node \(A\). Children of the same parent-for example, \(B\) and \(C\)-are called siblings. The three children \(D, E\), and \(F\) of node \(B\), for example, are siblings. The leftmost child \(D\) is called the oldest child, or first child, of \(B\).

FIGURE 15-1 \(\quad\) (a) A tree; (b) a subtree of the tree in part \(a\)
(a)

(b)


Each node in a tree has at most one parent, and exactly one node - called the root of the treehas no parent. Node \(A\) is the root of the tree in Figure 15-1a. A node that has no children is called a leaf of the tree. The leaves of the tree are \(C, D, E\), and \(F\).

The parent-child relationship between the nodes is generalized to the relationships ancestor and descendant. In Figure 15-1 a, \(A\) is an ancestor of \(D\), and thus \(D\) is a descendant of \(A\). Not all nodes are related by the ancestor or descendant relationship: \(B\) and \(C\), for instance, are not so related. However, the root of any tree is an ancestor of every node in that tree. A subtree in a tree is any node in the tree together with all of its descendants. A subtree of a node \(n\) is a subtree rooted at a child of \(n\). For example, Figure \(15-1 \mathrm{~b}\) shows a subtree of the tree in Figure \(15-1 \mathrm{a}\). This subtree has \(B\) as its root and is a subtree of the node \(A\).

Because trees are hierarchical in nature, you can use them to represent information that itself is hierarchical in nature-for example, organization charts and family trees, as Figure 15-2 depicts. It may be disconcerting to discover, however, that the nodes in the family tree in Figure 15-2b that represent Bart's parents (Homer and Marge) are the children of the node that represents Bart. That is, the nodes that represent Bart's ancestors in a family tree are the descendants of Bart's node in one of our trees!

FIGURE 15-2 (a) An organization chart; (b) a family tree


Formal definition of a binary tree

Intuitive definition of a binary tree

\subsection*{15.1.1 Kinds of Trees}

Formally, a general tree is a set \(T\) of one or more nodes such that \(T\) is partitioned into disjoint subsets:
- A single node \(r\), the root
- Sets that are general trees, called subtrees of \(r\)

Thus, the trees in Figures 15-1 and 15-2a are general trees.
An \(\boldsymbol{n}\)-ary tree is a set \(T\) of nodes that is either empty or partitioned into disjoint subsets:
- A single node \(r\), the root
- \(n\) possibly empty sets that are \(n\)-ary subtrees of \(r\)

Each node can have no more than \(n\) children. The tree in Figure \(15-1\) is an \(n\)-ary tree with \(n=3\). An \(n\)-ary tree is not a special kind of general tree, because an \(n\)-ary tree can be empty, whereas a general tree cannot.

If an \(n\)-ary tree has the restriction that every node has at most two children, it is a binary tree. The primary focus of this chapter will be on binary trees. Formally, a binary tree is a set \(T\) of nodes that is either empty or partitioned into disjoint subsets:
- A single node \(r\), the root
- Two possibly empty sets that are binary trees, called left and right subtrees of \(r\)

The trees in Figures 2-19 and 15-2b are binary trees. Notice that each node in a binary tree has no more than two children.

The following intuitive restatement of the definition of a binary tree is useful:
\(T\) is a binary tree if either
- Thas no nodes, or
- \(T\) is of the form

where \(r\) is a node and \(T_{L}\) and \(T_{R}\) are both binary trees.
Notice that the formal definition agrees with this intuitive one: If \(r\) is the root of \(T\), then the binary tree \(T_{L}\) is the left subtree of node \(r\) and \(T_{R}\) is the right subtree of node \(r\). If \(T_{L}\) is not empty, its root is the left child of \(r\), and if \(T_{R}\) is not empty, its root is the right child of \(r\). Notice that if both subtrees of a node are empty, that node is a leaf.
Example: Algebraic expressions. As an example of how you can use a binary tree to represent data in a hierarchical form, consider Figure 15-3. The binary trees in this figure represent algebraic expressions that involve the binary operators,,\(+- \times\), and /. To represent an expression such as \(a-b\), you place the operator in the root node and the operands \(a\) and \(b\) into left and right children, respectively, of the root. (See Figure 15-3a.) Figure 15-3b represents the expression \(a-b / c\); a subtree represents the subexpression \(b / c\). A similar situation exists in Figure 15-3c, which represents \((a-b) \times c\). The leaves of these trees contain the expression's operands, while other tree nodes contain the operators. Parentheses do not appear in these trees. The binary tree provides a hierarchy for the operations-that is, the tree specifies an unambiguous order for evaluating an expression. Operators lower in the tree are evaluated first.

Note: Although various kinds of trees are possible, they are either general trees or \(n\)-ary trees.

FIGURE 15-3 Binary trees that represent algebraic expressions

(a)
\(a-b / c\)

(b)
\((a-b) \times c\)

The nodes of a tree typically contain values. A binary search tree is a binary tree that is, in a sense, sorted according to the values in its nodes. For each node \(n\), a binary search tree satisfies the following three properties:
- \(n\) 's value is greater than all values in its left subtree \(T_{L}\).
- \(n\) 's value is less than all values in its right subtree \(T_{R}\).
- Both \(T_{L}\) and \(T_{R}\) are binary search trees.

Figure 15-4 is an example of a binary search tree. As its name suggests, a binary search tree organizes data in a way that facilitates searching it for a particular data item. Chapter 16 will discuss binary search trees in detail, and Chapter 19 will present several other trees that improve searching data.

FIGURE 15-4 A binary search tree of names


Question 1 What kind of tree is the tree in Figure 15-1a?
Question 2 Repeat the previous question, but use the tree in Figure 15-3c instead.
Question 3 Given the tree in Figure 15-3c, what node or nodes are
a. Ancestors of \(b\) ?
b. Descendants of \(\times\) ?
c. Leaves?

Properties of a binary search tree terms of levels

Question 4 Given the tree in Figure 15-4, what node or nodes are
a. The tree's root?
b. Parents?
c. Children of the parents in part \(b\) of this question?
d. Siblings?

\subsection*{15.1.2 The Height of Trees}

Trees come in many shapes. For example, although the binary trees in Figure 15-5 all contain the same nodes, their structures are quite different. Although each of these trees has seven nodes, some are "taller" than others. The height of a tree is the number of nodes on the longest path from the root to a leaf. For example, the trees in Figure 15-5 have respective heights of 3, 5, 7, and 7. Many people's intuitive notion of height would lead them to say that these trees have heights of \(2,4,6\), and 6. Indeed, many authors define height to agree with this intuition. However, the definition of height used in this book leads to a cleaner statement of many algorithms and properties of trees.

There are other equivalent ways to define the height of a tree \(T\). One way uses the following definition of the level of a node \(\boldsymbol{n}\) :
- If \(n\) is the root of \(T\), it is at level 1 .
- If \(n\) is not the root of \(T\), its level is 1 greater than the level of its parent.

For example, in Figure \(15-5\) a, node \(A\) is at level 1 , node \(B\) is at level 2, and node \(D\) is at level 3 .
The height of a tree \(T\) in terms of the levels of its nodes is defined as follows:
- If \(T\) is empty, its height is 0 .
- If \(T\) is not empty, its height is equal to the maximum level of its nodes.

By applying this definition to the trees in Figure 15-5, you will find that their heights are, respectively, 3,5 , and 7 , as was stated earlier.

FIGURE 15-5 Binary trees with the same nodes but different heights
(a)

(b)

(c)

(d)


For binary trees, it is often convenient to use an equivalent recursive definition of height:

Recursive definition of height
- If \(T\) is empty, its height is 0 .
- If \(T\) is a nonempty binary tree, then because \(T\) is of the form

the height of \(T\) is 1 greater than the height of its root's taller subtree; that is, \(\operatorname{height}(T)=1+\max \left\{\operatorname{height}\left(T_{\mathrm{L}}\right), \operatorname{height}\left(T_{\mathrm{R}}\right)\right\}\)

Question 5 What are the levels of all nodes in the trees in parts \(b, c\), and \(d\) of Figure 15-5?
Question 6 What is the height of the tree in Figure 15-4?

\subsection*{15.1.3 Full, Complete, and Balanced Binary Trees}

In a full binary tree of height \(h\), all nodes that are at a level less than \(h\) have two children each. Figure 15-6 depicts a full binary tree of height 3 . Each node in a full binary tree has left and right subtrees of the same height. Among binary trees of height \(h\), a full binary tree has as many leaves as possible, and they all are at level \(h\). Intuitively, a full binary tree has no missing nodes.

FIGURE 15-6 A full binary tree of height 3


When proving properties about full binary trees-such as how many nodes they have-the following recursive definition of a full binary tree is convenient:
- If \(T\) is empty, \(T\) is a full binary tree of height 0 .
- If \(T\) is not empty and has height \(h>0, T\) is a full binary tree if its root's subtrees are both full binary trees of height \(h-1\).

This definition closely reflects the recursive nature of a binary tree.
A complete binary tree of height \(h\) is a binary tree that is full down to level \(h-1\), with level \(h\) filled in from left to right, as Figure 15-7 illustrates. More formally, a binary tree \(T\) of height \(h\) is complete if

FIGURE 15-7 A complete binary tree


A complete binary tree

Full binary trees are complete

Complete binary trees are balanced
1. All nodes at level \(h-2\) and above have two children each, and
2. When a node at level \(h-1\) has children, all nodes to its left at the same level have two children each, and
3. When a node at level \(h-1\) has one child, it is a left child

Parts 2 and 3 of this definition formalize the requirement that level \(h\) be filled in from left to right. Note that a full binary tree is complete.

Finally, a binary tree is height balanced, or simply balanced, if the height of any node's right subtree differs from the height of the node's left subtree by no more than 1. The binary trees in Figures 15-7 and 15-5a are balanced, but the other trees in Figure 15-5 are not balanced. A complete binary tree is balanced.

\section*{Note: Summary of tree terminology}
\begin{tabular}{|c|c|}
\hline General tree & A set of one or more nodes, partitioned into a root node and subsets that are general subtrees of the root. \\
\hline Parent of node \(n\) & The node directly above node \(n\) in the tree. \\
\hline Child of node \(n\) & A node directly below node \(n\) in the tree. \\
\hline Root & The only node in the tree with no parent. \\
\hline Leaf & A node with no children. \\
\hline Height & The number of nodes on the longest path from the root to a leaf. \\
\hline Siblings & Nodes with a common parent. \\
\hline Ancestor of node \(n\) & A node on the path from the root to \(n\). \\
\hline Descendant of node \(n\) & A node on a path from \(n\) to a leaf. \\
\hline Subtree of node \(n\) & A tree that consists of a child (if any) of \(n\) and the child's descendants. \\
\hline Left (right) child of node \(n\) & A node directly below and to the left (right) of node \(n\) in a binary tree. \\
\hline Left (right) subtree of node \(n\) & In a binary tree, the left (right) child (if any) of node \(n\) plus its descendants. \\
\hline \(n\)-ary tree & A set of nodes that is either empty or partitioned into a root node and at most \(n\) subsets that are \(n\)-ary subtrees of the root. Each node has at most \(n\) children. \\
\hline Binary tree & A set of nodes that is either empty or partitioned into a root node and one or two subsets that are binary subtrees of the root. Each node has at most two children, the left child and the right child. \\
\hline Binary search tree & A binary tree in which the value in any node \(n\) is greater than the value in every node in \(n\) 's left subtree but less than the value in every node in \(n\) 's right subtree. \\
\hline
\end{tabular}
\begin{tabular}{ll} 
Full binary tree & \begin{tabular}{l} 
A binary tree of height \(h\) with no missing nodes. All \\
leaves are at level \(h\), and all other nodes each have \\
two children.
\end{tabular} \\
Complete binary tree & \begin{tabular}{l} 
A binary tree of height \(h\) that is full to level \(h-1\) and \\
has level \(h\) filled in from left to right.
\end{tabular} \\
Balanced binary tree & \begin{tabular}{l} 
A binary tree in which the left and right subtrees of \\
any node have heights that differ by at most 1.
\end{tabular}
\end{tabular}

\subsection*{15.1.4 The Maximum and Minimum Heights of a Binary Tree}

You can maximize the height of a binary tree with \(n\) nodes simply by giving each internal node (nonleaf ) exactly one child, as shown earlier in parts \(c\) and \(d\) of Figure 15-5. This process will result in a tree of height \(n\). The \(n\)-node tree with height \(n\) in Figure 15-5d strikingly resembles a chain of linked nodes.

Note: The maximum height of an \(n\)-node binary tree is \(n\).

A minimum-height binary tree with \(n\) nodes is a bit more difficult to obtain. As a first step, consider the number of nodes that a binary tree with a given height \(h\) can have. For example, if \(h=3\), the possible binary trees include those in Figure 15-8. Thus, binary trees of height 3 can have between three and seven nodes. In addition, Figure \(15-8\) shows that 3 is the minimum height for a binary tree with four, five, six, or seven nodes. Similarly, binary trees with more than seven nodes require a height greater than 3 .

Intuitively, to minimize the height of a binary tree given \(n\) nodes, you must fill each level of the tree as completely as possible. A complete tree meets this requirement, although it does not matter whether the nodes on the last level are filled from left to right. In fact, except for the tree in part \(a\), the trees in Figure 15-8 are complete trees. If a complete binary tree of a given height \(h\) is to have the maximum possible number of nodes, it should be full, as in part \(e\) of the figure.

FIGURE 15-8 Binary trees of height 3

Except for the last level, each level of a minimum-height binary tree must contain as many nodes as possible
(a)


(b)

(c)
(d)


(e)

FIGURE 15-9 Counting the nodes in a full binary tree of height \(h\)


Figure 15-9 counts the nodes of a full tree by level and demonstrates the following facts, whose proof is left as an exercise:


\section*{Note: Facts about full binary trees}
- A full binary tree of height \(h \geq 0\) has \(2^{h}-1\) nodes.
- You cannot add nodes to a full binary tree without increasing its height.
- The maximum number of nodes that a binary tree of height \(h\) can have is \(2^{h}-1\).

We now can determine the minimum height of an \(n\)-node binary tree.


Note: The minimum height of a binary tree with \(n\) nodes is \(\left\lceil\log _{2}(n+1)\right\rceil .^{1}\)

To prove the previous fact, we let \(h\) be the smallest integer such that \(n \leq 2^{h}-1\) and establish the following assertions:
1. A binary tree whose height is \(\leq h-1\) has \(<n\) nodes.

We know that a binary tree of height \(h-1\) has at most \(2^{h-1}-1\) nodes. If it is possible that \(n \leq 2^{h-1}-1\), then \(h\) is not the smallest integer such that \(n \leq 2^{h}-1\). Therefore, \(n\) must be greater than \(2^{h-1}-1\) or, equivalently, \(2^{h-1}-1<n\). Because a binary tree of height \(h-1\) has at most \(2^{h-1}-1\) nodes, it must have fewer than \(n\) nodes.
2. An n-node complete binary tree whose height is \(h\) exists.

Consider a full binary tree of height \(h-1\). It must have \(2^{h-1}-1\) nodes. As you just saw, \(n>2^{h-1}-1\) because \(h\) was selected so that \(n \leq 2^{h}-1\). You can thus add nodes to the full tree

from left to right until you have \(n\) nodes, as Figure 15-10 illustrates. Because \(n \leq 2^{h}-1\) and a binary tree of height \(h\) cannot have more than \(2^{h}-1\) nodes, you will reach \(n\) nodes by the time level \(h\) is filled up.
3. The minimum height of a binary tree with \(n\) nodes is the smallest integer \(h\) such that \(n \leq 2^{h}-1\).
If \(h\) is the smallest integer such that \(n \leq 2^{h}-1\), and if a binary tree has height \(\leq h-1\), then by fact 1 it has fewer than \(n\) nodes. Because by fact 2 there is a binary tree of height \(h\) that has exactly \(n\) nodes, \(h\) must be as small as possible.

The previous discussion implies that
\[
\begin{aligned}
& 2^{h-1}-1<n \leq 2^{h}-1 \\
& 2^{h-1}<n+1 \leq 2^{h} \\
& h-1<\log _{2}(n+1) \leq h
\end{aligned}
\]

Now \(h\) is either \(\log _{2}(n+1)\) or satisfies the inequality \(h-1<\log _{2}(n+1)<h\). In the latter case, \(\log _{2}(n+1)\) cannot be an integer, so we round it up to get \(h\). In either case, we have that \(h\) is \(\left\lceil\log _{2}(n+1)\right\rceil\). Thus, we have established that \(\left\lceil\log _{2}(n+1)\right\rceil\) is the minimum height of a binary tree with \(n\) nodes.

Note: Complete trees and full trees with \(n\) nodes have heights of \(\left\lceil\log _{2}(n+1)\right\rceil\), which, as you just saw, is the theoretical minimum.

Complete trees and full trees have minimum height

\section*{Question 7 Consider the binary trees in Figure 15-8.}
a. Which are complete?
b. Which are full?
c. Which are balanced?
d. Which have minimum height?
e. Which have maximum height?

\subsection*{15.2 The ADT Binary Tree}

As an abstract data type, the binary tree has operations that add and remove nodes, set or retrieve the data in the root of the tree, and test whether the tree is empty. By using these basic operations, you can build a binary tree.

Traversal operations that visit every node in a binary tree are typical. Visiting a node means "doing something with or to" the node. Chapter 4 introduced the concept of traversal for a linear

The general form of a recursive traversal algorithm

Three choices for when to visit the root

Traversal is \(\mathrm{O}(n)\)

Preorder traversal
chain of linked nodes: Beginning with the chain's first node, you visit each node sequentially until you reach the end of the chain. Traversal of a binary tree, however, can visit the tree's nodes in one of several different orders. We examine the details of the traversal operations next.

\subsection*{15.2.1 Traversals of a Binary Tree}

A traversal algorithm for a binary tree visits each node in the tree. While visiting a node, you do something with or to the node, such as display or modify the contents of the node. For the purpose of this discussion, assume that visiting a node simply means displaying the data portion of the node.

With the recursive definition of a binary tree in mind, you can construct a recursive traversal algorithm as follows. According to the definition, the binary tree \(T\) is either empty or is of the form


If \(T\) is empty, the traversal algorithm takes no action-an empty tree is the base case. If \(T\) is not empty, the traversal algorithm must perform three tasks: It must display the data in the root \(r\), and it must traverse the two subtrees \(T_{L}\) and \(T_{R}\), each of which is a binary tree smaller than \(T\). We can summarize this logic using the following pseudocode:
```

if (T is not empty)
{
Display the data in T's root
Traverse T's left subtree
Traverse T's right subtree
}

```

Although we arbitrarily visited the tree's root before traversing its subtrees, the algorithm actually has three choices of when to visit \(r\) : It can visit \(r\)
- Before it traverses both of \(r\) 's subtrees-as we just did
- After it has traversed \(r\) 's left subtree \(T_{L}\) but before it traverses \(r\) 's right subtree \(T_{R}\)
- After it has traversed both of \(r\) 's subtrees

These choices result in preorder, inorder, and postorder traversals, respectively. Figure 15-11 shows the results of these traversals for a given binary tree.

Each of these traversals visits every node in a binary tree exactly once. Thus, \(n\) visits occur for a tree of \(n\) nodes. Each visit performs the same operations on each node, independently of \(n\), so it must be \(\mathrm{O}(1)\). Thus, each traversal is \(\mathrm{O}(n)\).

Preorder traversal. The recursive preorder traversal algorithm is as follows:
// Traverses the given binary tree in preorder.
// Assumes that "visit a node" means to process the node's data item.
preorder(binTree: BinaryTree): void
if (binTree is not empty)
\{
Visit the root of binTree preorder(Left subtree of binTree's root)
preorder (Right subtree of binTree's root)

FIGURE 15-11 Three traversals of a binary tree

(Numbers beside nodes indicate traversal order.)

The preorder traversal of the tree in Figure 15-11a visits the nodes in this order: 60, 20, 10, 40, \(30,50,70\). If you apply preorder traversal to a binary tree that represents an algebraic expression, such as any tree in Figure 15-3, and display the nodes as you visit them, you will obtain the prefix form of the expression. \({ }^{2}\)
Inorder traversal. The recursive inorder traversal algorithm is as follows:
```

// Traverses the given binary tree in inorder.
// Assumes that "visit a node" means to process the node's data item.
inorder(binTree: BinaryTree): void
if (binTree is not empty)
{
inorder(Left subtree of binTree's root)
Visit the root of binTree
inorder(Right subtree of binTree's root)
}

```

The result of the inorder traversal of the tree in Figure \(15-11 \mathrm{~b}\) is \(10,20,30,40,50,60,70\). If you apply inorder traversal to a binary search tree, you will visit the nodes in order according to their data values. Such is the case for the tree in Figure 15-11b.

Postorder traversal. Finally, the recursive postorder traversal algorithm is as follows:
```

// Traverses the given binary tree in postorder.
// Assumes that "visit a node" means to process the node's data item.
postorder(binTree: BinaryTree): void
if (binTree is not empty)
{
postorder(Left subtree of binTree's root)
postorder(Right subtree of binTree's root)
Visit the root of binTree
}

```

2 The prefix expressions are (a) \(-a b\); (b) \(-a / b c\); (c) \(\times-a b c\).

A client-defined function, which the traversal methods call, defines the meaning of "visit"

The result of the postorder traversal of the tree in Figure 15-11c is 10, 30, 50, 40, 20, 70, 60. If you apply postorder traversal to a binary tree that represents an algebraic expression, such as any tree in Figure 15-3, and display the nodes as you visit them, you will obtain the postfix form of the expression. \({ }^{3}\)

)
Note: Although traversal means to visit each item in the ADT, traversal can be more difficult than you might imagine if you do more than simply display each item when you visit it. For example, you might copy the item into another data structure or even alter it. The details of traversal are thus quite application dependent, which makes traversal a difficult operation to define within the framework of an ADT.

Visiting a node. You could have a different traversal operation for each desired task during a visit to a node, such as preorderTraverseAndDisplay, preorderTraverseAndCopy, and so on. Or the traversal methods for the ADT binary tree could call a function, which the client defines and passes as an argument, when it visits a node. We will take this latter approach, as it is more general.

For example, if the ADT binary tree has the operation preorderTraverse, you can invoke it for the binary tree binTree as follows:
```

bintree.preorderTraverse(display);

```
where display is a client function that displays the data passed to it as an argument. Suppose the binary tree binTree stores string objects. You would define the function display to display data of type string:
```

void display(string someItem)
{
cout << someItem << endl;
} // end display

```

Now the call bintree. preorderTraverse(display) will call display each time it visits a node in the tree.

Despite the fact that the ADT operation preorderTraverse calls a client-supplied function, the wall between the program and the implementation of the ADT has not been violated. Because display is on the client's side of the wall, the function can access the data only by using the ADT operations.

The next chapter will discuss the implementation details of the traversal operations.


Question 8 What are the preorder, inorder, and postorder traversals of the binary trees in parts \(a, b\), and \(c\) of Figure 15-5?

\subsection*{15.2.2 Binary Tree Operations}

The ADT binary tree has the following operations:


Note: ADT binary tree operations
- Test whether a binary tree is empty.
- Get the height of a binary tree.

\footnotetext{
3 The postfix expressions are (a) \(a b-\); (b) \(a b c /-\); (c) \(a b-c \times\).
}
- Get the number of nodes in a binary tree.
- Get the data in a binary tree's root.
- Set the data in a binary tree's root.
- Add a new node containing a given data item to a binary tree.
- Remove the node containing a given data item from a binary tree.
- Remove all nodes from a binary tree.
- Retrieve a specific entry in a binary tree.
- Test whether a binary tree contains a specific entry.
- Traverse the nodes in a binary tree in preorder, inorder, or postorder.

The following contract specifies these operations in more detail, and a UML diagram for a class of binary trees appears in Figure 15-12.

\section*{Abstract Data Type: Binary Tree}

\section*{Data}
- A finite number of objects in hierarchical order.

\section*{Operations}
\begin{tabular}{|c|c|}
\hline Pseudocode & Description \\
\hline isEmpty () & \begin{tabular}{l}
Task: Tests whether this binary tree is empty. \\
Input: None. \\
Output: True if the binary tree is empty; otherwise false.
\end{tabular} \\
\hline getHeight() & \begin{tabular}{l}
Task: Gets the height of this binary tree. Input: None. \\
Output: The height of the binary tree.
\end{tabular} \\
\hline getNumberOfNodes() & \begin{tabular}{l}
Task: Gets the number of nodes in this binary tree. Input: None. \\
Output: The number of nodes in the binary tree.
\end{tabular} \\
\hline getRootData() & Task: Gets the data that is in the root of this binary tree. Input: None. Assumes the tree is not empty. Output: The root's data. \\
\hline setRootData (newData) & \begin{tabular}{l}
Task: Replaces the data item in the root of this binary tree with newData, if the tree is not empty. However, if the tree is empty, inserts a new root node whose data item is newData into the tree. \\
Input: newData is the data item. \\
Output: None.
\end{tabular} \\
\hline add(newData) & \begin{tabular}{l}
Task: Adds a new node containing a given data item to this binary tree. Input: newData is the data item. \\
Output: True if the addition is successful, or false if not.
\end{tabular} \\
\hline remove(data) & Task: Removes the node containing the given data item from this binary tree. Input: data is the data item. Output: True if the removal is successful, or false if not. \\
\hline
\end{tabular}
\begin{tabular}{|c|c|}
\hline clear () & \begin{tabular}{l}
Task: Removes all nodes from this binary tree. Input: None. \\
Output: None. (The binary tree is empty.)
\end{tabular} \\
\hline getEntry (anEntry) & \begin{tabular}{l}
Task: Gets a specific entry in this binary tree. \\
Input: anEntry is the desired data item. \\
Output: The entry in the binary tree that matches anEntry. Throws an exception if the entry is not found.
\end{tabular} \\
\hline contains(data) & \begin{tabular}{l}
Task: Tests whether the given data item occurs in this binary tree. \\
Input: data is the data item. \\
Output: True if the binary tree contains the given data item, or false if not.
\end{tabular} \\
\hline preorderTraverse(visit) & \begin{tabular}{l}
Task: Traverses this binary tree in preorder and calls the function visit once for each node. \\
Input: visit is a client-defined function that performs an operation on or with the data in each visited node. \\
Output: None.
\end{tabular} \\
\hline inorderTraverse(visit) & \begin{tabular}{l}
Task: Traverses this binary tree in inorder and calls the function visit once for each node. \\
Input: visit is a client-defined function that performs an operation on or with the data in each visited node. \\
Output: None.
\end{tabular} \\
\hline postorderTraverse(visit) & \begin{tabular}{l}
Task: Traverses this binary tree in postorder and calls the function visit once for each node. \\
Input: visit is a client-defined function that performs an operation on or with the data in each visited node. \\
Output: None.
\end{tabular} \\
\hline
\end{tabular}

FIGURE 15-12 UML diagram for the class BinaryTree
\begin{tabular}{|l|}
\hline \multicolumn{1}{|c|}{ BinaryTree } \\
\hline +isEmpty(): boolean \\
+getHeight(): integer \\
+getNumberOfNodes(): integer \\
+getRootData(): ItemType \\
+setRootData(newData: ItemType): void \\
+add(newData: ItemType): boolean \\
+remove(data: ItemType): boolean \\
+clear(): void \\
+getEntry(anEntry: ItemType): ItemType \\
+contains(data: ItemType): boolean \\
+preorderTraverse(visit(item: ItemType) : void) : void \\
+inorderTraverse(visit(item: ItemType): void): void \\
+postorderTraverse(visit(item: ItemType): void): void
\end{tabular}

Example. You can use the add operation to build the binary tree in Figure 15-5a, where the node labels represent character data:
```

tree = a new empty binary tree
tree.add('A')
tree.add('B')
tree.add('C')
tree.add('D')
tree.add('E')
tree.add('F')
tree.add('G')

```

Given the specification of the add method, you have no reason to expect that the previous statements, in fact, produce the binary tree in the figure. Assuming that the operations are successful, you will know only that the binary tree has seven nodes containing the given data. Our specification of add is intentionally vague and gives the programmer who implements the class BinaryTree flexibility. It also makes BinaryTree more useful as a base class, as you will see in the next chapter.

\subsection*{15.2.3 An Interface Template for the ADT Binary Tree}

We formalize our specifications for the ADT binary tree by writing the interface template given in Listing 15-1.

\section*{LISTING 15-1 An interface template for the ADT binary tree}
```

/** Interface for the ADT binary tree.
@file BinaryTreeInterface.h */
\#ifndef _BINARY_TREE_INTERFACE
\#define _BINARY_TREE_INTERFACE
template<class ItemType>
class BinaryTreeInterface
{
public:
/** Tests whether this binary tree is empty.
@return True if the binary tree is empty, or false if not. */
virtual bool isEmpty() const = 0;
/** Gets the height of this binary tree.
@return The height of the binary tree. */
virtual int getHeight() const = 0;
/** Gets the number of nodes in this binary tree.
@return The number of nodes in the binary tree. */
virtual int getNumberOfNodes() const = 0;
/** Gets the data that is in the root of this binary tree.
@pre The binary tree is not empty.
@post The root's data has been returned, and the binary tree is unchanged.
@return The data in the root of the binary tree. */
virtual ItemType getRootData() const = 0;
/** Replaces the data item in the root of this binary tree
with the given data, if the tree is not empty. However, if

```

Using ADT binary tree operations to build a binary tree

This organization of data enables you to search a binary search tree for a particular data item, given its value instead of its position.

Example. Let's consider a binary search tree whose nodes contain people's names. Suppose that these objects are Alan, Bob, Elisa, Jane, Nancy, Tom, and Wendy. Figure 15-13 illustrates one binary search tree that we can form with these names.

FIGURE 15-13 A binary search tree of names


Many different binary search trees can contain the same data, however. For example, in addition to the tree in Figure 15-13, each tree in Figure 15-14 is a valid binary search tree for the names Alan, Bob, Elisa, Jane, Nancy, Tom, and Wendy. Although these trees have different shapes, each one satis-

Several different binary search trees are possible for the same data fies the requirements of a binary search tree.

FIGURE 15-14 Binary search trees with the same data as in Figure 15-13


\subsection*{15.3.1 Binary Search Tree Operations}

As an ADT, the binary search tree has operations that are like the operations for the ADTs you studied in previous chapters, in that they involve inserting, removing, and retrieving data. Unlike the position-oriented ADTs stack, list, and queue, but like the ADT sorted list, the insertion, removal, and retrieval operations are by value, not by position. On the other hand, the other operations - including
the traversal operations-that you just saw for a binary tree apply to a binary search tree without change, because a binary search tree is a binary tree.

The operations that define the ADT binary search tree are as follows:


For simplicity, we will insist that a binary search tree contain unique data. The following contract specifies the insertion and removal operations in more detail. The UML diagram for a class of binary search trees will be almost the same as the one for a class of binary trees, so we do not write one here.

\section*{Abstract Data Type: Binary Search Tree}

\section*{Data}
- A finite number of objects in hierarchical order.
\begin{tabular}{l} 
OPERATIONS \\
\multicolumn{1}{c}{ Pseudocode } \\
add (newEntry) \\
\(\quad\)\begin{tabular}{l} 
Task: Inserts newEntry into this binary search tree such that the properties of a binary \\
search tree are maintained.
\end{tabular} \\
Input: newEntry is the data item to be inserted. Assumes the entries in the tree are \\
distinct and differ from newEntry. \\
Output: True if the insertion is successful, or false if not. \\
Task: Removes the given entry from this binary search tree such that the properties of a \\
binary search tree are maintained. \\
Input: anEntry is the entry to remove. \\
Output: True if the removal is successful, or false if not.
\end{tabular}

The methods isEmpty, getHeight, getNumberOfNodes, getRootData, clear, getEntry, contains, preorderTraverse, inorderTraverse, and postorderTraverse have the same specifications as for a binary tree.


Example. If nameTree is any one of the binary search trees in Figures 15-13 and 15-14, and if their nodes contain strings, the statement
```

string name = nameTree.getEntry(nancy);

```
where nancy is a string variable containing "Nancy", locates the node containing "Nancy" and assigns its data to name. It doesn't matter which one of the four trees getEntry searches; the result will be the
same. The time it takes for getEntry to find nancy, however, does depend on the shape of the tree, as you will see.

If the string variable hal contains " Ha 7 " and you insert it into nameTree by invoking
```

nameTree.add(ha1)

```
you will be able to retrieve hal later and still be able to retrieve nancy. If you remove jane by using
```

nameTree.remove(jane)

```
you will still be able to retrieve the objects nancy and ha1. Finally, if displayName is a client function that displays its string argument,
```

nameTree.inorderTraverse(displayName)

```
will display in alphabetical order the names of the people that nameTree represents. Again, it does not matter which one of the four trees you traverse.

Question 9 Show that each tree in Figures 15-13 and 15-14 is a binary search tree.
Question 10 Show that the inorder traversals of each binary search tree in Figures 15-13 and 15-14 are the same.

Question 11 What are the preorder and postorder traversals of each binary search tree in Figures 15-13 and 15-14? Are the preorder traversals the same? Are the postorder traversals the same?

\subsection*{15.3.2 Searching a Binary Search Tree}

Consider again the binary search tree in Figure 15-13. Each node in the tree contains a person's name. Suppose that you want to locate Elisa in the binary search tree. The root node of the tree contains Jane, so if Elisa is present in the tree, it must be in Jane's left subtree, because the string "Elisa" is before the string "Jane" alphabetically. Because a binary search tree is recursive by nature, it is natural to formulate recursive algorithms for operations on the tree. Thus, you know that Jane's left subtree is also a binary search tree, so you use exactly the same strategy to search this subtree for Elisa. The root of this binary search tree contains Bob, and, because the string "Elisa" is greater than the string "Bob", Elisa must be in Bob's right subtree. That right subtree is also a binary search tree, and it happens that Elisa is in the root node of this tree. Thus, the search has located Elisa.

The following pseudocode summarizes this search strategy:
```

// Searches the binary search tree for a given target value.
search(bstTree: BinarySearchTree, target: ItemType)
if (bstTree is empty)
The desired item is not found
else if (target == data item in the root of bstTree)
The desired item is found
else if (target < data item in the root of bstTree)
search(Left subtree of bstTree, target)
else
search(Right subtree of bstTree, target)

```

As you will see, this search algorithm is the basis of the other operations on a binary search tree.
The shape of the tree in no way affects the validity of the search algorithm. The algorithm requires only that a tree be a binary search tree. However, the search algorithm works more efficiently on some

A search algorithm for a binary search tree

Use search to determine the insertion point
trees than on others. For example, with the tree in Figure 15-14b, search inspects every node before locating Wendy. In fact, this binary search tree really has the same structure as a link-based representation of a sorted list and offers no advantage in efficiency. In contrast, with the full tree in Figure 15-13, the search algorithm inspects only the nodes that contain the names Jane, Tom, and Wendy. These names are exactly the names that a binary search of the sorted array in Figure 15-15 would inspect.

FIGURE 15-15 An array of names in sorted order


The shape of a binary search tree affects the efficiency of its operations. The more balanced a binary search tree is, the farther it is from a linear structure and the closer the behavior of the search algorithm will be to a binary search of an array (and the farther it will be from the behavior of a linear search). Later in this chapter, you will learn more about how the shape of a binary search tree affects search's efficiency and how the insertion and removal operations affect this shape.

Question 12 Using the tree in Figure 15-14c, trace the algorithm that searches a binary search tree for
a. Elisa
b. Kyle

In each case, list the nodes in the order in which the search visits them.

\subsection*{15.3.3 Creating a Binary Search Tree}

Suppose that you want to insert a record for Frank into the binary search tree of Figure 15-13. As a first step, imagine that you instead want to search for the string "Frank". The search algorithm first searches the tree rooted at Jane, then the tree rooted at Bob, and then the tree rooted at Elisa. It then searches the tree rooted at the right child of Elisa. Because this tree is empty, as Figure 15-16 illustrates, the search algorithm has reached a base case and will terminate with the report that Frank is not present. What does it mean that search looked for Frank in the right subtree of Elisa? For one thing, it means that if Frank were the right child of Elisa, search would have found Frank there.

This observation indicates that a good place to insert Frank is as the right child of Elisa. Because Elisa has no right child, the insertion is simple, requiring only the addition of a new leaf node. More important, Frank belongs in this location-search will look for Frank here. Specifically, inserting Frank as the right child of Elisa will preserve the tree's binary search tree property. Because search, when searching for Frank, would follow a path that leads to the right child of Elisa, you are assured that Frank is in the proper relation to the names above it in the tree.


Note: Using search to determine where in the tree to insert a new name always leads to an easy insertion. No matter what new item you insert into the tree, search will always terminate at an empty subtree. Thus, search always tells you to insert the item as a new leaf.

FIGURE 15-16 Empty subtree where the search algorithm terminates when looking for Frank


Example. Let's create the binary search tree shown in Figure 15-13. The following pseudocode statements begin with an empty binary search tree and add entries one at time until we get the desired tree:
```

nameTree = a new, empty binary search tree
nameTree.insert("Jane")
nameTree.insert("Bob")
nameTree.insert("Alan")
nameTree.insert("Elisa")
nameTree.insert("Tom")
nameTree.insert("Nancy")
nameTree.insert("Wendy")

```

Once we insert Jane into the empty tree, Jane becomes the tree's root. Inserting Bob into the tree leaves the root untouched, and so the only binary search tree that is possible places Bob as the left child of the root. Why? Because "Bob" is less than "Jane". When we insert the next name, Alan, it becomes the left child of Bob. Note that each insertion creates a new leaf of the tree at that point in time. However, it might not remain a leaf in the final tree.

Question 13 Beginning with an empty binary search tree, what binary search tree is formed when you insert the following letters in the order given? J, N, B, A, W, E, T

Question 14 Arrange nodes that contain the letters A, C, E, F, L, V, and Z into two binary search trees: one that has maximum height and one that has minimum height.

\subsection*{15.3.4 Traversals of a Binary Search Tree}

The traversals of a binary search tree are the same as the traversals of a binary tree. The inorder traversal of a binary search tree, however, is of special note. Recall its algorithm from earlier in this chapter:
// Traverses the given binary tree in inorder.
// Assumes that "visit a node" means to process the node's data item.
inorder(binTree: BinaryTree): void
```

if (binTree is not empty)
{
inorder(Left subtree of binTree's root)
Visit the root of binTree
inorder(Right subtree of binTree's root)
}

```

Note: The inorder traversal of a binary search tree visits the tree's nodes in sorted search-key order.

Use inorder traversal to visit nodes of a binary search tree in search-key order

The maximum number of comparisons for a retrieval, insertion, or removal is the height of the tree

The proof of the previous fact is by induction on \(h\), the height of the tree. Let \(T\) represent the tree. Basis: \(h=0\). When \(T\) is empty, the algorithm does not visit any nodes. This is the proper sorted order for the zero names that are in the tree!
Inductive hypothesis: Assume that the theorem is true for all \(k, 0<k<h\). That is, assume for all \(k\) ( \(0<\) \(k<h\) ) that the inorder traversal visits the nodes in sorted search-key order. Inductive conclusion: You must show that the theorem is true for \(k=h>0\). \(T\) has the form


Because \(T\) is a binary search tree, all the entries in the left subtree \(T_{L}\) are less than the entry in the root \(r\), and all the entries in the right subtree \(T_{R}\) are greater than the entries in \(r\). The inorder algorithm will visit all the nodes in \(T_{L}\), then visit \(r\), and then visit all the nodes in \(T_{R}\). Thus, the only concern is that inorder visit the nodes within each of the subtrees \(T_{L}\) and \(T_{R}\) in the correct sorted order. But because \(T\) is a binary search tree of height \(h\), each subtree is a binary search tree of height less than \(h\). Therefore, by the inductive hypothesis, inorder visits the nodes in each subtree \(T_{L}\) and \(T_{R}\) in the correct sorted search-key order. The proof is now complete.

\subsection*{15.3.5 The Efficiency of Binary Search Tree Operations}

You have seen binary search trees in many shapes. For example, even though the binary search trees in Figures 15-13 and 15-14 have seven nodes each, they have different shapes and heights. You saw that to locate Wendy in Figure 15-14b, you would have to inspect all seven nodes, but you can locate Wendy in Figure 15-13 by inspecting only three nodes (Jane, Tom, and Wendy). Consider now the relationship between the height of a binary search tree and the efficiency of the retrieval, insertion, and removal operations.

Each of these operations compares a specified value \(v\) to the entries in the nodes along a path through the tree. This path always starts at the root of the tree and, at each node \(n\), follows the left or right branch, depending on the comparison of \(v\) to the value in \(n\). The path terminates at the node that contains \(v\) or, if \(v\) is not present, at an empty subtree. Thus, each retrieval, insertion, or removal operation requires a number of comparisons equal to the number of nodes along this path. This means that the maximum number of comparisons that each operation can require is the number of nodes on the longest path through the tree. In other words, the maximum number of comparisons that these operations can require is equal to the height of the binary search tree.

Note: The height of a binary tree
Recall that the height of an \(n\)-node binary tree ranges from \(\left\lceil\log _{2}(n+1)\right\rceil\) to \(n\).

If the height of the binary search tree is \(\log _{2}(n+1)\), the efficiency of its operations is \(\mathrm{O}(\log n)\). However, what will the height of a binary search tree actually be? The factor that determines the height of a binary search tree is the order in which you perform insertion and removal operations on the tree. Recall that, starting with an empty tree, if you insert names in the order Alan, Bob, Elisa, Jane, Nancy, Tom, Wendy, you would obtain a binary search tree of maximum height, as shown in Figure 15-14b. On the other hand, if you insert names in the order Jane, Bob, Tom, Alan, Elisa, Nancy, Wendy, you would obtain a binary search tree of minimum height, as shown in Figure 15-13.

Which of these situations should you expect to encounter in the course of a real application? It can be proven mathematically that if the insertion and removal operations occur in a random order, the height of the binary search tree will be quite close to \(\log _{2} n\). Thus, in this sense, the previous analysis is not unduly optimistic. However, in a real-world application, is it realistic to expect the insertion and removal operations to occur in random order? In many applications, the answer is yes. There are, however, applications in which this assumption would be dubious. For example, the person preparing the previous sequence of names for the insertion operations might well decide to "help you out" by arranging the names to be inserted into sorted order. This arrangement, as has been mentioned, would lead to a tree of maximum height. Thus, while in many applications you can expect the behavior of a binary search tree to be excellent, you should be wary of the possibility of poor performance due to some characteristic of a given application.

Is there anything you can do if you suspect that the operations might not occur in a random order? Similarly, is there anything you can do if you have an enormous number of items and need to ensure that the height of the tree is close to \(\log _{2} n\) ? Chapter 19 presents variations of the basic binary search tree that are guaranteed always to remain balanced and therefore be of minimum height.

Figure \(15-17\) summarizes the order of the retrieval, insertion, removal, and traversal operations for the ADT binary search tree.

FIGURE 15-17 The Big O for the retrieval, insertion, removal, and traversal operations of the ADT binary search tree
\begin{tabular}{ccccc}
\hline & \(\frac{O p e r a t i o n}{}\) & & Average case & Worst case \\
& Retrieval & & \(\mathrm{O}(\log n)\) & \\
\(\mathrm{O}(n)\) \\
& Insertion & & \(\mathrm{O}(\log n)\) & \\
\(\mathrm{O}(n)\) \\
& Removal & \(\mathrm{O}(\log n)\) & & \(\mathrm{O}(n)\) \\
& Traversal & \(\mathrm{O}(n)\) & \(\mathrm{O}(n)\) \\
\hline
\end{tabular}

\section*{Summary}
1. Binary trees provide a hierarchical organization of data, which is important in many applications.
2. Traversing a tree is a useful operation. Intuitively, traversing a tree means to visit every node in the tree. Because the meaning of "visit" is application dependent, you can pass a client-defined visit function to the traversal operation.
3. The binary search tree allows you to use a binary search-like algorithm to search for an item with a specified value.
4. Binary search trees come in many shapes. The height of a binary search tree with \(n\) nodes can range from a minimum
of \(\left\lceil\log _{2}(n+1)\right\rceil\) to a maximum of \(n\). The shape of a binary search tree determines the efficiency of its operations. The closer a binary search tree is to a balanced tree (and the farther it is from a linear structure), the closer the behavior of the search algorithm will be to a binary search (and the farther it will be from the behavior of a linear search).
5. An inorder traversal of a binary search tree visits the tree's nodes in sorted search-key order.

Insertion in sorted order produces a maximum-height binary search tree

Insertion in random order produces a near-minimumheight binary search tree

- Arer

\section*{Exercises}
1. Consider the tree in Figure 15-18. What node or nodes are
a. The tree's root?
b. Parents?
c. Children of the parents in part \(b\) ?
d. Siblings?
e. Ancestors of 50?
f. Descendants of 20?
g. Leaves?
2. What is the height of the tree in Figure 15-18?
3. Write preconditions and postconditions for the ADT binary search tree operations.

FIGURE 15-18 A tree for Exercises 1, 2, 6, 7, 11, and 23

4. What are the preorder, inorder, and postorder traversals of the binary tree in Figure 15-19?
5. Consider a method isLeaf that returns true if a binary tree is a one-node tree-that is, if it consists of only a leaf-and returns fal se otherwise.
a. Specify the method isLeaf.
b. If isLeaf were not a method of a class of binary trees, would a client of the class be able to implement isLeaf? Explain.
6. Starting with an empty binary search tree, in what order should you insert items to get the binary search tree in Figure 15-18?
7. Using the binary search tree in Figure 15-18, trace the search algorithm when it searches for
a. 30
b. 15

In each case, list the nodes in the order in which the search visits them.
8. Is the tree in Figure 15-19 a binary search tree? Explain.

FIGURE 15-19 A tree for Exercises 4 and 8

9. Beginning with an empty binary search tree, what binary search tree is formed when you insert the following values in the order given?
a. W, T, N, J, E, B, A
b. W, T, N, A, B, E, J
c. A, B, W, J, N, T, E
d. \(\mathrm{B}, \mathrm{T}, \mathrm{E}, \mathrm{A}, \mathrm{N}, \mathrm{W}, \mathrm{J}\)
10. Consider the binary search tree in Figure 15-20. The numbers simply label the nodes so that you can reference them; they do not indicate the contents of the nodes.
a. Without performing an inorder traversal, which node must contain the value that comes immediately after the value in the root? Explain.
b. In what order will an inorder traversal visit the nodes of this tree? Indicate this order by listing the labels of the nodes in the order that they are visited.
11. Consider the binary search tree in Figure \(15-18\). What tree results after you insert the nodes \(80,65,75,45,35\), and 25 , in that order?

FIGURE 15-20 A binary search tree for Exercise 10

12. If duplicates are allowed in a binary search tree, it is important to have a convention that determines the relationship between the duplicates. Items that duplicate the root of a tree should either all be in the left subtree or all be in the right subtree, and, of course, this property must hold for every subtree. Why is this convention critical to the effective use of the binary search tree?
13. Consider a nonempty binary tree with two types of nodes: min nodes and max nodes. Each node has an integer value initially associated with it. This tree is a minimax tree and has a value, which we define as follows:
- If the root is a min node, the value of the tree is equal to the minimum of
- The integer stored in the root
- The value of the left subtree, but only if it is nonempty
- The value of the right subtree, but only if it is nonempty
- If the root is a max node, the value of the tree is equal to the maximum of the above three values.
a. Compute the value of the minimax tree in Figure 15-21. Each node is labeled with its initial value.
b. Design a general solution for representing and evaluating these trees.
*14. A binary search tree with a given set of data items can have several different structures that conform to the definition of a binary search tree. If you are given a list of data items, does at least one binary search tree whose preorder traversal matches the order of the items on your list always exist? Is there ever more than one binary search tree that has the given preorder traversal?
*15. How many differently shaped \(n\)-node binary trees are possible? How many differently shaped \(n\)-node binary search trees are possible? (Write recursive definitions.)
16. Write pseudocode for a binary search tree method that visits all nodes whose data lies within a given range of values (such as all values between 100 and 1,000 ).
17. By using mathematical induction, prove that a full binary tree of height \(h \geq 0\) has \(2^{h}-1\) nodes.
18. By using mathematical induction, prove that the maximum number of nodes in a binary tree of height \(h\) is \(2^{h}-1\).

FIGURE 15-21 A minimax tree for Exercise 13

19. What is the maximum number of nodes that a binary tree can have at level \(n\) ? Prove your answer by using mathematical induction. Use this fact to do the following:
a. Rewrite the formal definition of a complete tree of height \(h\).
b. Derive a closed form for the formula
\[
\sum_{i=1}^{h} 2^{i-1}
\]

What is the significance of this sum?
20. Prove by mathematical induction that a binary tree with \(n\) nodes has exactly \(n+1\) empty subtrees.
21. A binary tree is strictly binary if every nonleaf node has exactly two children. Prove by mathematical induction on the number of leaves that a strictly binary tree with \(n\) leaves has exactly \(2 n-1\) nodes.
22. Consider two algorithms for traversing a binary tree. Both are nonrecursive algorithms that use an extra container \(C\) for bookkeeping. Both algorithms have the following basic form:
```

Put the root of the tree in C
while (C is not empty)
{
Remove a node from C and name it N
VisitN
if (N has a left child)
Put the child in C
if (Nhas a right child)
Put the child in C
}

```

The difference between the two algorithms is the approach for choosing a node \(N\) to remove from the container \(C\) :
- Algorithm 1: Remove the newest (most recently added) node from \(C\).
- Algorithm 2: Remove the oldest (earliest added) node from C.
a. In what order would each algorithm visit the nodes of the tree in Figure 15-13?
b. For each algorithm, describe an appropriate ADT for the bookkeeping container \(C\). What data should the ADT have? Be conservative with the amount of memory needed for the ADT. Also, note that the traversal of a tree should not alter the tree in any way.
23. Suppose that you traverse the binary search tree in Figure 15-18 and write the data item in each node visited to a file. You plan to read this file later and create a new binary search tree by using the ADT binary search tree operation add. In creating the file, in what order should you traverse the tree so that the new tree will have exactly the same shape and nodes as the original tree?

\section*{Programming Problems}
1. Develop a program that can be used to test an implementation of BinaryTreeInterface.
2. Repeat the previous problem, but assume that the implementation represents a binary search tree.
3. Develop an interface template for a general tree. What methods are reasonable for a general tree?
4. Repeat the previous problem, but consider an \(n\)-ary tree instead.

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\section*{Tree Implementations}

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A free list keeps track of available nodes

The previous chapter discussed various kinds of trees, but it focused on the ADTs binary tree and binary search tree. This chapter will show you how to implement these two ADTs by using the constructs of \(\mathrm{C}++\). In each case, the data structures will be private data members of a class of trees.

\subsection*{16.1 The Nodes in a Binary Tree}

The first step in implementing a tree is to choose a data structure to represent its nodes. Since each node must contain both data and "pointers" to the node's children-which are other nodes in the tree-it is natural to make each node an object. Thus, we will use a C++ class to define the nodes in
the tree. If we place these nodes in an array, the "pointers" in the nodes are array indices. However, if the nodes are a part of a linked chain, we use \(\mathrm{C}++\) pointers to link them together.

\subsection*{16.1.1 An Array-Based Representation}

Let's name our class of nodes TreeNode. An array-based implementation of a tree uses an array of nodes, so a class of such trees could have the following data members:
\begin{tabular}{lll} 
TreeNode<ItemType> & tree[MAX_NODES]; // Array of tree nodes \\
int & root; & // Index of root \\
int & free; & \(/ /\) Index of free 7 ist
\end{tabular}

The variable root is an index to the tree's root node within the array tree. If the tree is empty, root is -1 .

As the tree changes due to insertions and removals, its nodes may not be in contiguous elements of the array. Therefore, this implementation requires you to establish a collection of available nodes, which is called a free list. To insert a new node into the tree, you first obtain an available node from the free list. If you remove a node from the tree, you place it into the free list so that you can reuse the node at a later time. The data member free is the index to the first node in the free list. We will talk about the free list in a moment.

Although these data members are appropriate for any tree, we want to restrict our conversation to binary trees and consider a class of its nodes. The class TreeNode, as given in Listing 16-1, is such a class.
```

LISTING 16-1 The class TreeNode for an array-based implementation
of the ADT binary tree
temp1ate<c1ass ItemType>
class TreeNode
{
private:
ItemType item; // Data portion
int leftChild; // Index to left child
int rightChild; // Index to right child
public:
TreeNode();
TreeNode(const ItemType\& nodeItem, int left, int right);
// Declarations of the methods setItem, getItem, setLeft, getLeft,
// setRight, and getRight are here.
}; // end TreeNode

```

Each node in the array-based binary tree has a data item and two array indices, one to each child. Both 7 eftChild and rightChild within a node are indices to the children of that node. If a node has no left child, 1 eftChild is -1 ; if a node has no right child, rightChild is -1 . Since root is the index of the root \(r\) of a binary tree, tree[root].getLeft() is leftChild, the index of the root of the left subtree of \(r\); tree[root].getRight() is rightChild, the index of the root of the right subtree of \(r\).

The free list. Even though the data member free is the index of the first node in the free list, the next available node is not necessarily at index free +1 . When a node is removed from a tree and returned to the free list, it could be anywhere in the array. Thus, we "link" the available nodes together by arbitrarily making the rightChild member of each node be the index of the next node in the free list. Thus, both the free list and the tree itself are array-based, but the nodes are linked by array indices instead of \(\mathrm{C}++\) pointers. While the free list is a linear data structure, the tree is not.

Figure 16-1 contains a binary tree and its data members for this array-based implementation. Note how the nodes in the tree are linked. For example, Bob in tree [1] has Elisa as its right child. Elisa is in tree[4], and tree[1]. rightChild is 4. Since Elisa is in a leaf of the tree, its links in tree[4]. leftChild and tree[4]. rightChild are each - L. Likewise, Alan and Nancy are in leaf nodes. The nodes in the free list are linked, but since this tree has not undergone removals, the free list happens to be in contiguous array elements.

An array-based implementation of a binary tree is much more attractive when the tree is complete. In such cases, indices that link parents and their children are not stored, so the data structure is

FIGURE 16-1 (a) A binary tree of names; (b) its implementation using the array tree
(a)


simpler than if the tree is not complete. In Chapter 17, you will see that an array-based representation of a complete binary tree is useful in the implementation of the ADT priority queue. We will not use any other array-based implementation of a tree.

Note: If you use an array-based implementation of a complete binary tree, you must be sure that the tree remains complete as a result of insertions or removals.

Question 1 Represent the binary tree in Figure 15-18 of Chapter 15 with an array.

\subsection*{16.1.2 A Link-Based Representation}

You can use C++ pointers to link the nodes in the tree. Doing so is the most common way of implementing a tree. Listing 16-2 shows the class BinaryNode for a link-based implementation of the ADT binary tree.

\section*{LISTING 16-2 The header file containing the class BinaryNode for a link-based implementation of the ADT binary tree}
```

/** A class of nodes for a link-based binary tree.
@file BinaryNode.h */
\#ifndef _BINARY_NODE
\#define _BINARY_NODE
template<class ItemType>
class BinaryNode
{
private:
ItemType item; // Data portion
BinaryNode<ItemType>* 1eftChildPtr; // Pointer to left child
BinaryNode<ItemType>* rightChildPtr; // Pointer to right child
public:
BinaryNode();
BinaryNode(const ItemType\& anItem);
BinaryNode(const ItemType\& anItem,
BinaryNode<ItemType>* 1eftPtr,
BinaryNode<ItemType>* rightPtr);
void setItem(const ItemType\& anItem);
ItemType getItem() const;
bool isLeaf() const;
BinaryNode<ItemType>* getLeftChildPtr() const;
BinaryNode<ItemType>* getRightChildPtr() const;
void setLeftChildPtr(BinaryNode<ItemType>* 1eftPtr);
void setRightChildPtr(BinaryNode<ItemType>* rightPtr);
}; // end BinaryNode
\#include "BinaryNode.cpp"
\#endif

```

FIGURE 16-2 A link-based implementation of a binary tree


Given the class BinaryNode in Listing 16-2, a class of link-based binary trees will declare one data member-a pointer rootPtr-to point to the tree's root node. If the tree is empty, rootPtr contains nullptr. For a binary tree that is not empty, rootPtr->getLeftChildPtr() points to the root's left subtree, and rootPtr->getRightChildPtr() points to the root's right subtree. If either of these subtrees is empty, the pointer to it would be nul1ptr. Figure 16-2 illustrates the root of a binary tree and its two children.

\subsection*{16.2 A Link-Based Implementation of the ADT Binary Tree}

We will derive our class of link-based binary trees from BinaryTreeInterface, which specifies the public methods for the ADT binary tree and appears in Listing 15-1 of Chapter 15. The nodes in the tree will be instances of the class BinaryNode as given previously in Listing 16-2.

\subsection*{16.2.1 The Header File}

The header file in Listing 16-3 declares the class BinaryNodeTree in anticipation of its link-based implementation. Note the protected methods that the public methods will call to perform their operations recursively. These methods require pointers as arguments. As such, they should not be public and available to clients of the class. After all, clients should not have access to node pointers. Although these methods could be private, they are protected so that a derived class can use them.

The public section declares more constructors than we have in the past, allowing a client to define binary trees in a variety of circumstances. For example, you can construct a binary tree
- That is empty
- From data for its root, which is its only node
- From data for its root and from its two subtrees

For example, the following statements invoke these three constructors:
```

BinaryNodeTree<string> tree1;
BinaryNodeTree<string>* tree2Ptr = new BinaryNodeTree<string>("A");
BinaryNodeTree<string>* tree3Ptr = new BinaryNodeTree<string>("B");
BinaryNodeTree<string>* tree4Ptr =
new BinaryNodeTree<string>("C", tree2Ptr, tree3Ptr);

```

Some methods should not be public

Sample uses of public constructors

Here tree1 is an empty binary tree; tree2Ptr and tree3Ptr each point to binary trees that have only a root node. These two roots contain the strings " A " and " B ", respectively; and tree4Ptr points to a binary tree whose root contains " C " and has subtrees pointed to by tree2Ptr and tree3Ptr.

!
Programming Tip: In the previous example, if you were to use BinaryTreeInterface instead of BinaryNodeTree in the data type of either tree2Ptr or tree3Ptr, you would have to cast the variable to BinaryNodeTree before you could call the third constructor. For example, suppose you defined treePtr2 as follows:

BinaryTreeInterface<string>* tree2Ptr = new BinaryNodeTree<string>("A");
You would then define treePtr4 as follows:
BinaryNodeTree<string>* tree4Ptr = new BinaryNodeTree<string>("C", static_cast<BinaryNodeTree<string>*>(tree2Ptr), tree3Ptr);

BinaryNodeTree also declares an overloaded assignment operator. You could also, of course, overload other operators.

\section*{LISTING 16-3 A header file for the link-based implementation of the class BinaryNodeTree}
```

/** ADT binary tree: Link-based implementation.
@file BinaryNodeTree.h */
\#ifndef _BINARY_NODE_TREE
\#define _BINARY_NODE_TREE
\#include "BinaryTreeInterface.h"
\#include "BinaryNode.h"
\#include "PrecondViolatedExcep.h"
\#include "NotFoundException.h"
temp1ate<c1ass ItemType>
void BinaryNodeTree : public BinaryTreeInterface<ItemType>
{
private:
BinaryNode<ItemType>* rootPtr;
protected:
//-------------------------------------------
// Recursive helper methods for the public methods.
//--
int getHeightHelper(BinaryNode<ItemType>* subTreePtr) const;
int getNumberOfNodesHe1per(BinaryNode<ItemType>* subTreePtr) const;
// Recursively deletes al1 nodes from the tree.
void destroyTree(BinaryNode<ItemType>* subTreePtr);
// Recursively adds a new node to the tree in a left/right fashion to
// keep the tree balanced.

```
```

    BinaryNode<ItemType>* balancedAdd(BinaryNode<ItemType>* subTreePtr,
                BinaryNode<ItemType>* newNodePtr);
    // Removes the target value from the tree by calling moveValuesUpTree
// to overwrite value with value from child.
BinaryNode<ItemType>* removeValue(BinaryNode<ItemType>* subTreePtr,
const ItemType target, bool\& success);
// Copies values up the tree to overwrite value in current node until
// a leaf is reached; the leaf is then removed, since its value is
// stored in the parent.
BinaryNode<ItemType>* moveValuesUpTree(BinaryNode<ItemType>* subTreePtr);
// Recursively searches for target value in the tree by using a
// preorder traversal.
BinaryNode<ItemType>* findNode(BinaryNode<ItemType>* treePtr,
const ItemType\& target,
bool\& success) const;
// Copies the tree rooted at treePtr and returns a pointer to
// the copy.
BinaryNode<ItemType>*
copyTree(const BinaryNode<ItemType>* treePtr) const;
// Recursive traversal helper methods:
void preorder(void visit(ItemType\&),
BinaryNode<ItemType>* treePtr) const;
void inorder(void visit(ItemType\&),
BinaryNode<ItemType>* treePtr) const;
void postorder(void visit(ItemType\&),
BinaryNode<ItemType>* treePtr) const;
public:
//---------------------------------------------
//-------------------
BinaryNodeTree(const ItemType\& rootItem);
BinaryNodeTree(const ItemType\& rootItem,
const BinaryNodeTree<ItemType>* 1eftTreePtr,
const BinaryNodeTree<ItemType>* rightTreePtr);
BinaryNodeTree(const BinaryNodeTree<ItemType>\& tree);
virtual ~BinaryNodeTree();
//
// Public BinaryTreeInterface Methods Section.
//------------------------------------------------------------------
bool isEmpty() const;
int getHeight() const;
int getNumberOfNodes() const;
ItemType getRootData() const throw(PrecondViolatedExcep);
void setRootData(const ItemType\& newData);
bool add(const ItemType\& newData); // Adds a node
bool remove(const ItemType\& data); // Removes a node
void clear();

```

An overloaded assignment operator
```

    ItemType getEntry(const ItemType& anEntry) const
                                    throw(NotFoundException);
    bool contains(const ItemType& anEntry) const;
    //
// Public Traversals Section.
//-
void preorderTraverse(void visit(ItemType\&)) const;
void inorderTraverse(void visit(ItemType\&)) const;
void postorderTraverse(void visit(ItemType\&)) const;
//-------------------------------------
//--}\mathrm{ BinaryNodeTree\& operator=(const BinaryNodeTree\& rightHandSide);
}; // end BinaryNodeTree
\#include "BinaryNodeTree.cpp"
\#endif

```

\subsection*{16.2.2 The Implementation}

Although we will not provide a complete implementation file here, we will examine its most significant portions.

The constructors. The public constructors have the following definitions in the implementation file:
```

template<class ItemType>
BinaryNodeTree<ItemType>::BinaryNodeTree() : rootPtr(nul1ptr)
{
} // end default constructor
template<class ItemType>
BinaryNodeTree<ItemType>::BinaryNodeTree(const ItemType\& rootItem)
{
rootPtr = new BinaryNode<ItemType>(rootItem, nu11ptr, nu11ptr);
} // end constructor
template<class ItemType>
BinaryNodeTree<ItemType>::BinaryNodeTree(const ItemType\& rootItem,
const BinaryNodeTree<ItemType>* 1eftTreePtr,
const BinaryNodeTree<ItemType>* rightTreePtr)
{
rootPtr = new BinaryNode<ItemType>(rootItem,
copyTree(leftTreePtr->rootPtr),
copyTree(rightTreePtr->rootPtr));
} // end constructor

```

The default constructor uses an initializer to set the value of rootPtr to nul1ptr. The second constructor sets rootPtr to point to a new node containing the given data item. The third constructor also makes rootPtr point to a new node, but the node contains the given data item and pointers to copies of the given subtrees. We talk about the method copyTree next during our discussion of the copy constructor.

The copy constructor and destructor. The copy constructor and the destructor implicitly use traversal. Since traversal is a recursive operation, the copy constructor and destructor each call a recursive method. As we have mentioned before, public ADT methods usually are not themselves recursive, but rather call a recursive method that is either private or protected. We do this to hide the underlying data structure from the client.

The protected method copyTree, which the copy constructor calls, uses a recursive preorder traversal to copy each node in the tree. By copying each node as soon as the traversal visits it, copyTree can make an exact copy of the original tree. To make the copy distinct from the original tree, the new nodes must be linked together by using new pointers. That is, you cannot simply copy the pointers in the nodes of the original tree. The result is a deep copy of the tree.

Thus, copyTree has the following definition:
```

temp7ate<class ItemType>
BinaryNode<ItemType>* BinaryNodeTree<ItemType>::
copyTree(const BinaryNode<ItemType>* treePtr) const
{
BinaryNode<ItemType>* newTreePtr = nul1ptr;
// Copy tree nodes during a preorder traversal
if (treePtr != nul1ptr)
{
// Copy node
newTreePtr = new BinaryNode<ItemType>(treePtr->getItem(),
nu17ptr, nu17ptr);
newTreePtr->setLeftChildPtr(copyTree(treePtr->getLeftChildPtr()));
newTreePtr->setRightChildPtr(copyTree(treePtr->getRightChildPtr()));
} // end if
// Else tree is empty (newTreePtr is nul7ptr)
return newTreePtr;
} // end copyTree

```

The copy constructor then looks like this:
```

temp7ate<class ItemType>
BinaryNodeTree<ItemType>::
BinaryNodeTree(const BinaryNodeTree<ItemType>\& treePtr)
{
rootPtr = copyTree(treePtr.rootPtr);
} // end copy constructor

```

Similarly, the protected method destroyTree, which the destructor calls, uses a recursive postorder traversal to delete each node in the tree. A postorder traversal is appropriate here because you can delete a node only after you have first traversed and deleted both of its subtrees. Thus, destroyTree has the following definition:
```

temp7ate<class ItemType>
void BinaryNodeTree<ItemType>::
destroyTree(BinaryNode<ItemType>* subTreePtr)
{
if (subTreePtr != nul7ptr)
{
destroyTree(subTreePtr->getLeftChildPtr());
destroyTree(subTreePtr->getRightChildPtr());
delete subTreePtr;
} // end if
} // end destroyTree

```

The destructor then only needs to make the call destroyTree(rootPtr).

\section*{To copy a tree,} traverse it in preorder and insert each item visited into a new node

The method getHeight. The public method getHeight calls the protected, recursive method getHeightHelper. The height of a subtree rooted at a particular node is 1 -for the node itself—plus the height of the node's tallest subtree. Thus, we define the protected method getHeightHelper as follows:
```

temp1ate<class ItemType>
int BinaryNodeTree<ItemType>::
getHeightHelper(BinaryNode<ItemType>* subTreePtr) const
{
if (subTreePtr == nu11ptr)
return 0;
else
return 1 + max(getHeightHe1per(subTreePtr->getLeftChildPtr()),
getHeightHelper(subTreePtr->getRightChildPtr()));
} // end getHeightHe7per

```

The public method getHeight contains just the statement
return getHeightHelper (rootPtr);
The method getNumberOfNodes has a similar definition.

Question 2 What are the definitions of the public method getNumberOfNodes and the protected helper method getNumberOfNodesHelper?
Question 3 What is the definition of the public method setRootData?
Question 4 What is the definition of the public method getRootData? Recall that this method has a precondition.

The method add. The specification of the public method add says to add a new node containing given data to a binary tree. It does not indicate where that new node should be in the tree. Thus, we have flexibility in how we define the method. Moreover, derived classes of BinaryNodeTree can both override add to change its behavior and adhere to its specification.

Let's add the new node so that the resulting tree is balanced. Once again, we can use recursion to achieve our goal. The public method add can create a new node containing the data passed to it as an argument. Then the method can pass this node to a protected, recursive method that adds it in an appropriate place in the tree. The protected method must then return a pointer to the tree's root. If we name the protected method balancedAdd, the public add method has the following definition:
```

template<class ItemType>
boo1 BinaryNodeTree<ItemType>::add(const ItemType\& newData)
{
BinaryNode<ItemType>* newNodePtr = new BinaryNode<ItemType> (newData);
rootPtr = balancedAdd(rootPtr, newNodePtr);
return true;
} // end add

```

Now we need to define the protected method balancedAdd. Notice that the add method passes two arguments to balancedAdd: a pointer to the root of the tree - or subtree - to which we will add a node and a pointer to the new node. To add a node to the tree, we add the node to the root's shorter
subtree. This is the recursive step, since the subtree is a binary tree. The base case occurs when a subtree is empty.

Here is the definition of balancedAdd:
```

temp7ate<class ItemType>
BinaryNode<ItemType>* BinaryNodeTree<ItemType>::
balancedAdd(BinaryNode<ItemType>* subTreePtr,
BinaryNode<ItemType>* newNodePtr)
{
if (subTreePtr == nu17ptr)
return newNodePtr;
else
{
BinaryNode<ItemType>* 1eftPtr = subTreePtr->getLeftChi1dPtr();
BinaryNode<ItemType>* rightPtr = subTreePtr->getRightChildPtr();
if (getHeightHelper(leftPtr) > getHeightHelper(rightPtr))
{
rightPtr = balancedAdd(rightPtr, newNodePtr);
subTreePtr->setRightChildPtr(rightPtr);
}
else
{
1eftPtr = balancedAdd(1eftPtr, newNodePtr);
subTreePtr->setLeftChildPtr(1eftPtr);
} // end if
return subTreePtr;
} // end if
} // end balancedAdd

```

For example, the following statements in bal ancedAdd add a new node to a right subtree:
```

rightPtr = balancedAdd(rightPtr, newNodePtr);
subTreePtr->setRightChildPtr(rightPtr);

```

The recursive call to balancedAdd adds the new node and returns a pointer to the revised subtree. However, we need to link this subtree to the rest of the tree. The call to setRightChildPtr accomplishes this. The addition of a new node to a left subtree is performed in a similar manner. Note that such insertions are analogous to the recursive addition of a new node to a linear chain of linked nodes, as discussed in Section 9.2.3 of Chapter 9. Reviewing that section will give you insight into the process here.

Figure 16-3 shows the effect of a sequence of add operations on an initially empty binary tree.

Question 5 Where would a new node be placed next in the binary tree shown in Figure 16-3?

The traversals. Since the traversals are recursive, the public traversal methods each call a protected method that performs the actual recursion. For example, the public method inorderTraverse calls the protected method inorder. Like inorderTraverse, inorder has the function visit as a parameter. This function specifies the tree item as a reference parameter, which enables the client not only to examine the item but also to modify it. The second parameter of inorder is

Implement traversals so that visit remains on the client's side of the wall

\section*{FIGURE 16-3 Adding nodes to an initially empty binary tree}









the pointer treePtr, which, due to the recursive calls, eventually points to every node in the tree. The definition of inorder follows:

Protected methods that enable recursive traversals
```

temp1ate<c1ass ItemType>
void BinaryNodeTree<ItemType>::
inorder(void visit(ItemType\&), BinaryNode<ItemType>* treePtr) const

```
```

{
if (treePtr != nul1ptr)
{
inorder(visit, treePtr->getLeftChildPtr());
ItemType theItem = treePtr->getItem();
visit(theItem);
inorder(visit, treePtr->getRightChildPtr());
} // end if
} // end inorder

```

Now the definition of the public method inorderTraverse contains only the call
```

inorder(visit, rootPtr);

```

The other traversals have similar definitions.


Programming Tip: Why did we write
ItemType theItem = treePtr->getItem(); visit(theItem);
instead of
```

    visit(treePtr->getItem());
    ```
in the definition of inorder? Because visit has a reference parameter, we need an lvalue to reference. The variable theItem serves this purpose.

Question 6 Define the protected method postorder.

Nonrecursive traversal (optional). Before leaving the topic of traversals, let's develop a nonrecursive traversal algorithm to illustrate further the relationship between stacks and recursion that was discussed in Chapter 6. In particular, we will develop a nonrecursive inorder traversal for the linkbased implementation of a binary tree.

The conceptually difficult part of a nonrecursive traversal is determining where to go next after a particular node has been visited. To gain some insight into this problem, consider how the recursive inorder method works:
```

if (treePtr != nul1ptr)
{
inorder(visit, treePtr->getLeftChildPtr()); // Point 1
ItemType theItem = treePtr->getItem();
visit(theItem);
inorder(visit, treePtr->getRightChildPtr()); // Point 2
} // end if

```

The method has its recursive calls marked as points 1 and 2.
During the course of the method's execution, the value of the pointer treePtr actually marks the current position in the tree. Each time inorder makes a recursive call, the traversal moves to another node. In terms of the stack that is implicit to recursive methods, a call to inorder pushes the new

Recursive calls from points 1 and 2

FIGURE 16-4 Contents of the implicit stack as treePtr progresses through a given tree during a recursive inorder traversal
(The notation \(\rightarrow 60\) means "a pointer to the node containing 60.")


Recursive inorder's implicit stack gives insight into a nonrecursive traversal
value of treePtr - that is, a pointer to the new current node-onto the stack. At any given time, the stack contains pointers to the nodes along the path from the tree's root to the current node \(n\), with the pointer to \(n\) at the top of the stack and the pointer to the root at the bottom. Note that \(n\) is possibly "empty"-that is, it may be indicated by a nullptr value for treePtr at the top of the stack. Such an occurrence corresponds to the base case of the recursion.

Figure 16-4 partially traces the execution of inorder and shows the contents of the implicit stack. The first four steps of the trace show the stack as treePtr points first to 60 , then to 20 , then to 10 , and then becomes nul1ptr. The recursive calls for these four steps are from point 1 in inorder.

Now consider what happens when inorder returns from a recursive call. The traversal retraces its steps by backing up the tree from a node \(n\) to its parent \(p\), from which the recursive call to \(n\) was made. Thus, the pointer to \(n\) is popped from the stack and the pointer to \(p\) comes to the top of the stack, as occurs in step 5 of the trace in Figure 16-4. In this case, \(n\) happens to be empty, so nul1ptr is popped from the stack.

What happens next depends on which subtree of \(p\) has just been traversed. If you have just finished traversing \(p\) 's left subtree-that is, if \(n\) is the left child of \(p\) and thus the return is made to point 1 in inorder-control is returned to the statement that visits node \(p\). Such is the case for steps 6 and 10 of the trace in Figure 16-4. Figure 16-5a illustrates steps 9 and 10 in more detail.

After node \(p\) has been visited, a recursive call is made from point 2 and the right subtree of \(p\) is traversed. However, if, as Figure 16-5b illustrates, you have just traversed \(p\) 's right subtree-that is, if \(n\) is the right child of \(p\) and thus the return is made to point 2 -control is returned to the end of the method. As a consequence another return is made, the pointer to \(p\) is popped off the stack, and you go back up the tree to \(p\) 's parent, from which the recursive call to \(p\) was made. In this latter case, node \(p\) is not visited - it was visited before the recursive call to \(n\) was made from point 2 .

FIGURE 16-5 Traversing (a) the left subtree (steps 9 and 10 in Figure 16-4) and (b) the right subtree of 20


Thus, two facts emerge from the recursive version of inorder when a return is made from a recursive call:
- The implicit recursive stack of pointers is used to find the node \(p\) that the traversal must go back to.
- Once the traversal backs up to node \(p\), it either visits \(p\) (for example, displays its data) or backs farther up the tree. It visits \(p\) if \(p\) 's left subtree has just been traversed; it backs up if its right subtree has just been traversed. The appropriate action is taken simply as a consequence of the point-1 or 2-to which control is returned.

You could directly mimic this action by using an iterative method and an explicit stack, as long as some bookkeeping device kept track of which subtree of a node had just been traversed. However, you can use the following observation both to eliminate the need for the bookkeeping device and to speed up the traversal somewhat. Consider the tree in Figure 16-6. After you have finished traversing the subtree rooted at node \(R\), there is no need to return to nodes \(C\) and \(B\), because the right subtrees of these nodes have already been traversed. You can instead return directly to node \(A\), which is the nearest ancestor of \(R\) whose right subtree has not yet been traversed.

This strategy of not returning to a node after its right subtree has been traversed is simple to implement: You place a pointer to a node in the stack only before the node's left subtree is traversed, but not before its right subtree is traversed. Thus, in Figure 16-6, when you are at node \(R\), the stack contains \(A\) and \(R\), with \(R\) on top. Nodes \(B\) and \(C\) are not in the stack, because you have visited them already and are currently traversing their right subtrees. On the other hand, \(A\) is in the stack because you are currently traversing its left subtree. When you return from node \(R\), nodes \(B\) and \(C\) are thus bypassed because you have finished with their right subtrees and do not need to return to these nodes. Thus, you pop \(R\) 's pointer from the stack and go directly to node \(A\), whose left subtree has just been traversed. You then visit \(A\), pop its pointer from the stack, and traverse \(A\) 's right subtree.

Actions upon a return from a recursive call to inorder

FIGURE 16-6 Avoiding returns to nodes \(B\) and \(C\)

nodeStack


Nonrecursive inorder traversal

This nonrecursive traversal strategy is captured by the following pseudocode, assuming a linkbased implementation. Exercise 9 at the end of this chapter asks you to trace this algorithm for the tree in Figure 16-4.
```

// Nonrecursively traverses a binary tree in inorder.
traverse(visit(item: ItemType): void): void
// Initialize
nodeStack = A new, empty stack
curPtr = rootPtr() // Start at root
done = false
while (!done)
{
if (curPtr != nul1ptr)
{
// Place pointer to node on stack before traversing the node's left subtree
nodeStack.push(curPtr)
// Traverse the left subtree
curPtr = curPtr->getLeftChildPtr()
}
e1se // Backtrack from the empty subtree and visit the node at the top of
// the stack; however, if the stack is empty, you are done
{
if (!nodeStack.isEmpty())
{
nodeStack.peek(curPtr)

```
```

    visit(curPtr->getItem())
    nodeStack.pop()
    // Traverse the right subtree of the node just visited
    curPtr = curPtr ->getRightChildPtr()
    }
        else
    done = true
    }
    }

```

Eliminating recursion can be more complicated than the example given here. However, the general case is beyond the scope of this book.

\subsection*{16.3 A Link-Based Implementation of the ADT Binary Search Tree}

Since a binary search tree is a binary tree, its implementation can use the same node objects as for a binary-tree implementation. As we plan a link-based implementation, we will use the class BinaryNode, as given earlier in Listing 16-2.

The recursive search algorithm that we presented in Section 15.3.2 of Chapter 15 is the basis of the insertion, removal, and retrieval operations on a binary search tree. The algorithms that follow for the binary search tree operations assume the link-based implementation of a binary tree that was discussed earlier in this chapter. Also keep in mind the assumption that the data items in the binary search tree are unique.

\subsection*{16.3.1 Algorithms for the ADT Binary Search Tree Operations}

Adding a new entry. As Section 15.3.3 of Chapter 15 demonstrated, you insert a new entry into a binary search tree in the same place that the search algorithm would look for it. For example, if you want to insert Kody into the binary search tree shown in Figure 16-7a, you would first search the tree

Use search to determine the insertion point

FIGURE 16-7 \(\quad\) (a) A binary search tree; (b) empty subtree where the search algorithm terminates when looking for Kody; (c) the tree after Kody is inserted as a new leaf


First draft of the insertion algorithm

Refinement of the insertion algorithm
for Kody. The search algorithm would terminate at Nancy's empty left subtree, as Figure 16-7b illustrates. You would then insert Kody as Nancy's left child, as Figure 16-7c shows.

Because Nancy has no left child, the insertion is simple, requiring only that Nancy's left-child pointer point to Kody. Because searching for an entry that is not in the binary search tree always ends at an empty subtree, you always insert a new item as a new leaf. Because adding a leaf requires only a change of the appropriate pointer in the parent, the work required for an insertion is virtually the same as that for the corresponding search.

We begin by defining the public method add as follows:
```

temp1ate<class ItemType>
boo1 BinarySearchTree<ItemType>::add(const ItemType\& newData)
{
BinaryNode<ItemType>* newNodePtr = new BinaryNode<ItemType>(newData);
rootPtr = insertInorder(rootPtr, newNodePtr );
return true;
} // end add

```

The method creates a new node and passes pointers to the tree and the new node to a recursive method that actually performs the insertion. The following high-level pseudocode describes this insertion process:
```

// Inserts a new node into a binary search tree.
insertInorder(subTreePtr: BinaryNodePointer,
newNodePtr: BinaryNodePointer): BinaryNodePointer
Search the tree pointed to by subTreePtr for the item in the node pointed
to by newNodePtr
if (the search terminates at parentNode's left subtree)
Set leftChildPtr of parentNode to newNodePtr
else
Set rightChildPtr of parentNode to newNodePtr

```

The appropriate pointer-1eftChildPtr or rightChildPtr—of node parentNode must be set to point to the new node. The recursive nature of the search algorithm provides an elegant means of setting the pointer, provided that you return this possibly changed pointer. The situation is quite similar to the recursive insertion method for the ADT list that you saw in Chapter 9. If the tree was empty before the insertion, the external pointer to the root of the tree would be nul1ptr and the method would not make a recursive call. Thus, you would return a pointer to the new node. Our pseudocode is refined as follows:
```

// Inserts a new item into the binary search tree to which subTreePtr points.
insertInorder(subTreePtr: BinaryNodePointer,
newNodePtr: BinaryNodePointer): BinaryNodePointer
if (subTreePtr is nul1ptr)
return newNodePtr
else if (subTreePtr->getItem() > newNodePtr->getItem())
{
tempPtr = insertInorder(subTreePtr->getLeftChildPtr(), newNodePtr)
subTreePtr->setLeftChildPtr(tempPtr)
}
e1se
{

```

FIGURE 16-8 (a) Insertion into an empty tree; (b) search for Frank terminates at a leaf;
(c) insertion at a leaf

```

    tempPtr = insertInorder(subTreePtr->getRightChildPtr(), newNodePtr)
    subTreePtr->setRightChildPtr(tempPtr)
    }
return subTreePtr

```

When subTreePtr is nul1ptr, insertInorder returns a pointer to the new node. For an initially empty tree, add assigns the returned pointer to rootPtr, as Figure 16-8a illustrates. For a tree that is not empty, add assigns the returned pointer to the pointer leftChildPtr or rightChildPtr-whose current value is nullptr-in the parent of the empty subtree. Thus, the appropriate pointer within the parent is set to point to the new node. Parts \(b\) and \(c\) of Figure 16-8 illustrate the general case of insertion.

You can use the public method add to create a binary search tree. For example, beginning with an empty tree, if you insert the names Jane, Bob, Alan, Elisa, Tom, Nancy, and Wendy in order, you will get the binary search tree in Figure 15-13 of Chapter 15. It is interesting to note that the names Jane, Bob, Alan, Elisa, Tom, Nancy, and Wendy constitute the preorder traversal of the tree in Figure 15-13. Thus, if you take the output of a preorder traversal of a binary search tree and use it with insertInorder to create a binary search tree, you will obtain a duplicate tree. This result should not surprise you, as the copy constructor for the ADT binary tree used a preorder traversal to copy the tree.

By inserting the previous names in a different order, you will get a different binary search tree. For example, by inserting the previous names in alphabetical order, you will get the binary search tree in Figure 15-14b of Chapter 15.

Question 7 Starting with an empty binary search tree, in what order should you insert items to get the binary search tree in Figure 15-18 of Chapter 15?

First draft of the removal algorithm

Three cases for the node \(N\) containing the item to be removed

Case 1: Set the pointer in a leaf's parent to nu11ptr

Case 2: Two possibilities for a node with one child

Let N's parent adopt \(N\) 's child

Case 3: \(N\) has two children

Removing an item whose node has two children

Removing an entry. Removing an entry from a binary search tree is a bit more involved than adding one. First, you use the search algorithm to locate the specified item and then, if it is found, you must remove it from the tree. A first draft of the algorithm follows:
```

// Removes the given target from a binary search tree.
// Returns true if the removal is successful or false otherwise.
removeValue(target: ItemType): boolean
Locate the target by using the search algorithm
if (target is found)
{
Remove target from the tree
return true
}
else
return false

```

The essential task here is to remove the target from the tree. Assuming that removeValue locates the target in a particular node \(N\), there are three cases to consider:
- \(N\) is a leaf
- \(N\) has only one child
- \(N\) has two children

Case 1. The first case is the easiest. To remove the leaf containing the target, you need only set the pointer in its parent to nullptr.

Case 2. The second case is a bit more involved. If \(N\) has only one child, you have two possibilities:
- \(N\) has only a left child
- \(N\) has only a right child

The two possibilities are symmetrical, so it is sufficient to illustrate the solution for a left child. In Figure \(16-9 \mathrm{a}, L\) is the left child of \(N\) and \(P\) is the parent of \(N . N\) can be either the left or right child of \(P\). If you deleted \(N\) from the tree, \(L\) would be without a parent and \(P\) would be without one of its children. Suppose you let \(L\) take the place of \(N\) as one of \(P\) 's children, as in Figure 16-9b. Does this adoption preserve the binary search tree property?

If \(N\) is the left child of \(P\), for example, all data items in the subtree rooted at \(N\) are less than the item in \(P\). Thus, all data items in the subtree rooted at \(L\) are less than the item in \(P\). Therefore, after \(N\) is removed and \(L\) is adopted by \(P\), all items in \(P\) 's left subtree are still less than the item in \(P\). This strategy thus preserves the binary search tree property. A parallel argument holds if \(N\) is a right child of \(P\), and therefore the binary search tree property is preserved in either case.

Case 3. The most difficult of the three cases occurs when the item to be removed is in a node \(N\) that has two children, as in Figure 16-10. As you just saw, when \(N\) has only one child, the child replaces \(N\). However, when \(N\) has two children, these children cannot both replace \(N\) : \(N\) 's parent has room for only one of \(N\) 's children as a replacement for \(N\). A different strategy is necessary.

In fact, you will not delete \(N\) at all. You can find another node that is easier to delete and delete it instead of \(N\). This strategy may sound like cheating, but remember that the client expects only a certain entry to be removed from the ADT. It has no right, because of the wall between the program and the ADT implementation, to expect a particular node in the tree to be deleted.

Consider, then, an alternate strategy. To remove from a binary search tree an item that resides in a node \(N\) that has two children, take the following steps:
1. Locate another node \(M\) that is easier to remove from the tree than the node \(N\).
2. Copy the item that is in \(M\) to \(N\), thus effectively removing from the tree the item originally in \(N\).
3. Remove the node \(M\) from the tree.

FIGURE 16-9 (a) \(N\) with only a left child \(-N\) can be either the left child or right child of \(P\);
(b) after removing node \(N\)


What kind of node \(M\) is easier to remove than the node \(N\) ? Because you know how to delete a node that has either no children or one child, \(M\) could be such a node. You have to be careful, though. Can you choose any node and copy its data into \(N\) ? No, because you must preserve the tree's status as a binary search tree. For example, if in the tree of Figure 16-11a, you copied the data from \(M\) to \(N\), the result in Figure 16-11b would no longer be a binary search tree.

What data item, when copied into the node \(N\), will preserve the tree's status as a binary search tree? All of the entries in the left subtree of \(N\) are less than the entry in \(N\), and all of the entries in the right subtree of \(N\) are greater than the entry in \(N\). You must retain this property when you replace the

FIGURE 16-10 \(N\) with two children


The inorder successor of N's entry is in the leftmost node in \(N\) 's right subtree

FIGURE 16-11 (a) Not any node will do; (b) no longer a binary search tree
(a)

(b)

entry \(x\) in node \(N\) with the entry \(y\). There are two suitable possibilities for the value \(y\) : It can come immediately after or immediately before \(x\) in the sorted order of entries. If \(y\) comes immediately after \(x\), clearly all entries in the left subtree of \(N\) are smaller than \(y\), because they are all smaller than \(x\), as Figure 16-12 illustrates. Further, all entries in the right subtree of \(N\) are greater than or equal to \(y\), because they are greater than \(x\) and, by assumption, there are no entries in the tree between \(x\) and \(y\). A similar argument illustrates that if \(y\) comes immediately before \(x\) in the sorted order, it is greater than or equal to all entries in the left subtree of \(N\) and smaller than all entries in the right subtree of \(N\).

You can thus copy into \(N\) either the item that is immediately after \(N\) 's entry or the item that is immediately before it. Suppose that, arbitrarily, you decide to use the node whose entry \(y\) comes immediately after \(N\) 's entry \(x\). This entry is called \(x\) 's inorder successor. \({ }^{1}\) How can you locate this node? Because \(N\) has two children, the inorder successor of its data item is in the leftmost node of \(N\) 's right subtree. That is, to find the node that contains \(y\), you follow \(N\) 's right-child pointer to its right child \(C\), which must be present because \(N\) has two children. You then descend the tree rooted at \(C\) by taking left branches at each node until you encounter a node \(S\) with no left child. You copy the item in

FIGURE 16-12 Search key \(x\) can be replaced by \(y\)


\footnotetext{
1 We also will use the term \(N\) 's inorder successor to mean the inorder successor of \(N\) 's entry.
}

FIGURE 16-13 Copying the item whose search key is the inorder successor of \(N\) 's search key

this node \(S\) into node \(N\) and then, because \(S\) has no left child, you can remove \(S\) from the tree as one of the two easy cases, as Figure 16-13 illustrates.

A more detailed high-level description of the removal algorithm follows:
```

// Removes the given target from a binary search tree.
// Returns true if the removal is successful, or false otherwise.

```
removeValue(subTreePtr: BinaryNodePointer, target: ItemType): boolean
    Locate the target by using the search algorithm; it occurs in node \(N\)
    if (target is found in node \(N\) )
    \{
        removeNode(N) // Defined next
        return true
    \}
    e1se
        return false
// Removes the item in a given node of a binary search tree.
removeNode(N: BinaryNode)
    if ( N is a leaf)
        Remove N from the tree
    else if (N has only one child C)
    \{
        if (N was a left child of its parent P)
            Make C the left child of P
        else
            Make C the right child of P
    \}
    e1se // Node has two children
    \{
        Find S, the node that contains N's inorder successor
        Copy the item from node S into node N
        Remove S from the tree by using the previous technique for a leaf or a node
        with one child
    \}

In the following refinement, search's algorithm is adapted and inserted directly into removeValue. Also, the method removeNode uses the method findSuccesorNode to find the node, \(S\), that contains the inorder successor of node \(N\). The method findSuccesorNode returns the item in \(S\) and then deletes node \(S\) from the tree. The returned item then replaces the item in node \(N\), thus removing it from the binary search tree.

Second draft of the removal algorithm
```

// Removes the given target from the binary search tree to which subTreePtr points.
// Returns a pointer to the node at this tree location after the value is removed.
// Sets success to true if the removal is successful, or false otherwise.
removeValue(subTreePtr: BinaryNodePointer, target: ItemType,
success: boolean\&): BinaryNodePointer
if (subTreePtr == nul1ptr)
{
success = false
return nul1ptr
}
else if (subTreePtr->getItem() == target)
{
// Item is in the root of some subtree
subTreePtr = removeNode(subTreePtr) // Remove the item
success = true
return subTreePtr
}
else if (subTreePtr->getItem() > target)
{
// Search the left subtree
tempPtr = removeValue(subTreePtr->getLeftChildPtr(), target, success)
subTreePtr->setLeftChildPtr(tempPtr)
return subTreePtr
}
else
{
// Search the right subtree
tempPtr = removeValue(subTreePtr->getRightChildPtr(), target, success)
subTreePtr->setRightChildPtr(tempPtr)
return subTreePtr
}
// Removes the item in the node, N, to which nodePtr points.
// Returns a pointer to the node at this tree location after node N is deleted.
removeNode(nodePtr: BinaryNodePointer): BinaryNodePointer
if (N is a leaf)
{
// Remove leaf from the tree
delete nodePtr
nodePtr = nul1ptr
return nodePtr
}
else if (N has only one child C)
{
// C replaces N as the child of N's parent
if (C is a left child)
nodeToConnectPtr = nodePtr->getLeftChildPtr()
else
nodeToConnectPtr = nodePtr->getRightChildPtr()
delete nodePtr
nodePtr = nullptr
return nodeToConnectPtr
}
else // Nhas two children
{
// Find the inorder successor of the entry in N: it is in the left subtree rooted
// at N's right child
tempPtr = removeLeftmostNode(nodePtr->getRightChildPtr(), newNodeValue)
nodePtr->setRightChildPtr(tempPtr)

```
```

        nodePtr->setItem(newNodeValue) // Put replacement value in node N
        return nodePtr
        }
    //Removes the leftmost node in the left subtree of the node pointed to by nodePtr.
// Sets inorderSuccessor to the value in this node.
// Returns a pointer to the revised subtree.
removeLeftmostNode(nodePtr: BinaryNodePointer,
inorderSuccesssor: ItemType\&): BinaryNodePointer
if (nodePtr->getLeftChildPtr() == nullptr)
{
// This is the node you want; it has no left child, but it might have a right subtree
inorderSuccesssor = nodePtr->getItem()
return removeNode(nodePtr)
}
else
return removeLeftmostNode(nodePtr->getLeftChildPtr(), inorderSuccesssor)

```

Now the public method remove calls removeValue and assigns the pointer it receives back from this call to rootPtr, as the following pseudocode indicates:
```

// Removes the given data from this binary search tree.
remove(target: ItemType): boolean
success $=$ false
rootPtr $=$ removeValue(rootPtr, target, success)
return success

```

Let's see how the recursive calls work. Suppose that we have the tree in Figure 16-14a, and we want to remove Dave. We call remove, which calls removeValue. Assuming that removeValue recursively locates Dave in Node \(N\), it calls removeNode to remove \(N\). Any change that removeNode makes to its parameter nodePtr, is returned by removeNode (Figure 16-14b) and then by removeValue. Because Dave isn't in the tree's root, the return from removeValue is due to a recursive call. The method now assigns the pointer returned by removeNode - and the recursive call to removeValue-to the left child pointer of Node \(N\) 's parent. The result is shown in Figure 16-14c.

Observe that in the case of the removeNode method, the argument that corresponds to its parameter nodePtr is either one of the pointers of the parent of \(N\), as Figure 16-14 depicts, or the external pointer to the root, in the case where \(N\) is the root of the original tree. In either case, nodePtr points to \(N\). Thus, any change you make to nodePtr by calling the method removeNode with the argument nodePtr must be returned so you can change either a pointer in the parent of node \(N\) or the root pointer.

Note: The methods removeValue, removeNode, removeLeftmostNode are behind the wall of abstraction, since they are protected. Therefore they can use pointers as parameters, and if necessary, a reference parameter (success or inorderSuccessor). Each method receives as an argument a pointer to the root of the subtree on which it is to perform its process. When that process is complete, the method reflects-like a mirrorback to the parent the pointer to the node that is now the root of that subtree.

Note that what we have done is safe programming:
- Only one reference parameter is used in a method, and the argument passed to it is local to the calling method.
- The calling method has a choice of whether or not to replace the pointer it has passed as an argument with the one returned-which it should do here.
- Each of the three methods performs a single, clearly defined task.

The recursive method removeLeftmostNode, which is called by removeNode if \(N\) has two children, also uses this strategy to remove the inorder successor of the node containing the item to be removed.

Exercise 12 at the end of this chapter describes an easier removal algorithm. However, that algorithm tends to increase the height of the tree, and, as you will see later, an increase in height can decrease the efficiency of searching the tree.

FIGURE 16-14 Recursive deletion of node \(N\)

(c)

Pointer returned by removeNode/removeValue is assigned to left child pointer of Node N's parent


Question 8 Given the binary search tree in Figure 15-18 of Chapter 15, trace the removal algorithms when removing each of the following values from the tree. Begin with the original tree each time.
a. 70
b. 20
c. 60

Retrieving an entry. By refining the search algorithm, you can implement the operation getEntry. Recall from Section 15.3.2 of Chapter 15 that the search algorithm is
```

// Searches a binary search tree for a given target value.
search(binTree: BinarySearchTree, target: ItemType)
if (binTree is empty)
The desired item is not found
else if (target == data item in the root of binTree)
The desired item is found
else if (target < data item in the root of binTree)
search(Left subtree of binTree, target)
else
search(Right subtree of binTree, target)

```

We can refine this algorithm as follows:
```

// Locates the node in the binary search tree to which subTreePtr points that contains
// the value target. Returns either a pointer to the located node or nu17ptr if such a
// node is not found.
findNode(subTreePtr: BinaryNodePointer, target: ItemType): BinaryNodePointer
if (subTreePtr == nul1ptr)
return nul1ptr // Not found
else if (subTreePtr->getItem() == target)
return subTreePtr; // Found
else if (subTreePtr->getItem() > target)
// Search left subtree
return findNode(subTreePtr->getLeftChildPtr(), target)
e1se
// Search right subtree
return findNode(subTreePtr->getRightChildPtr(), target)

```

The operation getEntry must return the item with the desired value if it exists; otherwise it must throw an exception NotFoundException. The method, therefore, calls findNode and checks its return value. If the desired target is found, getEntry returns it. If findNode returns nul1ptr, getEntry throws an exception.

\section*{Note: Traversals}

The traversals of a binary search tree are the same as the traversals of a binary tree. Recall from Chapter 15 that an inorder traversal of a binary search tree will visit the tree's nodes in sorted order according to their entries.

\subsection*{16.3.2 The Class BinarySearchTree}

A C++ link-based definition of the class BinarySearchTree is given in Listing 16-4. Notice the protected methods that implement the recursive algorithms. These methods are not public, because
findNode is a refinement of search

\[
1
\]
clients do not have access to node pointers. The methods could be private instead, but making them protected enables a derived class to use them directly.

The ADT binary tree and the ADT binary search tree have many methods in common, as Section 15.3.1 of Chapter 15 noted. This observation should not surprise you, as a binary search tree is a binary tree. For this reason, we have derived BinarySearchTree from BinaryNodeTree.

\section*{LISTING 16-4 A header file for the link-based implementation of the class BinarySearchTree}
```

/** Link-based implementation of the ADT binary search tree.
@file BinarySearchTree.h */
\#ifndef _BINARY_SEARCH_TREE
\#define _BINARY_SEARCH_TREE
\#include "BinaryTreeInterface.h"
\#include "BinaryNode.h"
\#include "BinaryNodeTree.h"
\#include "NotFoundException.h"
\#include "PrecondViolatedExcep.h"
template<class ItemType>
class BinarySearchTree : public BinaryNodeTree<ItemType>
{
private:
BinaryNode<ItemType>* rootPtr;
protected:

```

    // Protected Utility Methods Section:
    // Recursive he1per methods for the pub1ic methods.
    //-----------------------------------------------------------------1
    // Recursively finds where the given node should be placed and
    // inserts it in a leaf at that point.
    BinaryNode<ItemType>* insertInorder(BinaryNode<ItemType>* subTreePtr,
                BinaryNode<ItemType>* newNode);
    // Removes the given target value from the tree while maintaining a
    // binary search tree.
    BinaryNode<ItemType>* removeValue(BinaryNode<ItemType>* subTreePtr,
                                    const ItemType target,
                                    bool\& success);
    // Removes a given node from a tree while maintaining a
    // binary search tree.
    BinaryNode<ItemType>* removeNode(BinaryNode<ItemType>* nodePtr);
    // Removes the 1eftmost node in the left subtree of the node
    // pointed to by nodePtr.
    // Sets inorderSuccessor to the value in this node.
    // Returns a pointer to the revised subtree.
    BinaryNode<ItemType>* removeLeftmostNode(
                        BinaryNode<ItemType>* subTreePtr, ItemType\& inorderSuccessor);
```

    // Returns a pointer to the node containing the given value,
    // or nullptr if not found.
    BinaryNode<ItemType>* findNode(BinaryNode<ItemType>* treePtr,
                                    const ItemType& target) const;
    public:
//--
// Constructor and Destructor Section.
//-----------------
BinarySearchTree(const ItemType\& rootItem);
BinarySearchTree(const BinarySearchTree<ItemType>\& tree);
virtual ~BinarySearchTree();
//----------------------------
//----------------
int getHeight() const;
int getNumberOfNodes() const;
ItemType getRootData() const throw(PrecondViolatedExcep);
void setRootData(const ItemType\& newData) const
throw(PrecondViolatedExcep);
bool add(const ItemType\& newEntry);
bool remove(const ItemType\& anEntry);
void clear();
ItemType getEntry(const ItemType\& anEntry) const
throw(NotFoundException);
bool contains(const ItemType\& anEntry) const;
//-------------------------------------------------------------------
// Public Traversals Section.
//------------------------------------------------------------------
void preorderTraverse(void visit(ItemType\&)) const;
void inorderTraverse(void visit(ItemType\&)) const;
void postorderTraverse(void visit(ItemType\&)) const;
//------------------------------------
// O
BinarySearchTree<ItemType>\&
operator=(const BinarySearchTree<ItemType>\& rightHandSide);
}; // end BinarySearchTree
\#include "BinarySearchTree.cpp"
\#endif

```

\subsection*{16.4 Saving a Binary Search Tree in a File}

Imagine a program that maintains the names, addresses, and telephone numbers of your friends and relatives. While the program is running, you can enter a name and get the person's address and phone number. If you terminate program execution, the program must save its database of people in a form that it can recover at a later time.

Use a preorder traversal to save a binary search tree in a file; restore it to its original shape by using add

A balanced binary search tree increases the efficiency of the ADT operations

Building a full binary search tree

FIGURE 16-15 An initially empty binary search tree after the insertion of \(60,20,10,40,30,50\), and 70


If the program uses a binary search tree to represent the database, it must save the tree's data in a file so that it can later restore the tree. Two different algorithms for saving and restoring a binary search tree will be considered here. The first algorithm restores a binary search tree to its original shape. The second restores a binary search tree to a shape that is balanced.

Saving a binary search tree and then restoring it to its original shape. The first algorithm restores a binary search tree to exactly the same shape it had before it was saved. For example, consider the tree in Figure 16-15. If you save the tree in preorder, you get the sequence \(60,20,10,40,30,50,70\). If you then use the add method to insert these values into a binary search tree that is initially empty, you will get the original tree.

Saving a binary search tree and then restoring it to a balanced shape. After you save a binary search tree in a file, do you necessarily want the restored tree to have its original shape? Recall that you can use a given set of data items to create several binary search trees with different shapes. Although the shape of a binary search tree has no effect whatsoever on the correctness of the ADT operations, it will affect the efficiency of those operations. Efficient operations are ensured if the binary search tree is balanced.

The algorithm that restores a binary search tree to a balanced shape is surprisingly simple. In fact, you can even guarantee a restored tree of minimum height-a condition stronger than balanced. To gain some insight into the solution, consider a full tree, because it is balanced. If you save a full tree in a file by using an inorder traversal, the file will be in sorted order, as Figure 16-16 illustrates.

A full tree with exactly \(n=2^{h}-1\) nodes for some height \(h\) has the exact middle of the data items in its root. The left and right subtrees of the root are full trees of \(2^{h-1}-1\) nodes each (that is, half of \(n-1\), as \(n\) is odd or, equivalently, \(n / 2\) ). Thus, you can use the following recursive algorithm to create a full binary search tree with \(n\) nodes, provided you either know or can determine \(n\) beforehand.
```

// Builds a full binary search tree from $n$ sorted values in a file.
// Returns a pointer to the tree's root.
readFullTree(n: integer): BinaryNodePointer
if ( $\mathrm{n}>0$ )
\{
// Get the root
treePtr $=$ pointer to new node with nu11ptr as its child pointers
rootItem = next item from file
treePtr->setItem(rootItem)

```

FIGURE 16-16 A full tree saved in a file by using inorder traversal

\[
\begin{array}{lllllll}
10 & 20 & 25 & 30 & 40 & 50 & 60
\end{array}
\]
```

    // Construct the left subtree
    leftPtr = readFul1Tree(treePtr->getLeftChildPtr(), n / 2)
    treePtr->setLeftChildPtr(7eftPtr)
    // Construct the right subtree
    rightPtr = readFul1Tree(treePtr->getRightChildPtr(), n / 2)
    treePtr->setRightChildPtr(rightPtr)
    return treePtr
    }
else
return nul1ptr

```

Surprisingly, you can construct the tree directly by reading the sorted data sequentially from the file.

This algorithm for building a full binary search tree is simple, but what can you do if the tree to be restored is not full-that is, if it does not have \(n=2^{h}-1\) nodes for some \(h\) ? The first thing that comes to mind is that the restored tree should be complete-full up to the last level, with the last level filled in from left to right. Actually, because you care only about minimizing the height of the restored tree, it does not matter where the nodes on the last level go, as Figure 16-17 shows.

The method readFul1Tree is essentially correct even if the tree is not full. However, you do have to be a bit careful when computing the sizes of the left and right subtrees of the tree's root. If \(n\) is odd, both subtrees are of size \(n / 2\), as before, and the root is automatically accounted for. If \(n\) is even, however, you have to deal with the root and the fact that one of the root's subtrees will have

FIGURE 16-17 A tree of minimum height that is not complete


Building a minimumheight binary search tree
one more node than the other. In this case, you can arbitrarily choose to put the extra node in the left subtree. The following algorithm makes these compensations:
```

// Builds a minimum-height binary search tree from n sorted values in a file
// Returns a pointer to the tree's root.
readTree(n: integer): BinaryNodePointer
if (n > 0)
{
// Get the root
treePtr = pointer to new node with nul1ptr as its child pointers
rootItem = next item from file
treePtr->setItem(rootItem)
// Construct the left subtree
1eftPtr = readFul1Tree(treePtr->getLeftChildPtr(), n / 2)
treePtr->setLeftChildPtr(1eftPtr)
// Construct the right subtree
rightPtr = readFul1Tree(treePtr->getRightChildPtr(), (n - 1) / 2)
treePtr->setRightChildPtr(rightPtr )
return treePtr
}
else
return nullptr

```

You should trace this algorithm and convince yourself that it is correct for both even and odd values of \(n\).

To summarize, you can easily restore a tree as a balanced binary search tree if the data is sorted-that is, if it has been produced from the inorder traversal-and you know the number \(n\) of nodes in the tree. You need \(n\) so that you can determine the middle item and, in turn, the number of nodes in the left and right subtrees of the tree's root. Knowing these numbers is a simple matter of counting nodes as you traverse the tree and then saving the number in a file that the restore operation can read.

Note that readTree would be an appropriate protected method of BinarySearchTree, if you also had a public method to call it.

Note: You can use the readTree algorithm to restore an out-of-balance binary search tree. Exercise 18 asks you to try it using an array instead of a file.

Question 9 Consider the pseudocode operation readTree.
a. What binary search tree results when you execute readTree with a file of the six integers \(2,4,6,8,10,12\) ?
b. Is the resulting tree's height a minimum? Is the tree complete? Is it full?

\subsection*{16.5 Tree Sort}

You can use the ADT binary search tree to sort an array efficiently. To simplify the discussion, we will sort an array of integers into ascending order, as we did with the sorting algorithms in Chapter 11.

The basic idea of the algorithm is simple:
> // Sorts the integers in an array into ascending order.
> treeSort(anArray: array, n: integer)
> Insert anArray's entries into a binary search tree bst
> Traverse bst in inorder. As you visit bst's nodes, copy their data items into successive locations of anArray

An inorder traversal of the binary search tree bst visits the integers in bst's nodes in ascending order.

A tree sort can be quite efficient. As Figure 15-17 in Chapter 15 indicates, each insertion into a binary search tree requires \(\mathrm{O}(\log n)\) operations in the average case and \(\mathrm{O}(n)\) operations in the worst case. Thus, tree sort's \(n\) insertions require \(\mathrm{O}(n \times \log n)\) operations in the average case and \(\mathrm{O}\left(n^{2}\right)\) operations in the worst case. The traversal of the tree involves one copy operation for each of the \(n\) entries and so is \(\mathrm{O}(n)\). Because \(\mathrm{O}(n)\) is less than \(\mathrm{O}(n \times \log n)\) and \(\mathrm{O}\left(n^{2}\right)\), tree sort in the average case is \(\mathrm{O}(n \times\) \(\log n)\) and \(\mathrm{O}\left(n^{2}\right)\) in the worst case.

2
Question 10 Trace the tree sort algorithm as it sorts the following array into ascending order: 208040256030 .

\subsection*{16.6 General Trees}

This chapter ends with a brief discussion of general trees and their relationship to binary trees. Consider the general tree in Figure 16-18. The three children \(B, C\), and \(D\) of node \(A\), for example, are siblings. The leftmost child \(B\) is the oldest child, or first child, of \(A\), as mentioned in Section 15.1 of Chapter 15. One way to implement this tree uses the BinaryNode objects that we used for a link-based binary tree. That is, each node has two pointers: The left pointer points to the node's oldest child and the right pointer points to the node's next sibling. Thus, you can use the data structure in Figure 16-19a to implement the tree in Figure 16-18. Notice that the structure in Figure 16-19a also represents the binary tree pictured in Figure 16-19b.

Recall from Chapter 15 that an \(n\)-ary tree is a generalization of a binary tree whose nodes each can have no more than \(n\) children. The tree in Figure 16-18 could be an \(n\)-ary tree with \(n=3\) instead of a general tree. You can, of course, use the implementation just described for an \(n\)-ary tree. However, because you know the maximum number of children for each node, you can let each node point directly to its children. Figure 16-20 illustrates such a representation for the tree in Figure 16-18. This tree is shorter than the tree in Figure 16-19b.

FIGURE 16-18 A general tree


FIGURE 16-19 (a) A link-based implementation of the general tree in Figure 16-18; (b) the binary tree that part \(a\) represents
(a)

(b)


FIGURE 16-20 An implementation of the \(n\)-ary tree in Figure 16-18


Exercise 17 discusses general trees further.

\section*{Summary}
1. The implementation of a binary tree is usually link based. However, if the binary tree is complete, an efficient array-based implementation is possible.
2. A node in a binary tree is an object that references a data entry and two child nodes.
3. Because a binary tree has a recursive nature, recursion is useful in the implementation of its operations. As is typical, the recursive methods are not public, since they require parameters that are a part of the underlying data structure. Those methods can be either private or protected, but as protected methods, they are available to classes derived from the class of binary trees.
4. The ADT binary search tree has several methods that are the same as those in the ADT binary tree. However, its insertion, removal, and retrieval operations have different definitions than those for the ADT binary tree.
5. Operations on a binary search tree can be quite efficient. In the worst case, however-when the tree approaches a linear shape - the performance of its operations degrades and is comparable to that of a linear linked chain. If you must avoid such a situation for a given application, you should use the balancing methods presented in Chapter 19.
6. The tree sort algorithm efficiently sorts an array by using the binary search tree's insertion and traversal operations.
7. If you save a binary search tree's data in a file while performing an inorder traversal of its nodes, you can restore the tree as a binary search tree of minimum height. If you save a binary search tree's data in a file while performing a preorder traversal of its nodes, you can restore the tree to its original form
8. You can use a binary tree to represent a general tree or an \(n\)-ary tree. However, an \(n\)-ary tree can have nodes that point to all of its children.

\section*{Exercises}
1. Consider the binary search tree in Figure \(15-20\) of Chapter 15 . The numbers simply label the nodes so that you can reference them; they do not indicate the contents of the nodes.
a. Which node must contain the inorder successor of the value in the root? Explain.
b. Which node must contain the inorder predecessor of the value in the root? Explain.
2. Arrange nodes that contain the letters \(A, C, E, F, L, V\), and \(Z\) into two binary search trees: one that has maximum height and one that has minimum height.
3. Consider the binary search tree in Figure 15-18 of Chapter 15.
a. What tree results after you insert the entries \(80,65,75,45,35\), and 25 , in that order?
b. After inserting the nodes mentioned in part \(a\), what tree results when you remove the entries 50 and 20 ?
4. Consider the binary search tree in Figure 15-19 of Chapter 15. What does the tree look like after you remove \(M\), \(D, G\), and \(T\), in that order?
5. If you remove an item from a binary search tree and then insert it back into the tree, will you ever change the shape of the tree?
6. Suppose that the ADT binary tree has the operation
replace(item: ItemType, replacementItem: ItemType): boolean
It locates, if possible, the node in a binary tree that contains item and replaces item with replacementItem.
a. Add the operation replace to the link-based implementation of the ADT binary tree given in this chapter. The operation should replace an item without altering the tree structure.
b. Add the operation replace to the link-based implementation of the ADT binary search tree. Be sure that the tree remains a binary search tree.
c. Implement a method replace within a client of BinarySearchTree.
7. Consider an array-based implementation of a binary search tree bst. Figure 16-1 presents such a representation for a particular binary search tree.
a. Depict the array in an array-based implementation for the binary search tree in Figure 15-14a of Chapter 15. Assume that tree items are strings.
b. Show the effect of each of the following sequential operations on the array in part \(a\) of this exercise.
```

bst.add("Doug");
bst.add("Nancy");
bst.add("Bob");
bst.add("Sarah");

```
c. Repeat parts \(a\) and \(b\) of this exercise for the tree in Figure 15-14b.
d. Write an inorder traversal algorithm for this array-based implementation.
8. Duplicates in an ADT could mean either identical items or, more subtly, objects that match because their class overloads the equality operator so that it compares only certain fields, while differences exist in other fields. If duplicates are allowed in a binary search tree, it is important to have a convention that determines the relationship between the duplicates. Items that duplicate the root of a tree should either all be in the left subtree or all be in the right subtree, and, of course, this property must hold for every subtree.
a. Why is this convention critical to the effective use of the binary search tree?
b. This chapter stated that you can remove an item from a binary search tree by replacing it with the item that either immediately follows or immediately precedes the item to be deleted. If duplicates are allowed, however, the choice between inorder successor and inorder predecessor is no longer arbitrary. How does the convention of putting duplicates in either the left or right subtree affect this choice?
9. Complete the trace of the nonrecursive inorder traversal algorithm that Figure \(16-4\) began. Show the contents of the implicit stack as the traversal progresses.
10. Implement in C++ the nonrecursive inorder traversal algorithm for a binary tree that was presented in this chapter.
11. Exercise 13 in Chapter 15 introduced the minimax tree. It has min nodes and max nodes. Part \(b\) of the exercise asked you to design a solution for representing and evaluating these trees. Implement your design in \(\mathrm{C}++\).
12. Design another algorithm to remove items from a binary search tree. This algorithm differs from the one described in this chapter when a node \(N\) has two children. First let \(N\) 's right child take the place of the deleted node \(N\) in the same manner in which you delete a node with one child. Next reconnect \(N\) 's left child (along with its subtree, if any) to the left side of the node containing the inorder successor of the value in \(N\).
13. Write iterative methods to perform insertion and removal operations on a binary search tree.
14. A level-order traversal of a tree processes (visits) nodes one level at a time, from left to right, beginning with the root. Design an algorithm that performs a level-order traversal of a binary tree.
15. If you know in advance that you often access a given item in a binary search tree several times in succession before accessing a different item, you will end up searching for the same item repeatedly. One way to avoid this problem is to add an extra bookkeeping component to your implementation. That is, you can maintain a lastaccessed pointer that will always reference the last item that any binary search tree operation accessed. Whenever you perform such an operation, you can check the search key of the item most recently accessed before performing the operation.

Revise the implementation of the ADT binary search tree to add this new feature by adding the data member lastAccessed to the class.
16. Exercise 10 in Chapter 4 introduced the doubly linked chain. The analogy for a binary search tree is to maintain parent pointers in each binary node in addition to the pointers to the node's children. That is, every node except the root will have a pointer to its parent in the tree. This type of binary tree is called a doubly linked binary tree. Write insertion and removal operations for this tree.
17. A node in a general tree can have an arbitrary number of children.
a. Describe a C++ implementation of a general tree in which every node contains an array of child pointers. Write a recursive preorder traversal method for this implementation. What are the advantages and disadvantages of this implementation?
b. Consider the implementation of a general tree that is illustrated in Figure 16-19. Each node has two pointers: The left pointer points to the node's oldest child and the right pointer points to the node's next sibling. Write a recursive preorder traversal method for this implementation.
c. Every node in a binary tree \(T\) has at most two children. Compare the oldest-child/next-sibling representation of \(T\), as part \(b\) describes, to the left-child/right-child representation of a binary tree, as this chapter describes. Does one representation simplify the implementation of the ADT operations? Are the two representations ever the same?
18. Given an unbalanced binary search tree, use an inorder traversal to copy its data to an array. Then create a balanced binary search tree using the readTree algorithm given in Section 16.4, but use your array instead of a file.

Add an overloaded \(==\) operator to the class BinaryNodeTree.

\section*{Programming Problems}
1. Complete the implementation of the class BinaryNodeTree that was begun in Section 16.2.2 of this chapter.
2. Implement the class BinarySearchTree, as given in Listing 16-4.
3. Write an array-based implementation of the ADT binary tree that uses dynamic memory allocation. Use a data structure like the one in Figure 16-1.
4. Repeat the previous problem, but define a binary search tree instead.
5. Write a program that maintains a database containing data, such as name and birthday, about your friends and relatives. You should be able to enter, remove, modify, or search this data. Initially, you can assume that the names are unique. The program should be able to save the data in a file for use later.

Design a class to represent the database and another class to represent the people. Use a binary search tree of people as a data member of the database class.

You can enhance this problem by adding an operation that lists everyone who satisfies a given criterion. For example, you could list people born in a given month. You should also be able to list everyone in the database.
*6. Implement the ADT queue operations as well as a sorted traversal operation for a queue that points into a doubly linked binary search tree, as shown in Figure 16-21. Doubly linked binary trees are explained in Exercise 16. You will need the insertion and removal operations for a binary search tree that contains parent pointers.

FIGURE 16-21 A queue that points into a doubly linked binary search tree


\section*{C++}

\section*{Iterators}

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\section*{Prerequisites}
\begin{tabular}{ll} 
Chapter 9 & List Implementations \\
C++ Interlude 5 & Overloaded Operators and Friend Access
\end{tabular}

\(\mathrm{A}_{\mathrm{n}}\)n iterator is an object that traverses a collection of data. During the traversal, you can look at the data entries and process them. You can compare iterators to see if they point to the same entry in a collection, and C++ provides special functions that let you perform other actions with collections, such as searching and counting, when using iterators. Iterators are commonly implemented by overloading operators that correspond to the functions an iterator performs. A container can instantiate an iterator and give it to a client to use to traverse the contained data collection.


C++ iterators

An iterator is an object that traverses a collection of like objects

\section*{C6.1 Iterators}

How would you count the number of lines on this page? You could use your finger to point to each line as you counted it. Your finger would keep your place on the page. If you paused at a particular line, your finger would be on the current line, and there would be a previous line and a next line. If you think of this page as a list of lines, you would be traversing the list as you counted the lines.

An iterator is a program component that enables you to traverse a collection of data, such as the data in a list, beginning with the first entry. During one complete traversal, or iteration, each data item is considered once. You control the progress of the iteration by repeatedly asking the iterator to give you a reference to the next entry in the
collection. You also can modify the collection as you traverse it by adding, removing, or simply changing entries.

You are familiar with iteration because you have written loops. For example, if nameList is a list of strings, you can write the following for loop to display the entire list:
```

int listSize = nameList.getLength();
for (int position = 1; position <= listSize; position++)
cout << nameList.getEntry(position) << endl;

```

Here the loop traverses, or iterates, through the entries in the list. Instead of simply displaying each entry, we could do other things to or with it.

Notice that the previous loop is at the client level, since it uses the ADT operation getEntry to access the list. For an array-based implementation of the list, getEntry can retrieve the desired array entry directly and quickly. But if a chain of linked nodes represents the list's entries, getEntry must move from node to node until it locates the desired one. For example, to retrieve the \(n^{\text {th }}\) entry in the list, getEntry would begin at the first node in the chain and then move to the second node, the third node, and so on until it reached the \(n^{\text {th }}\) node. At the next repetition of the loop, getEntry would retrieve the \(n+1^{\text {st }}\) entry in the list by beginning again at the first node in the chain and stepping from node to node until it reached the \(n+1^{\text {st }}\) node. This wastes time.

\section*{Note: Iterators}

An iterator is a program component that steps through, or traverses, a collection of data. The iterator keeps track of its progress during the traversal, or iteration. It can tell you whether a next entry exists and, if so, return a reference to it. During one cycle of the iteration, each data item is considered once.

Iteration is such a common operation that we could include it as part of the ADT list. Doing so would enable a more efficient implementation than you could achieve at the client level. Notice that the operation toVector of the ADT bag in Chapters 3 and 4 performs a traversal and is an example of a traversal controlled by the ADT. A client can invoke toVector but cannot control its traversal once it begins.

But toVector only returns the list's entries. What if we want to do something else with them as we traverse them? We do not want to add another operation to the ADT each time we think of another way to use an iteration. We need a way for a client to step through a collection of data and retrieve or modify the entries. The traversal should keep track of its progress; that is, it should know where it is in the collection and whether it has accessed each entry. An iterator provides such a traversal.

C++ provides a rich library of tools to build iterators, but in this interlude we focus on the capabilities necessary to implement simple iterators that traverse a collection of items, retrieve an item in a collection, and compare two iterators to determine whether they access the same entry in the collection. Such iterators are known as input iterators.

\section*{C6.1.1 Common Iterator Operations}

Though not required, iterators in \(\mathrm{C}++\) typically provide basic functionality by using overloaded operators. An iterator operation that accesses the item it currently references is usually implemented by overloading the \(\mathrm{C}++\) dereferencing operator *. For example, you can define the * operator for an iterator \(i\) so that \(* i\) is the item that \(i\) references.

Iterators also have operations that move the iterator forward through the collection. Often, these operations take the form of the overloaded operator ++ . Bidirectional iterators also overload the --
operator to allow you to move backward through the collection. Lastly, the operators \(==\) and \(!=\) are usually overloaded to compare iterators for equality.

\[
\begin{array}{ll}
\text { Note: Common iterator operations } \\
\text { Operation } & \begin{array}{l}
\text { Description } \\
\text { * }
\end{array} \\
++ & \begin{array}{l}
\text { Return the item that the iterator currently references } \\
\text { Move the iterator to the next item in the collection }
\end{array} \\
-- & \begin{array}{l}
\text { Move the iterator to the previous item in the collection (used only } \\
\text { for bidirectional or random iterators) }
\end{array} \\
== & \begin{array}{l}
\text { Compare two iterators for equality }
\end{array} \\
!= & \text { Compare two iterators for inequality }
\end{array}
\]

By enabling you to overload these operators for your iterator class, \(\mathrm{C}++\) gives you the advanced iterator operations covered in the next section without further programming. An additional step you must take to enable this functionality, however, is to derive your iterator class from the C ++ template class iterator. This template does not have any methods. Instead, it is used to identify the category of iterator you are creating by using an iterator category tag as the template type.

For example, to declare an input iterator for the class LinkedList, we would use the lines:
```

template <class ItemType>
class LinkedIterator : public iterator<input_iterator_tag, int>

```
where input_iterator_tag indicates that this iterator implements input-iterator functionality. The second template type specified, int, identifies the type of value used to measure the distance between two iterators. These values are normally integers. The distance between two iterators is the number of elements or positions between the current positions of the two iterators. For example, if an iterator points to the first item in a list and another iterator points to the fifth item in the same list, their distance is 4.

\section*{Note: C++ iterator categories}

All of the following iterators provide operations that copy or assign (=) and increment (++).
\begin{tabular}{|c|c|c|}
\hline Category & Tag & Operation \\
\hline Input iterator & input_iterator_tag & Equality/inequality (==, !=), access collection entry (*) \\
\hline Output iterator & output_iterator_tag & Change a collection entry (*) \\
\hline Forward iterator & forward_iterator_tag & Same as the input and output iterators and has a default constructor \\
\hline Bidirectional iterator & bidirectional_iterator_tag & Same as the forward iterator, but also can traverse the collection backward (--) \\
\hline Random-access iterator & random_iterator_tag & Same as the bidirectional iterator and adds support for arithmetic (+, -, +=, \(-=\) ) and relational ( \(<,<=,>,>=\) ) operations between iterators. Supports the [] operator to directly access collection entries. \\
\hline
\end{tabular}

\section*{C6.1.2 Using Iterator Operations}

Before discussing the details of a class of iterators and its implementation, we should examine how clients of LinkedList would use these iterators. Standard containers in C++ implement two special methods, begin and end, that return an iterator to the first entry and last entry respectively. These iterators should have an order of magnitude in performance of \(\mathrm{O}(1)\) as they move from entry to entry. Let's see how a client could use these methods if we implemented them in our LinkedList class.

Up to this point, if we want to display all the entries in a list myList, we need to perform this \(\mathrm{O}\left(n^{2}\right)\) action:
```

int currentPosition = 1;
while (currentPosition <= myList.getLength())
{
cout << myList.getEntry(currentPosition); // O(n) operation
currentPosition++;
} // end while

```

Now assume that LinkedIterator is a class of iterators for LinkedList, and LinkedList defines the methods begin and end and declares them as follows:

LinkedIterator<ItemType> begin() const;
LinkedIterator<ItemType> end() const;
We can rewrite the above snippet of code using LinkedIterator objects:
```

LinkedIterator<ItemType> currentIterator = myList.begin();
while (currentIterator != myList.end())
{
cout << *currentIterator // O(1) operation
++currentIterator;
} // end while

```

The first line creates a LinkedIterator object, currentIterator, that points to the first entry in the list. The while-loop test checks to see whether currentIterator points to the entry that is at the end of the list. Then, for each entry in the list, we dereference the iterator using the * operator to access and display the entry at the iterator's current position. The iterator is then incremented to reference the next entry in the list for the following iteration.

Notice that the ++ operator precedes the name of the iterator. We use this prefix operator to differentiate it from an arithmetic increment of an integer, which uses the postfix ++ operator. Iterators may implement the postfix ++ operator, but we have chosen not to do this to avoid possible confusion in reading the code.

Note: Using the ++ and -- operators
In C++, the operators ++ and -- can be used in either prefix or postfix form to represent incrementing or decrementing the corresponding variable. To differentiate between incrementing or decrementing arithmetic variables from iterator positions, we use ++ and -- as postfix operators for the former operations and as prefix operators for the latter.

The implementation of the methods begin and end in the LinkedList class are straightforward:
```

template <class ItemType>
LinkedIterator<ItemType> LinkedList<ItemType>::begin()
{
return LinkedIterator<ItemType>(this, headPtr);
} // end begin

```
```

template <class ItemType>
LinkedIterator<ItemType> LinkedList<ItemType>::end()
{
return LinkedIterator<ItemType>(this, nu11ptr);
} // end end

```

In the begin method, we send the iterator references to the list object and the first node in the chain that contains the list items. Giving the iterator direct access to LinkedList's data member violates the wall of abstraction and must be done with care. Well-designed iterators only traverse the collection and access the items stored in the collection; they should not change the structure of a collection.

\section*{C6.1.3 Implementing an Iterator}

Listing C6-1 is the header file for the class, LinkedIterator, of input iterators designed to work with our LinkedList class. It is a distinct class separate from LinkedList. The constructor has two parameters, the list traversed by the iterator and an initial node for the iterator to reference. The second parameter is used by the LinkedList class methods begin and end to initialize the iterator's position.

\section*{LISTING C6-1 The header file for the class LinkedIterator}
```

\#ifndef _LINKED_ITERATOR
\#define _LINKED_ITERATOR

```
\#include <iterator>
\#include "Node.h"
temp1ate<c1ass ItemType>
class LinkedList;
template <class ItemType>
class LinkedIterator : pub7ic iterator<input_iterator_tag, int>
\{
private:
    // ADT associated with iterator
    const LinkedList<ItemType>* containerPtr;
    // Current location in collection
    Node<ItemType>* currentItemPtr;
public:
    LinkedIterator (const LinkedList<ItemType>* someList,
                    Node<ItemType>* nodePtr);
    /** Dereferencing operator overload.
        @return The item at the position referenced by iterator. */
    const ItemType operator*() ;
    /** Prefix increment operator overload.
        @return The iterator referencing the next position in
            the 1ist. */
    LinkedIterator<ItemType> operator++();
    /** Equality operator overload.
    @param LinkedList The iterator for comparison.
    @return True if this iterator references the same list and
            the same position as rightHandSide, false otherwise. * /
```

    bool operator==(const LinkedIterator<ItemType>& rightHandSide) const;
    /** Inequality operator overload.
    @param LinkedList The iterator for comparison.
    @return True if this iterator does not reference the same
        list and the same position as rightHandSide,
        false otherwise. */
    bool operator!=(const LinkedIterator<ItemType>& rightHandSide) const;
    }; // end LinkedIterator
\#include "LinkedIterator.cpp"
\#endif

```

In addition to the base class specification and class constructor, there are two other lines in this header file that differ from previous header files. Before the LinkedIterator class is declared, the following two lines provide a forward declaration of the class LinkedList:
```

template<class ItemType>
class LinkedList;

```

These lines are necessary because LinkedList must instantiate an object of the class LinkedIterator<ItemType> in the begin and end methods, but the LinkedIterator class must have a complete declaration of LinkedList<ItemType> to instantiate itself. To resolve these circular references, we include this forward declaration. When the compiler begins to compile the LinkedIterator class, it will read these statements and know that LinkedList<ItemType> is also a class. The compiler can then use placeholders for the details of LinkedList.

The implementation of LinkedIterator is shown in Listing C6-2.

\section*{LISTING C6-2 The implementation file for the class LinkedIterator}
```

\#include "LinkedIterator.h"
template <class ItemType>
LinkedIterator<ItemType>::
LinkedIterator(const LinkedList<ItemType>* someList,
Node<ItemType>* nodePtr):
containerPtr(someList), currentItemPtr(nodePtr)
{
} // end constructor
template <class ItemType>
const ItemType LinkedIterator<ItemType>::operator*()
{
return currentItemPtr->getItem();
} // end operator*
template <class ItemType>
LinkedIterator<ItemType> LinkedIterator<ItemType>: :operator++()
{

```
```

    currentItemPtr = currentItemPtr->getNext();
    return *this;
    } // end prefix operator++
template <class ItemType>
bool LinkedIterator<ItemType>::operator==(const
LinkedIterator<ItemType>\& rightHandSide) const
{
return ((containerPtr == rightHandSide.containerPtr) \&\&
(currentItemPtr == rightHandSide.currentItemPtr));
} // end operator==
template <class ItemType>
bool LinkedIterator<ItemType>::operator!=(const
LinkedIterator<ItemType>\& rightHandSide) const
{
return ((containerPtr != rightHandSide.containerPtr) ||
(currentItemPtr != rightHandSide.currentItemPtr));
} // end operator!=

```

The implementations of the iterator methods use the methods of the Node class. For example, the dereferencing * operator returns the result of calling getItem on the current node, and the ++ operator uses getNext to advance the iterator to the next position. Notice that the ++ operator returns the current object to the caller. This is part of the \(\mathrm{C}++\) standard for the ++ operator, but clients seldom use the return value.

An alternative is to use only iterators that have access to public methods of your data structure, but then you lose the efficiencies gained by directly accessing the structure. The best approach is to design an iterator for your class at the same time that you design your ADT, so that you can coordinate the features and ensure that the iterator does not change the class's structure.

\section*{C6.2 Advanced Iterator Functionality}

Deriving our iterator from the C++ template class iterator enables the iterator to be used in a number of standard functions provided by \(\mathrm{C}++\) that simplify the processing of many common algorithms, such as displaying the items in a collection, searching a collection, and counting the number of occurrences of an item in the collection.
The function for_each. Earlier in this interlude, we presented a whi 1 e loop that used iterators to access and display each item in a list of strings. C++ provides a function for_each that can accomplish this in a simpler manner. The syntax of the for_each function is:
```

for_each(start_iterator, end_iterator, function_to_perform);

```
where start_iterator and end_iterator are iterators to the first and last position that should be processed in the collection, and function_to_perform is a function with a single parameter of the same type as the entries stored in the collection. This is similar to the visit function discussed with respect to ADT binary tree traversals in Chapter 15. This function is applied to each of the items in the collection from start_iterator to end_iterator.

To use the for_each function, we must create a function to be applied to each of the items in the collection. In this case, the items are strings, and we want to display each one:
```

void displayOneItem(string itemToDisplay)
{
cout << itemToDisplay << endl;
} // end displayOneItem

```

We can then call the for_each function to display all items in the collection as follows:
```

for_each(myList.begin(), myList.end(), displayOneItem);

```

With this one statement, \(\mathrm{C}++\) takes care of all the details. The compiler will insert code that begins at the first position of myList and sends each item to the function displayOneItem to be displayed until all items have been processed.

(
Note: The syntax used to call most of the special functions that use iterators to process collections follows the form
```

function_name(start_iterator, end_iterator, special_parameters);

```
where special_parameters represents one or more parameters that the function needs to perform its task.

The function find. To search a collection for a specified target entry, you use the function find. This function returns an iterator to the target if it exists in the collection. If the target is not in the collection, find returns an iterator to the last position:
```

// Find the entry "Ace" in the col1ection myList
LinkedIterator<string> myCard = find(myList.begin(), myList.end(), "Ace");

```

The function count. Another useful function is count. This function returns an integer that represents the number of times a specified value occurs in the collection between the starting iterator and ending iterator:
```

// Count the number of occurrences of "Ace" in collection myList
int aceCount = count(myList.begin(), myList.end(), "Ace");

```

The function advance. To move an iterator forward a specific number of positions, you can use the advance function. It has a slightly different form:
```

advance(someIterator, distanceToAdvance);

```

For example, if we needed an iterator to the fifth element of myList, we would use the following statements:
```

// Iterator to first position in myList
LinkedIterator<string> myIterator = myList.begin();
// Advance from first position to fifth (4 positions ahead)
advance(myIterator, 4);

```

The function distance. To determine how many entries remain to be processed, we can determine the distance from our current position to the end of the collection, using the distance function:
```

int numberRemaining = distance(myIterator, myList.end());

```

The function equa1. If you want to determine whether the elements in two lists are the same, either for the entire first list or for a subrange within the lists, you can use the equal function. Suppose myList and yourList are two instances of LinkedList<string>. To determine if all entries in myList are equal to the first entries in yourList, you would write:
```

bool same = equa1(myList.begin(), myList.end(), yourList.begin());

```

This function is concerned only with a range of positions. The number of entries to compare, \(n\), is determined by the distance between the first two iterator parameters. The third iterator parameter specifies a starting location in the second collection for the comparison. The number of entries in the second collection can be greater than the number in the first collection, but only \(n\) entries are compared.

\section*{Note: Some useful C++ functions that use iterators}
- Process entries in a collection from start_iterator position to end_iterator position using the function function_to_perform:
```

for_each(start_iterator, end_iterator, function_to_perform);

```
- Return an iterator to the position of the first occurrence of target between start_ iterator and end_iterator:
iteratorType someIterator \(=\) find(start_iterator, end_iterator, target);
- Return the number of occurrences of target between start_iterator and end_iterator:
```

int numberOccur = count(start_iterator, end_iterator, target);

```
- Compare entries in collection 1 from startl_iterator through endl_iterator to those in collection 2 beginning at start2_iterator:
```

boo1 result = equa1(start1_iterator, end1_iterator, start2_iterator);

```
- Move someIterator from its current position forward distanceToAdvance positions: advance(someIterator, distanceToAdvance);
- Determine the distance or number of positions from someIterator to anotherIterator: int theDistance \(=\) distance \((\) someIterator, anotherIterator \()\);

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\section*{Heaps}

\section*{Chapter}

\section*{Contents}

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\section*{Prerequisites}

Chapter 13 Queues and Priority Queues
Chapter 15 Trees
Chapter 16 Tree Implementations

Chapter 15 introduced the ADTs binary tree and binary search tree. This chapter will present the ADT heap, which is a special kind of complete binary tree that provides an efficient implementation of the ADT priority queue. You can also use a heap to sort an array as efficiently as you can by using either a quick sort or a merge sort.

\subsection*{17.1 The ADT Heap}

A heap is a complete binary tree that either is empty or whose root

A heap is a special complete binary tree
- Contains a value greater than or equal to the value in each of its children, and
- Has heaps as its subtrees

VideoNote
The ADT heap

A heap is similar to a binary search tree, although it differs from a binary search tree in two significant ways:
- While you can view a binary search tree as sorted, a heap is ordered in a much weaker sense.
- While binary search trees come in many different shapes, heaps are always complete binary trees.

In our definition of a heap, the root contains the item with the largest value. Such a heap is also known as a maxheap. A minheap, on the other hand, places the item with the smallest value in its root. An example of a maxheap and a minheap are given in Figure 17-1. Exercises 1 and 8 and Programming Problem 3 consider the minheap further.

FIGURE 17-1 (a) A maxheap and (b) a minheap


The ADT heap has the following operations:

Note: ADT heap operations
- Test whether a heap is empty
- Get the number of nodes in a heap
- Get the height of a heap
- Get the item in the heap's root
- Insert a new item into the heap
- Remove the item in the heap's root
- Remove all nodes from the heap

Depending on whether the heap is a maxheap or a minheap, the removal and retrieval operations respectively involve either the largest or the smallest item in the heap

FIGURE 17-2 UML diagram for the class Heap
\begin{tabular}{|l|}
\hline \multicolumn{1}{|c|}{ Heap } \\
\hline \\
\hline +isEmpty(): boolean \\
+getNumberOfNodes(): integer \\
+getHeight(): integer \\
+peekTop(): ItemType \\
+add(newData: ItemType) : boolean \\
+remove(): boolean \\
+clear(): void
\end{tabular}

The following contract specifies these operations in more detail, and a UML diagram for a class of heaps appears in Figure 17-2.

\section*{Abstract Data Type: Heap}

\section*{Data}
- A finite number of objects in hierarchical order.

\section*{Operations}
\begin{tabular}{|c|c|}
\hline Pseudocode & Description \\
\hline isEmpty () & \begin{tabular}{l}
Task: Sees whether this heap is empty. \\
Input: None. \\
Output: True if the heap is empty; otherwise false.
\end{tabular} \\
\hline getNumberOfNodes() & \begin{tabular}{l}
Task: Gets the number of nodes in this heap. Input: None. \\
Output: The number of nodes in the heap.
\end{tabular} \\
\hline getHeight() & \begin{tabular}{l}
Task: Gets the height of this heap. Input: None. \\
Output: The height of the heap.
\end{tabular} \\
\hline peekTop() & \begin{tabular}{l}
Task: Gets the data that is in the root (top) of this heap. \\
Input: None. Assumes the heap is not empty. \\
Output: The item in the root of the heap. If the heap is a maxheap, this item is the largest value in the heap. For a minheap, the item is the smallest value.
\end{tabular} \\
\hline add (newData) & \begin{tabular}{l}
Task: Inserts newData into this heap. \\
Input: newData is the data item to be inserted. \\
Output: True if the insertion is successful, or false if not.
\end{tabular} \\
\hline remove() & \begin{tabular}{l}
Task: Removes the item in the root of this heap. \\
Input: None. \\
Output: True if the removal is successful, or false if not.
\end{tabular} \\
\hline clear () & \begin{tabular}{l}
Task: Removes all nodes from this heap. Input: None. \\
Output: The heap is empty.
\end{tabular} \\
\hline
\end{tabular}

An interface that finalizes the operation contract for the ADT heap appears in Listing 17-1.

LISTING 17-1 An interface for the ADT heap
```

/** Interface for the ADT heap.
@file HeapInterface.h */
\#ifndef _HEAP_INTERFACE
\#define _HEAP_INTERFACE
template<class ItemType>
class HeapInterface
{
public:

```
    /** Sees whether this heap is empty.
    @return True if the heap is empty, or false if not. */
    virtual bool isEmpty() const = 0;
    /** Gets the number of nodes in this heap.
    @return The number of nodes in the heap. */
    virtual int getNumberOfNodes() const = 0;
    /** Gets the height of this heap.
    @return The height of the heap. */
    virtual int getHeight() const = 0;
    /** Gets the data that is in the root (top) of this heap.
        For a maxheap, the data is the largest value in the heap;
        for a minheap, the data is the smallest value in the heap.
    @pre The heap is not empty.
    @post The root's data has been returned, and the heap is unchanged.
    @return The data in the root of the heap. */
    virtual ItemType peekTop() const = 0;
    /** Adds a new node containing the given data to this heap.
    @param newData The data for the new node.
    @post The heap contains a new node.
    @return True if the addition is successful, or false if not. */
    virtual bool add(const ItemType& newData) = 0;
    /** Removes the root node from this heap.
    @return True if the removal is successful, or false if not. */
    virtual bool remove() = 0;
    /** Removes all nodes from this heap. */
    virtual void clear() = 0;
}; // end HeapInterface
#endif
```



Note: Do not confuse the ADT heap with the collection of memory cells known as a heap. This memory is available for allocation to your program when you use the new operator. The heap that contains this available memory is not an instance of the ADT heap.

Question 1 Is the full binary tree in Figure 16-16 of Chapter 16 a heap? Why?

### 17.2 An Array-Based Implementation of a Heap

VideoNote
Overview of ArrayHeap

Because a heap is a binary tree, you can use the array-based implementation of a binary tree discussed in Section 16.1.1 of Chapter 16, if you know the maximum size of the heap. However, because a heap is a complete binary tree, you can use a simpler array-based implementation that saves

FIGURE 17-3 (a) Level-by-level numbering of a complete binary tree; (b) its array-based implementation

memory. As you saw in Chapter 15, a complete tree of height $h$ is full to level $h-1$ and has level $h$ filled from left to right.

Figure 17-3a shows a complete binary tree with its nodes numbered according to a level-by-level scheme. The root is numbered 0 , and the children of the root-which are at the next level of the treeare numbered, left to right, 1 and 2 . The nodes at the next level are numbered, left to right, 3,4 , and 5 . You place these nodes into the array items in numeric order. That is, items [i] contains the node numbered $i$, as Figure 17-3b illustrates. Now, given any node items [i], you can easily locate both of its children and its parent:

- Its left child, if it exists, is items [2 * i + 1]
- Its right child, if it exists, is items [2 * $\mathbf{i}+2$ ]
- Its parent, if it exists, is items [(i-1) / 2]

Remember that only the root in items [0] does not have a parent.
This array-based representation requires a complete binary tree. If nodes were missing from the middle of the tree, the numbering scheme would be thrown off, and the parent-child relationship among nodes would be ambiguous. This requirement implies that any changes to the tree must maintain its completeness.

As you will see, an array-based representation of a heap is useful in the implementation of the ADT priority queue.

If the binary tree is complete and remains complete, you can use a memory-efficient array-based implementation

Question 2 What array represents the maxheap shown in Figure 17-1a?
Question 3 What array represents the minheap shown in Figure 17-1b?
Question 4 What criterion can you use to tell whether the node in items [i] is a leaf?

## Question 5 What complete binary tree does the following array represent?

| 5 | 1 | 2 | 8 | 6 | 10 | 3 | 9 | 4 | 7 |
| :--- | :--- | :--- | :--- | :--- | :--- | :--- | :--- | :--- | :--- |
| 0 |  |  |  |  |  |  |  |  |  |

## Question 6 Does the array in the previous question represent a heap?

The data members for an array-based implementation of a heap
peekTop's logic
remove's first step produces a semiheap

### 17.2.1 Algorithms for the Array-Based Heap Operations

Let's assume that our class of heaps has the following private data members:

- items: an array of heap items
- itemCount: an integer equal to the number of items in the heap
- maxItems: an integer equal to the maximum capacity of the heap

The array items corresponds to the array-based representation of a complete binary tree, such as the array shown in Figure 17-3b. We will assume that we are working with a maxheap of integers.

Retrieving an item from a heap. First consider the heap's peekTop operation. Where is the largest item in the heap? Because the item in every tree node is greater than or equal to the item in either of its children, the largest item must be in the root of the tree-that is, at the top of the heap. Thus, the peekTop operation needs to take only the following step:

```
// Return the item in the root
return items[0]
```

Removing an item from a heap. While the peekTop operation returns the largest item in the heapwhich we know is in its root-the heap operation remove must remove it. Removing the root of the heap leaves two disjoint heaps, as Figure 17-4a indicates. Therefore, you do not want to actually remove the root. Instead, you remove the last node of the tree and place its item in the root, as Figure $17-4 \mathrm{~b}$ shows. The result of this step is not necessarily a heap. It is, however, a complete binary tree whose left and right subtrees are both heaps. The only problem is that the item in the root usually is out of place. Such a structure is called a semiheap.

You thus need a way to transform a semiheap into a heap. One strategy allows the item in the root to trickle down the tree until it reaches a node in which it will not be out of place; that is, the item will come to rest in the first node where it would be greater than or equal to the item in each of its children. To accomplish this, you first compare the item in the root of the semiheap to the items in its children. If the root's item is smaller than the larger of the items in its children, you swap the item in the root with that larger item. For convenience, we will refer to the child containing the larger item as the larger child. The result of this step is shown in Figure 17-4c. Although the value 5 trickles down in this example to its correct position after only one swap, in general more swaps may be necessary.

We need to remember that an array represents the original heap before we remove its root. The remove operation will simply move items within this array. Figure 17-5 shows the array for the heaps and semiheap illustrated in Figure 17-4. The transformation from the heap in Figure 17-5a to the semiheap in Figure $17-5 b$ is accomplished by the following steps:

[^12]FIGURE 17-4 (a) Disjoint heaps after removing the heap's root; (b) a semiheap; (c) the restored heap
(a)

(b)

(c)


The transformation from the semiheap in Figure $17-5$ b to the heap in Figure $17-5$ c is accomplished by the following recursive algorithm. Note that the algorithm will not work if it does not begin with a semiheap.

```
// Converts a semiheap rooted at index root into a heap. remove's final step
heapRebuild(root: integer, items: ArrayType, itemCount: integer)
    // Recursively trickle the item at index root down to its proper position by
    // swapping it with its larger child, if the child is larger than the item.
    // If the item is at a leaf, nothing needs to be done.
    if (the root is not a leaf)
    {
        // The root must have a left child; assume it is the larger child
        largerChildIndex = 2* rootIndex + 1 // Left child index
        if (the root has a right child)
    {
        rightChildIndex = largerChildIndex + 1 // Right child index
        if (items[rightChi1dIndex] > items[largerChildIndex ])
            7argerChildIndex = rightChildIndex // Larger child index
    }
    // If the item in the root is smaller than the item in the larger child, swap items
    if (items[rootIndex] < items[largerChildIndex])
    {
        Swap items[rootIndex] and items[largerChildIndex]
```

remove's final step transforms the semiheap into a heap

Every nonleaf in a complete tree has a left child

```
    // Transform the semiheap rooted at largerChi1dIndex into a heap
    heapRebuild(largerChildIndex, items, itemCount)
    }
}
// Else root is a leaf, so you are done
```

Figure 17-6 illustrates heapRebuild's recursive calls.

FIGURE 17-5 The array representation of (a) the heap in Figure 17-4a; (b) the semiheap in Figure 17-4b; (c) the restored heap in Figure 17-4c
(a)


| 0 | 10 |
| :---: | :---: |
| 1 | 9 |
| 2 | 6 |
| 3 | 3 |
| 4 | 2 |
| 5 | 5 |
|  |  |
|  |  |
|  |  |

Item in root is (b) smaller than items



| 0 | 9 |
| :---: | :---: |
| 1 | 5 |
| 2 | 6 |
| 3 | 3 |
| 4 | 2 |
|  |  |
|  |  |
|  |  |
|  |  |

FIGURE 17-6 Recursive calls to heapRebuild


First semiheap passed to heapRebuild


Second semiheap passed to heapRebuild

Now the heap's remove operation uses heapRebuild as follows:
// Copy the item from the last node into the root
items[0] = items[itemCount -1]
// Remove the last node
itemCount--
// Transform the semiheap back into a heap
heapRebuild(0, items, itemCount)
Consider briefly the efficiency of remove. Because the tree is stored in an array, the removal of a node requires you to swap array entries rather than simply to change a few pointers. These swaps may concern you, but they do not necessarily indicate that the algorithm is inefficient. At most, how many array entries will you have to swap? After remove copies the item in the last node of the tree into the root, heapRebui 1d trickles this item down the tree until its appropriate place is found. This item travels down a single path from the root to, at worst, a leaf. Therefore, the number of array items that heapRebuild must swap is no greater than the height of the tree. The height of a complete binary tree with $n$ nodes is always $\left\lceil\log _{2}(n+1)\right\rceil$, as you know from Chapter 15 . Because each swap requires three data moves, remove requires

$$
3 \times\left\lceil\log _{2}(n+1)\right\rceil+1
$$

remove's efficiency
remove is $\mathrm{O}(\log n)$
data moves. Thus, remove is $\mathrm{O}(\log n)$, which is in fact quite efficient.
Adding an item to a heap. The strategy for the add algorithm is the opposite of that for remove. A new item is inserted at the bottom of the tree, and it trickles up to its proper place, as Figure 17-7 illustrates. It is easy to trickle up a node, because the parent of the node in items [i] is always stored in items [(i-1)/2]-unless, of course, the node is the root.

## FIGURE 17-7 Insertion into a heap



The pseudocode for add follows:

```
// Insert newData into the bottom of the tree
items[itemCount] = newData
// Trickle new item up to the appropriate spot in the tree
newDataIndex = itemCount
inPlace = false
while ( (newDataIndex >= 0) and !inPlace)
{
    parentIndex = (newDataIndex - 1) / 2
    if (items[newDataIndex] < items[parentIndex])
```

```
        inPlace = true
        else
        {
        Swap items[newDataIndex ] and items[parentIndex]
        newDataIndex = parentIndex
    }
}
itemCount++
```

The efficiency of insert is like that of remove. At worst, insert has to swap array entries on a path from a leaf to the root. The number of swaps, therefore, cannot exceed the height of the tree. Because the height of the tree, which is complete, is always $\left\lceil\log _{2}(n+1)\right\rceil$, insert is also $\mathrm{O}(\log n)$.

Question 7 Is the full binary tree in Figure 16-16 of Chapter 16 a semiheap?
Question 8 Consider the maxheap in Figure 17-1a. Draw the heap after you insert 12 and then remove 12 .

Question 9 What does the initially empty heap myHeap contain after the following sequence of pseudocode operations?

```
myHeap.add(2)
myHeap.add(3)
myHeap.add(4)
myHeap.add(1)
myHeap.add(9)
myHeap.remove()
myHeap.add(7)
myHeap.add(6)
myHeap.remove()
myHeap.add(5)
```


### 17.2.2 The Implementation

The header file for the class ArrayMaxHeap, which is an array-based implementation of the ADT heap, is in Listing 17-2. This heap is a maxheap.

## LISTING 17-2 The header file for the class ArrayMaxHeap

```
/** Array-based implementation of the ADT heap.
    @file ArrayMaxHeap.h */
#ifndef _ARRAY_MAX_HEAP
#define _ARRAY_MAX_HEAP
#include "HeapInterface.h"
#include "PrecondViolatedExcep.h"
temp1ate<c1ass ItemType>
class ArrayMaxHeap : public HeapInterface<ItemType>
{
```

```
private:
    static const int ROOT_INDEX = 0; // Helps with readability
    static const int DEFAULT_CAPACITY = 21; // Smal7 capacity to test for a ful7 heap
    ItemType* items; // Array of heap items
    int itemCount; // Current count of heap items
    int maxItems; // Maximum capacity of the heap
    // Most of the private utility methods use an array index as a parameter
    // and in calculations. This should be safe, even though the array is an
    // implementation detail, since the methods are private.
    //
    // Returns the array index of the left child (if it exists).
    int getLeftChildIndex(const int nodeIndex) const;
    // Returns the array index of the right child (if it exists).
    int getRightChildIndex(int nodeIndex) const;
    // Returns the array index of the parent node.
    int getParentIndex(int nodeIndex) const;
    // Tests whether this node is a leaf.
    bool isLeaf(int nodeIndex) const;
    // Converts a semiheap to a heap.
    void heapRebuild(int subTreeRootIndex);
    // Creates a heap from an unordered array.
    void heapCreate();
public:
    ArrayMaxHeap();
    ArrayMaxHeap(const ItemType someArray[], const int arraySize);
    virtual ~ArrayMaxHeap();
    // HeapInterface Public Methods:
    bool isEmpty() const;
    int getNumberOfNodes() const;
    int getHeight() const;
    ItemType peekTop() const throw(PrecondViolatedExcep);
    bool add(const ItemType& newData);
    bool remove();
    void clear();
}; // end ArrayMaxHeap
#include "ArrayMaxHeap.cpp"
#endif
```

Some method definitions. Let's examine the definitions of some of the methods in ArrayMaxHeap, beginning with the private ones. To make our code more readable, we established the private methods getLeftChildIndex, getRightChildIndex, and getParentIndex. The implementations of these methods follow from our observations at the beginning of Section 17.2. For example, the left child of node items[i] is items [2 * $i+1$ ], so the method getLeftChildIndex has the following definition:

```
template<class ItemType>
int ArrayMaxHeap<ItemType>::getLeftChildIndex(const int nodeIndex) const
{
```

```
    return (2 * nodeIndex) + 1;
} // end getLeftChildIndex
```

The recursive private method heapRebuild closely follows the pseudocode given earlier in this chapter, so let's assume that we have coded it and look at a method that uses it. The second constructor creates a heap from an array of entries. To do so, this constructor can call the private method heapCreate, which in turn will need to call heapRebuild. Consider the details.

Notice that heapCreate has no parameters, but it does have access to our class's private data. In particular, heapCreate can use the array items and the counter itemCount. These two data members will serve as both the input to and the output from heapCreate. So if the constructor places values into itemCount and the array items, it can call heapCreate to rearrange the values in the array into a heap.

Thus, we can define the constructor as follows:

```
template<class ItemType>
ArrayMaxHeap<ItemType>: :
ArrayMaxHeap(const ItemType someArray[], const int arraySize):
    itemCount(arraySize), maxItems(2 * arraySize)
{
    // Allocate the array
    items = new ItemType[2 * arraySize];
    // Copy given values into the array
    for (int i = 0; i < itemCount; i++)
        items[i] = someArray[i];
    // Reorganize the array into a heap
    heapCreate();
} // end constructor
```

We use initializers to set the initial values of itemCount and maxItems. After allocating the array items, we copy the values from the given array someArray into items. Finally, we call heapCreate.

Now heapCreate must form a heap from the values in the array items. One way to accomplish this transformation is to use the heap's add method to insert the items into the heap one by one. However, a more efficient technique of building a heap out of the items in an array is possible.

For example, assume that the initial contents of an array are as shown in Figure 17-8a. First you imagine the array as a complete binary tree by assigning the array's items to the tree's nodes, beginning with the root and proceeding left to right down the tree. Figure $17-8 b$ shows the resulting tree. Next, you transform this tree into a heap by calling heapRebuild repeatedly. Each call to heapRebuild

FIGURE 17-8 (a) The initial contents of an array; (b) the array's corresponding complete binary tree
(a)
(b)

transforms a semiheap-a tree whose subtrees are both heaps but whose root may be out of placeinto a heap. But are there any semiheaps in the tree for heapRebuild to work on? Although the tree in Figure $17-8 \mathrm{~b}$ is not a semiheap, if you look at its leaves you will find semiheaps-that is, each leaf is a semiheap. In fact, each leaf is a heap, but for the sake of simplicity, ignore this fact.

You first call heapRebui 1d on the leaves from right to left. You then move up the tree, knowing that by the time you reach a node $N$, its subtrees are heaps, and thus heapRebuild will transform the semiheap rooted at $N$ into a heap. The following pseudocode steps transform the array items of itemCount items into a heap:

```
for (index = itemCount - 1 down to 0)
{
    // Assertion: The tree rooted at index is a semiheap
    heapRebuild(index)
    // Assertion: The tree rooted at index is a heap
}
```

Actually, you can replace itemCount - 1 with itemCount / 2 in the previous for statement. Thus, our pseudocode becomes

```
for (index = itemCount / 2 down to 0)
{
    // Assertion: The tree rooted at index is a semiheap
    heapRebuild(index)
    // Assertion: The tree rooted at index is a heap
}
```

Exercise 11 at the end of this chapter asks you to explain why this improvement is possible. Figure 17-9 traces this algorithm for the array in Figure 17-8a. Since the array contains six items, index in the for statement begins at $5 / 2$, or 2 .

FIGURE 17-9 Transforming an array into a heap


We now translate this pseudocode into $\mathrm{C}++$ to get the method heapCreate:

```
template<class ItemType>
void ArrayMaxHeap<ItemType>::heapCreate()
{
    for (int index = itemCount / 2; index >= 0; index--)
        heapRebuild(index);
} // end heapCreate
```

Programming Problem 1 asks you to complete the implementation file for the class ArrayMaxHeap.
Finally, we will define the method peekTop, since it must throw an exception if the heap is empty. Although you might be tempted to simply return items [0], since it always contains the top of the heap, if you do, peekTop will return a value even when the heap might be empty. Thus, you need to have peekTop test for an empty heap, as in the following definition:

```
temp1ate<class ItemType>
ItemType ArrayMaxHeap<ItemType>::peekTop() const throw(PrecondViolatedExcep)
{
        if (isEmpty())
        throw PrecondViolatedExcep("Attempted peek into an empty heap.");
    return items[0];
} // end peekTop
```

Question 10 Execute the following pseudocode statements on the array shown in Checkpoint Question 5.

```
for (index = n - 1 down to 0)
heapRebuild(index)
```


### 17.3 A Heap Implementation of the ADT Priority Queue

Priority-queue operations and heap operations are analogous

Section 13.3 of Chapter 13 introduced the ADT priority queue, and Section 14.2 of Chapter 14 offered an implementation based on a sorted list. At that time, we mentioned that using a heap to define a priority queue results in a more time-efficient implementation.

Once you have implemented the ADT heap, the implementation of the ADT priority queue is straightforward, because priority queue operations are exactly analogous to heap operations. The priority value in a priority queue item corresponds to an item in a heap. Thus, the implementation of the priority queue can reuse ArrayMaxHeap. To do so, we could use an instance of ArrayMaxHeap as a data member of the class of priority queues, or we could use inheritance. Although a heap provides an excellent implementation of a priority queue, a priority queue is not a heap. Since an is- $a$ relationship does not exist between ArrayMaxHeap and the class of priority queues, public inheritance is not appropriate. But we can use private inheritance, and that is what we will do. Let's assume that we have an interface for the ADT priority queue, as Exercise 8 in Chapter 13 asked you to write, so that we can derive our new class from it as well.

Listing 17-3 contains a header file for a class of priority queues.

## LISTING 17-3 A header file for the class Heap_PriorityQueue

Header file

```
/** ADT priority queue: Heap-based implementation.
    @file Heap_PriorityQueue.h */
#ifndef _HEAP_PRIORITY_QUEUE
#define _HEAP_PRIORITY_QUEUE
```

```
#include "ArrayMaxHeap.h"
#include "PriorityQueueInterface.h"
template<class ItemType>
class Heap_PriorityQueue : public PriorityQueueInterface<ItemType>,
    private ArrayMaxHeap<ItemType>
{
public:
    Heap_PriorityQueue();
    bool isEmpty() const;
    bool add(const ItemType& newEntry);
    bool remove();
    /** @pre The priority queue is not empty. */
    ItemType peek() const throw(PrecondViolatedExcep);
}; // end Heap_PriorityQueue
#include "Heap_PriorityQueue.cpp"
#endif
```

Each of the method definitions in Listing 17-4 calls the corresponding method in ArrayMaxHeap. Notice how the method peek handles the exception thrown by the heap's peekTop method, so that the message given mentions the priority queue and not the underlying heap.

## LISTING 17-4 An implementation of the class Heap_PriorityQueue

```
/** Heap-based implementation of the ADT priority queue.
```

    @file Heap_PriorityQueue.cpp */
    \#include "Heap_PriorityQueue.h"
temp1ate<c1ass ItemType>
Heap_PriorityQueue<ItemType>: :Heap_PriorityQueue()
\{
ArrayMaxHeap<ItemType>();
\} // end constructor
template<class ItemType>
boo1 Heap_PriorityQueue<ItemType>::isEmpty() const
\{
return ArrayMaxHeap<ItemType>::isEmpty();
\} // end isEmpty
temp1ate<c1ass ItemType>
boo1 Heap_PriorityQueue<ItemType>: :add(const ItemType\& newEntry)
\{
return ArrayMaxHeap<ItemType>: :add(newEntry);
\} // end add
temp1ate<c1ass ItemType>
bool Heap_PriorityQueue<ItemType>: :remove()
\{
return ArrayMaxHeap<ItemType>: :remove();
\} // end remove

The heap implementation requires knowledge of the priority queue's maximum size

A heap is always balanced

A heap of queues
template<c1ass ItemType>
ItemType Heap_PriorityQueue<ItemType>: :peek() const throw(PrecondViolatedExcep) \{
try
\{
return ArrayMaxHeap<ItemType>: :peekTop();
\}
catch (PrecondViolatedExcep e)
\{
throw PrecondViolatedExcep("Attempted peek into an empty priority queue.");
\} // end try/catch
\} // end peek

A heap versus a binary search tree as a priority queue. How does a heap compare to a binary search tree as an implementation of a priority queue? If you know the maximum number of items in the priority queue, the heap is the better implementation.

Because a heap is complete, it is always balanced, which is its major advantage. If the binary search tree is balanced, both implementations will have the same average performance for $n$ items: They both will be $\mathrm{O}(\log n)$. The height of a binary search tree, however, can increase during the insertion and removal of entries, greatly exceeding $\log _{2} n$ and degrading the implementation's efficiency to $\mathrm{O}(n)$ in the worst case. The heap implementation avoids this decrease in performance. In Chapter 19, you will see how to keep a search tree balanced, but the operations that do this are far more complex than the heap operations.

Finite, distinct priority values. If you have a finite number of distinct priority values, such as the integers 1 through 20, many items will likely have the same priority value. You could place items whose priority values are the same in the order in which you encounter them.

A heap of queues accommodates this situation, one queue for each distinct priority value. To insert an item into the priority queue, you add a queue for the item's priority value to the heap, if it is not already there. Then you insert the item into the corresponding queue. To remove an item from a priority queue, you remove the item at the front of the queue that corresponds to the highest priority value in the heap. If this removal leaves the queue empty, you remove the queue from the heap. Programming Problem 7 at the end of this chapter treats distinct priority values further.

Question 11 Consider a heap-based implementation of the ADT priority queue. What does the underlying heap contain after the following sequence of pseudocode operations, assuming that pQueue is an initially empty priority queue?

```
pQueue.add(5)
pQueue.add(9)
pQueue.add(6)
pQueue.add(7)
pQueue.add(3)
pQueue.add(4)
pQueue.remove()
pQueue.add(9)
pQueue.add(2)
pQueue.remove()
```


### 17.4 Heap Sort

As its name implies, the heap sort algorithm uses a heap to sort an array of items that are in no particular order. Suppose that we have a heap and an empty array whose size is the number of items in the heap. A call to peekTop gets the largest item in the heap, and we can place that item at the end of the array. A call to remove then removes that item from the heap. By repeatedly calling peekTop and remove, we can move the items from the heap in descending order and place them into sequentially decreasing positions in the array. The result is an array sorted into ascending order.

While this approach would work, it uses more memory and time than is necessary. Let's start over with an array of unsorted values. Our first step would be to transform the array into a heap. Recall the private method heapCreate from our implementation of the ADT heap in Section 17.2.2. This method transformed the array items, which is a data member of the class of heaps, into a heap. The body of heapCreate is simply the following loop:

```
for (int index = itemCount / 2; index >= 0; index--)
    heapRebuild(index);
```

As members of the class ArrayMaxHeap, both heapCreate and heapRebui 1 d have access to the class's data members, including the array items and its number of entries, itemCount. To use heapRebuild in a heap sort, we must revise it so that it has the array and its size as parameters. Let's declare heapRebuild as a function outside of ArrayMaxHeap, as follows:
void heapRebuild(int index, ItemType\& anArray[], int n)
After transforming the array into a heap, heap sort partitions the array into two regions-the Heap region and the Sorted region-as Figure 17-10 illustrates. The Heap region is in anArray [0.. last], and the Sorted region is in anArray[1ast $+1 . . n-1]$. Initially, the Heap region is all of anArray and the Sorted region is empty.

FIGURE 17-10 Heap sort partitions an array into two regions


Each step of the algorithm moves an item $I$ from the Heap region to the Sorted region. During this process, the following statements are true - they are the algorithm's invariant:

- The Sorted region contains the largest values in anArray, and they are in sorted order-that is, anArray $[n-1]$ is the largest item, anArray [n-2] is the second largest, and so on.
- The items in the Heap region form a heap.

So that the invariant holds, $I$ must be the item that has the largest value in the Heap region, and therefore $I$ must be in the root of the heap. To accomplish the move, you exchange the item in the root of the heap with the last item in the heap-that is, you exchange anArray[0] with anArray[last]-and then decrement the value of 1 ast. As a result, the item just swapped from the root into anArray [last]

Building a heap from an array of items

This version of heapRebuild is not in ArrayMaxHeap

Invariant for heap sort
becomes the smallest item in the Sorted region (and is in the first position of the Sorted region). After the move, you must transform the Heap region back into a heap because the new root may be out of place. You can accomplish this transformation by using heapRebuild to trickle down the item now in the root so that the Heap region is once again a heap.

The following algorithm summarizes these steps:

```
// Sorts anArray[0..n-1].
heapSort(anArray: ArrayType, n: integer)
    // Build initial heap
    for (index = n / 2 down to 0)
    {
        // Assertion:The tree rooted at index is a semiheap
        heapRebuild(index, anArray, n)
        // Assertion: The tree rooted at index is a heap
    }
    // Assertion: anArray [0] is the largest item in heap anArray [0..n-1]
    // Move the largest item in the Heap region -the root anArray [0] - to the beginning
    // of the Sorted region by swapping items and then adjusting the size of the regions
    Swap anArray[0] and anArray[n - 1]
    heapSize = n - 1 // Decrease the size of the Heap region, expand the Sorted region
    while (heapSize > 1)
    {
        // Make the Heap region a heap again
        heapRebuild(0, anArray, heapSize)
        // Move the largest item in the Heap region - the root anArray [0] - to the beginning
        // of the Sorted region by swapping items and then adjusting the size of the regions
        Swap anArray[0] and anArray[heapSize - 1]
        heapSize-- // Decrease the size of the Heap region, expand the Sorted region
    }
```

If we begin with the array shown in Figure 17-9, the first steps of heap sort are the same as those shown in the figure, and they transform the array into a heap. Figure 17-11 traces the heap sort from that point on. The C++ implementation of heap sort is left as an exercise.

The analysis of the efficiency of heap sort is similar to that of merge sort, as given in Chapter 11. Both algorithms are $\mathrm{O}(n \times \log n)$ in both the worst and average cases. Heap sort has an advantage over merge sort in that it does not require a second array. Quick sort is also $\mathrm{O}(n \times \log n)$ in the average case but is $\mathrm{O}\left(n^{2}\right)$ in the worst case. Even though quick sort has poor worst-case efficiency, it is generally the preferred sorting algorithm.

Question 12 Trace the heap sort as it sorts the following array into ascending order: 2530 20804060

FIGURE 17-11 A trace of heap sort, beginning with the heap in Figure 17-9

After making anAr ray a heap

After swapping anArray [0] with anArray [5] and decreasing the size of the Heap region

After heapRebui1d(0, anArray, 4)

After swapping anArray [0] with anArray[4] and decreasing the size of the Heap region

After heapRebui1d(0, anArray, 3)


$5^{1} \quad 2$ anArray [3] and decreasing the size of the Heap region

After rebuildHeap (0, anArray, 2)

anArray [2] and decreasing the size of the Heap region

After heapRebui1d(0, anArray, 1)


After swapping anArray [0] with anArray[1] and decreasing the size of the Heap region


Array is sorted


## Summary

1. Since a heap is a complete binary tree, it has an efficient array-based implementation.
2. The root of a maxheap contains the heap's largest value. The root of a minheap contains the heap's smallest value.
3. To remove the value in the root of a heap, you first replace the value with the value in the last node of the heap. This step likely results in a semiheap. After deleting the last node, you convert the semiheap to a heap by a trickle-down step.
4. To add an item to a heap, you insert a new leaf into the heap so that it remains a complete binary tree. You then use a trickle-up step to reposition the new value.
5. A heap that uses an array-based representation of a complete binary tree is a good implementation of a priority queue when you know the maximum number of items that will be stored at any one time.
6. Heap sort converts an array into a heap to locate the array's largest item. This step enables the heap sort to sort an array in an efficient manner.
7. Heap sort, like merge sort, has good worst-case and average-case behaviors, but neither algorithm is as good in the average case as quick sort. Heap sort has an advantage over merge sort in that it does not require a second array.

## Exercises

1. Given the minheap myHeap in Figure 17-12a, show what it would look like after each of the following pseudocode operations:
a. myHeap.add (8)
b. myHeap.add(5)
c. myHeap. remove()

FIGURE 17-12 (a) Minheap for Exercise 1; (b) maxheap for Exercise 2

2. Given the maxheap myHeap in Figure $17-12$ b, show what it would look like after each of the following pseudocode operations:
a. myHeap.add(16)
b. myHeap.add(14)
c. myHeap. remove()
3. Repeat Checkpoint Question 10, but instead use the array $9,12,4,8,3,11,6,15$.
4. Prove that the root of a maxheap contains the largest value in the tree.
5. Does the order in which you insert items into a heap affect the heap that results? Explain.
6. Revise the pseudocode for the ADT heap methods add and heapRebuild so that they do not swap items.
7. Suppose that you have two items with the same priority value. How does the order in which you insert these items into a priority queue affect the order in which they will be removed? What can you do if you need entries whose priority values are equal to be served on a first-come, first-served basis?
8. Suppose that you wanted the remove operation of a priority queue to remove the entry whose priority value is smallest instead of largest. You would then use a minheap. Convert the maxheap implementation to a minheap implementation.
9. Suppose that you wanted to maintain the index of the item with the smallest value in a maxheap. That is, in addition to a getTop operation, you might want to support a getMin operation. How difficult would it be to maintain this index within the add and remove operations?
10. Suppose that after you have placed several items into a priority queue, you need to adjust one of their priority values. For example, a particular task in a priority queue of tasks could become either more or less urgent. How can you adjust a heap if a single value changes?
11. Show that within the pseudocode for the method heapCreate you can replace the statement

```
for (index = n - 1 down to 0)
```

with

```
for (index = n / 2 down to 0)
```

12. Trace the action of heapSort on the array given in Checkpoint Question 5.
13. Implement heapSort in $\mathrm{C}++$.
14. Revise heapSort so that it sorts an array into descending order.

## Programming Problems

1. Complete the implementation file for the class ArrayMaxHeap that is described in Section 17.2.2.
2. Use a binary search tree in the implementation of HeapInterface. Where in the tree will the largest entry occur? How efficient is this implementation?
3. Implement a class of minheaps using an array. Name your class ArrayMinHeap.
4. Consider the problem of combining two heaps together into a single heap.
a. Write an efficient algorithm for combining two heaps, one with size $n$ and the other with size 1 . What is the Big O performance of your algorithm?
b. Write an efficient algorithm for combining two heaps of equal size $n$. What is the Big O performance of your algorithm?
c. Write an efficient algorithm for combining two arbitrary-sized heaps into one heap. What is the Big O performance of your algorithm?
d. Implement the algorithm that you wrote in part $c$.
5. Section 2.4.4 of Chapter 2 discussed the problem of finding the $k^{\text {th }}$ smallest value in an array of $n$ values. Design an algorithm that uses a minheap to solve this problem. Using the class ArrayMinHeap defined in Programming Problem 3, implement your algorithm as a function at the client level.
6. Implement the to-do list described at the beginning of Section 13.3 in Chapter 13. Use the class Heap_PriorityQueue as given in Listings 17-3 and 17-4.
7. Suppose that you wanted to implement a priority queue whose priority values are integers 1 through 20.
a. Implement the priority queue as a heap of queues, as described in this chapter.
b. Another solution uses an array of 20 queues, one for each priority value. Use this approach to implement the priority queue.
8. Write an interactive program that will monitor the flow of patients in a large hospital. The program should account for patients who check in and out of the hospital and should allow access to information about a given patient. In addition, the program should manage the scheduling of three operating rooms. Doctors make a request that includes a patient's name and a priority value between 1 and 10 that reflects the urgency of the operation. Patients are chosen for the operating room by priority value, and patients with the same priority are served on a first-come, first-served basis.

The user should use either one-letter or one-word commands to control the program. As you design your solution, try to identify the essential operations (excuse the pun) that you must perform on the data, and only then choose an appropriate data structure for implementation. This approach will allow you to maintain the wall between the main part of the program and the implementations.
9. Implement a class of priority queues using a binary search tree to contain its items. Discuss the appropriateness and efficiency of this implementation.

## Dictionaries and Their Implementations

## Chapter

 18
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The ADT dictionary

Thhis chapter considers the ADT dictionary, which is appropriate for problems that must manage data by value. Several dictionary implementations-which use arrays, linked chains, binary search trees, and a new technique known as hashing-will be presented, along with their advantages and disadvantages.

To make an intelligent choice among the various possible dictionary implementations, you must analyze the efficiency with which each of the implementations supports the dictionary operations. For example, we analyze the time-efficiency of array-based and link-based dictionary implementations and conclude that, in many applications, the implementations do not support the dictionary operations as efficiently as possible. This conclusion motivates the use of more sophisticated dictionary implementations.

### 18.1 The ADT Dictionary

When we began our discussion of sorting a collection of data in Chapter 11, we introduced the notion of a sort key. For example, you might want to sort a collection of data about people based on their names, their ages, or their zip codes. The criterion you choose is known as the sort key. Similarly, you might want to search the same collection of data for a name, an address, or a phone number. The criterion you choose for this search is known as a search key.

Applications that require value-oriented operations are extremely prevalent, as you might imagine. For example, the tasks

- Find the phone number of John Smith
- Delete all the information about the employee with ID number 12908
involve values instead of positions. We first encountered a value-oriented ADT, the sorted list, in Chapter 12. Chapter 15 presented the binary search tree, which is also a value-oriented ADT. This section describes yet another value-oriented ADT, the dictionary.

Consider the data given in Figure 18-1 for some major cities in the world. Each city has certain information listed, and the design of the data collection enables you to look up this information. For

FIGURE 18-1 A collection of data about certain cities

| City | Country | Population |
| :---: | :---: | :---: |
| Buenos Aires | Argentina | 13,170,000 |
| Cairo | Egypt | 14,450,000 |
| Cape Town | South Africa | 3,092,000 |
| London | England | 12,875,000 |
| Madrid | Spain | 4,072,000 |
| Mexico City | Mexico | 20,450,000 |
| Mumbai | India | 19,200,000 |
| New York City | U.S.A. | 19,750,000 |
| Paris | France | 9,638,000 |
| Sydney | Australia | 3,665,000 |
| Tokyo | Japan | 32,450,000 |
| Toronto | Canada | 4,657,000 |

example, if you wanted to know the population of London, you could scan the column of city names, starting at the top, until you came to London. Because the cities are listed in alphabetical order, you could also mimic a binary search. You could begin the search near the middle of the column determine in which half London lies, and recursively apply the binary search to the appropriate half. As you know, a binary search is far more efficient than scanning the entire column of city names from the beginning.

If, however, you wanted to find out which cities in the data collection are in Spain, you would have no choice but to scan the entire data collection. The alphabetical order of the city names does not help you for this problem at all. The data arrangement facilitates the search for a given city, but other types of questions require a complete scan of the data.

The ADT dictionary, or map, or table, also allows you to look up information easily and has a special operation for this purpose. Typically, the items in the ADT dictionary are objects that contain several pieces of data. You can facilitate the retrieval of an item by basing the search on a specified search key. In the dictionary of cities, for example, you could designate City as the search key if you often needed to retrieve the information about a city. You can devise implementations of a dictionary that allow the rapid retrieval of the item(s) whose search key matches some specified value. However, if you need to retrieve item(s) based on a value that is not a search key, you will have to inspect the entire dictionary. Therefore, the choice of a search key sends the ADT implementer the following message:

Arrange the data to facilitate the search for an item, given the value of its search key.
The basic operations that define the ADT dictionary are as follows:

## Note: ADT dictionary operations

- Test whether a dictionary is empty.
- Get the number of items in a dictionary.
- Insert a new item into a dictionary.
- Remove the item with a given search key from a dictionary.
- Remove all items from a dictionary.
- Get the item with a given search key from a dictionary.
- Test whether a dictionary contains an item with a given search key.
- Traverse the items in a dictionary in sorted search-key order.

For simplicity, we will assume that all items in the dictionary have distinct search keys. This assumption can affect our design in one of in several ways:

- Distinct search keys can be a precondition of the insertion operation
- The insertion operation can deny an attempt to insert a new entry whose search key already exists in the dictionary
- The insertion operation can replace an existing entry whose search key matches the search key of a new entry with the new entry
We will choose the first option.
The following operation contract specifies in more detail an ADT dictionary of items with distinct search keys. Figure 18-2 shows a UML diagram for a class of dictionaries.

The ADT dictionary uses a search key to identify its items

## Abstract Data Type: Dictionary

DATA

- A finite number of objects, each associated with a search key.

Operations

| Pseudocode | Description |
| :---: | :---: |
| isEmpty () | Task: Sees whether this dictionary is empty. <br> Input: None. <br> Output: True if the dictionary is empty; otherwise false |
| getNumberOfItems() | Task: Gets the number of items in this dictionary. Input: None. <br> Output: The number of items in the dictionary. |
| add(searchKey, newItem) | Task: Inserts an item into this dictionary according to the item's search key. Input: newItem is the data item to be inserted; searchKey is the item's associated search key and differs from all search keys presently in the dictionary. Output: True if the insertion is successful, or false if not. |
| remove(searchKey) | Task: Removes the item with the given search key from this dictionary. Input: searchKey is the search key of the item to be removed. Output: True if the removal is successful, or false if not. |
| clear () | Task: Removes all entries from this dictionary. Input: None. Output: None. |
| getItem(searchKey) | Task: Gets an item with a given search key from this dictionary. Input: searchKey is the search key of the item to be retrieved. Output: The item associated with the given search key. |
| contains(searchKey) | Task: Sees whether this dictionary contains an item with a given search key. Input: searchKey is the search key of the desired item. <br> Output: True if the dictionary contains the designated item, or false if not. |
| traverse(visit) | Task: Traverses this dictionary and calls a given client function once for each item. Input: The client function visit, which exists outside of the ADT implementation. Output: visit's action occurs once for each item in the dictionary and possibly alters the item. |

You should realize that these operations are only one possible set of dictionary operations. The

Various sets of dictionary operations are possible

Our dictionary assumes distinct search keys client may require either a subset of these operations or other operations not listed here to fit the application at hand. It may also be convenient to modify the definitions of some of the operations. For example, these operations assume that no two dictionary items have the same search keys. However, in many applications it is quite reasonable to expect duplicate search keys. If this is the case, you must redefine several of the operations to eliminate the ambiguity that would arise from duplicate search keys. For example, which item should getItem return if several items have the specified search key? You should tailor your definition of the ADT dictionary to the problem at hand.

| Dictionary |
| :--- |
|  |
| +isEmpty(): boolean |
| +getNumber0fItems(): integer |
| +add(itemKey: KeyType, newItem: ItemType): boolean |
| +remove(itemKey: KeyType): boolean |
| +clear() : void |
| +geItem(itemKey: KeyType): ItemType |
| +contains(itemKey: KeyType): boolean |
| +traverse(visit(item: ItemType): void) : void |

Although just the operations add, remove, and getItem in the previous set of operations are sufficient for some applications, you cannot do several significant things without additional operations. For instance, you cannot display all the dictionary items, because you cannot retrieve a data item unless you know the value of its search key. Thus, you cannot display the entire dictionary unless you can traverse the dictionary.

The traverse operation visits each item in the dictionary once. In defining this operation, you must specify the order in which the traversal should visit the items. One common specification is to visit the items in sorted order by search key, but perhaps you do not care about the order in which the items are visited. As you will see, the way you define traverse-if you request it at all-can affect the way you implement the dictionary.

## Note: Search keys

The concept of a search key for the dictionary items is essential to the implementation of the dictionary. It is important that the value of the search key remain the same as long as the item is stored in the dictionary. Changing the search key of an existing entry in the dictionary could make that entry or other dictionary entries impossible to find. Thus, the client should not be able to modify the value of an entry's search key once that entry is in the dictionary.

Like the traversal operations for the ADT binary tree, traverse has a visit function as its argument. Because visit can do any number of things-including access the dictionary via the ADT operations-traverse is a versatile operation. We illustrate this versatility with three brief examples, in which the dictionary contains the data for the cities listed in Figure 18-1.

Examples: Each of the following tasks will use the city's name as the search key. The class City contains all the information for a city, including its name, country, and population. The class has accessor and mutator methods for each these pieces of data.

- Display, in alphabetical order, the name of each city and its population.This task requires you to write the city names in alphabetical order. Since the city names are the search keys, traverse must visit items alphabetically by search key. You pass the method the name of the function displayItem, which appears in pseudocode as follows:

```
displayItem(anItem: ItemType): void
Display anItem.getName()
Display anItem.getPopulation()
```

Other dictionaries could allow duplicate search keys
traverse visits all dictionary items

Tasks that use the city's name as the search key

Although the visitation order of traverse is significant for this first task, it is immaterial for the next one.

- Increase the population of each city by 10 percent. To perform this task, you pass to traverse the name of the function updatePopulation, which we define in pseudocode as follows:

```
updatePopulation(anItem: ItemType): void
    anItem.setPopulation(1.1 * anItem.getPopulation())
```


### 18.1.1 An Interface for the ADT Dictionary

To complete our specification of the ADT dictionary, we write the interface shown in Listing 18-1. Note that the interface specifies two data-type parameters-KeyType and ItemType-in the statement

```
template<class KeyType, class ItemType>
```

These parameters make it possible for the data type of the dictionary's data items to differ from the data type of their search keys.

## LISTING 18-1 An interface for the ADT dictionary

```
/** An interface for the ADT dictionary.
@file DictionaryInterface.h */
#ifndef _DICTIONARY_INTERFACE
#define _DICTIONARY_INTERFACE
#include "NotFoundException.h"
template<class KeyType, class ItemType>
class DictionaryInterface
{
public:
    /** Sees whether this dictionary is empty.
        @return True if the dictionary is empty;
            otherwise returns false. */
    virtual bool isEmpty() const = 0;
    /** Gets the number of items in this dictionary.
        @return The number of items in the dictionary. */
    virtual int getNumberOfItems() const = 0;
    /** Inserts an item into this dictionary according to the item's
            search key.
        @pre The search key of the new item differs from all search
            keys presently in the dictionary.
        @post If the insertion is successful, newItem is in its
            proper position within the dictionary.
        @param searchKey The search key associated with the item to be inserted.
        @param newItem The item to add to the dictionary.
        @return True if item was successfully added, or false if not. */
        virtua1 boo1 add(const KeyType& searchKey, const ItemType& newItem) = 0;
```

```
    /** Removes an item with the given search key from this dictionary.
    @post If the item whose search key equals searchKey
        existed in the dictionary, the item was removed.
    @param searchKey The search key of the item to be removed.
    @return True if the item was successfully removed, or false if not. */
    virtual bool remove(const KeyType& searchKey) = 0;
    /** Removes all entries from this dictionary. */
    virtual void clear() = 0;
    /** Retrieves an item with a given search key from a dictionary.
    @post If the retrieval is successful, the item is returned.
    @param searchKey The search key of the item to be retrieved.
    areturn The item associated with the search key.
    @throw NotFoundException if the item does not exist. */
    virtual ItemType getItem(const KeyType& searchKey) const
                        throw (NotFoundException) = 0;
        /** Sees whether this dictionary contains an item with a given
        search key.
    @post The dictionary is unchanged.
    @param searchKey The search key of the item to be retrieved.
    @return True if an item with the given search key exists in the
        dictionary. */
    virtual bool contains(const KeyType& searchKey) const = 0;
    /** Traverses this dictionary and calls a given client function once
        for each item.
    @post The given function's action occurs once for each item in the
        dictionary and possibly alters the item.
    @param visit A client function. */
    virtual void traverse(void visit(ItemType&)) const = 0;
}; // end DictionaryInterface
#endif
```

Question 1 Using the ADT dictionary operations, write pseudocode for a replace function at the client level that replaces the dictionary item whose search key is $x$ with another item whose search key is also $x$.

### 18.2 Possible Implementations

In previous chapters, ADT implementations often were either array based or link based. That is, you used either an array or a chain of linked nodes to store the ADT's items. Such implementations are called linear because they represent items one after another in a data structure and thus mirror the flat, list-like appearance of the city data given in Figure 18-1.

Linear implementations of a dictionary are certainly possible and fall into four categories:

- Sorted (by search key), array-based
- Sorted (by search key), link-based
- Unsorted, array-based
- Unsorted, link-based

The unsorted implementations store the items in no particular order; they can insert a new item into any convenient location. Notice that the add operation has two parameters: a data item and the item's search key. The sorted implementations must insert a new item into its proper position as determined by the value of its search key. Whether sorted or unsorted, the dictionary must not only store both pieces of data-search key and data item - but also form an association between the two. To do so, we will encapsulate each item with its search key into an object of a new class, Entry, as pictured in Figure 18-3.

## FIGURE 18-3 A dictionary entry



The array-based and link-based linear implementations have the basic structures shown in Figure 18-4. Both implementations maintain a count of the present number of items in the dictionary. The two implementations in this figure happen to be sorted. As you will see, the unsorted and sorted implementations have their relative advantages and disadvantages.

FIGURE 18-4 The data members for two sorted linear implementations of the ADT dictionary for the data in Figure 18-1: (a) array based; (b) link based


Note: A linear array-based implementation of the ADT dictionary must shift data during an insertion in sorted order and when removing an entry. These shifts can be expensive, particularly for large dictionaries.

Although a linear link-based implementation of the ADT dictionary eliminates the need to shift data, it does not support the insertion and removal operations any more efficiently than does an array-based implementation, because you cannot perform a binary search in a reasonable fashion.

The header file for the class Entry is in Listing 18-2. Notice that, like DictionaryInterface, the class Entry specifies two data-type parameters. We leave Entry's implementation for you as an exercise.

## LISTING 18-2 A header file for a class of dictionary entries

```
/** A class of entry objects for an array-based implementation of the
    ADT dictionary.
    @file Entry.h */
#ifndef _ENTRY
#define _ENTRY
template <class KeyType, class ItemType>
class Entry
{
private:
    ItemType item;
    KeyType searchKey;
protected:
    void setKey(const KeyType& searchKey);
public:
    Entry();
    Entry(ItemType newEntry, KeyType searchKey);
    ItemType getItem() const;
    KeyType getKey() const;
    void setItem(const ItemType& newEntry);
    bool operator==(const Entry<KeyType, ItemType>& rightHandItem) const;
    bool operator>(const Entry<KeyType, ItemType>& rightHandItem) const;
}; // end Entry
#include "Entry.cpp"
#endif
```


## Note: Observations about the class Entry

The class Entry has two data members, a dictionary item and an associated search key. It also has a set method and a get method for the item portion, but only a public get method for the search key. This design prevents us from accidentally changing an Entry object's search key, as doing so would destroy the integrity of the dictionary. However, Entry does have a protected method setKey. This method is necessary so that the constructor of any derived class of Entry can initialize the search key.

Notice that the class overloads the operators $==$ and $>$. While we will not need these overloaded operators for the array-based dictionary, we will use them when we store the dictionary's entries in a binary search tree.

At this point in your study of ADTs, you have other choices for a dictionary implementation. For instance, you can implement the ADT dictionary by using the ADT list, sorted list, or binary search tree. The binary search tree implementation, as illustrated in Figure 18-5, is an example of a nonlinear implementation and offers several advantages over linear implementations. Among these advantages is the opportunity to reuse the implementation of the ADT binary search tree discussed in Chapter 16. Implementations based on the ADTs list and sorted list also share this advantage, and they are left for you to consider as exercises.

A binary search tree implementation is nonlinear

FIGURE 18-5 The data members for a binary search tree implementation of the ADT dictionary for the data in Figure 18-1


We also will consider an important unsorted implementation that uses a technique known as hashing. Although based on an array, this implementation is unlike any you have seen in this book.

### 18.2.1 A Sorted Array-Based Implementation of the ADT Dictionary

Whether a dictionary organizes its entries in sorted order by search key or leaves them unsorted, an array-based implementation has data members that are similar to those in array-based implementations of other ADTs that you have seen, namely an array of data items, a count of the items, and a maximum number of items. We have already seen that the data items will be Entry objects that contain both a search key and some other data.

For a sorted dictionary, the operation traverse should visit the dictionary's data items in an order such that their search keys are in sorted order. Note that we have specified this behavior in the header file for the class of dictionaries given in Listing 18-3. Notice also that this dictionary requires that its search keys be unique. This requirement will simplify our implementation, but Exercises 7 and 8 at the end of this chapter ask you to remove this restriction.

The declaration of the class ArrayDictionary contains nothing unusual, but it does declare two private methods-destroyDictionary and findEntryIndex. The former method can be called by both the destructor and the public method clear. The latter method locates the entry in the array items that has a given search key. This method is useful for the public methods remove, getItem, and contains. Notice that both private methods perform operations needed by more than one public method, enabling us to avoid repetitive code in the definitions of the public methods.

## LISTING 18-3 A header file for the class ArrayDictionary

```
/** An array-based implementation of the ADT dictionary
    that organizes its data items in sorted search-key order.
    Search keys in the dictionary are unique.
    @file ArrayDictionary.h */
#ifndef _ARRAY_DICTIONARY
#define _ARRAY_DICTIONARY
#include "DictionaryInterface.h"
#include "Entry.h"
#include "NotFoundException.h"
template <class KeyType, class ItemType>
class ArrayDictionary : public DictionaryInterface<KeyType, ItemType>
{
private:
    static const int DEFAULT_CAPACITY = 21; // Small capacity to test for
                                    // a full dictionary
    Entry<KeyType, ItemType>* items; // Array of dictionary entries
    int itemCount; // Current count of dictionary items
    int maxItems; // Maximum capacity of the dictionary
```

    void destroyDictionary();
    int findEntryIndex(int firstIndex, int lastIndex,
                const KeyType\& searchKey) const;
    pub7ic:
ArrayDictionary();
ArrayDictionary (int maxNumberOfEntries);
ArrayDictionary (const ArrayDictionary<KeyType, ItemType>\& dict);
virtual ~ArrayDictionary();
bool isEmpty() const;
int getNumberOfItems() const;
bool add(const KeyType\& searchKey, const ItemType\& newItem);
bool remove(const KeyType\& searchKey);
void clear();
ItemType getItem(const KeyType\& searchKey) const
throw (NotFoundException);
bool contains (const KeyType\& searchKey) const;
/** Traverses the items in this dictionary in sorted search-key order
and calls a given client function once for each item. */
void traverse(void visit(ItemType\&)) const;
\}; // end ArrayDictionary
\#include "ArrayDictionary.cpp"
\#endif

The requirement that we traverse the dictionary in sorted search-key order provides one reason to sort the array items according to the search key. While we could sort items each time traverse is called, maintaining this array in sorted order has advantages other than avoiding a sorting step. One important advantage of a sorted array is that the private method findEntryIndex can perform a recursive binary search when looking for a particular search key. Notice that the
declaration of this method in Listing 18-3 provides two indices as parameters to indicate the portion of the array to search.

The public method add bears the responsibility for keeping the array items sorted. Let's examine its definition:

```
template <class KeyType, class ItemType>
bool ArrayDictionary<KeyType, ItemType>::add(const ItemType& newItem,
                                    const KeyType& searchKey)
{
    bool ableToInsert = (itemCount < maxItems);
    if (ableToInsert)
    {
        // Make room for new entry by shifting al1 entries at
        // positions >= newPosition toward the end of the array
        // (no shift if newPosition == itemCount + 1). Performing
        // a binary search doesn't help here, because we need to
        // shift the entries while looking for the insertion location.
        int index = itemCount;
        // Short-circuit evaluation is important
        while ( (index > 0) && (searchKey < items[index-1].getKey()) )
        {
            items[index] = items[index-1];
            index--;
        } // end while
        // Insert new entry
        items[index] = Entry<KeyType, ItemType>(newItem, searchKey);
        itemCount++; // Increase count of entries
    } // end if
    return ab7eToInsert;
} // end add
```

Programming Problem 1 at the end of this chapter asks you to complete the implementation of ArrayDictionary.

Question 2 Explain how the while loop in the previous definition of the method add locates the insertion point for the new entry in the array items.

Question 3 Why is short-circuit evaluation important in the while loop of the previous definition of the method add?

Question 4 We mentioned that the remove method calls the private method findEntryIndex to locate the entry to remove. Assuming that the entry is located, what does remove need to do after it gets the index of this entry?

Question 5 What is the definition of the method traverse for ArrayDictionary?

### 18.2.2 A Binary Search Tree Implementation of the ADT Dictionary

This dictionary implementation has a binary search tree as a data member

Although linear implementations of the ADT dictionary are fine for specific applications, they are not suitable as general-purpose classes. The following nonlinear link-based implementation uses a binary search tree to represent the items in the ADT dictionary. That is, the dictionary class will use composition and thus will have a binary search tree as one of its data members. In this way, our class reuses the class BinarySearchTree from Chapter 16.

The header file. The header file in Listing 18-4 declares the class TreeDictionary. Like ArrayDictionary, TreeDictionary stores Entry objects that encapsulate data items with their corresponding search keys. Much of this class is the same as or is similar to ArrayDictionary, except for the private section. It is there that we declare the binary search tree, itemTree, along with a method to help with the traversal operation.

## LISTING 18-4 A header file for the class TreeDictionary

```
/** A binary search tree implementation of the ADT dictionary
    that organizes its data items in sorted search-key order.
    Search keys in the dictionary are unique.
    @file TreeDictionary.h */
#ifndef _TREE_DICTIONARY
#define _TREE_DICTIONARY
#include "DictionaryInterface.h"
#include "BinarySearchTree.h"
#include "Entry.h"
#include "NotFoundException.h"
template <class KeyType, class ItemType>
class TreeDictionary : public DictionaryInterface<KeyType, ItemType>
{
private:
    // Binary search tree of dictionary entries
    BinarySearchTree<Entry<KeyType, ItemType> > itemTree;
    void traversa1He1per(Entry<KeyType, ItemType>& theEntry);
public:
    TreeDictionary();
    TreeDictionary(int maxNumberOfEntries);
    TreeDictionary(const TreeDictionary<KeyType, ItemType>& dict);
    virtual ~TreeDictionary();
    // The declarations of the public methods appear here and are the
    // same as given in Listing 18-3 for the class ArrayDictionary.
    . . .
}; // end TreeDictionary
#include "TreeDictionary.cpp"
#endif
```

The method implementations. Several of the methods in TreeDictionary simply call the analogous method in BinarySearchTree. For example, the method add creates an Entry object from the item and search key passed to it and then passes that object to the add method in BinarySearchTree. The resulting definition follows:

```
template <class KeyType, class ItemType>
boo1 TreeDictionary<KeyType, ItemType>::add(const ItemType& newItem,
                                    const KeyType& searchKey)
{
    return itemTree.add(Entry<KeyType, ItemType>(newItem, searchKey));
} // end add
```

Each of the methods remove, getItem, and contains is passed a search key as its only argument. Each method must search the binary search tree for an Entry object that has a particular search key. How will TreeDictionary's method remove, for example, call BinarySearchTree's method remove? After all, it requires an Entry object as its argument.

The methods of BinarySearchTree compare Entry objects by using either of the operators $==$ or $>$. These operators are the same ones that Entry overloads. Thus, we can control how BinarySearchTree will make these comparisons. Because we are searching for a specific search key, we overload $==$ and $>$ to examine only the search-key portion of an Entry object. For example, the definition of the overloaded operator $==$ in Entry is

```
template <class KeyType, class ItemType>
boo1 Entry<KeyType, ItemType>::operator==(
    const Entry<KeyType, ItemType>& rightHandItem) const
{
    return (searchKey == rightHandItem.getKey());
} // end operator==
```

Now that BinarySearchTree can compare Entry objects based only on their search keys, how does TreeDictionary's method remove call BinarySearchTree's remove? Because BinarySearchTree's remove expects an Entry object as an argument, TreeDictionary's remove creates an Entry object that has a specific search key and any item portion, and passes it to the invoked method. To create this Entry object, you can write the following expression:

```
Entry<KeyType, ItemType>(ItemType(), searchKey)
```

The result is a call to Entry's second constructor with these two arguments: the result of a call to ItemType's default constructor and the search key that was passed to TreeDictionary's remove as an argument. Thus, the remove method in TreeDictionary has the following definition:

```
template <class KeyType, class ItemType>
boo1 TreeDictionary<KeyType, ItemType>::remove(const KeyType& searchKey)
{
    return itemTree.remove(Entry<KeyType, ItemType>(ItemType(), searchKey));
} // end remove
```

We leave the rest of this implementation for you to complete as an exercise.


Note: Usually a binary search tree can support the ADT dictionary operations quite efficiently. However, in the worst case, when the tree approaches a linear shape, the performance of the dictionary operations is comparable to that of a linear link-based implementation. If a given application cannot tolerate poor performance, you should use the dictionary implementations presented either in Section 18.4 of this chapter or in Chapter 19.

### 18.3 Selecting an Implementation

Perspective, efficiency, and motivation are reasons for studying the linear implementations

A major goal of this chapter is to indicate how the requirements of a particular application influence the selection of an implementation. The discussion here elaborates on the comments made in Section 10.2.4 of Chapter 10. Some applications require all of the ADT dictionary operations given earlier; others require either a subset of them or additional operations. Before choosing an implementation of the ADT dictionary, you as problem solver should carefully analyze which operations
you really need for the application at hand. It is tempting to want all possible operations, but this strategy is a poor one, because often one implementation supports some of the operations more efficiently than another implementation does. Therefore, if you include an operation that you never use, you might end up with an implementation of the ADT that does not best suit your purpose.

In addition to knowing what operations are needed for a given application, the ADT implementer should know approximately how often the application will perform each operation. Although some applications may require many occurrences of every operation, other applications may not. For example, if you maintained a dictionary of major cities - such as those in Figure 18-1-you would expect to perform many more retrieval operations than additions or removals. Thus, if you seldom add items to a dictionary, you can tolerate an implementation that results in an inefficient add operation, as long as frequently used operations are efficient. Of course, as Chapter 10 mentioned, if an ADT operation is to be used in a life-or-death situation, that operation must be efficient even if you rarely need it. The necessary operations, their expected frequency of occurrence, and their required response times are therefore some factors that influence which implementation of an ADT you should select for a particular application. You should, however, remain conscious of factors other than efficiency, as discussed in Chapter 10.

### 18.3.1 Three Scenarios

Consider now several different application scenarios, each of which requires a particular mix of the dictionary operations. The analysis of various implementations of the ADT dictionary will illustrate some of the basic concerns of the analysis of algorithms. You will see, given an application, how to select an implementation that supports in a reasonably efficient manner the required mix of dictionary operations.

Scenario A: Insertion and traversal in no particular order. Mary's sorority plans to raise money for a local charity. Tired of previous fund-raisers, Mary suggests a brainstorming session to discover a new money-making strategy. As sorority members voice their ideas, Mary records them by adding each new thought to a dictionary. Later, she will print a report of all the ideas currently in the dictionary. Assume that the organization of the report is irrelevant - the items can be sorted or unsorted. Also assume that operations such as retrieval, removal, or traversal in sorted order either do not occur or occur so infrequently that they do not influence your choice of an implementation.

For this application, maintaining the items in a sorted order has no advantage. In fact, by not maintaining a sorted order, the add operation can be quite efficient. For either unsorted linear implementation, you can insert a new item into any convenient location. For an unsorted array-based implementation, you can easily insert a new item after the last item in the array-that is, at location items [itemCount]. Figure 18-6a shows the result of this insertion after itemCount has been updated. For a link-based implementation, you can simply insert a new item at the beginning of the linked chain. As Figure 18-6b illustrates, the head pointer points to the new item, and the new item points to the item that was previously first in the chain. Thus, you can insert a new item quickly into either unsorted implementation of a dictionary; in fact, the add operation is $\mathrm{O}(1)$ : It requires a constant time for either implementation regardless of the dictionary size.

Should you choose the array-based or the link-based implementation? As you have seen with other ADTs, an implementation that uses dynamically allocated memory is appropriate if you do not have a good estimate of the maximum possible size of the dictionary. Mary's brainstorming session likely falls into this category. On the other hand, if you know that the dictionary's maximum size is not drastically larger than its expected size, ${ }^{1}$ the choice is mostly a matter of style. An array-based

What operations are needed?

How often is each operation required?

An unsorted order is efficient

Array-based versus pointer-based

[^13]A sorted arraybased implementation can use a binary search

Questions

implementation requires less space than a link-based implementation, because no explicit pointer is stored. The extra cost of this pointer relative to the size of the data items, however, is insignificant in most situations because large data items are typical.

Should you use a binary search tree in the dictionary implementation for this application? Because such an implementation orders the dictionary items, it does more work than the application requires. In fact, as you saw in Chapter 15, insertion into a binary search tree is $\mathrm{O}(\log n)$ in the average case; it is slower than the $\mathrm{O}(1)$ linear implementations.
Scenario B: Retrieval. When you use a word processor's thesaurus to look up synonyms for a word, you use a retrieval operation. If an ADT dictionary represents the thesaurus, each dictionary item is a record that contains both the word-which is the search key-and the word's synonyms. Frequent retrieval operations require a dictionary implementation that allows you to search efficiently for an item, given its search key. Typically, you cannot alter the thesaurus, so no insertion or removal operations are necessary.

For an array-based implementation, you can use a binary search to retrieve a particular word's synonyms, if the array is sorted. On the other hand, for a link-based implementation, you must traverse the linked chain from its beginning until you encounter the word. The binary search performs this retrieval in significantly less time than is required to traverse a linked chain. Two questions come to mind at this point:

- Is a binary search of a linked chain possible?
- How much more efficient is a binary search of an array than a sequential search of a linked chain?

Can you perform a binary search of a linked chain? Yes, but too inefficiently to be practical. Consider the very first step of the binary search algorithm:

Look at the "middle" item in the dictionary

If $n$ items are in a linked chain, how can you possibly get to the middle one? You can traverse the chain from its beginning until you have visited $n / 2$ items. But, as you will see in the answer to the second question posed before, just this first step will often take longer than the entire binary search of an array. Further, you would have the same problem of finding the "middle" item at each recursive step. It is thus not practical to perform a binary search for the linear link-based implementation. This observation is extremely significant.

On the other hand, if $n$ items are in an array items, the middle item is at location $n / 2$ and can be accessed directly. Thus, a binary search of an array requires considerably less time than an algorithm that must inspect every item in the dictionary. What does "considerably less time" mean? As you know, without the ability to perform a binary search, you may have to inspect every item in the dictionary, either to locate an item with a particular search key or to detect that such an item is not present. In other words, if a dictionary has size $n$, you will have to inspect as many as $n$ items; thus, such a search is $\mathrm{O}(n)$. How much better can you do with a binary search? Recall from Chapter 10 that a binary search is $\mathrm{O}\left(\log _{2} n\right)$ in its worst case and that an $\mathrm{O}\left(\log _{2} n\right)$ algorithm is substantially more efficient than an $\mathrm{O}(n)$ algorithm. For example, $\log _{2} 1,024=10$ and $\log _{2} 1,048,576=20$. For a large dictionary, the binary search has an enormous advantage.

Because a thesaurus is probably large, you must choose an implementation for which a binary search is practical. As you have just seen, this observation eliminates the linear link-based implementations. The sorted array-based implementation is fine here, because you know the size of the thesaurus.

An implementation using a binary search tree is also a good choice for retrieval-dominated applications. As you saw in Chapter 15, searching a binary search tree is $\mathrm{O}(\log n)$ if the tree is balanced. Because the thesaurus does not change, you can create a balanced tree that remains balanced and be assured of an efficient search. Although the pointers in a binary search tree add a space cost, as scenario A mentioned, this cost is relatively insignificant when the data items are large.

Scenario C: Insertion, removal, retrieval, and traversal in sorted order. If your local library has computerized its catalog of books, you perform a retrieval operation when you access this catalog. The library staff uses insertion and removal operations to update the catalog and a traversal to save the entire catalog in a file. Presumably, retrieval is the most frequent operation, but the other operations are not infrequent enough to ignore. If they were, this scenario would be the same as scenario B!

To insert into a dictionary an item whose search key is $X$, you must first determine where the item belongs in the dictionary's sorted order. Similarly, to remove from the dictionary an item that has $X$ as its search key, you must first locate the item. Thus, both the add and remove operations perform the following steps:

1. Find the appropriate position in the dictionary.
2. Insert into (or remove from) this position.

Step 1 is far more efficient if the dictionary implementation is array based instead of link based. For an array-based implementation, you can use a binary search to determine-in the case of insertionwhere the new item $X$ belongs and-in the case of removal-where the item is located. On the other hand, for a link-based implementation, you know from the discussion in scenario B that a binary search is impractical, and so you must traverse the chain from its beginning. You also saw in scenario B that it takes significantly less time to perform a binary search of an array than it does to traverse a linked chain.

Thus, because it facilitates a binary search, the array-based implementation is superior with respect to step 1 of add and remove. However, as you may have guessed, the link-based implementation is better for step 2, the actual insertion or removal of the item. Under the array-based implementation, add must shift array entries to make room for the new item, as Figure 18-7a illustrates. The worst case would require that every array entry be shifted. On the other hand, under the link-based

A binary search is impractical with a link-based implementation

If you know the dictionary's maximum size, a sorted array-based implementation is appropriate for frequent retrievals

If you do not know the dictionary's maximum size, use a binary search tree in the implementation

Both add and remove perform these two steps

Use an array-based implementation for step 1

Use a link-based implementation for step 2

A sorted array-based implementation shifts data during insertions and removals

The sorted linear implementations are comparable here, but none is suitable

FIGURE 18-7 Insertion for sorted linear implementations: (a) array based; (b) pointer based

implementation, you can accomplish this second step simply by changing at most two pointers, as Figure 18-7b indicates. Similar comments are true for the remove operation. It must shift array entries to fill in the gap created when an item is removed, but it needs to change only two pointers at most to remove an item from a linked chain.

When you take steps 1 and 2 together, you will find that the sorted array-based and sorted linkbased implementations of add or remove both require roughly the same amount of time - they are both $\mathrm{O}(n)$. Neither implementation supports these two operations particularly well. The binary search tree implementation, however, combines the best features of the two linear implementations. Because it is link based, you avoid shifting data, and the dictionary can grow dynamically as needed. You can also retrieve items from a binary search tree efficiently.

D
Note: When designing an ADT to solve a particular problem, do not include unnecessary operations. The proper choice of an implementation depends on the mix of requested operations, and if you request an operation that you do not need, you might get an implementation that does not best support what you are really doing.

Despite certain difficulties, linear implementations of a dictionary can be appropriate

Summary. Although linear implementations of a dictionary are less sophisticated and generally require more time to perform their operations than a binary search tree implementation, they are nevertheless useful for many applications. Because linear implementations are easy to understand conceptually, they are appropriate for dictionaries that will contain only a small number of entries. In such cases efficiency is not as great a concern as are simplicity and clarity. Even when a dictionary is large, a linear implementation may still be appropriate for applications that can use an unsorted dictionary and have few removals.

## Note: Comparing linear implementations of the ADT dictionary

- An unsorted array-based implementation of the ADT dictionary can efficiently insert an item at the end of an array. A removal, however, will usually require shifting data so that no gap remains in the array. Because the items are unsorted, retrieval will require a sequential search.
- A sorted array-based implementation usually requires shifting data during both insertions and removals. Retrieval, however, can use an efficient binary search because the items are sorted.
- An unsorted link-based implementation can efficiently insert an item at the beginning of a linked chain. A removal will require a sequential search but no data shifts. Retrieval will also require a sequential search.
- A sorted link-based implementation requires a sequential search but no data shifts during both insertions and removals. Retrieval will also require a sequential search.

The nonlinear, binary search tree implementation of the ADT dictionary can be a better choice than a linear implementation, in general. If an $n$-node binary search tree has minimum height-that is, has height $\left\lceil\log _{2}(n+1)\right\rceil$-the binary search tree implementation of the ADT dictionary certainly succeeds where the linear implementations failed: You can, with efficiency comparable to that of a binary search, locate an item in both the retrieval operation and the first steps of the add and remove operations. In addition, the link-based implementation of the binary search tree permits dynamic allocation of its nodes, so that it can handle a dictionary whose maximum size is unknown.

This implementation also efficiently performs the second step of the insertion and removal operations: The actual insertion or removal of a node requires only a few pointer changes - plus a short traversal to the inorder successor if the node to be removed has two children - rather than the possible shifting of all the dictionary entries, as the array-based implementations require. The binary search tree implementation therefore combines the best aspects of the two linear implementations, yet avoids their disadvantages.

As Chapter 15 showed, however, the height of a binary search tree depends on the order in which you perform the add and remove operations on the tree and can be as large as $n$. If the insertion and removal operations occur in a random order, the height of the binary search tree will be quite close to its minimum value. You do need to watch for a possible increase in the tree's height, however, and the resulting decrease in performance. If instead you use a variation of the binary search tree that remains balanced-as the next chapter will describe-you can keep the height of the tree near $\log _{2} n$.

Figure 18-8 summarizes the order of the insertion, removal, retrieval, and traversal operations for the dictionary implementations discussed so far in this chapter.

FIGURE 18-8 The average-case order of the ADT dictionary operations for various implementations

|  | Insertion | Removal | Retrieval | Traversal |
| :---: | :---: | :---: | :---: | :---: |
| Unsorted array-based | O(1) | $\mathrm{O}(n)$ | $\mathrm{O}(n)$ | $\mathrm{O}(\mathrm{n})$ |
| Unsorted link-based | $\mathrm{O}(1)$ | $\mathrm{O}(\mathrm{n})$ | $\mathrm{O}(\mathrm{n})$ | $\mathrm{O}(\mathrm{n})$ |
| Sorted array-based | $\mathrm{O}(\mathrm{n})$ | $\mathrm{O}(\mathrm{n})$ | $\mathrm{O}(\log n)$ | $\mathrm{O}(\mathrm{n})$ |
| Sorted link-based | $\mathrm{O}(\mathrm{n})$ | $\mathrm{O}(\mathrm{n})$ | $\mathrm{O}(\mathrm{n})$ | $\mathrm{O}(\mathrm{n})$ |
| Binary search tree | $\mathrm{O}(\log n)$ | $\mathrm{O}(\log n)$ | $\mathrm{O}(\log n)$ | $\mathrm{O}(\mathrm{n})$ |

A binary search tree implementation is a better choice than a linear implementation, in general

A balanced binary search tree increases the efficiency of the ADT dictionary operations

## Note: Perspective

You might wonder why, if the binary search tree implementation of the ADT dictionary is so good, we studied the linear implementations at all. There are three reasons. The first and foremost reason is perspective. Section 10.2.4 of Chapter 10 spoke of the dangers of overanalyzing a problem. If the size of the problem is small, the difference in efficiency among the possible solutions is likely insignificant. In particular, if the size of the dictionary is small, a linear implementation is adequate and simple to understand.

The second reason is efficiency: A linear implementation can be quite efficient for certain situations. For example, a linear implementation was best for scenario A, where the predominant operations are insertion and traversal in no particular order. For scenario B , where the predominant operation is retrieval, the sorted array-based implementation is adequate, if the maximum number of items is known. For these situations, a concern for simplicity suggests that you use a linear implementation and not a binary search tree, even for large dictionaries.

The third reason is motivation. By seeing scenarios for which the linear implementations are not adequate, you are forced to look beyond arrays and consider other implementations, such as the binary search tree. Actually looking at both a linear implementation and a binary search tree implementation allows you to see these inadequacies more clearly.

Dictionary operations without searches

### 18.4 Hashing

VideoNote Hashing

The binary search tree provides an excellent implementation of the ADT dictionary, as do the balanced search trees discussed in Chapter 19. They allow you to perform all of the dictionary operations quite efficiently. If, for example, a dictionary contains 10,000 items, the operations getItem, add, and remove each require approximately $\log _{2} 10,000 \approx 13$ steps. As impressive as this efficiency may be, situations do occur for which the search-tree implementations are not adequate.

As you know, time can be vital. For example, when a person calls the 911 emergency system, the system detects the caller's telephone number and searches a database for the caller's address. Similarly, an air traffic control system searches a database of flight information, given a flight number. Clearly these searches must be rapid.

A radically different strategy is necessary to locate (and insert or remove) an item virtually instantaneously. Imagine an array table of $N$ items-with each array slot capable of holding a single dictionary item - and a seemingly magical box called an "address calculator." Whenever you have a new item that you want to insert into the dictionary, the address calculator will tell you where you should place it in the array. Figure 18-9 illustrates this scenario.

You can thus easily perform an insertion into the dictionary as follows:

```
add(newItem: ItemType): boolean
    \(i=\) the array index that the address calculator gives you for newItem's search key
    table[i] = newItem
```

This add operation is $\mathrm{O}(1)$; that is, it requires constant time.
You also use the address calculator for the getItem and remove operations. If you want to retrieve an item that has a particular search key, you simply ask the address calculator to tell you where it would insert such an item. Because you would have inserted the item earlier by using the add algorithm just given, if the desired item is present in the dictionary, it will be in the array location that the address calculator specifies.


Thus, the getItem operation appears in pseudocode as follows:

```
getItem(searchKey: KeyType): ItemType throw NotFoundException
    i = the array index that the address calculator gives you for an item whose search
        keyequals searchKey
    if (table[i].getKey() == searchKey)
        return tab7e[i]
    else
        Throw NotFoundException
```

Similarly, the pseudocode for the remove operation is

```
remove(searchKey: KeyType): boolean
    i = the array index that the address calculator gives you for an item whose search
        keyequals searchKey
    if (tab7e[i].getKey() == searchKey)
    {
        Remove the item from table[i]
        isSuccessful = true
    }
    else
        isSuccessful = false
    return isSuccessful
```

It thus appears that you can perform the operations getItem, add, and remove virtually instantaneously. You never have to search for an item; instead, you simply let the address calculator determine where the item should be. The amount of time required to carry out the operations is $\mathrm{O}(1)$ and depends only on how quickly the address calculator can perform this computation.

If you are to implement such a scheme, you must, of course, be able to construct an address calculator that can, with very little work, tell you where a given item should be. Address calculators are actually not as mysterious as they seem; in fact, many exist that can approximate the idealized behavior just described. Such an address calculator is usually referred to as a hash function. The scheme just described is an idealized description of a technique known as hashing, and the array table is called the hash table.

To understand how a hash function works, consider the 911 emergency system mentioned earlier. If, for each person, the system had a record whose search key was the person's telephone number, it could store these records in a search tree. Although searching a tree would be fast, faster access to a particular record would be possible by storing the records in an array table, as follows. You store the record for a person whose telephone number is $t$ into table[t]. Retrieval of the record, then, is

A hash function tells you where to place an item in an array called a hash table

A hash function maps an integer into an array index

A perfect hash function maps each search key into a unique location of the hash table

A perfect hash function is possible if you know all the search keys
almost instantaneous given its search key t. For example, you can store the record for the telephone number 123-4567 in tab7e [1234567]. If you can spare 10 million memory locations for table, this approach is fine. You need not use memory so extravagantly, however, because 911 systems are regional. If you consider only one telephone exchange, for example, you can store the record for the number 123-4567 in table [4567] and get by with an array table of 10,000 locations.

The transformation of 1234567 into an array index 4567 is a simple example of a hash function. A hash function $h$ must take an arbitrary integer $x$ and map it into an integer that you can use as an array index. In our example, such indices would be in the range 0 through 9999. That is, $h$ is a function such that for any integer $x$,
$h(x)=i$, where $i$ is an integer in the range 0 through 9999
Because the database contains entries for every telephone number in a particular exchange, the array table is completely full. In this sense, our example is not typical of hashing applications and serves only to illustrate the idea of a hash function. What if many fewer entries were in the array? Consider, for example, an air traffic control system that stores an entry for each current flight according to its four-digit flight number. You could store an entry for Flight 4567 in table [4567], but you still would need an array of 10,000 locations, even if only 50 flights were current.

A different hash function would save memory. If you allow space for a maximum of 101 flights, for example, so that the array table has indices 0 through 100, the necessary hash function $h$ should map any four-digit flight number into an integer in the range 0 through 100.

If you have such a hash function $h$-and you will see several suggestions for hash functions later-the dictionary operations are easy to write. For example, in the getItem algorithm, the step
$\begin{aligned} & i= \text { the array index that the address calculator gives you for an item whose search } \\ & \text { key equals searchKey }\end{aligned}$
is implemented simply as
i $=$ h (searchKey)
In the previous example, searchKey would be the flight number.
Although the dictionary operations appear to be virtually instantaneous, is hashing really as good as it sounds? If it really was this good, there would have been little reason for developing all those other dictionary implementations. Hashing would beat them hands down!

Why is hashing not quite as simple as it seems? You might first notice that since the hashing scheme stores the items in an array, it would appear to suffer from the familiar problems associated with a fixed-size implementation. Obviously, the hash table must be large enough to contain all of the items that you want to store. This requirement is not the crux of the implementation's difficulty, however, for-as you will see later-there are ways to allow the hash table to grow dynamically. But the implementation does have a major pitfall, even given the assumption that the number of items to be stored will never exceed the size of the hash table.

Ideally, you want the hash function to map each search key $x$ into a unique integer $i$. The hash function in the ideal situation is called a perfect hash function. In fact, it is possible to construct perfect hash functions if you know all of the possible search keys that actually occur in the dictionary. You have this knowledge for the 911 example, since everyone is in the database, but not for the air traffic control example. Usually, you will not know the values of the search keys in advance.

In practice, a hash function can map two or more search keys $x$ and $y$ into the same integer. That is, the hash function tells you to store two or more items in the same array location table[i]. This occurrence is called a collision. Thus, even if fewer than 101 items were present in the hash table table [0..100], $h$ could very well tell you to place more than one item into the same array location. For example, if two items have search keys 4567 and 7597, and if

$$
h(4567)=h(7597)=22
$$

$h$ will tell you to place the two items into the same array location, tab7e [22]. That is, the search keys 4567 and 7597 have collided.

Even if the number of items that can be in the array at any one time is small, the only way to avoid collisions completely is for the hash table to be large enough that each possible search-key value can have its own location. If, for example, Social Security numbers were the search keys, you would need an array location for each integer in the range 000000000 through 999999999 . This situation would certainly require a good deal of storage! Because reserving vast amounts of storage is usually not practical, collision-resolution schemes are necessary to make hashing feasible. Such schemes usually require that the hash function place items evenly throughout the hash table.

To summarize, a typical hash function must

- Be easy and fast to compute
- Place items evenly throughout the hash table

Note that the size of the hash table affects the ability of the hash function to distribute the items evenly throughout the table. The requirements of a hash function will be discussed in more detail later in this chapter.

Consider now several hash functions and collision-resolution schemes.

### 18.4.1 Hash Functions

It is sufficient to consider hash functions that have an arbitrary integer as an argument. Why? If a search key is not an integer, you can simply map the search key into an integer, which you then hash. At the end of this section, you will see one way to convert a string into an integer.

There are many ways to convert an arbitrary integer into an integer within a certain range, such as 0 through 100. Thus, there are many ways to construct a hash function. Many of these functions, however, will not be suitable. Here are several simple hash functions that operate on positive integers.
Selecting digits. If your search key is the nine-digit employee ID number 001364825, you could select the fourth digit and the last digit, to obtain 35 as the index to the hash table. That is,
$h(001364825)=35($ select the fourth and last digits $)$
Therefore, by using digit selection to define your hash function, you would store the item whose search key is 001364825 in tab1e [35].

You do need to be careful about which digits you choose in a particular situation. For example, the first three digits of a Social Security number are based on the geographic region in which the number was assigned. If you select only these digits, you will map all people from the same state into the same location of the hash table.

Digit-selection hash functions are simple and fast, but generally they do not evenly distribute the items in the hash table. A hash function really should utilize the entire search key.

Folding. One way to improve on the previous approach of selecting digits is to add the digits. The resulting process is known as folding. For example, you can add all of the digits in 001364825 to obtain

$$
0+0+1+3+6+4+8+2+5=29 \quad(\text { add the digits })
$$

Therefore, you would store the item whose search key is 001364825 in table[29]. Notice that if you add all of the digits from a nine-digit search key,
$0 \leq h($ search key $) \leq 81$

Collisions occur when the hash function maps more than one item into the same array location

Requirements for a hash function

It is sufficient for hash functions to operate on integers

Digit selection does not distribute items evenly in the hash table

Applying more than one hash function to a single search key

The dictionary size should be prime

That is, you would use only table [0] through table [81] of the hash table. To change this situation or to increase the size of the hash table, you can group the digits in the search key and add the groups. For example, you could form three groups of three digits from the search key 001364825 and add them as follows:

$$
001+364+825=1,190
$$

For this hash function,

$$
0 \leq h(\text { search key }) \leq 3 \times 999=2,997
$$

Clearly, if 2,997 is larger than the size of the hash table that you want, you can alter the groups that you choose. Perhaps not as obvious is that you can apply more than one hash function to a search key. For example, you could select some of the digits from the search key before adding them, or you could either select digits from the previous result 2,997 or apply folding to it once again by adding 29 and 97.

Modulo arithmetic. Modulo arithmetic provides a simple and effective hash function that we will use in the rest of this chapter. For example, consider the function ${ }^{2}$

$$
h(x)=x \bmod \text { tableSize }
$$

where the hash table has tableSize elements. In particular, if tableSize is $101, h(x)=x \bmod 101$ maps any integer $x$ into the range 0 through 100. For example, $h$ maps 001364825 into 12.

For $h(x)=x \bmod$ tableSize, many $x$ 's map into table [0], many $x$ 's map into table [1], and so on. That is, collisions occur. However, you can distribute the dictionary items evenly over all of tablethus reducing collisions-by choosing a prime number as tableSize. For instance, 101 in the previous example is prime. The choice of table size will be discussed in more detail later in this chapter. For now, realize that 101 is used here as a simple example of a prime table size. For the typical hash table, 101 is much too small.

Converting a character string to an integer. If your search key is a character string-such as a name-you could convert it into an integer before applying the hash function $h(x)$. To do so, you could first assign an integer value to each character in the string. For example, for the word "NOTE" you could assign the ASCII values $78,79,84$, and 69 , to the letters $N, O, T$, and $E$, respectively. Or, if you assign the values 1 through 26 to the letters $A$ through $Z$, you could assign 14 to $N, 15$ to $O, 20$ to $T$, and 5 to $E$.

If you now simply add these numbers, you will get an integer, but it will not be unique to the character string. For example, the word "TONE" will give you the same result. Instead, you can write the numeric value for each character in binary and concatenate the results. If you assign the values 1 through 26 to the letters $A$ through $Z$, you obtain the following result for the word "NOTE":
$N$ is 14 , or 01110 in binary
$O$ is 15 , or 01111 in binary
$T$ is 20, or 10100 in binary
$E$ is 5 , or 00101 in binary
Concatenating the binary values gives you the binary integer
01110011111010000101
which is 474,757 in decimal. You can apply the hash function $x \bmod$ tableSize for $x=474,757$.

2 Remember that this book uses "mod" as an abbreviation for the mathematical operation modulo. In $\mathrm{C}++$, the modulo operator is $\%$.

Now consider a more efficient way to compute 474,757. Rather than converting the previous binary number to decimal, you can evaluate the expression

$$
14 \times 32^{3}+15 \times 32^{2}+20 \times 32^{1}+5 \times 32^{0}
$$

This computation is possible because we have represented each character as a 5-bit binary number, and $2^{5}$ is 32 .

By factoring this expression, you can minimize the number of arithmetic operations. This technique is called Horner's rule and results in

$$
((14 \times 32+15) \times 32+20) \times 32+5
$$

Horner's rule minimizes the number of computations

Although both of these expressions have the same value, the result in either case could very well be larger than a typical computer can represent; that is, an overflow can occur. If you use the hash function

$$
h(x)+x \bmod \text { tableSize }
$$

you can prevent an overflow by applying the modulo operator after computing each parenthesized expression in Horner's rule. The implementation of this algorithm is left as an exercise.

```
!
Programming Tip: You can use the following hash function-named getHashIndex-in your programs. C++ Interlude 7 will discuss the Standard Template Library (STL) and make sense of the way that this function uses the STL.
```

```
template <class KeyType, class ItemType>
```

template <class KeyType, class ItemType>
int HashedDictionary<KeyType, ItemType>::
int HashedDictionary<KeyType, ItemType>::
getHashIndex(const KeyType\& key) const
getHashIndex(const KeyType\& key) const
{
{
// We are creating a hash function type called hashFunction that hashes
// We are creating a hash function type called hashFunction that hashes
// a search key. First we create an unordered_map object for our KeyType
// a search key. First we create an unordered_map object for our KeyType
// and ItemType.
// and ItemType.
std::tr1::unordered_map<KeyType, ItemType> mapper;
std::tr1::unordered_map<KeyType, ItemType> mapper;
// Then we invoke the method hash_function to return the hash function
// Then we invoke the method hash_function to return the hash function
// for the KeyType and assign it to 'hashFunction'
// for the KeyType and assign it to 'hashFunction'
typename std::tr1::unordered_map<KeyType, ItemType> ::
typename std::tr1::unordered_map<KeyType, ItemType> ::
hasher hashFunction = mapper.hash_function();
hasher hashFunction = mapper.hash_function();
// Need static_cast because hashFunction returns an unsigned long.
// Need static_cast because hashFunction returns an unsigned long.
return static_cast<int>(hashFunction(searchKey) % hashTableSize);
return static_cast<int>(hashFunction(searchKey) % hashTableSize);
} // end getHashIndex

```
} // end getHashIndex
```


### 18.4.2 Resolving Collisions

Consider the problems caused by a collision. Suppose that you want to insert an item whose search key is 4567 into the hash table table, as was described previously. The hash function $h(x)=x \bmod$ 101 tells you to place the new item in table[22], because $4567 \bmod 101$ is 22 . Suppose, however, that table[22] already contains an item, as Figure 18-10 illustrates. If earlier you had placed 7597 into table [22] because $7597 \bmod 101$ equals 22, where do you place the new item? You certainly do not want to disallow the insertion on the grounds that the dictionary is full: You could have a collision even when inserting into a dictionary that contains only one item!

Two approaches to collision resolution

Begin at the hash location and search the dictionary sequentially

Three states: occupied, empty, removed

## FIGURE 18-10 A collision



Two general approaches to collision resolution are common. One approach places the new item into another location within the hash table. A second approach changes the structure of the hash table so that each location table[i] can accommodate more than one item. The collision-resolution schemes described next exemplify these two approaches.
Approach 1: Open addressing. During an attempt to insert a new item into a dictionary, if the hash function indicates a location in the hash table that is already occupied, you look-or probe-for some other empty, or open, location in which to place the item. The sequence of locations that you examine is called the probe sequence.

Such schemes are said to use open addressing. The concern, of course, is that you must be able to find a dictionary item efficiently after you have inserted it. That is, the remove and getItem operations must be able to reproduce the probe sequence that add used and must do so efficiently.

The difference among the various open-addressing schemes is the technique used to probe for an empty location. We briefly describe three such techniques.
Linear probing. In this simple scheme to resolve a collision, you search the hash table sequentially, starting from the original hash location. More specifically, if table [h(searchKey)] is occupied, you check the dictionary locations table[h(searchKey) +1$]$, table[h(searchKey) +2 ], and so on until you find an available location. Figure 18-11 illustrates the placement of four items that all hash into the same location table [22] of the hash table, assuming a hash function $h(x)=x \bmod 101$. Typically, you wrap around from the last array location to the first array location if necessary.

In the absence of removals, the implementation of getItem under this scheme is straightforward. You need only follow the same probe sequence that add used until you either find the item you are searching for; reach an empty location, which indicates that the item is not present; or visit every table location.

Removals, however, complicate matters slightly. The remove operation itself is no problem. You merely find the desired item, as in getItem, and remove it from the hash table, making the location empty. But now what happens when getItem needs to locate an item? The new empty locations that remove created along a probe sequence could cause getItem to stop prematurely, incorrectly indicating a failure. You can solve this problem by placing a table location into one of three states: occupied

FIGURE 18-11 Linear probing with $h(x)=x \bmod 101$

(currently in use), empty (has not been used), or removed (was once occupied but is now available). You then modify the getItem operation to continue probing when it encounters a location in the removed state. Similarly, you modify add to insert into locations that are in either the empty or removed states.

One of the problems with the linear-probing scheme is that items tend to cluster together in the hash table. That is, the table contains groups of consecutively occupied locations. This phenomenon is called primary clustering. Clusters can get close to one another and, in fact, merge into a larger cluster. Large clusters tend to get even larger. Thus, one part of the hash table might be quite densely populated, even though another part has relatively few items.

Note: Primary clustering causes long probe searches and therefore decreases the overall efficiency of hashing.

Question 6 Write the pseudocode for the remove operation when linear probing is used to implement the hash table.

Quadratic probing. You can virtually eliminate primary clusters simply by adjusting the linear probing scheme just described. Instead of probing consecutive table locations from the original hash locationtable[h(searchKey)], youchecklocationstable[h(searchKey) $\left.+1^{2}\right]$, table[h(searchKey) $\left.+2^{2}\right]$, table $\left[h\right.$ (searchKey) $\left.+3^{2}\right]$, and so on until you find an available location. Figure 18-12 illustrates this open-addressing scheme-which is called quadratic probing-for the same items that appear in Figure 18-11.

Unfortunately, when two items hash into the same location, quadratic probing uses the same probe sequence for each item. The resulting phenomenon-called secondary clustering-delays the resolution of the collision. Although the research of quadratic probing remains incomplete, it appears that secondary clustering is not a problem.

Clustering can be a problem

A hash address and a step size determine the probe sequence

Guidelines for the step-size function $\mathrm{h}_{2}$

Primary and secondary hash functions

FIGURE 18-12 Quadratic probing with $h(x)=x \bmod 101$


Doub7e hashing. Double hashing, which is yet another open-addressing scheme, drastically reduces clustering. The probe sequences that both linear probing and quadratic probing use are key independent. For example, linear probing inspects the locations in the hash table sequentially no matter what the hash key is. In contrast, double hashing defines key-dependent probe sequences. In this scheme the probe sequence still searches the hash table in a linear order, starting at the location $h_{1}$ (key), but a second hash function $h_{2}$ determines the size of the steps taken.

Although you choose $h_{1}$ as usual, you must follow these guidelines for $h_{2}$ :

$$
\begin{aligned}
& h_{2}(k e y) \neq 0 \\
& h_{2} \neq h_{1}
\end{aligned}
$$

Clearly, you need a nonzero step size $h_{2}(\mathrm{key})$ to define the probe sequence. In addition, $h_{2}$ must differ from $h_{1}$ to avoid clustering.

For example, let $h_{1}$ and $h_{2}$ be the primary and secondary hash functions defined as

$$
\begin{aligned}
& h_{1}(\text { key })=k e y \bmod 11 \\
& h_{2}(\text { key })=7-(\text { key } \bmod 7)
\end{aligned}
$$

where a hash table of only 11 items is assumed, so that you can readily see the effect of these functions on the hash table. If key $=58, h_{1}$ hashes key to dictionary location $3(58 \bmod 11)$, and $h_{2}$ indicates that the probe sequence should take steps of size $5(7-58 \bmod 7)$. In other words, the probe sequence will be $3,8,2$ (wraps around), 7,1 (wraps around), $6,0,5,10,4,9$. On the other hand, if key $=14, h_{1}$ hashes key to table location $3(14 \bmod 11)$, and $h_{2}$ indicates that the probe sequence should take steps of size $7(7-14 \bmod 7)$, and so the probe sequence would be $3,10,6,2,9,5,1,8,4,0$.

Each of these probe sequences visits all the table locations. This phenomenon always occurs if the size of the hash table and the size of the probe step are relatively prime, that is, if their greatest common divisor is 1 . Because the size of a hash table is commonly a prime number, it will be relatively prime to all step sizes.

Figure 18-13 illustrates the insertion of 58,14 , and 91 into an initially empty hash table. Because $h_{1}(58)$ is 3 , you place 58 into table [3]. You then find that $h_{1}(14)$ is also 3 , so to avoid a collision, you step by $h_{2}(14)=7$ and place 14 into table [3+7], or table[10]. Finally, $h_{1}(91)$ is 3 and $h_{2}(91)$ is 7 .

FIGURE 18-13 Double hashing during the insertion of 58 , 14, and 91


Because table[3] is occupied, you probe tab7e[10] and find that it, too, is occupied. You finally store 91 in table [ $(10+7) \% 11]$, or table [6].

While more than two hash functions can be desirable, such schemes are difficult to implement.

Question 7 What is the probe sequence that double hashing uses when
$h_{1}($ key $)=$ key $\bmod 11, h_{2}($ key $)=7-($ key $\bmod 7)$, and $k e y=19 ?$

Increasing the size of the hash table. With any of the open-addressing schemes, as the hash table fills, the probability of a collision increases. At some point, a larger hash table becomes desirable. If you use a dynamically allocated array for the hash table, you can increase its size whenever the dictionary becomes too full.

You cannot simply double the size of the array, as we did in earlier chapters, because the size of the hash table must remain prime. In addition, you do not simply move the items from the original hash table to the new hash table. If your hash function is $x \bmod$ tableSize, it changes as tableSize changes. Thus, you need to apply your new hash function to every item in the old hash table before placing it into the new hash table.

## Note: Rehashing

After creating a new, larger hash table of an appropriate size, you use the dictionary method add to add each item in the original hash table to the new table. The method computes the hash index using the size of the new table and handles any collisions. This process of enlarging a hash table and computing new hash indices for its contents is called rehashing. You can see that increasing the size of a hash table requires considerably more work than increasing the size of an ordinary array. Rehashing is a task that you should not do often.

Each hash-table location can accommodate more than one item

A bucket is an element of a hash table that is itself an array

Each hashdictionary location is a linked chain

Separate chaining successfully resolves collisions

The load factor measures how full a hash table is

Approach 2: Restructuring the hash table. Another way to resolve collisions is to change the structure of the array table - the hash table-so that it can accommodate more than one item in the same location. We describe two such ways to alter the hash table.
Buckets. If you define the hash table so that each location table[i] is itself an array-called a bucket-you can store the items that hash into table[i] in this array. The problem with this approach, of course, is choosing the size $b$ of each bucket. If $b$ is too small, you will only have postponed the problem of collisions until $b+1$ items map into some array location. If you attempt to make $b$ large enough so that each array location can accommodate the largest number of items that might map into it, you are likely to waste a good deal of storage.
Separate chaining. A better approach than using buckets is to design the hash table as an array of linked chains. In this collision-resolution technique, known as separate chaining, each entry table[i] is a pointer to a chain of linked nodes containing the items that the hash function has mapped into location i, as Figure 18-14 illustrates. Separate chaining provides a successful approach to resolving collisions. With separate chaining, the size of the dictionary is dynamic and can exceed the size of the hash table, because each linked chain can be as long as necessary. As you will see in the next section, the length of these chains affects the efficiency of retrievals and removals. Even so, separate chaining is the most time-efficient collision-resolution scheme.

We will discuss an implementation of the ADT dictionary using hashing and separate chaining in Section 18.4.6 of this chapter.

FIGURE 18-14 Separate chaining



Question 8 If $h(x)=x \bmod 7$ and separate chaining resolves collisions, what does the hash table look like after the following insertions occur: $8,10,24,15,32,17$ ? Assume that each item contains only a search key.

### 18.4.3 The Efficiency of Hashing

An analysis of the average-case efficiency of hashing involves the load factor alpha $(\boldsymbol{\alpha})$, which is the ratio of the current number of items in the dictionary to the maximum size of the array dictionary. That is,

$$
\alpha=\frac{\text { Current number of table items }}{\text { tableSize }}
$$

$\alpha$ is a measure of how full the hash table is. As the table fills, $\alpha$ increases and the chance of collision increases, so search times increase. Thus, hashing efficiency decreases as $\alpha$ increases.

Unlike the efficiency of earlier dictionary implementations, the efficiency of hashing does not depend solely on the number $n$ of items in the dictionary. While it is true that for a fixed tableSize efficiency decreases as $n$ increases, for a given $n$ you can choose tableSize to increase efficiency. Thus, when determining the size of the hash table, you should estimate the largest possible $n$ and select tableSize so that $\alpha$ is small. As you will see shortly, $\alpha$ should not exceed $2 / 3$.

Hashing efficiency for a particular search also depends on whether the search is successful. An unsuccessful search requires more time in general than a successful search. The following analyses ${ }^{3}$ enable a comparison of collision-resolution techniques.
Linear probing. For linear probing, the approximate average number of comparisons that a search requires is

$$
\begin{array}{ll}
\frac{1}{2}\left[1+\frac{1}{1-\alpha}\right] & \text { for a successful search, and } \\
\frac{1}{2}\left[1+\frac{1}{(1-\alpha)^{2}}\right] & \text { for an unsuccessful search }
\end{array}
$$

As collisions increase, the probe sequences increase in length, causing increased search times. For example, for a dictionary that is two-thirds full $(\alpha=2 / 3)$, an average unsuccessful search might require at most five comparisons, or probes, while an average successful search might require at most two comparisons. To maintain efficiency, it is important to prevent the hash table from filling up.
Quadratic probing and double hashing. The efficiency of both quadratic probing and double hashing is given by

$$
\begin{array}{ll}
\frac{-\log _{e}(1-\alpha)}{\alpha} & \text { for a successful search, and } \\
\frac{1}{1-\alpha} & \text { for an unsuccessful search }
\end{array}
$$

On average, both techniques require fewer comparisons than linear probing. For example, for a dictionary that is two-thirds full, an average unsuccessful search might require at most three comparisons, or probes, while an average successful search might require at most two comparisons. As a result, you can use a smaller hash table for both quadratic probing and double hashing than you can for linear probing. However, because they are open-addressing schemes, all three approaches suffer when you are unable to predict the number of insertions and removals that will occur. If your hash table is too small, it will fill up, and search efficiency will decrease.

Separate chaining. Because the dictionary's add operation places the new item at the beginning of a linked chain within the hash table, it is $\mathrm{O}(1)$. The getItem and remove operations, however, are not as fast. They each require a search of the linked chain of items, so ideally you would like for these chains to be short.

For separate chaining, tableSize is the number of chains, not the maximum number of dictionary items. Thus, it is entirely possible, and even likely, that the current number of dictionary items $n$ exceeds tableSize. That is, the load factor $\alpha$, or $n /$ tableSize, can exceed 1 . Because tableSize is the number of linked chains, $n /$ tableSize -that is, $\alpha$-is the average length of each linked chain.

[^14]Unsuccessful searches generally require more time than successful searches

Do not let the hash table get too full

Open-addressing schemes require a good estimate of the number of insertions and removals

Insertion is instantaneous

Average-case efficiency of retrievals and removals

Some searches of the hash table are unsuccessful because the relevant chain is empty. Such searches are virtually instantaneous. For an unsuccessful search of a nonempty chain, however, getItem and remove must examine the entire chain, or $\alpha$ items in the average case. On the other hand, a successful search must examine a nonempty chain. In the average case, the search will locate the item in the middle of the chain. That is, after determining that the linked chain is not empty, the search will examine $\alpha$ / 2 items.

Thus, the efficiency of the retrieval and removal operations under the separate-chaining approach is

$$
1+\frac{\alpha}{2} \quad \text { for a successful search, and }
$$

$\alpha \quad$ for an unsuccessful search
Even if the linked chains typically are short, you should still estimate the worst case. If you seriously underestimate tableSize, or if most of the dictionary items happen to hash into the same location, the number of items in a chain could be quite large. In fact, in the worst case, all $n$ items in the dictionary could be in the same linked chain!

As you can see, the time that a retrieval or removal operation requires can range from almost nothing-if the linked chain to be searched either is empty or has only a couple of items in it - to the time required to search a chain that contains all the items in the dictionary, if all the items hashed into the same location.

Comparing techniques. Figure $18-15$ plots the relative efficiency of the collision-resolution schemes just discussed. When the hash table is about half full-that is, when $\alpha$ is 0.5 -the techniques are nearly equal in efficiency. As the dictionary fills and $\alpha$ approaches 1 , separate chaining is the most efficient technique. Does this mean that we should discard all other search algorithms in favor of hashing with separate chaining?

FIGURE 18-15 The relative efficiency of four collision-resolution methods


No. The analyses here are average-case analyses. Although an implementation of the ADT dictionary that uses hashing might often be faster than one that uses a search tree, in the worst case it can be much slower. If you can afford both an occasional slow search and a large tableSize-that is, a small $\alpha$-then hashing can be an attractive dictionary implementation. However, if you are performing a life-and-death search for your city's poison control center, a search-tree implementation would at least provide you with a guaranteed limit on its worst-case behavior.

Furthermore, while separate chaining is the most time-efficient collision-resolution scheme, you do have the storage overhead of the pointers in the linked chain. If the data entries in the dictionary are small, the pointers add a significant overhead in storage, and you may want to consider a simpler collision-resolution scheme. On the other hand, if the entries are large, the addition of a pointer is insignificant, so separate chaining is a good choice.

## Note: Maintaining the performance of hashing

Collisions and their resolution typically cause the load factor $\alpha$ to increase and the efficiency of the dictionary operations to decrease. To maintain efficiency, you should restrict the size of $\alpha$ as follows:

- $\alpha<0.5$ for open addressing
- $\alpha<1.0$ for separate chaining

Should the load factor exceed these bounds, you must increase the size of the hash table using rehashing.

### 18.4.4 What Constitutes a Good Hash Function?

Before we conclude this introduction to hashing, consider in more detail the issue of choosing a hash function to perform the address calculations for a given application. A great deal has been written on this subject, most of which is beyond the mathematical level of this book. However, this section will present a brief summary of the major concerns.

- Is the hash function easy and fast to compute? If a hashing scheme is to perform dictionary operations almost instantaneously and in constant time, you certainly must be able to calculate the hash function rapidly. Most of the common hash functions require only a single division (like the modulo operation), a single multiplication, or some kind of "bit-level" operation on the internal representation of the search key. In all these cases, the requirement that the hash function be easy and fast to compute is satisfied.
- Does the hash function scatter the data evenly throughout the hash table? Unless you use a perfect hash function-which is usually impractical to construct-you typically cannot avoid collisions entirely. For example, to achieve the best performance from a separatechaining scheme, each entry table[i] should point to a chain containing approximately the same number of items; that is, each chain should contain approximately $n$ / tableSize items (and thus no chain should contain significantly more than $n$ / tableSize items). To accomplish this goal, your hash function should scatter the search keys evenly throughout the hash table.
- How well does the hash function scatter random data? If every search-key value is equally likely, will the hash function scatter the search keys evenly? For example, consider nine-digit employee ID numbers as the search keys, a hash table table[0..39], and a hash function $h(x)=($ first two digits of $x) \bmod 40$. If all employee ID numbers are equally likely, does a given ID number $x$ have equal probability of hashing into any one of the 40 array locations? For this hash function, the answer is no. Only ID numbers that start with 19,59 , and 99 map into

In the worst case, a hashing implementation of a dictionary can be much slower than other implementations

You cannot avoid collisions entirely

A function that does not scatter random data evenly

A function that does scatter random data evenly

General requirements of a hash function
table[19], while only ID numbers that start with 20 and 60 map into tab7e[20]. In general, three different ID prefixes - that is, the first two digits of an ID number-map into each array location 0 through 19, while only two different prefixes map into each array location 20 through 39. Because all ID numbers are equally likely—and thus all prefixes 00 through 99 are equally likely - a given ID number is 50 percent more likely to hash into one of the locations 0 through 19 than it is to hash into one of the locations 20 through 39. As a result, each array location 0 through 19 would contain, on average, 50 percent more items than each location 20 through 39.

Thus, the hash function

$$
h(x)=(\text { first two digits of } x) \bmod 40
$$

does not scatter random data evenly throughout the array table[0..39]. On the other hand, it can be shown that the hash function

$$
h(x)=x \bmod 101
$$

does, in fact, scatter random data evenly throughout the array table[0. .100].

- How well does the hash function scatter nonrandom data? Even if a hash function scatters random data evenly, it may have trouble with nonrandom data. In general, no matter what hash function you select, it is always possible that the data will have some unlucky pattern that will result in uneven scattering. Although there is no way to guarantee that a hash function will scatter all data evenly, you can greatly increase the likelihood of this behavior.

As an example, consider a hash table table[0..39] and a hash function $h(x)=$ first two digits of $x$. If every ID number is equally likely, $h$ will scatter the search keys evenly throughout the array. But what if every ID number is not equally likely? For instance, a company might assign employee IDs according to department, as follows:

```
10xxxxx Sales
20xxxxx Customer Relations
90xxxxx Data Processing
```

Under this assignment, only 9 out of the 100 array locations would contain any items at all. Further, those locations corresponding to the largest departments (Sales, for example, which corresponds to table[10]) would contain more items than those locations corresponding to the smallest departments. This scheme certainly does not scatter the data evenly.

Much research has been done into the types of hash functions you should use to guard against various types of patterns in the data. The results of this research are really in the province of more advanced courses, but two general principles can be noted here:

- The calculation of the hash function should involve the entire search key. Thus, for example, computing a modulo of the entire ID number is much safer than using only its first two digits.
- If a hash function uses modulo arithmetic, the base should be prime; that is, if $h$ is of the form
$h(x)=x \bmod$ tableSize
then tableSize should be a prime number. This selection of tableSize is a safeguard against many subtle kinds of patterns in the data (for example, search keys whose digits are likely to be multiples of one another). Although each application can have its own particular kind of patterns and thus should be analyzed on an individual basis, choosing a prime number for tableSize is an easy way to safeguard against some common types of patterns in the data.


### 18.4.5 Dictionary Traversal: An Inefficient Operation Under Hashing

For many applications, hashing provides the most efficient implementation of the ADT dictionary. One important dictionary operation-traversal in sorted order-performs poorly when hashing implements the dictionary. As was mentioned previously, a good hash function scatters items as randomly as possible throughout the array, so that no ordering relationship exists between a search key that hashes into table[i] and one that hashes into table[i+1]. As a consequence, if you must traverse the dictionary in sorted order, you first would have to sort the items. If sorting were required frequently, hashing would be a far less attractive implementation than a search tree.

Traversing a dictionary in sorted order is really just one example of a whole class of operations that hashing does not support well. Many similar operations that you often wish to perform on a dictionary require that the items be ordered. For example, consider an operation that must find the dictionary item whose search key is the smallest or largest. If you use a search-tree implementation, these items are in the leftmost and rightmost nodes of the tree, respectively. If you use a hashing implementation, however, you do not know where these items are-you would have to search the entire dictionary. A similar type of operation is a range query, which requires that you retrieve all items whose search keys fall into a given range of values. For example, you might want to retrieve all items whose search keys are in the range 129 to 755 . This task is relatively easy to perform by using a search tree (see Exercise 3 in the next chapter), but if you use hashing, there is no efficient way to answer the range query.

In general, if an application requires any of these ordered operations, you should probably use a search tree. Although the getItem, add, and remove operations are somewhat more efficient when you use hashing to implement the dictionary instead of a balanced search tree, the balanced search tree supports these operations so efficiently itself that, in most contexts, the difference in speed for these operations is negligible (whereas the advantage of the search tree over hashing for the ordered operations is significant).

In the context of external storage, however, the story is different. For data that is stored externally, the difference in speed between hashing's implementation of getItem and a search tree's implementation may well be significant, as you will see in Chapter 21. In an external setting, it is not uncommon to see a hashing implementation of the getItem operation and a search-tree implementation of the ordered operations used simultaneously.

### 18.4.6 Using Hashing and Separate Chaining to Implement the ADT Dictionary

Let's now use hashing to implement the ADT dictionary. We will resolve collisions by using separate chaining.

The hash table. Recall from Figure 18-14 that to use separate chaining, you create a hash table of pointers to chains of linked nodes that represent entries in the dictionary. Although each node in a chain could contain an object of the class Entry, as given earlier in Listing 18-2-and as you might do in any of the linked-based implementations mentioned earlier in this chapter-we will derive a new class, HashedEntry, from Entry and add a pointer field, as Figure 18-16 illustrates. Listing 18-5 shows the declaration of this new class. Notice that HashedEntry is like a combination of the classes Node-which we have used in previous link-based implementations-and Entry. In particular, it has the data members and methods of Entry, as well as the pointer data field and the methods getNext and setNext of Node.

Items hashed into table[i] and table[i+1] have no ordering relationship

Hashing versus balanced search trees

FIGURE 18-16 A dictionary entry when separate chaining is used


## LISTING 18-5 The class HashedEntry

```
/** A class of entry objects for a hashing implementation of the
    ADT dictionary.
    @file HashedEntry.h */
#ifndef _HASHED_ENTRY
#define _HASHED_ENTRY
#include "Entry.h"
template<class KeyType, class ItemType>
class HashedEntry : public Entry<KeyType, ItemType>
{
private:
    HashedEntry<KeyType, ItemType>* nextPtr;
public:
    HashedEntry();
    HashedEntry(ItemType newEntry, KeyType searchKey);
    HashedEntry(ItemType newEntry, KeyType searchKey,
                    HashedEntry<KeyType, ItemType>* nextEntryPtr);
    void setNext(HashedEntry<KeyType, ItemType>* nextEntryPtr);
    HashedEntry<KeyType, ItemType>* getNext() const;
}; // end HashedEntry
#include "HashedEntry.cpp"
#endif
```

Now the data members of our dictionary class, HashedDictionary, are
HashedEntry<KeyType, ItemType>** hashTable; // Array of pointers to entries
int itemCount; // Count of dictionary entries
int hashTableSize; // Table size must be prime
static const int DEFAULT_SIZE = 101;
The array hashTable contains pointers to the itemCount entries in the dictionary. The field hashTableSize is the size of this array. The constructors will set the value of this field, either to DEFAULT_SIZE or to a value given by the client. We note again that this value must be a prime number. Any value given to the constructor will need to be checked to determine whether it is prime and, if not, replaced by a prime number.
The method add. To insert a new entry into the dictionary, you place it at the beginning of the chain that the hash function indicates. Assuming that the initial values of the pointers in the hash table are each nul1ptr, the following pseudocode describes the algorithm for the add operation:

```
add(newItem: ItemType, searchKey: KeyType): boolean
    p = pointer to a new HashedEntry object containing newItem and searchKey
```

```
i = h(searchKey)
if (hashTab7e[i] == nu11ptr)
    hashTab7e[i] = p
else
{
    // Add new entry to beginning of chain
    p->setNext(hashTab1e[i])
    hashTab7e[i] = p
}
return true
```

Recall that we assume that the search key given to add is not already in the dictionary.
The definition of add in the implementation file for HashedDictionary is

```
template <class KeyType, class ItemType>
bool HashedDictionary<KeyType, ItemType>::add(const ItemType& newItem,
                                    const KeyType& searchKey)
{
    // Create entry to add to dictionary
    HashedEntry<KeyType, ItemType>* entryToAddPtr =
                        new HashedEntry<KeyType, ItemType>(newItem, searchKey);
    // Compute the hashed index into the array
    int itemHashIndex = getHashIndex(searchKey);
    // Add the entry to the chain at itemHashIndex
    if (hashTable[itemHashIndex] == nul1ptr)
    {
        hashTab1e[itemHashIndex] = entryToAddPtr;
    }
    else
    {
            entryToAddPtr->setNext(hashTab1e[itemHashIndex]);
            hashTable[itemHashIndex] = entryToAddPtr;
    } // end if
    return true;
} // end add
```

The method remove. To remove an entry, given its search key, from the dictionary, you proceed as the following pseudocode indicates:

```
remove(searchKey: KeyType): boolean
    i = h(searchKey)
    if (hashTab1e[i] != nu11ptr)
    {
        Search the chain pointed to by hashTab7e[i] for searchKey
        if (searchKey is in the first node)
        {
            p = hashTab7e[i]
            hashTab7e[i] = hashTab1e[i]->getNext()
            de1ete p
            return true
        }
        else
        {
            curPtr = pointer to the entry containing searchKey
            prevPtr = pointer to the previous entry in the chain
```

```
        prevPtr->setNext(curPtr->getNext())
        delete curPtr
        return true
    }
}
e7se
    return false
HashedDictionary's definition of remove follows:
template <class KeyType, class ItemType>
boo1 HashedDictionary<KeyType, ItemType>::remove(const KeyType& searchKey)
{
    bool itemFound = false;
    // Compute the hashed index into the array
    int itemHashIndex = getHashIndex(searchKey);
    if (hashTable[itemHashIndex] != nul1ptr)
    {
        // Special case - first node has target
        if (searchKey == hashTable[itemHashIndex]->getKey())
        {
            HashedEntry<KeyType, ItemType>* entryToRemovePtr =
                                    hashTab7e[itemHashIndex];
            hashTab1e[itemHashIndex] = hashTab1e[itemHashIndex]->getNext();
            deTete entryToRemovePtr;
            entryToRemovePtr = nul1ptr; // For safety
            itemFound = true;
        }
        else // Search the rest of the chain
        {
            HashedEntry<KeyType, ItemType>* prevPtr = hashTab7e[itemHashIndex];
            HashedEntry<KeyType, ItemType>* curPtr = prevPtr->getNext();
            while ((curPtr != nul1ptr) && !itemFound )
            {
                // Found item in chain so remove that node
                    if (searchKey == curPtr->getKey())
                {
                        prevPtr->setNext(curPtr->getNext());
                            delete curPtr;
                        curPtr = nul1ptr; // For safety
                            itemFound = true;
                }
                e1se // Look at next entry in chain
                {
                        prevPtr = curPtr;
                        curPtr = curPtr->getNext();
            } // end if
            } // end while
        } // end if
    } // end if
    return itemFound;
} // end remove
```

The rest of the implementation of the class HashedDictionary is left as an exercise.

## Summary

1. The ADT dictionary supports value-oriented operations, such as "Retrieve all the information about John Smith."
2. The linear implementations (array-based and link-based) of a dictionary are adequate only in limited situations, such as when the dictionary is small or for certain operations. In those situations, the simplicity of a linear implementation may be an advantage. A linear implementation of a dictionary, however, is not suitable as a general-purpose, reusable class.
3. A nonlinear link-based (binary search tree) implementation of the ADT dictionary provides the best aspects of the two linear implementations. The link-based implementation allows the dictionary to grow dynamically and allows insertions and removals of data to occur through pointer changes instead of data movement. In addition, the binary search tree allows you to use a binary-search-like algorithm when searching for an item with a specified value. These characteristics make a nonlinear dictionary implementation far superior to the linear implementations in many applications.
4. Hashing as a dictionary implementation calculates where a data item should be rather than searching for it. Hashing allows for very efficient retrievals, insertions, and removals.
5. The hash function should be extremely easy to compute-it should require only a few operations-and it should scatter the search keys evenly throughout the hash table.
6. A collision occurs when two different search keys hash into the same array location. Two ways to resolve collisions are through probing and chaining.
7. Separate chaining is the most time-efficient collision-resolution scheme.
8. Hashing does not efficiently support operations that require the table items to be ordered-for example, traversing the table in sorted order.
9. When dictionary operations such as traversal are not important to a particular application, if you know the maximum number of dictionary items, and if you have ample storage, hashing is an implementation for a dictionary that is simpler and faster than balanced search tree implementations. Tree implementations, however, are dynamic and do not require you to estimate the maximum number of table items.

## Exercises

1. Implement the class Entry, as given in Listing 18-2, and then complete the sorted array-based implementation of the ADT dictionary.
2. Consider an operation replace (replacementItem, searchKey) that locates, if possible, the item in a dictionary with the given search key. If the dictionary contains such an item, the method replaces it with replacementItem.
a. Write implementations of replace for the classes ArrayDictionary, TreeDictionary, and HashedDictionary, as described in this chapter.
b. For TreeDictionary, under what circumstances can replace replace an item without altering the structure of the binary search tree? (See Exercise 6 in Chapter 16.)
3. Imagine an application program that behaves like an English dictionary. The user types a word and the program provides the word's definition. Thus, the dictionary needs only a retrieval operation. Which implementation of the ADT dictionary would be most efficient as an English dictionary?
4. When you use a word processor's spell checker, it compares the words in your document with words in a dictionary. You can add new words to the dictionary as necessary. Thus, this dictionary needs frequent retrievals and occasional insertions. Which implementation of the ADT dictionary would be most efficient as a spell checker's dictionary?
5. A C++ compiler uses a symbol table to keep track of the identifiers that a program uses. When the compiler encounters an identifier, it searches the symbol table to see whether that identifier has already been encountered. If the identifier is new, it is inserted into the table. Thus, the symbol table needs only insertion and retrieval operations. Which implementation of the ADT dictionary would be most efficient as a symbol table?
6. Consider adding operations to the ADT dictionary to form the union and intersection of two given dictionaries. Each operation returns a new dictionary. The union should combine all of the entries in both dictionaries into a third dictionary. The intersection should be a dictionary of the entries common to both of the two dictionaries.

Within each given dictionary, search keys are not repeated. However, an entry in one dictionary could have the same search key as an entry in the second dictionary. Propose and discuss ways to specify these two operations for this case.
7. The implementations of the ADT dictionary given in this chapter make the following assumption: At any time, a dictionary contains at most one item with a given search key. Although the definition of the ADT required for a specific application may not allow duplicates, it is probably wise to test for them rather than simply to assume that they will not occur. Why?

Modify the dictionary implementations so that they test for-and disallow-any duplicates. What dictionary operations are affected? What are the implications for the unsorted linear implementations?
8. Although disallowing duplicate search keys in the ADT dictionary is reasonable for some applications, it is just as reasonable to have an application that will allow duplicates.
a. What are the implications of inserting identical items with the same search keys? What are the implications for the removal and retrieval operations?
b. What are the implications of inserting items that are not identical but have the same search key? Specifically, what would the implementations of add, remove, and getItem do?
9. Suppose that you want to support two remove operations for the ADT dictionary-for example, removeByName and removeByID. Describe an efficient implementation for these operations.
10. Repeat Exercise 9, but instead use a binary search tree to organize one of the search keys and a sorted linked chain to organize the second search key.
11. Can you use a heap instead of a binary search tree as an implementation of the ADT dictionary?
12. Write pseudocode for the dictionary operations add, remove, and getItem when the implementation uses hashing and linear probing to resolve collisions.
13. Write the pseudocode for the remove operation when the implementation uses hashing and separate chaining to resolve collisions.
14. The success of a hash-table implementation of the ADT dictionary is related to the choice of a good hash function. A good hash function is one that is easy to compute and will evenly distribute the possible data. Comment on the appropriateness of the following hash functions. What patterns would hash to the same location?
a. The hash table has size 2,048. The search keys are English words. The hash function is

$$
h(k e y)=(\text { Sum of positions in alphabet of key's letters }) \bmod 2048
$$

b. The hash table has size 2,048. The keys are strings that begin with a letter. The hash function is

$$
h(k e y)=(\text { position in alphabet of first letters key }) \bmod 2048
$$

Thus, "BUT" maps to 2 . How appropriate is this hash function if the strings are random? What if the strings are English words?
c. The hash table is 10,000 entries long. The search keys are integers in the range 0 through 9999 . The hash function is

$$
h(\text { key })=(\text { key } * \text { random }) \text { truncated to an integer }
$$

where random represents a sophisticated random-number generator that returns a real value between 0 and 1.
d. The hash table is 10,000 entries long (HASH_TABLE_SIZE is 10000). The search keys are integers in the range 0 through 9999 . The hash function is given by the following $\mathrm{C}++$ function:

```
int hashIndex(int x)
{
    for (int i = 1; i <= 1000000; i++)
        x = (x * x) % HASH_TABLE_SIZE;
    return x;
} // end hashIndex
```


## Programming Problems

1. Complete the implementation of the class ArrayDictionary as given in Listing 18-3.
2. Complete the implementation of the class TreeDictionary as given in Listing 18-4.
3. Write the sorted link-based, unsorted array-based, and unsorted link-based implementations of the ADT dictionary described in this chapter.
4. Write unsorted and sorted implementations of the ADT dictionary that use, respectively, the ADTs list and sorted list, which Chapters 8 and 12 described.
5. Repeat Programming Problem 5 of Chapter 16 , using the ADT dictionary as the database.
6. Develop a program that can be used to test an implementation of the ADT dictionary.
7. Implement the symbol table described in Exercise 5 by reusing the class TreeDictionary, as described in Section 18.2.2 of this chapter.
8. Consider any collection of data that you can organize in at least two ways. For example, you can order employees by name or by Social Security number and books by title or by author. Note that other information about the employees or books is present in the database but is not used to organize these items. This program assumes that the search keys (for example, book title or book author) are unique and are strings. Thus, in the previous examples, the Social Security number must be a string instead of an integer, and only one book per author is permitted. Choose any set of data that conforms to these requirements, and create a text file.

Program behavior. When your program begins execution, it should read your text file. It then should provide some typical database management operations, all under user control via an interface of your design. For example, you should be able to add an item, remove an item, display (that is, retrieve) an item, and display all of the items in search-key order. You should be able to use either of two search keys to designate the item to be removed or displayed.

Implementation notes. The items in the database should be objects that contain two search keys and additional data, all of which appear in the text file. Thus, you need to design and implement a class of these objects.

Although your program could create two dictionaries from these objects-one organized by one search key (such as the employee name) and the other organized by another search key (such as the Social Security number)-this approach could waste a substantial amount of memory due to the duplication of all of the data in both dictionaries.

A better approach revises the ADT dictionary to provide operations according to two search keys. For example, you want to be able to remove by name and by Social Security number. The underlying data structure for the dictionary's implementation should be a binary search tree. Actually, you will want two binary search trees so that you can organize the data in two ways: by name and by Social Security number, for example.

To avoid duplicated data, store the data in an ADT list and let each node in the binary search trees contain the position of the actual data item in the list, instead of the data item itself.

Your program can be specific to the type of database (employees, books, and so on), or it can be more general. For example, you could determine the search-key descriptions that the user interface displays by requiring that they be in the text file.
9. Repeat Programming Problem 7, but use the class HashedDictionary, as described in Section 18.4.6, which uses separate chaining to resolve collisions. Use the hash function $h(x)=x \bmod$ tableSize and the algorithm that involves Horner's rule, as described in Section 18.4.1 about hash functions, to convert a variable into an integer $x$.

Because you add an item to the dictionary only if its search key is not already present, does the time required for an insertion increase?
10. Repeat Programming Problem 9, but this time
a. Use linear probing as the collision-resolution scheme.
b. Use double hashing as the collision-resolution scheme.
c. Use quadratic probing as the collision-resolution scheme.
11. Repeat Programming Problem 9, but allocate the hash table dynamically. If the hash table becomes more than half full, increase its size to the first prime number greater than $2 \times$ tableSize.
12. Repeat Programming Problem 9, but experiment with variations of chaining. For example, the hash table could point to binary search trees instead of chains of linked entries.

## Balanced Search Trees

## Chapter

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lthough Chapter 18 described the advantages of using the binary search tree to implement the ADT dictionary, the efficiency of this implementation suffers when the tree loses its balance. This chapter introduces various other search trees, which remain balanced in all situations and thus enable dictionary operations whose efficiency is comparable to a binary search.

The height of a binary search tree is sensitive to the order of insertions and removals

Various search trees can retain their balance despite insertions and removals

### 19.1 Balanced Search Trees

As you saw in the previous chapter, the efficiency of the binary search tree implementation of the ADT dictionary is related to the tree's height. The operations add, remove, and getItem follow a path from the root of the tree to the node that contains the desired item (or, in the case of the add operation, to the node that is to become the parent of the new item). At each node along the path, you compare a given value to the search key in the node and determine which subtree to search next. Because the maximum number of nodes on such a path is equal to the height of the tree, the maximum number of comparisons that the dictionary operations can require is also equal to this height.

As you know, the height of a binary search tree of $n$ nodes ranges from a maximum of $n$ to a minimum of $\left\lceil\log _{2}(n+1)\right\rceil$ As a consequence, locating a particular item in a binary search tree requires between $n$ and $\left\lceil\log _{2}(n+1)\right\rceil$ comparisons. Thus, a search of a binary search tree can be as inefficient as a sequential search of a chain of linked nodes or as efficient as a binary search of a sorted array. Efficiency was the primary reason for developing the binary search tree implementation of the dictionary: We wanted to perform a search of a linked structure as efficiently as we could perform a binary search of an array. Thus, we certainly want the optimum behavior of the binary search tree.

What affects the height of a binary search tree? As you learned in Chapters 15 and 16, the height of the tree is quite sensitive to the order in which you insert or remove items. For example, consider a binary search tree that contains the items $10,20,30,40,50,60$, and 70 . If you inserted the items into the tree in ascending order, you would obtain a binary search tree of maximum height, as shown in Figure 19-1a. If, on the other hand, you inserted the items in the order $40,20,60,10,30,50,70$, you would obtain a balanced binary search tree of minimum height, as shown in Figure 19-1b.

As you can see, if you use the algorithms in Chapter 16 to maintain a binary search tree, insertions and removals can cause the tree to lose its balance and approach a linear shape. Such a tree is no better than a linear chain of linked nodes. For this reason, it is desirable in many applications to use one of

FIGURE 19-1 $\quad$ (a) A binary search tree of maximum height; (b) a binary search tree of minimum height

several variations of the basic binary search tree. Such trees can absorb insertions and removals without a deterioration of their balance and are easier to maintain than a minimum-height binary search tree. In addition, you can search these trees almost as efficiently as you can search a minimum-height binary search tree. This chapter discusses the better-known search trees to give you a sense of the possibilities. We continue to assume that the entries in a tree are unique-that is, that there are no duplicates.

### 19.2 2-3 Trees

A 2-3 tree is a tree in which each internal node (nonleaf) has either two or three children and all leaves are at the same level. For example, Figure 19-2 shows a 2-3 tree of height 3. A node with two children is called a 2 -node - the nodes in a binary tree are all 2 -nodes-and a node with three children is called a 3-node.

## FIGURE 19-2 A 2-3 tree of height 3



A 2-3 tree is not a binary tree, because a node can have three children; nevertheless, a 2-3 tree does resemble a full binary tree. If a particular 2-3 tree contains only 2 -nodes-a possibility, according to the definition-it is like a full binary tree, because all of its internal nodes have two children and all of its leaves are at the same level. If, on the other hand, some of the internal nodes of a 2-3 tree do have three children, the tree will contain more nodes than a full binary tree of the same height. Therefore, a 2-3 tree of height $h$ always has at least as many nodes as a full binary tree of height $h$; that is, it always has at least $2^{h}-1$ nodes. To put this another way, a 2-3 tree with $n$ nodes never has height greater than $\left\lceil\log _{2}(n+1)\right\rceil$, the minimum height of a binary tree with $n$ nodes.

Given these observations, a 2-3 tree might be useful as an implementation of the ADT dictionary. Indeed, this is the case if the 2-3 tree orders its nodes to make it useful as a search tree. Just as we distinguish between a binary tree and a binary search tree, we could distinguish between a " $2-3$ tree" and a "2-3 search tree." The previous description would define a $2-3$ tree, and the recursive definition given next would define a $2-3$ search tree. Most people, however, do not make such a distinction and use the term " $2-3$ tree" to mean " $2-3$ search tree;" we will do so as well. The following recursive definition of a 2-3 tree specifies the order of its nodes, making it a search tree:

A 2-3 tree is not a
binary tree

A 2-3 tree is never taller than a minimum-height binary tree

## Note: 2-3 trees

$T$ is a 2-3 tree of height $h$ if one of the following is true:

- $T$ is empty, in which case $h$ is 0 .
- $T$ is of the form


VideoNote 2-3 trees
where $r$ is a node that contains one data item and $T_{\mathrm{L}}$ and $T_{\mathrm{R}}$ are both 2-3 trees, each of height $h-1$. In this case, the item in $r$ must be greater than each item in the left subtree $T_{\mathrm{L}}$ and smaller than each item in the right subtree $T_{\mathrm{R}}$.

- $T$ is of the form

where $r$ is a node that contains two data items and $T_{\mathrm{L}}, T_{\mathrm{M}}$, and $T_{\mathrm{R}}$ are 2-3 trees, each of height $h-1$. In this case, the smaller item in $r$ must be greater than each item in the left subtree $T_{\mathrm{L}}$ and smaller than each item in the middle subtree $T_{\mathrm{M}}$. The larger item in $r$ must be greater than each item in the middle subtree $T_{\mathrm{M}}$ and smaller than each item in the right subtree $T_{\mathrm{R}}$.


## Note: Rules for placing data items in the nodes of a 2-3 tree

The previous definition of a 2-3 tree implies the following rules for how you may place data items in its nodes:

- A 2-node, which has two children, must contain a single data item that is greater than the left child's item(s) and less than the right child's item(s), as Figure 19-3a illustrates.
- A 3-node, which has three children, must contain two data items, $S$ and $L$, that satisfy the following relationships, as Figure 19-3b illustrates: $S$ is greater than the left child's item(s) and less than the middle child's item(s); $L$ is greater than the middle child's item(s) and less than the right child's item(s).
- A leaf may contain either one or two data items.

FIGURE 19-3 Nodes in a 2-3 tree: (a) a 2-node; (b) a 3-node


Items in a 2-3 tree are ordered

Thus, the items in a 2-3 tree are ordered. For example, the tree in Figure 19-4 is a 2-3 tree.

FIGURE 19-4 A 2-3 tree


Listing 19-1 declares a class of nodes for a 2-3 tree. When a node contains only one data item, you can place it in smal1Item and use leftChildPtr and midChildPtr to point to the node's children. To be safe, you can place nullptr in rightChildPtr.

## LISTING 19-1 A header file for a class of nodes for a 2-3 tree

```
/** A class of nodes for a link-based 2-3 tree.
    @file TriNode.h */
#ifndef _TRI_NODE
#define _TRI_NODE
temp1ate<c1ass ItemType>
class TriNode
{
private:
    ItemType sma11Item, largeItem; // Data portion
    TriNode<ItemType>* leftChildPtr; // Left-child pointer
    TriNode<ItemType>* midChildPtr; // Middle-child pointer
    TriNode<ItemType>* rightChildPtr; // Right-child pointer
public:
    TriNode();
    TriNode(const ItemType& anItem);
    TriNode(const ItemType& anItem, TriNode<ItemType>* 1eftPtr,
                TriNode<ItemType>* midPtr, TriNode<ItemType>* rightPtr);
    bool isLeaf() const;
    bool isTwoNode() const;
    bool isThreeNode() const;
    ItemType getSmal1Item() const;
    ItemType getLargeItem() const;
    void setSma11Item(const ItemType& anItem);
    void setLargeItem(const ItemType& anItem);
```

Inorder traversal

Searching a 2-3 tree is efficient

```
        TriNode<ItemType>* getLeftChildPtr() const;
    TriNode<ItemType>* getMidChildPtr() const;
    TriNode<ItemType>* getRightChildPtr() const;
    void setLeftChildPtr(TriNode<ItemType>* 1eftPtr);
    void setMidChildPtr(TriNode<ItemType>* midPtr);
    void setRightChildPtr(TriNode<ItemType>* rightPtr);
}; // end TriNode
#include "TriNode"
#endif
```


### 19.2.1 Traversing a 2-3 Tree

The algorithms for the traversal, retrieval, insertion, and removal operations on a 2-3 tree are recursive. You can avoid distracting implementation details by defining the base case for these recursive algorithms to be a leaf rather than an empty subtree. As a result, the algorithms must assume that they are not passed an empty tree as an argument.

You can traverse a 2-3 tree in sorted order by performing the analogue of an inorder traversal on a binary tree:

```
// Traverses a nonempty 2-3 tree in sorted order.
inorder(23Tree: TwoThreeTree): void
    if (23Tree's root node r is a leaf)
        Visit the data item(s)
    else if (r has two data items)
    {
        inorder(left subtree of 23Tree's root)
        Visit the first data item
        inorder (middle subtree of 23Tree's root)
        Visit the second data item
        inorder(right subtree of 23Tree's root)
}
    else // r has one data item
    {
        inorder(left subtree of 23Tree's root)
        Visit the data item
        inorder(right subtree of 23Tree's root)
    }
```


### 19.2.2 Searching a 2-3 Tree

The ordering of items in a $2-3$ tree is analogous to the ordering for a binary search tree and allows you to search a 2-3 tree efficiently for a particular item. In fact, the retrieval operation for a 2-3 tree is quite similar to the retrieval operation for a binary search tree, as you can see from the following pseudocode:

```
// Locates the value target in a nonempty 2-3 tree. Returns either the located
// entry or throws an exception if such a node is not found.
findItem(23Tree: TwoThreeTree, target: ItemType): ItemType
    if (target is in 23Tree's rootnode r)
    \{ // The item has been found
        treeItem = the data portion of r
        return treeItem // Success
    \}
```

```
else if (r is a leaf)
    throw NotFoundException // Failure
// Else search the appropriate subtree
else if (r has two data items)
{
    if (target < smaller item in r)
        return findItem(r's left subtree, target)
    else if (target < larger item in r)
        return findItem(r's middle subtree, target)
    else
        return findItem(r's right subtree, target)
}
e1se // r has one data item
{
    if (target < r's data item)
        return findItem(r's left subtree, target)
    else
        return findItem(r's right subtree, target)
}
```

You can search the 2-3 tree and the shortest binary search tree with approximately the same efficiency, because

- A binary search tree with $n$ nodes cannot be shorter than $\left\lceil\log _{2}(n+1)\right\rceil$
- A 2-3 tree with $n$ nodes cannot be taller than $\left\lceil\log _{2}(n+1)\right\rceil$
- A node in a 2-3 tree has at most two items

Searching a 2-3 tree is not more efficient than searching a binary search tree, however. This observation may surprise you because, after all, the nodes of a 2-3 tree can have three children, and hence a 2-3 tree might indeed be shorter than the shortest possible binary search tree. Although true, this advantage in height is offset by the extra time required to compare a given value with two values instead of only one. In other words, although you might visit fewer nodes when searching a 2-3 tree, you might have to make more comparisons at each node. As a consequence, the number of comparisons that you need to search a 2-3 tree for a given item is approximately equal to the number of comparisons required to search a binary search tree that is as balanced as possible. This number is approximately $\log _{2} n$.

If you can search a 2-3 tree and a balanced binary search tree with approximately the same efficiency, why then should you use a 2-3 tree? Because although maintaining the balance of a binary search tree is difficult in the face of insertion and removal operations, maintaining the shape of a 2-3 tree is relatively simple. For example, consider the two trees in Figure 19-5. The first tree is a binary

FIGURE 19-5 (a) A balanced binary search tree; (b) a 2-3 tree with the same entries


Searching a 2-3 tree is $\mathrm{O}(\log n)$

FIGURE 19-6 (a) The binary search tree of Figure 19-5a after inserting the sequence of values 32 through 39 ; (b) the 2-3 tree of Figure 19-5b after the same insertions


Maintaining the shape of a 2-3 tree is relatively easy
search tree and the second is a 2-3 tree. Both trees contain the same data items. The binary search tree is as balanced as possible, and thus you can search both it and the 2-3 tree for an item with approximately the same efficiency. If, however, you perform a sequence of insertions on the binary search tree-by using the insertion algorithm of Chapter 16-the tree can quickly lose its balance, as Figure 19-6a indicates. In this example, we have inserted the values 32 through 39 in numerical order. As you soon will see, you can perform the same sequence of insertions on the 2-3 tree without a degradation in the tree's shape-it will retain its structure, as Figure 19-6b shows.

The new values ( 32 through 39) that were inserted into the binary search tree of Figure 19-5a appear along a single path in Figure 19-6a. The insertions increased the height of the binary search tree from 4 to 12 -an increase of 8 . On the other hand, the new values have been spread throughout the 2-3 tree in Figure 19-6b. As a consequence, the height of the resulting tree is only 1 greater than the height of the original 2-3 tree in Figure 19-5b. We demonstrate these insertions into the original 2-3 tree next.

### 19.2.3 Inserting Data into a 2-3 Tree

Because the nodes of a 2-3 tree can have either two or three children and can contain one or two values, you can insert items into the tree while maintaining its shape. The following paragraphs informally describe the sequence of insertions that produced the 2-3 tree shown in Figure 19-6b from the original tree in Figure 19-5b.

Insert 39. As is true with a binary search tree, the first step in inserting a node into a 2-3 tree is to locate the node at which the search for the value to be inserted would terminate. To do this, you can use the search strategy of the findItem algorithm given previously; an unsuccessful search will always terminate at a leaf. With the tree in Figure 19-5b, the search for 39 terminates at the leaf $<40>$. (We will use this angle-bracket notation to denote a node and its contents.) Because this node contains only one item, you can simply insert the new item into this node. The result is the 2-3 tree in Figure 19-7.

FIGURE 19-7 After inserting 39 into the tree in Figure 19-5b


Insert 38. In a similar manner, you would search the tree in Figure 19-7 for 38 and find that the search terminates at the node $<3940>$. You cannot place 38 in this node, as Figure 19-8a illustrates, because a node cannot contain three values. You arrange these three values, however, into the smallest (38), middle (39), and largest (40) values. You can move the middle value (39) up to the node's parent $p$ and separate the remaining values, 38 and 40 , into two nodes that you attach to $p$ as children, as Figure 19-8b indicates. Because you chose to move up the middle value of $<383940>$, the parent correctly separates the values of its children; that is, 38 is less than 39 , which is less than 40 . The result of the insertion is the 2-3 tree in Figure 19-8c.

Insertion into a 2 -node leaf is simple

Insertion into a 3-node causes it to divide

FIGURE 19-8 The steps for inserting 38 into the tree in Figure 19-7: (a) The located node has no room; (b) the node splits; (c) the resulting tree


Insert 37. The insertion of 37 into the tree in Figure 19-8c is easy because 37 belongs in a leaf that currently contains only one value, 38 . The result of this insertion is the 2-3 tree in Figure 19-9.

FIGURE 19-9 After inserting 37 into the tree in Figure 19-8c


Insert 36. The search for 36 terminates at the node $<3738>$ of the tree in Figure 19-9, but the node has no room for 36 , as Figure 19-10a indicates. Thus, you arrange the three values 36,37 , and 38 into the smallest (36), middle (37), and largest (38) values, as you did previously. You then try to move the middle value (37) up to the node's parent $<3039>$ and give the parent children containing the smallest (36) and largest (38) values, as Figure 19-10b illustrates. However, the node $<3039>$ cannot contain three values and have four children. This situation is familiar, with the slight difference that the overcrowded node is not a leaf but rather has four children. As you did before, you divide the node into the smallest (30), middle (37), and largest (39) values and then move the middle value up to the node's parent.

Because you are splitting an internal node, you now must account for its four children-that is, what happens to nodes $<1020>,<36>,<38>$, and $<40>$ ? The solution is to attach the left pair of children-<10 $20>$ and $<36>$-to the smallest value (30) and attach the right pair of children-<38> and $<40>$ - to the largest value (39), as shown in Figure 19-10c. The final result of this insertion is the 2-3 tree in Figure 19-10d.

FIGURE 19-10 (a), (b), (c) The steps for inserting 36 into the tree in Figure 19-9; (d) the resulting tree


Insert 35, 34, and 33. Each of these insertions is similar to the previous ones. Figure 19-11 shows the tree after the three insertions.

FIGURE 19-11 The tree after the insertion of 35, 34, and 33 into the tree in Figure 19-10d


Question 1 To be sure that you fully understand the insertion algorithm, insert 32 into the 2-3 tree in Figure 19-11. The result should be the tree shown in Figure 19-6b. Once again, compare this tree with the binary search tree in Figure 19-6a and notice the dramatic advantage of the 2-3 tree's insertion strategy.

The insertion algorithm. Let's consider the general strategy for inserting an item into a 2-3 tree. To begin, you locate the leaf at which the search for the new item would terminate. If the leaf does not contain two items, you insert the new item into the leaf, and you are done. However, if the leaf already contains two items, you must split it into two nodes, $n_{1}$ and $n_{2}$. As Figure 19-12 illustrates, you place the smallest item $S$ into $n_{1}$, place the largest item $L$ into $n_{2}$, and try to move the middle item $M$ up to the original leaf's parent. If the parent has room for a new item-as is true here - nodes $n_{1}$ and $n_{2}$ become children of the parent, and you are finished. The parent now has only three children and contains two items.

FIGURE 19-12 Splitting a leaf in a 2-3 tree when the leaf is a (a) left child; (b) right child
(a)

(b)



When an internal node would contain three items, split it into two nodes and accommodate its children

When the root would contain three items, split it into two nodes and create a new root node

On the other hand, if the parent cannot accommodate the item moving up, you must split the parent, as follows. You split an internal node $n$ that contains three items by using the process just described for a leaf, except that you must also take care of $n$ 's four children. As Figure 19-13 illustrates, you split $n$ into $n_{1}$ and $n_{2}$, place $n$ 's smallest item $S$ into $n_{1}$, attach $n$ 's two leftmost children to $n_{1}$, place $n$ 's largest item $L$ into $n_{2}$, attach $n$ 's two rightmost children to $n_{2}$, and move $n$ 's middle item $M$ up to $n$ 's parent.

FIGURE 19-13 Splitting an internal node in a 2-3 tree when the node is a (a) left child; (b) right child


After this, the process of splitting a node and moving an item up to the parent continues recursively until a node is reached that had only one item before the insertion and thus has only two items after it takes on a new item. Notice in the previous sequence of insertions that the tree's height never increased from its original value of 3. In general, an insertion will not increase the height of the tree as long as there is at least one node containing only one item on the path from the root to the leaf into which the new item is inserted. The insertion strategy of a $2-3$ tree has thus postponed the growth of the tree's height much more effectively than the strategy of a basic binary search tree did.

When the height of a 2-3 tree does grow, it does so from the top. An increase in the height of a 2-3 tree will occur if every node on the path from the root of the tree to the leaf into which the new item is inserted contains two items. In this case, the recursive process of splitting a node and moving an item up to the node's parent will eventually reach the root $r$. When this occurs you must split $r$ into $r_{1}$ and $r_{2}$ exactly as you would any other internal node. However, you must create a new node that contains the middle item of $r$ and becomes the parent of $r_{1}$ and $r_{2}$. Thus, the new node is the new root of the tree, as Figure 19-14 illustrates.

FIGURE 19-14 Splitting the root of a 2-3 tree


The following high-level algorithm summarizes the entire insertion strategy:

```
// Inserts a new item into a 2-3 tree whose items are distinct and differ from the
// new item.
insertItem(23Tree: TwoThreeTree, newItem: ItemType)
    Locate the leaf, 1eafNode, in which newItem belongs
    Add newItem to leafNode
    if (7eafNode has three items)
        split(1eafNode)
// Splits node n, which contains two items. Note: If n is
// not a leaf, it has four children.
split(n: TwoThreeNode)
    if ( n is the root)
        Create a new node p
    e1se
        Let p be the parent of n
    Replace node n with two nodes, n 1 and n 2 , so that \(p\) is their parent
    Give n 1 the item in n with the smallest value
    Give n 2 the item in n with the largest value
    if ( n is not a leaf)
    \{
        n 1 becomes the parent of n's two leftmost children
        n 2 becomes the parent of n 's two rightmost children
    \}
    Move the item in n that has the middle value up to p
    if (p now has three items)
        split(p)
```

Question 2 What is the result of inserting 5, 40, 10, 20, 15, and 30-in the order giveninto an initially empty 2-3 tree? Note that insertion of one item into an empty 2-3 tree will create a single node that contains the inserted item.

Question 3 What is the result of inserting 3 and 4 into the 2-3 tree that you created in the previous question?

### 19.2.4 Removing Data from a 2-3 Tree

The removal strategy for a 2-3 tree is the inverse of its insertion strategy. Just as a 2-3 tree spreads insertions throughout the tree by splitting nodes when they would become too full, it spreads removals throughout the tree by merging nodes when they become empty. As an illustration of the 2-3 tree's removal strategy, consider the removal of 70, 100, and 80 from the tree in Figure 19-15a.

Remove 70. By searching the tree in Figure 19-15a, you discover that 70 is in the node $<7090>$. Because you always want to begin the removal process at a leaf, the first step is to swap 70 with its inorder successor-the value that follows it in the sorted order. Because 70 is the smaller of the two values in the node, its inorder successor (80) is the smallest value in the node's middle subtree. Note that the inorder successor of an item in an internal node will always be in a leaf. After the swap, the tree appears as shown in Figure 19-15b. The value 80 is in a legal position of the search tree because it is larger than all the values in its node's left subtree and smaller than all the values in its node's right subtree. The value 70 is not in a legal position, but this is of no concern, because the next step is to remove this value from the leaf.

2-3 tree insertion algorithm

FIGURE 19-15 (a) A 2-3 tree; (b), (c), (d), (e) the steps for removing 70; (f) the resulting tree


In general, after you remove a value from a leaf, another value may remain in the leaf (because the leaf contained two values before the removal). If this is the case, you are done, because a leaf of a 2-3 tree can contain a single value. In this example, however, once you remove 70 from the leaf, the node is left without a value, as Figure 19-15c indicates.
Merge nodes You then delete the node, as Figure 19-15d illustrates. At this point you see that the parent of the deleted node contains two values ( 80 and 90 ) but has two children ( 60 and 100). This situation is not allowed in a 2-3 tree. You can remedy the problem by moving the smaller value (80) down from the parent into the left child, as Figure 19-15e illustrates. Deleting the leaf node and moving a value down to a sibling of the leaf is called merging the leaf with its sibling. The 2-3 tree that results from this removal operation is shown in Figure 19-15f.

Redistribute values
Remove 100. The search strategy discovers that 100 is in the leaf $<100\rangle$ of the tree in Figure 19-15f. When you remove the value from this leaf, the node becomes empty, as Figure 19-16a indicates. In this case, however, no merging of nodes is required, because the sibling $<6080>$ can spare a value. That is, the sibling has two values, whereas a 2-3 tree requires only that it have at least one value. However, if you simply move the value 80 into the empty node-as Figure 19-16b illustrates-you find that the search-tree order is destroyed: The value in 90 's right child should be greater than 90 , but it is 80 . Instead, you move the larger value ( 80 ) from $<6080>$ into the parent and move the value

FIGURE 19-16 (a), (b), (c) The steps for removing 100 from the tree in Figure 19-15f;
(d) the resulting tree

(d)


90 down from the parent into the node that had been empty, as Figure 19-16c shows. This distribution preserves the search-tree order, and you have thus completed the removal. The resulting 2-3 tree is shown in Figure 19-16d.
Remove 80. The search strategy finds that 80 is in an internal node of the tree in Figure 19-16d. You thus must swap 80 with its inorder successor, 90, as Figure 19-17a illustrates. When you remove 80 from the leaf, the node becomes empty, as Figure 19-17b shows. Because the sibling of the empty node has only one value, you cannot redistribute as you did in the previous removal of 100. Instead you must merge the nodes, bringing the value 90 down from the parent and deleting the empty leaf, as Figure 19-17c indicates.

You are not yet finished, however, because the parent contains no data and has only one child. You must recursively apply the removal strategy to this internal node without a value. First, you should check to see whether the node's sibling can spare a value. Because the sibling $\langle 30\rangle$ contains only the single value 30 , you cannot redistribute-you must merge the nodes. The merging of two internal nodes is identical to the merging of leaves, except that the child $<6090>$ of the empty node must be adopted. Because the sibling of the empty node contains only one value - and hence can have only two children, as stated in the rule for 2-nodes-it can become the parent of $<6090>$ only if you bring the value 50 down from the sibling's parent. The tree now appears as shown in Figure 19-17d. Note that this operation preserves the search property of the tree.

Now the parent of the merged nodes is left without a data item and only a single child. Usually, you would apply the recursive removal strategy to this node, but this case is special because the node is the root. Because the root is empty and has only one child, you can simply delete it, allowing $<30$ $50>$ to become the root of the tree, as Figure 19-17e illustrates. This deletion has thus caused the height of the tree to shrink by 1 .

To summarize, we have removed 70, 100, and 80 from the 2-3 tree in Figure 19-15a and obtained the 2-3 tree in Figures 19-17e and 19-18a. In contrast, after removing 70, 100, and 80 from the balanced binary search tree in Figure 19-5a, you are left with the tree in Figure 19-18b. Notice that the

FIGURE 19-17 The steps for removing 80 from the tree in Figure 19-16d


FIGURE 19-18 Results of removing 70, 100, and 80 from (a) the 2-3 tree of Figure 19-15a and (b) the binary search tree of Figure 19-5a

removals affected only one part of the binary search tree, causing it to lose its balance. The left subtree has not been affected at all, and thus the overall height of the tree has not been diminished.

The removal algorithm. In summary, to remove an item $I$ from a 2-3 tree, you first locate the node $n$ that contains it. If $n$ is not a leaf, you find $l$ 's inorder successor and swap it with $I$. As a result of the swap, the removal always begins at a leaf. If the leaf contains an item in addition to $I$, you simply remove $I$ and you are done. On the other hand, if the leaf contains only $I$, removing $I$ would leave the leaf without a data item. In this case you must perform some additional work to complete the removal.

You first check the siblings of the now-empty leaf. If a sibling has two items, you redistribute the items among the sibling, the empty leaf, and the leaf's parent, as Figure 19-19a illustrates. If no sibling of the leaf has two items, you merge the leaf with an adjacent sibling by moving an item down from the leaf's parent into the sibling -it had only one item before, so it has room for another-and deleting the empty leaf. This case is shown in Figure 19-19b.

By moving an item down from a node $n$, as just described, you might cause $n$ to be left without a data item and with only one child. If so, you recursively apply the removal algorithm to $n$. Thus, if $n$ has a sibling with two items-and three children-you redistribute the items among $n$, the sibling, and $n$ 's parent. You also give $n$ one of its sibling's children, as Figure 19-19c indicates.

If $n$ has no sibling with two items, you merge $n$ with a sibling, as Figure 19-19d illustrates. That is, you move an item down from the parent and let the sibling adopt $n$ 's one child. (At this point you know that the sibling previously had only one item and two children.) You then delete the empty leaf. If the merge causes $n$ 's parent to be without an item, you recursively apply the removal process to it.

If the merging continues so that the root of the tree is without an item-and has only one child-you simply delete the root. When this step occurs, the height of the tree is reduced by 1 , as Figure 19-19e illustrates.

## A high-level statement of the algorithm for removing data from a 2-3 tree follows:

```
// Removes the given data item from a 2-3 tree. Returns true if successful
// or false if no such item exists.
removeItem(23Tree: TwoThreeTree, dataItem: ItemType): boolean
    Attempt to locate dataItem
    if (dataItem is found)
    {
        if (dataItem is not in a leaf)
            Swap dataItem with its inorder successor, which will be in a leaf 1eafNode
        // The removal always begins at a leaf
        Remove dataItem from leaf 1eafNode
        if (leafNode now has no items)
            fixTree(7eafNode)
        return true
    }
    e1se
        return false
// Completes the removal when node n is empty by either deleting the root,
// redistributing values, or merging nodes. Note: If n is internal, it has one child.
fixTree(n: TwoThreeNode)
    if (n is the root)
    Delete the root
    else
    {
        Let p be the parent of n
        if (some sibling of n has two items)
        {
Distribute items appropriately among n , the sibling, and p
```

FIGURE 19-19 (a) Redistributing values; (b) merging a leaf; (c) redistributing values and children; (d) merging internal nodes; (e) deleting the root
(a)

(b)


Sibling
Leaf

(c)

(d)

$\xrightarrow{\text { Merge }}$

(e)


```
    if (n is internal)
        Move the appropriate child from sibling to n
    }
    e1se // Merge the node
    {
    Choose an adjacent sibling s of n
    Bring the appropriate item down from p into s
    if (n is internal)
        Move n's child to s
    Remove node n
    if (p is now empty)
        fixTree(p)
    }
}
```

The details of the $\mathrm{C}++$ implementation of the preceding insertion and removal algorithms for 2-3 trees are rather involved. The implementation is left as a challenging exercise (Programming Problem 2).

You might be concerned about the overhead that the insertion and removal algorithms incur in the course of maintaining the 2-3 structure of the tree. That is, after the search strategy locates either the item or the position for the new item, the insertion and removal algorithms sometimes have to perform extra work, such as splitting and merging nodes. However, this extra work is not a real concern. A rigorous mathematical analysis would show that the extra work required to maintain the structure of a 2-3 tree after an insertion or a removal is not significant. In other words, when analyzing the efficiency of the insertItem and removeItem algorithms, it is sufficient to consider only the time required to locate the item (or the position for the insertion). Given that a 2-3 tree is always balanced, you can search a 2-3 tree in all situations with the logarithmic efficiency of a binary search.

Thus, a 2-3 tree implementation of the ADT dictionary is guaranteed to have efficient operations. Although a binary search tree that is as balanced as possible minimizes the amount of work required by the ADT dictionary operations, its balance is difficult to maintain. A 2-3 tree is a compromise: Although searching it may not be quite as efficient as searching a binary search tree of minimum height, it is relatively simple to maintain.


Question 4 What is the result of removing the 10 from the 2-3 tree that you created in Checkpoint Question 1?

### 19.3 2-3-4 Trees

If a 2-3 tree is so good, are trees whose nodes can have more than three children even better? To some extent, the answer is yes. A 2-3-4 tree is like a 2-3 tree, but it also allows 4-nodes, which are nodes that have four children and three data items. For example, Figure 19-20 shows a 2-3-4 tree of height 3

A 2-3 tree is always balanced

A 2-3 tree implementation of a dictionary is $\mathrm{O}(\log n)$ for all of its operations

VideoNote
2-3-4 and red-black trees

FIGURE 19-20 A 2-3-4 tree with the same data items as the 2-3 tree in Figure 19-6b

that has the same items as the 2-3 tree in Figure 19-6b. As you will see, you can perform insertions and removals on a 2-3-4 tree with fewer steps than a $2-3$ tree requires.

## Note: 2-3-4 trees

$T$ is a 2-3-4 tree of height $h$ if one of the following is true:

- $T$ is empty, in which case $h$ is 0 .
- $T$ is of the form

where $r$ is a node that contains one data item and $T_{\mathrm{L}}$ and $T_{\mathrm{R}}$ are both 2-3-4 trees, each of height $h-1$. In this case, the item in $r$ must be greater than each item in the left subtree $T_{\mathrm{L}}$ and smaller than each item in the right subtree $T_{\mathrm{R}}$.
- $T$ is of the form

where $r$ is a node that contains two data items and $T_{\mathrm{L}}, T_{\mathrm{M}}$, and $T_{\mathrm{R}}$ are 2-3-4 trees, each of height $h-1$. In this case, the smaller item in $r$ must be greater than each item in the left subtree $T_{\mathrm{L}}$ and smaller than each item in the middle subtree $T_{\mathrm{M}}$. The larger item in $r$ must be greater than each item in $T_{\mathrm{M}}$ and smaller than each item in the right subtree $T_{\mathrm{R}}$.
- $T$ is of the form

where $r$ is a node that contains three data items and $T_{\mathrm{L}}, T_{\mathrm{ML}}, T_{\mathrm{MR}}$, and $T_{\mathrm{R}}$ are 2-3-4 trees, each of height $h-1$. In this case, the smallest item in $r$ must be greater than each item in the left subtree $T_{\mathrm{L}}$ and smaller than each item in the middle-left subtree $T_{\mathrm{ML}}$. The middle item in $r$ must be greater than each item in $T_{\mathrm{ML}}$ and smaller than each item in the mid-dle-right subtree $T_{\mathrm{MR}}$. The largest item in $r$ must be greater than each item in $T_{\mathrm{MR}}$ and smaller than each item in the right subtree $T_{\mathrm{R}}$.


## Note: Rules for placing data items in the nodes of a 2-3-4 tree

The previous definition of a 2-3-4 tree implies the following rules for how you may place data items in its nodes:

- A 2-node, which has two children, must contain a single data item that satisfies the relationships pictured earlier in Figure 19-3a.
- A 3-node, which has three children, must contain two data items that satisfy the relationships pictured earlier in Figure 19-3b.
- A 4-node, which has four children, must contain three data items $S, M$, and $L$ that satisfy the following relationships, as Figure 19-21 illustrates: $S$ is greater than the left child's item(s) and less than the middle-left child's item(s); $M$ is greater than the middle-left child's item(s) and less than the middle-right child's item(s); $L$ is greater than the middle-right child's item(s) and less than the right child's item(s).
- A leaf may contain either one, two, or three data items.

FIGURE 19-21 A 4-node in a 2-3-4 tree


Although a 2-3-4 tree has more efficient insertion and removal operations than a 2-3 tree, a 2-3-4 tree has greater storage requirements due to the additional data members in its 4-nodes, as the beginning of the class QuadNode indicates:

```
template<class ItemType>
class QuadNode
{
private:
    ItemType smal1Item, middleItem, largeItem; // Data portion
    QuadNode<ItemType>* 1eftChildPtr; // Left-child pointer
    QuadNode<ItemType>* leftMidChildPtr; // Middle-7eft-child pointer
    QuadNode<ItemType>* rightMidChildPtr; // Middle-right-child pointer
    QuadNode<ItemType>* rightChildPtr; // Right-child pointer
    // Constructors, accessor methods, and mutator methods are here.
}; // end QuadNode
```

As you will see later, however, you can transform a 2-3-4 tree into a special binary tree that reduces the storage requirements.

### 19.3.1 Searching and Traversing a 2-3-4 Tree

The search algorithm and the traversal algorithm for a 2-3-4 tree are simple extensions of the corresponding algorithms for a 2-3 tree. For example, to search the tree in Figure 19-20 for 31, you would search the left subtree of the root, because 31 is less than 37 ; search the middle subtree of the node $<3035>$, because 31 is between 30 and 35 ; and terminate the search at the left child pointer of $<3233$ $34>$, because 31 is less than 32 -deducing that 31 is not in the tree. Exercise 8 asks you to complete the details of searching and traversing a 2-3-4 tree.

### 19.3.2 Inserting Data into a 2-3-4 Tree

The insertion algorithm for a 2-3-4 tree, like the insertion algorithm for a 2-3 tree, splits a node by moving one of its items up to its parent node. For a 2-3 tree, the search algorithm traces a path from the root to a leaf and then backs up from the leaf as it splits nodes. To avoid this return path after

A 2-3-4 tree requires more storage than a 2-3 tree

A node in a 2-3-4 tree

Split 4-nodes as they are encountered
reaching a leaf, the insertion algorithm for a 2-3-4 tree splits 4-nodes as soon as it encounters them on the way down the tree from the root to a leaf. As a result, when a 4 -node is split and an item is moved up to the node's parent, the parent cannot possibly be a 4-node, and so it can accommodate another item.

As an example of the algorithm, consider the tree in Figure 19-22a. This one-node tree is the result of inserting 60,30 , and 10 into an initially empty 2-3-4 tree. We will now insert more data into this tree.

Insert 20. While determining the insertion point, you begin at the root and encounter the 4 -node $<10$ $3060>$, which you split by moving the middle value 30 up. Because the node is the root, you create a new root, move 30 into it, and attach two children, as Figure 19-22b illustrates. You continue the search for 20 by examining the left subtree of the root, because 20 is less than 30 . The insertion results in the tree in Figure 19-22c.

FIGURE 19-22 Inserting 20 into a one-node 2-3-4 tree (a) the original tree; (b) after splitting the node; (c) after inserting 20
(a)


Insert 50 and 40. The insertions of 50 and 40 do not require split nodes and result in the tree in Figure 19-23.

FIGURE 19-23 After inserting 50 and 40 into the tree in Figure 19-22c


Insert 70. While searching Figure 19-23 for 70's insertion point, you encounter the 4-node $<4050$ $60>$, because 70 is greater than 30 . You split this 4 -node by moving 50 up to the node's parent, $<30>$, to get the tree in Figure 19-24a. You then insert 70 into the leaf $<60>$, as Figure 19-24b illustrates.

FIGURE 19-24 The steps for inserting 70 into the tree in Figure 19-23: (a) after splitting the 4-node; (b) after inserting 70
(a)

(b)


Insert 80 and 15. These insertions do not require split nodes and result in the tree in Figure 19-25.

FIGURE 19-25 After inserting 80 and 15 into the tree in Figure 19-24b


Insert 90. As you search Figure 19-25 for 90's insertion point, you traverse the root's right subtree, because 90 is greater than 50, and encounter the 4-node $<607080>$. You split this 4-node into two nodes and move 70 up to the root, as Figure 19-26a indicates. Finally, because 90 is greater than 70, you insert 90 into the leaf $<80>$ to get the tree in Figure 19-26b.

FIGURE 19-26 The steps for inserting 90 into the tree in Figure 19-25


Insert 100. As you begin to search Figure 19-26b, you immediately encounter a 4-node at the tree's root. You split this node into two nodes and move 50 up to a new root, as Figure 19-27a indicates. After continuing the search, you insert 100 into $<8090>$ to get the tree in Figure 19-27b.

FIGURE 19-27 The steps for inserting 100 into the tree in Figure 19-26b


Splitting 4-nodes during insertion. As you have just seen, you split each 4-node as soon as you encounter it during your search from the root to the leaf that will accommodate the new item to be inserted. As a result, each 4-node either will

- Be the root,
- Have a 2-node parent, or
- Have a 3-node parent

Figure 19-28 illustrates how to split a 4-node that is the tree's root. You have seen two previous examples of this: We split $<103060>$ in Figure 19-22a, resulting in the tree in Figure 19-22b. We also split $<305070>$ during the insertion of 100 into the tree in Figure 19-26b, giving us the tree in Figure 19-27a.

FIGURE 19-28 Splitting a 4-node root during insertion into a 2-3-4 tree


Figure 19-29 illustrates the two possible situations that can occur when you split a 4-node whose parent is a 2-node. For example, when you split <40 $5060>$ during the insertion of 70 into the tree in Figure 19-23, you get the tree in Figure 19-24a.

FIGURE 19-29 Splitting a 4-node whose parent is a 2 -node during insertion into a 2-3-4 tree, when the 4 -node is a (a) left child; (b) right child


Figure 19-30 illustrates the three possible situations that can occur when you split a 4-node whose parent is a 3-node. For example, when you split $<607080>$ during the insertion of 90 into the tree in Figure 19-25, you get the tree in Figure 19-26a.

Question 5 Insert 25 into the 2-3-4 tree in Figure 19-27b.
Question 6 Insert 3 and 4 into the 2-3-4 tree that you created in the previous question.

FIGURE 19-30 Splitting a 4-node whose parent is a 3-node during insertion into a 2-3-4 tree, when the 4 -node is a (a) left child; (b) middle child; (c) right child


### 19.3.3 Removing Data from a 2-3-4 Tree

The removal algorithm for a 2-3-4 tree has the same beginning as the removal algorithm for a 2-3 tree. You first locate the node $n$ that contains the item $I$ that you want to remove. You then find $I$ 's inorder successor and swap it with $I$ so that the removal will always be at a leaf. If that leaf is either a 3 -node or a 4 -node, you simply remove $I$. If you can ensure that $I$ does not occur in a 2 -node, you can perform the removal in one pass through the tree from root to leaf, unlike removal from a 2-3 tree. That is, you will not have to back away from the leaf and restructure the tree.

In fact, you can guarantee that $I$ does not occur in a 2-node by transforming each 2-node that you encounter during the search for $I$ into either a 3-node or a 4 -node. Several cases are possible, depending on the configuration of the 2 -node's parent and its nearest sibling. (Arbitrarily, a node's nearest sibling is its left sibling, unless the node is a left child, in which case its nearest sibling is to its right.) That is, either the parent or the sibling could be a 2 -node, a 3-node, or a 4-node. For example, if the next node that you encounter is a 2-node and both its parent and nearest sibling are 2 -nodes, apply the transformation that Figure 19-28 illustrates, but in reverse; however, if the parent is a 3-node, apply the transformation that Figure 19-29 illustrates, but in reverse; and if the parent is a 4-node, apply the transformation that Figure 19-30 illustrates, but in reverse.

The details of removal from a 2-3-4 tree are left to you as a challenging exercise (Exercise 8).

[^15]2-3 and 2-3-4 trees are attractive because their balance is easy to maintain

Insertion and removal algorithms for a 2-3-4 tree require fewer steps than those for a 2-3 tree

Allowing nodes with more than four children is counterproductive

Note: 2-3 trees versus 2-3-4 trees
The advantage of both 2-3 and 2-3-4 trees is their easy-to-maintain balance, not their shorter height. Even if a 2-3 tree is shorter than a balanced binary search tree, the reduction in height is offset by the increased number of comparisons that the search algorithm might require at each node. The situation is similar for a 2-3-4 tree, but its insertion and removal algorithms require only one pass through the tree and so are more efficient than those for a 2-3 tree. This decrease in effort makes the 2-3-4 tree more attractive than the 2-3 tree.

## Note: Nodes with more than four children?

Should we consider trees whose nodes have even more than four children? Although a tree whose nodes can each have 100 children would be shorter than a 2-3-4 tree, its search algorithm would require more comparisons at each node to determine which subtree to search. Thus, allowing the nodes of a tree to have many children is counterproductive. Such a search tree is appropriate, however, when it is implemented in external storage, because moving from node to node is far more expensive than comparing the data values in a node. In such cases, a search tree with the minimum possible height is desirable, even at the expense of additional comparisons at each node. Chapter 21 will discuss external search trees further.

### 19.4 Red-Black Trees

A 2-3-4 tree requires more storage than a binary search tree

A red-black tree has the advantages of a 2-3-4 tree but requires less storage

A 2-3-4 tree is appealing because it is balanced and because its insertion and removal operations use only one pass from root to leaf. On the other hand, a 2-3-4 tree requires more storage than a binary search tree that contains the same data because a 2-3-4 tree has nodes that must accommodate up to three data items. A typical binary search tree, however, might not be balanced.

You can use a special binary search tree-a red-black tree-to represent a 2-3-4 tree that retains the advantages of a 2-3-4 tree without the storage overhead. The idea is to represent each 3-node and 4-node in a 2-3-4 tree as an equivalent binary search tree. To distinguish between 2-nodes that appeared in the original 2-3-4 tree and 2-nodes that were generated from 3-nodes and 4-nodes, you use red and black child pointers. Let all the child pointers in the original 2-3-4 tree be black, and use red child pointers to link the 2 -nodes that result when you split 3-nodes and 4-nodes.


Note: In a red-black tree, red pointers link the 2-nodes that now contain the values that were in a 3-node or a 4-node.

Figure 19-31 indicates how to represent a 4-node and a 3-node as binary trees. Because there are two possible ways to represent a 3-node as a binary tree, a red-black representation of a 2-3-4 tree is not unique. Figure 19-32 gives a red-black representation for the 2-3-4 tree in Figure 19-20. In all of these figures, a dashed blue line represents a red pointer and a solid black line represents a black pointer.

FIGURE 19-31 Red-black representation of (a) a 4-node; (b) a 3-node
(a)

(b)


FIGURE 19-32 A red-black tree that represents the 2-3-4 tree in Figure 19-20


A node in a red-black tree is similar to a node in a binary search tree, but it must also store the pointer colors, as the following C++ statements indicate:

```
enum Color {RED, BLACK};
template<class ItemType>
class RedBlackNode : public BinaryNode<ItemType>
{
private:
    Color leftColor, rightColor;
public:
    // Get and set methods for 1eftColor and rightColor
    //
}; // end RedBTackNode
```

Note: Even with the pointer colors, a node in a red-black tree requires less storage than a node in a 2-3-4 tree. Keep in mind that the transformations in Figure 19-31 imply a change in the structure of the nodes.

A node in a red-black tree

Question 7 Why does a node in a red-black tree require less memory than a node in a 2-3-4 tree?

### 19.4.1 Searching and Traversing a Red-Black Tree

Because a red-black tree is a binary search tree, you can search and traverse it by using the algorithms for a binary search tree. You simply ignore the color of the pointers.

### 19.4.2 Inserting into and Removing from a Red-Black Tree

Because a red-black tree actually represents a 2-3-4 tree, you simply need to adjust the 2-3-4 insertion algorithms to accommodate the red-black representation. Recall that while searching a 2-3-4 tree, you split each 4-node that you encounter, so it is sufficient to reformulate that process in terms of the red-black representation. For example, Figure 19-31a shows the red-black representation of a 4-node. Thus, to identify a 4-node in its red-black form, you look for a node that has two red pointers.

Suppose that the 4-node is the root of the 2-3-4 tree. Figure 19-28 shows how to split the root into 2-nodes. By comparing this figure with Figure 19-31a, you see that to perform an equivalent operation on a red-black tree, you simply change the color of its root's pointers to black, as Figure 19-33 illustrates.

FIGURE 19-33 Splitting a red-black representation of a 4-node that is the root


Splitting the equivalent of a 4-node requires only simple color changes

Pointer changes called rotations result in a shorter tree

Figure 19-29 shows how to split a 4-node whose parent is a 2-node. If you reformulate this figure by using the red-black notation given in Figure 19-31, you get Figure 19-34. Notice that this case also requires only color changes within the red-black tree.

Finally, Figure 19-30 shows how to split a 4 -node whose parent is a 3 -node. Figure 19-35 shows this process for a red-black tree, using the transformations that Figure 19-31 describes. Note that each of the configurations before the split in Figure 19-30 has two red-black representations, as Figure 19-35 illustrates. As you can see, each pair of representations transforms into the same red-black configuration. Of the six possibilities given in Figure 19-35, only two require simple color changes. The others also require changes to the pointers themselves. These pointer changes, which are called rotations, result in a shorter tree.

The removal algorithm is derived in an analogous fashion from the 2-3-4 tree removal algorithm. Because insertion and removal operations on a red-black tree frequently require only color changes, they are more efficient than the corresponding operations on a 2-3-4 tree.

Exercise 11 asks you to complete the details of the insertion and removal algorithms.

Question 8 What red-black tree represents the 2-3-4 tree in Figure 19-27a?

FIGURE 19-34 Splitting a red-black representation of a 4-node whose parent is a 2 -node, when the 4 -node is a (a) left child; (b) right child
(a)

(b)


FIGURE 19-35 Splitting a red-black representation of a 4-node whose parent is a 2-node
(a)


FIGURE 19-35 Splitting a red-black representation of a 4-node whose parent is a 2-node (continued)
(b)


Rotation and color changes

(c)


Rotation and color changes



### 19.5 AVL Trees

An AVL tree-named for its inventors, Adel'son-Vel'skii and Landis-is a balanced binary search tree. Because the heights of the left and right subtrees of any node in a balanced binary tree differ by no more than 1 , you can search an AVL tree almost as efficiently as a minimum-height binary search tree. This section will simply introduce you to the notion of an AVL tree-which is the oldest form of balanced binary tree-and leave the details for another course.

It is, in fact, possible to rearrange any binary search tree of $n$ nodes to obtain a binary search tree with the minimum possible height $\left\lceil\log _{2}(n+1)\right\rceil$. Recall, for example, the algorithms developed in Section 16.4 of Chapter 16 that use a file to save and restore a binary search tree. You can start with an arbitrary binary search tree, save its values in a file, and then construct from these same values a new binary search tree of minimum height. Although this approach may be appropriate in the context of a dictionary that occasionally is saved and restored, it requires too much work to be performed every time an insertion or removal leaves the tree unbalanced. The cost of repeatedly rebuilding the tree could very well outweigh the benefit of searching a tree of minimum height.

The AVL algorithm is a compromise. It maintains a binary search tree with a height close to the minimum, but it is able to do so with far less work than would be necessary to keep the height of the tree exactly equal to the minimum. The basic strategy of the AVL algorithm is to monitor the shape of the binary search tree. You insert or delete nodes just as you would for any binary search tree, but after each insertion or deletion, you check that the tree is still an AVL tree. That is, you determine whether any node in the tree has left and right subtrees whose heights differ by more than 1 . For example, suppose that the binary search tree in Figure 19-36a is the result of a sequence of insertions and removals. The heights of the left and right subtrees of the root $<30>$ differ by 2 . You can restore this tree's AVL property-that is, its balance-by rearranging its nodes. For instance, you can rotate the tree so that the node $<20>$ becomes the root, with left child $<10>$ and right child $<30>$, as in Figure 19-36b. Notice that you cannot arbitrarily rearrange the tree's nodes, because you must take care not to destroy the search tree's ordering property in the course of the rebalancing.

Rotations are not necessary after every insertion or removal. For example, you can insert 40 into the AVL tree in Figure 19-36b and still have an AVL tree, as Figure 19-36c shows. However, when a rotation is necessary to restore a tree's AVL property, the rotation will be one of two possible types. Let's look at an example of each type.

Suppose that you insert a new node containing 60 into an AVL tree to get the tree in Figure 19-37a. An imbalance occurs at the node $\langle 20>$; that is, this node's left and right subtrees differ in height by more than 1. A single rotation to the left is necessary to obtain the balanced tree in Figure 19-37b: $<40>$ becomes the parent of $<20>$, which adopts $<30>$ as its right child. Figure 19-37c shows this rotation in a more general form. It shows, for example, that before the rotation the left and right

FIGURE 19-36 (a) An unbalanced binary search tree; (b) a balanced tree after rotation; (c) a balanced tree after insertion


An AVL tree is a balanced binary search tree

An AVL tree maintains a height close to the minimum

Rotations restore the balance

FIGURE 19-37 (a) Before; (b) and after a single left rotation that decreases the tree's height; (c) the rotation in general
(a)





FIGURE 19-38 (a) Before; (b) and after a single left rotation that does not affect the tree's height; (c) the rotation in general
(a)

(b)

(c)


subtrees of the node $<40>$ have heights $h$ and $h+1$, respectively. After the rotation, the tree is balanced and, in this particular case, has decreased in height from $h+3$ to $h+2$.

Figure 19-38 shows examples of a single left rotation that restores a tree's balance but does not affect its height. An analogous single right rotation would produce mirror images of these examples.

A more complex rotation might be necessary. For example, consider the tree in Figure 19-39a, which is the result of nodes being added to or deleted from an AVL tree. The left and right subtrees of the node $<40>$ differ in height by more than 1. A double rotation is necessary to restore this tree's balance. Figure $19-39$ b shows the result of a left rotation about $<20>$, and Figure 19-39c shows the result of a right rotation about $\langle 40\rangle$. Figure 19-39d illustrates this double rotation in a more general form. Mirror images of these figures would provide examples of other possible double rotations.

It can be proven that the height of an AVL tree with $n$ nodes will always be very close to the theoretical minimum of $\left\lceil\log _{2}(n+1)\right\rceil$. The AVL tree implementation of a dictionary is, therefore, one implementation that guarantees a binary search-like efficiency. Usually, however, implementations that use either a 2-3-4 tree or a red-black tree will be simpler.

An AVL tree implementation of a dictionary is more difficult than other implementations

FIGURE 19-39 (a) Before; (b) during; and (c) after a double rotation that decreases the tree's height; (d) the double rotation in general


## Summary

1. A 2-3 tree and a 2-3-4 tree are variants of a binary search tree. The internal nodes of a 2-3 tree can have either two or three children. The internal nodes of a 2-3-4 tree can have either two, three, or four children. Allowing the number of children to vary permits the insertion and removal algorithms to maintain the balance of the tree easily.
2. The insertion and removal algorithms for a 2-3-4 tree require only a single pass from root to leaf and, thus, are more efficient than the corresponding algorithms for a 2-3 tree.
3. A red-black tree is a binary tree representation of a 2-3-4 tree that requires less storage than a 2-3-4 tree. Insertions and removals for a red-black tree are more efficient than the corresponding operations on a 2-3-4 tree.
4. An AVL tree is a binary search tree that is guaranteed to remain balanced. The insertion and removal algorithms perform rotations in the event that the tree starts to stray from a balanced shape.

## Exercises

1. Consider the following sequence of operations on an initially empty search tree:

Insert 10
Insert 100
Insert 30
Insert 80
Insert 50
Remove 10
Insert 60
Insert 70
Insert 40
Remove 80
Insert 90
Insert 20
Remove 30
Remove 70
What does the tree look like after these operations execute if the tree is
a. A binary search tree?
d. A red-black tree?
b. A 2-3 tree?
e. An AVL tree?
c. A 2-3-4 tree?
2. What are the advantages of implementing the ADT dictionary with a $2-3$ tree instead of a binary search tree? Why do you not, in general, maintain a completely balanced binary search tree?
3. Write a pseudocode function that performs a range query for a 2-3 tree. That is, the function should visit all items that are within a given range of values. For example, your function should visit all values between 100 and 1,000 .

4. Given the 2-3 tree in Figure 19-40, draw the tree that results after inserting $k, b, c, y$, and $w$ into the tree.
5. Given the 2-3 tree in Figure 19-41, draw the tree that results after removing $t, e, k$, and $d$ from the tree.
6. Draw the 2-3-4 tree that results from inserting $o, d, j, h, s, g$, and $a$, in the order given, into a 2-3-4 tree that contains a single node whose value is $n$.
7. Assume that the tree in Figure 19-5b is a 2-3-4 tree, and insert $39,38,37,36,35,34,33$, and 32 into it. What 2-3-4 tree results?
*8. Write pseudocode for the insertion, removal, retrieval, and traversal operations for a 2-3-4 tree.
9. Figure 19-32 is a red-black tree that represents the 2-3-4 tree in Figure 19-20. Draw another red-black tree that also represents the same 2-3-4 tree.
10. What 2-3-4 tree does the red-black tree in Figure 19-42 represent?
11. Write pseudocode for the insertion, removal, retrieval, and traversal operations for a red-black tree.
12. Write a $\mathrm{C}++$ function that converts a $2-3-4$ tree to a red-black tree.

FIGURE 19-41 A 2-3 tree for Exercise 5


FIGURE 19-42 A red-black tree for Exercise 10


## Programming Problems

1. Implement the ADT dictionary by using a 2-3-4 tree.
*2. Implement the ADT dictionary by using a 2-3 tree. (This implementation is more difficult than the 2-3-4 implementation.)
2. Implement the ADT dictionary by using a red-black tree.
3. Repeat Programming Problem 5 of Chapter 16 using a dictionary implemented with a balanced search tree.

## Chapter

## Graphs

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## Prerequistes

Chapter 5 Recursion as a Problem-Solving Technique
Chapter 6 Stacks
Chapter 15 Trees

Uraphs are an important mathematical concept with significant applications not only in computer science, but also in many other fields. You can view a graph as a mathematical construct, a data structure, or an abstract data type. This chapter provides an introduction to graphs that allows you to view a graph in any of these three ways. It also presents the major operations and applications of graphs that are relevant to the computer scientist.
$G=\{V, E\}$; that is, a graph is a set of vertices and edges

### 20.1 Terminology

You are undoubtedly familiar with graphs: Line graphs, bar graphs, and pie charts are in common use. The simple line graph in Figure 20-1 is an example of the type of graph that this chapter considers: a set of points that are joined by lines. Clearly, graphs provide a way to illustrate data. However, graphs also represent the relationships among data items, and it is this feature of graphs that is important here.

A graph $G$ consists of two sets: a set $V$ of vertices, or nodes, and a set $E$ of edges that connect the vertices. For example, the campus map in Figure 20-2a is a graph whose vertices represent buildings and whose edges represent the sidewalks between the buildings. This definition of a graph is more general than the definition of a line graph. In fact, a line graph, with its points and lines, is a special case of the general definition of a graph.

FIGURE 20-1 An ordinary line graph

## FIGURE 20-2 $\quad$ (a) A campus map as a graph; (b) a subgraph

(a)

(b)


Adjacent vertices are joined by an edge

A path between two vertices is a sequence of edges

A subgraph consists of a subset of a graph's vertices and a subset of its edges. Figure 20-2b shows a subgraph of the graph in Figure 20-2a. Two vertices of a graph are adjacent if they are joined by an edge. In Figure 20-2b, the Library and the Student Union are adjacent. A path between two vertices is a sequence of edges that begins at one vertex and ends at another vertex. For example, there is a path in Figure 20-2a that begins at the Dormitory, leads first to the Library, then to the Student

FIGURE 20-3 Graphs that are (a) connected; (b) disconnected; and (c) complete


Union, and finally back to the Library. Although a path may pass through the same vertex more than once, as the path just described does, a simple path may not. The path Dormitory-Library-Student Union is a simple path. A cycle is a path that begins and ends at the same vertex; a simple cycle is a cycle that does not pass through other vertices more than once. The path Library-Student Union-Gymnasium-Dormitory-Library is a simple cycle in the graph in Figure 20-2a. A graph is connected if each pair of distinct vertices has a path between them. That is, in a connected graph you can get from any vertex to any other vertex by following a path. Figure 20-3a shows a connected graph. Notice that a connected graph does not necessarily have an edge between every pair of vertices. Figure 20-3b shows a disconnected graph.

In a complete graph, each pair of distinct vertices has an edge between them. The graph in Figure 20-3c is complete. Clearly, a complete graph is also connected, but the converse is not true; notice that the graph in Figure 20-3a is connected but is not complete.

Because a graph has a set of edges, a graph cannot have duplicate edges between vertices. However, a multigraph, as illustrated in Figure 20-4a, does allow multiple edges. Thus, a multigraph is not a graph. A graph's edges cannot begin and end at the same vertex. Figure 20-4b shows such an edge, which is called a self edge or loop.

You can label the edges of a graph. When these labels represent numeric values, the graph is called a weighted graph. The graph in Figure 20-5a is a weighted graph whose edges are labeled with the distances between cities.

All of the previous graphs are examples of undirected graphs, because the edges do not indicate a direction. That is, you can travel in either direction along the edges between the vertices of an undirected graph. In contrast, each edge in a directed graph, or digraph, has a direction and is called a

FIGURE 20-4 $\quad$ (a) A multigraph is not a graph; (b) a self edge is not allowed in a graph

A simple path passes through a vertex only once

A cycle is a path that begins and ends at the same vertex

A connected graph has a path between each pair of distinct vertices

A complete graph has an edge between each pair of distinct vertices

A complete graph is connected

A multigraph has multiple edges and so is not a graph

The edges of a weighted graph have numeric labels

Each edge in a directed graph has a direction

(a)

(b)

In a directed graph, vertex $y$ is adjacent to vertex $x$ if there is a directed edge from $x$ to $y$

Vertices can have values

FIGURE 20-5 $\quad$ (a) A weighted graph; (b) a directed graph
(a)

## 2,600

San Francisco

directed edge. Although each distinct pair of vertices in an undirected graph has only one edge between them, a directed graph can have two edges between a pair of vertices, one in each direction. For example, the airline flight map in Figure 20-5b is a directed graph. There are flights in both directions between Providence and New York, but, although there is a flight from San Francisco to Albuquerque, there is no flight from Albuquerque to San Francisco. You can convert an undirected graph to a directed graph by replacing each edge with two directed edges that point in opposite directions.

The definitions just given for undirected graphs apply also to directed graphs, with changes that account for direction. For example, a directed path is a sequence of directed edges between two vertices, such as the directed path in Figure 20-5b that begins in Providence, goes to New York, and ends in San Francisco. However, the definition of adjacent vertices is not quite as obvious for a digraph. If there is a directed edge from vertex $x$ to vertex $y$, then $y$ is adjacent to $x$. (Alternatively, $y$ is a successor of $x$, and $x$ is a predecessor of $y$.) It does not necessarily follow, however, that $x$ is adjacent to $y$. Thus, in Figure 20-5b, Albuquerque is adjacent to San Francisco, but San Francisco is not adjacent to Albuquerque.

### 20.2 Graphs as ADTs

You can treat graphs as abstract data types. Insertion and removal operations are somewhat different for graphs than for other ADTs that you have studied, in that they apply to either vertices or edges. You can define the ADT graph so that its vertices either do or do not contain values. A graph whose vertices do not contain values represents only the relationships among vertices. Such graphs are not unusual, because many problems have no need for vertex values. However, the following ADT graph operations do assume that the graph's vertices contain values.

## Note: ADT graph operations

- Test whether a graph is empty.
- Get the number of vertices in a graph.
- Get the number of edges in a graph.
- See whether an edge exists between two given vertices.
- Insert a vertex in a graph whose vertices have distinct values that differ from the new vertex's value.
- Insert an edge between two given vertices in a graph.
- Remove a particular vertex from a graph and any edges between the vertex and other vertices.
- Remove the edge between two given vertices in a graph.
- Retrieve from a graph the vertex that contains a given value.

Several variations of this ADT are possible. For example, if the graph is directed, you can replace occurrences of "edges" in the previous operations with "directed edges." You can also add traversal operations to the ADT. Graph-traversal algorithms are discussed in Section 20.3.

Listing 20-1 contains an interface that specifies in more detail the ADT operations for an undirected graph.

## LISTING 20-1 A C++ interface for undirected, connected graphs

```
/** An interface for the ADT undirected, connected graph.
    @file GraphInterface.h */
#ifndef _GRAPH_INTERFACE
#define _GRAPH_INTERFACE
template<c7ass Labe1Type>
class GraphInterface
{
public:
    /** Gets the number of vertices in this graph.
        @pre None.
        @return The number of vertices in the graph. */
    virtual int getNumVertices() const = 0;
    /** Gets the number of edges in this graph.
        @pre None.
        @return The number of edges in the graph. */
    virtual int getNumEdges() const = 0;
    /** Creates an undirected edge in this graph between two vertices
        that have the given labels. If such vertices do not exist, creates
        them and adds them to the graph before creating the edge.
        @param start A label for the first vertex.
        @param end A labe1 for the second vertex.
        @param edgeWeight The integer weight of the edge.
        @return True if the edge is created, or false otherwise. */
    virtual bool add(Labe1Type start, LabelType end, int edgeWeight) = 0;
    /** Removes an edge from this graph. If a vertex has no other edges,
        it is removed from the graph since this is a connected graph.
```

```
    @pre None.
    @param start A label for the first vertex.
    @param end A label for the second vertex.
    @return True if the edge is removed, or false otherwise. */
    virtual bool remove(Labe1Type start, Labe1Type end) = 0;
    /** Gets the weight of an edge in this graph.
    @return The weight of the specified edge.
        If no such edge exists, returns a negative integer. */
    virtual int getEdgeWeight(Labe1Type start, LabelType end) const = 0;
    /** Performs a depth-first search of this graph beginning at the given
        vertex and calls a given function once for each vertex visited.
    @param start A label for the first vertex.
    @param visit A client-defined function that performs an operation on
        or with each visited vertex. */
    virtual void depthFirstTraversa1(Labe1Type start, void visit(Labe1Type&)) = 0;
    /** Performs a breadth-first search of this graph beginning at the given
        vertex and calls a given function once for each vertex visited.
    @param start A label for the first vertex.
    @param visit A client-defined function that performs an operation on
        or with each visited vertex. */
    virtual void breadthFirstTraversal(Labe1Type start, void visit(Labe1Type&)) = 0;
}; // end GraphInterface
#endif
```


### 20.2.1 Implementing Graphs

The two most common implementations of a graph are the adjacency matrix and the adjacency list.

Adjacency matrix

Adjacency list An adjacency matrix for a graph with $n$ vertices numbered $0,1, \ldots, n-1$ is an $n$ by $n$ array matrix such that matrix [i] [j] is 1 (true) if there is an edge from vertex $i$ to vertex $j$, and 0 (false) otherwise. Figure $20-6$ shows a directed graph and its adjacency matrix. Notice that the diagonal entries matrix[i] [i] are 0 , although sometimes it can be useful to set these entries to 1 . You should choose the value that is most convenient for your application.

When the graph is weighted, you can let matrix[i][j] be the weight that labels the edge from vertex $i$ to vertex $j$, instead of simply 1 , and let matrix[i][j] equal $\infty$ instead of 0 when there is no edge from vertex $i$ to vertex $j$. For example, Figure 20-7 shows a weighted undirected graph and its adjacency matrix. Notice that the adjacency matrix for an undirected graph is symmetrical; that is, matrix[i][j] equals matrix[j][i].

Our definition of an adjacency matrix does not mention the value, if any, in a vertex. If you need to associate values with vertices, you can use a second array, values, to represent the $n$ vertex values. The array values is one-dimensional, and values [ $i$ ] is the value in vertex $i$.

An adjacency list for a graph with $n$ vertices numbered $0,1, \ldots, n-1$ consists of $n$ linked chains. The $i^{\text {th }}$ linked chain has a node for vertex $j$ if and only if the graph contains an edge from vertex $i$ to vertex $j$. This node can contain the vertex $j$ 's value, if any. If the vertex has no value, the node needs to contain some indication of the vertex's identity. Figure 20-8 shows a directed graph and its adjacency list. You can see, for example, that vertex $0(P)$ has edges to vertex $2(R)$ and vertex $5(W)$. Thus, the first linked chain in the adjacency chain contains nodes for $R$ and $W$.

Figure 20-9 shows an undirected graph and its adjacency list. The adjacency list for an undirected graph treats each edge as if it were two directed edges in opposite directions. Thus, the edge

FIGURE 20-6 (a) A directed graph and (b) its adjacency matrix


FIGURE 20-7 (a) A weighted undirected graph and (b) its adjacency matrix
(a)

(b)

between $A$ and $B$ in Figure 20-9a appears as edges from $A$ to $B$ and from $B$ to $A$ in Figure 20-9b. The graph in 20-9a happens to be weighted; you can include the edge weights in the nodes of the adjacency list, as shown in Figure 20-9b.

Which of these two implementations of a graph-the adjacency matrix or the adjacency list-is better? The answer depends on how your particular application uses the graph. For example, the two most commonly performed graph operations are

1. Determine whether there is an edge from vertex $i$ to vertex $j$
2. Find all vertices adjacent to a given vertex $i$

The adjacency matrix supports the first operation somewhat more efficiently than does the adjacency list. To determine whether there is an edge from $i$ to $j$ by using an adjacency matrix, you need

Two common operations on graphs

## FIGURE 20-8 $\quad$ (a) A directed graph and (b) its adjacency list



FIGURE 20-9 (a) A weighted undirected graph and (b) its adjacency list
(a)

(b)


An adjacency list often requires less space than an adjacency matrix

An adjacency matrix supports operation 1 more efficiently
only examine the value of matrix[i] [j]. If you use an adjacency list, however, you must traverse the $i^{\text {th }}$ linked chain to determine whether a vertex corresponding to vertex $j$ is present.

The second operation, on the other hand, is supported more efficiently by the adjacency list. To determine all vertices adjacent to a given vertex $i$, given the adjacency matrix, you must traverse the $i^{\text {th }}$ row of the array; however, given the adjacency list, you need only traverse the $i^{\text {th }}$ linked chain. For a graph with $n$ vertices, the $i^{\text {th }}$ row of the adjacency matrix always has $n$ entries, whereas the $i^{\text {th }}$ linked chain has only as many nodes as there are vertices adjacent to vertex $i$, a number typically far less than $n$.

Consider now the space requirements of the two implementations. On the surface it might appear that the adjacency matrix requires less memory than the adjacency list, because each entry in the
matrix is simply an integer, whereas each list node contains both a value to identify the vertex and a pointer. The adjacency matrix, however, always has $n^{2}$ entries, whereas the number of nodes in an adjacency list equals the number of edges in a directed graph or twice that number for an undirected graph. Even though the adjacency list also has $n$ head pointers, it often requires less storage than an adjacency matrix.

Thus, when choosing a graph implementation for a particular application, you must consider such factors as what operations you will perform most frequently on the graph and the number of edges that the graph is likely to contain. For example, Chapters 5 and 6 presented the HPAir problem, which was to determine whether an airline provided a sequence of flights from an origin city to a destination city. The flight map for that problem is in fact a directed graph and appeared earlier in this chapter in Figure 20-8a. Figures 20-6b and 20-8b show, respectively, the adjacency matrix and adjacency list for this graph. Because the most frequent operation was to find all cities (vertices) adjacent to a given city (vertex), the adjacency list would be the more efficient implementation of the flight map. The adjacency list also requires less storage than the adjacency matrix, which you can demonstrate as an exercise.

### 20.3 Graph Traversals

The solution to the HPAir problem in Chapter 6 involved an exhaustive search of the graph in Figure $20-8$ a to determine a directed path from the origin vertex (city) to the destination vertex (city). The algorithm searchS started at a given vertex and traversed edges to other vertices until it either found the desired vertex or determined that no (directed) path existed between the two vertices.

What distinguishes searchS from a standard graph traversal is that searchS stops when it first encounters the designated destination vertex. A graph-traversal algorithm, on the other hand, will not stop until it has visited all of the vertices that it can reach. That is, a graph traversal that starts at vertex $v$ will visit all vertices $w$ for which there is a path between $v$ and $w$. Unlike a tree traversal, which always visits all of the nodes in a tree, a graph traversal does not necessarily visit all of the vertices in the graph unless the graph is connected. In fact, a graph traversal visits every vertex in the graph if and only if the graph is connected, regardless of where the traversal starts. (See Exercise 18.) Thus, you can use a graph traversal to determine whether a graph is connected.

If a graph is not connected, a graph traversal that begins at vertex $v$ will visit only a subset of the graph's vertices. This subset is called the connected component containing $v$. You can determine all of the connected components of a graph by repeatedly starting a traversal at an unvisited vertex.

If a graph contains a cycle, a graph-traversal algorithm can loop indefinitely. To prevent such a misfortune, the algorithm must mark each vertex during a visit and must never visit a vertex more than once.

Two basic graph-traversal algorithms, which apply to either directed or undirected graphs, are presented next. These algorithms visit the vertices in different orders, but if they both start at the same vertex, they will visit the same set of vertices. Figure 20-10 shows the traversal order for the two algorithms when they begin at vertex $v$.

### 20.3.1 Depth-First Search

From a given vertex $v$, the depth-first search (DFS) strategy of graph traversal proceeds along a path from $v$ as deeply into the graph as possible before backing up. That is, after visiting a vertex, a DFS visits, if possible, an unvisited adjacent vertex.

An adjacency list supports operation 2 more efficiently

VideoNote
Graph operations

A graph traversal visits all of the vertices that it can reach

A graph traversal visits all vertices if and only if the graph is connected

A connected component is the subset of vertices visited during a traversal that begins at a given vertex

DFS traversal goes as far as possible from a vertex before backing up

## FIGURE 20-11 A connected graph with cycles



```
{
    if (no unvisited vertices are adjacent to the vertex on the top of the stack)
        s.pop() // Backtrack
    else
    {
        Select an unvisited vertex u adjacent to the vertex on the top of the stack
        s.push(u)
        Mark u as visited
    }
}
```

The dfs algorithm is similar to searchS of Chapter 6, but the while statement in searchS terminates when the top of the stack is destinationCity.

For another example of a DFS traversal, consider the graph in Figure 20-11. Figure 20-12 shows the contents of the stack as the previous function dfs visits vertices in this graph, beginning at vertex $a$. Because the graph is connected, a DFS traversal will visit every vertex. In fact, the traversal visits the vertices in this order: $a, b, c, d, g, e, f, h, i$.

The vertex from which a depth-first traversal embarks is the vertex that it visited most recently. This last visited, first explored strategy is reflected both in the explicit stack of vertices that the iterative dfs uses and in the implicit stack of vertices that the recursive dfs generates with its recursive calls.

### 20.3.2 Breadth-First Search

After visiting a given vertex $v$, the breadth-first search (BFS) strategy of graph traversal visits every vertex adjacent to $v$ that it can before visiting any other vertex. As Figure 20-10b illustrates, after marking and visiting $v$, the BFS traversal algorithm marks and then visits each of the vertices $u$, $w$, and $x$. Since no other vertices are adjacent to $v$, the BFS algorithm visits, if possible, all unvisited vertices adjacent to $u$. Thus, the traversal visits $q$ and $t$. Continuing in this manner, the traversal visits vertices in the order given in the figure.

A BFS traversal will not embark from any of the vertices adjacent to $v$ until it has visited all possible vertices adjacent to $v$. Whereas a DFS is a last visited, first explored strategy, a BFS is a first visited, first explored strategy. It is not surprising, then, that a breadth-first search uses a queue. An iterative version of this algorithm follows.

BFS traversal visits all vertices adjacent to a vertex before going forward

FIGURE 20-12 The results of a depth-first traversal, beginning at vertex $a$, of the graph in Figure 20-11

| Node visited | Stack (bottom to top) |
| :---: | :---: |
| a | a |
| b | $a \mathrm{~b}$ |
| c | $a b c$ |
| d | $a b c d$ |
| g | $a b c d g$ |
| e | $a b c d g e$ |
| (backtrack) | $a b c d g$ |
| f | $a b c d g f$ |
| (backtrack) | $a b c d g$ |
| (backtrack) | $a b c d$ |
| h | $a b c d h$ |
| (backtrack) | $a b c d$ |
| (backtrack) | $a b c$ |
| (backtrack) | $a \mathrm{~b}$ |
| (backtrack) | a |
| i | ai |
| (backtrack) | a |
| (backtrack) | (empty) |

An iterative BFS traversal algorithm uses a queue

```
// Traverses a graph beginning at vertex v by using a
// breadth-first search: Iterative version.
bfs(v: Vertex)
    q = a new empty queue
    // Addv to queue and mark it
    q.enqueue(v)
    Mark v as visited
    while (!q.isEmpty())
    {
        q.dequeue(w)
        // Loop invariant: there is a path from vertex w to every vertex in the queue q
        for (each unvisited vertex u adjacent to w)
        {
            Mark u as visited
            q.enqueue(u)
        }
    }
```

Figure 20-13 shows the contents of the queue as bfs visits vertices in the graph in Figure 20-11, beginning at vertex $a$. In general, a breadth-first search will visit the same vertices as a depth-first search, but in a different order. In this example, the BFS traversal visits all of the vertices in this order: $a, b, f, i, c, e, g, d, h$.

A recursive version of BFS traversal is not as simple as the recursive version of DFS traversal. Exercise 19 at the end of this chapter asks you to think about why this is so.

FIGURE 20-13 The results of a breadth-first traversal, beginning at vertex $a$, of the graph in Figure 20-11

| Node visited | Queue (front to back) |
| :---: | :---: |
| a | a (empty) |
| b | b |
| f | b f |
| i | $\underset{\mathrm{fi}}{\mathrm{bfi}}$ |
| c | fic |
| e | $\begin{aligned} & \text { fice } \\ & \text { ice } \end{aligned}$ |
| g | $\begin{aligned} & \text { iceg } \\ & \text { ceg } \end{aligned}$ e g |
| d | $\begin{aligned} & \text { e g d } \\ & \text { g d } \\ & \text { d } \\ & \text { (empty) } \end{aligned}$ |
| h | $\begin{aligned} & \text { h } \\ & \text { (empty) } \end{aligned}$ |

### 20.4 Applications of Graphs

There are many useful applications of graphs. This section surveys some of these common applications.

### 20.4.1 Topological Sorting

A directed graph without cycles, such as the one in Figure 20-14, has a natural order. For example, vertex $a$ precedes $b$, which precedes $c$. Such a graph has significance in ordinary life. If the vertices represent academic courses, the graph represents the prerequisite structure for the courses. For example, course $a$ is a prerequisite to course $b$, which is a prerequisite to both courses $c$ and $e$. In what order should you take all seven courses so that you will satisfy all prerequisites? There is a linear order, called a topological order, of the vertices in a directed graph without cycles that answers this question. In a list of vertices in topological order, vertex $x$ precedes vertex $y$ if there is a directed edge from $x$ to $y$ in the graph.

FIGURE 20-14 A directed graph without cycles


FIGURE 20-15 The graph in Figure 20-14 arranged according to the topological orders (a) $a, g$, $d, b, e, c, f$ and (b) $a, b, g, d, e, f, c$
(a)

(b)


The vertices in a given graph may have several topological orders. For example, two topological orders for the vertices in Figure 20-14 are

$$
a, g, d, b, e, c, f
$$

and

$$
a, b, g, d, e, f, c
$$

If you arrange the vertices of a directed graph linearly and in a topological order, the edges will all point in one direction. Figure 20-15 shows two versions of the graph in Figure 20-14 that correspond to the two topological orders just given.

Arranging the vertices into a topological order is called topological sorting. There are several simple algorithms for finding a topological order. First, you could find a vertex that has no successor. You remove from the graph this vertex and all edges that lead to it, and add it to the beginning of a list of vertices. You add each subsequent vertex that has no successor to the beginning of the list. When the graph is empty, the list of vertices will be in topological order. The following pseudocode describes this algorithm:

A simple topological sorting algorithm

The DFS topological sorting algorithm

```
// Arranges the vertices in graph theGraph into a
// topological order and places them in list aList.
topSort1(theGraph: Graph, aList: List)
    n = number of vertices in theGraph
    for (step = 1 through n)
    {
        Select a vertex v that has no successors
        aList.insert(1, v)
        Remove from theGraph vertex v and its edges
}
```

When the traversal ends, the list aList of vertices will be in topological order. Figure 20-16 traces this algorithm for the graph in Figure 20-14. The resulting topological order is the one that Figure 20-15a represents.

Another algorithm is a simple modification of the iterative depth-first search algorithm. First you push all vertices that have no predecessor onto a stack. Each time you pop a vertex from the stack, you add it to the beginning of a list of vertices. The pseudocode for this algorithm is
// Arranges the vertices in graph theGraph into a
// topological order and places them in list aList.
topSort2 (theGraph: Graph, aList: List)

FIGURE 20-16 A trace of topSort1 for the graph in Figure 20-14
Sraph theGraph

```
s = a new empty stack
for (all vertices v in the graph)
    if (v has no predecessors)
    {
        s.push(v)
        Mark v as visited
    }
while (!s.isEmpty())
{
    if (all vertices adjacent to the vertex on the top of the stack have been visited)
    {
        s.pop(v)
        aList.insert(1, v)
    }
    else
    {
```

```
        Select an unvisited vertex u adjacent to the vertex on the top of the stack
        s.push(u)
        Mark u as visited
    }
}
```

When the traversal ends, the list aList of vertices will be in topological order. Figure 20-17 traces this algorithm for the graph in Figure 20-14. The resulting topological order is the one that Figure 20-15b represents.

FIGURE 20-17 A trace of topSort2 for the graph in Figure 20-14

| Action | Stack s (bottom to top) | List aList (beginning to end) |
| :---: | :---: | :---: |
| Push a | a |  |
| Push g | a g |  |
| Push d | a g d |  |
| Push e | agde | c |
| Push c | agdec | c |
| Pop c, add c to aList | agde | $f \mathrm{c}$ |
| Push f | agdef | efc |
| Pop f, add f to aList | agde | defc |
| Pop e, add e to aList | agd | $g \mathrm{defc}$ |
| Pop d, add d to aList | a g | $g d e f c$ |
| Pop g, add g to aList | a | $b \mathrm{gdefc}$ |
| Push b | a b | $a b g d e f c$ |
| Pop b, add b to aList | a |  |
| Pop a, add a to aList | (empty) |  |

A tree is an undirected connected graph without cycles

Observations about undirected graphs that enable you to detect a cycle

### 20.4.2 Spanning Trees

A tree is a special kind of undirected graph, one that is connected but that has no cycles. Each vertex in the graph in Figure 20-3a could be the root of a different tree. Although all trees are graphs, not all graphs are trees. The nodes (vertices) of a tree have a hierarchical arrangement that is not required of all graphs.

A spanning tree of a connected undirected graph $G$ is a subgraph of $G$ that contains all of $G$ 's vertices and enough of its edges to form a tree. For example, Figure 20-18 shows a spanning tree for the graph in Figure 20-11. The dashed lines in Figure 20-18 indicate edges that were omitted from the graph to form the tree. There may be several spanning trees for a given graph.

If you have a connected undirected graph with cycles and you remove edges until there are no cycles, you will obtain a spanning tree for the graph. It is relatively simple to determine whether a graph contains a cycle. One way to make this determination is based on the following observations about undirected graphs:

1. A connected undirected graph that has $\boldsymbol{n}$ vertices must have at least $\boldsymbol{n} \boldsymbol{- 1}$ edges. To establish this fact, recall that a connected graph has a path between every pair of vertices. Suppose that, beginning with $n$ vertices, you choose one vertex and draw an edge between it and any other vertex. Next, draw an edge between this second vertex and any other unattached vertex. If you continue this process until you run out of unattached vertices, you will get a connected graph like the ones in Figure 20-19. If the graph has $n$ vertices, it has $n-1$ edges. In addition, if you remove an edge, the graph will not be connected.


FIGURE 20-19 Connected graphs that each have four vertices and three edges

2. A connected undirected graph that has $\boldsymbol{n}$ vertices and exactly $\boldsymbol{n}-\mathbf{1}$ edges cannot contain a cycle. To see this, begin with the previous observation: To be connected, a graph with $n$ vertices must have at least $n-1$ edges. If a connected graph did have a cycle, you could remove any edge along that cycle and still have a connected graph. Thus, if a connected graph with $n$ vertices and $n-1$ edges did contain a cycle, removing an edge along the cycle would leave you with a connected graph with only $n-2$ edges, which is impossible according to observation 1 .
3. A connected undirected graph that has $\boldsymbol{n}$ vertices and more than $\boldsymbol{n}-\mathbf{1}$ edges must contain at least one cycle. For example, if you add an edge to any of the graphs in Figure 20-19, you will create a cycle within the graph. This fact is harder to establish and is left as an exercise. (See Exercise 17 at the end of this chapter.)

Thus, you can determine whether a connected graph contains a cycle simply by counting its vertices and edges.

It follows, then, that a tree, which is a connected undirected graph without cycles, must connect its $n$ nodes with $n-1$ edges. Thus, to obtain the spanning tree of a connected graph of $n$ vertices, you must remove edges along cycles until $n-1$ edges are left.

Two algorithms for determining a spanning tree of a graph are based on the previous traversal algorithms and are presented next. In general, these algorithms will produce different spanning trees for any particular graph.
The DFS spanning tree. One way to determine a spanning tree for a connected undirected graph is to traverse the graph's vertices by using a depth-first search. As you traverse the graph, mark the edges that you follow. After the traversal is complete, the graph's vertices and marked edges form a spanning

Count a graph's vertices and edges to determine whether it contains a cycle

FIGURE 20-20 The DFS spanning tree rooted at vertex $a$ for the graph in Figure 20-11


DFS spanning tree algorithm
// Forms a spanning tree for a connected undirected graph
// beginning at vertex $v$ by using depth-first search:
// Recursive version.
dfsTree(v: Vertex)
Mark v as visited
for (each unvisited vertex u adjacent to v )
\{
Mark the edge from $u$ to $v$ dfsTree(u)
\}
When you apply this algorithm to the graph in Figure 20-11, you get the DFS spanning tree rooted at vertex $a$ shown in Figure 20-20. The figure indicates the order in which the algorithm visits vertices and marks edges. You should reproduce these results by tracing the algorithm.

The BFS spanning tree. Another way to determine a spanning tree for a connected undirected graph is to traverse the graph's vertices by using a breadth-first search. As you traverse the graph, mark the edges that you follow. After the traversal is complete, the graph's vertices and marked edges form a spanning tree, which is called the breadth-first search (BFS) spanning tree. (Alternatively, you can remove the unmarked edges from the graph to form the spanning tree.) You can modify the previous iterative version of bfs by marking the edge between $w$ and $u$ before you add $u$ to the queue. The result is the following iterative algorithm to create a BFS spanning tree.
> // Forms a spanning tree for a connected undirected graph
> // beginning at vertex $\vee$ by using breadth-first search:
> // Iterative version.
> bfsTree(v: Vertex)
> $\mathrm{q}=$ a new empty queue

// Add v to queue and mark it


```
q.enqueue(v)
Mark v as visited
while (!q.isEmpty())
{
    q.dequeue(w)
    // Loop invariant: there is a path from vertex w to
    // every vertex in the queue q
    for (each unvisited vertex u adjacent to w)
    {
        Mark u as visited
        Mark edge between w and u
        q.enqueue(u)
    }
}
```

When you apply this algorithm to the graph in Figure 20-11, you get the BFS spanning tree rooted at vertex $a$ shown in Figure 20-21. The figure indicates the order in which the algorithm visits vertices and marks edges. You should reproduce these results by tracing the algorithm.

### 20.4.3 Minimum Spanning Trees

Imagine that a developing country hires you to design its telephone system so that all the cities in the country can call one another. Obviously, one solution is to place telephone lines between every pair of cities. However, your engineering team has determined that due to the country's mountainous terrain, it is impossible to put lines between certain pairs of cities. The team's report contains the weighted undirected graph in Figure 20-22. The vertices in the graph represent $n$ cities. An edge between two vertices indicates that it is feasible to place a telephone line between the cities that the vertices represent, and each edge's weight represents the installation cost of the telephone line. Note that if this graph is not connected, you will be unable to link all of the cities with a network of telephone lines. The graph in Figure 20-22 is connected, however, making the problem feasible.

If you install a telephone line between each pair of cities that is connected by an edge in the graph, you will certainly solve the problem. However, this solution may be too costly. From observation 1 in

A minimum spanning tree of a connected undirected graph has a minimal edge-weight sum

Minimum spanning tree algorithm

FIGURE 20-22 A weighted, connected, undirected graph

the previous section, you know that $n-1$ is the minimum number of edges necessary for a graph of $n$ vertices to be connected. Thus, $n-1$ is the minimum number of lines that can connect $n$ cities.

If the cost of installing each line is the same, the problem is reduced to one of finding any spanning tree of the graph. The total installation cost-that is, the cost of the spanning tree-is the sum of the costs of the edges in the spanning tree. However, as the graph in Figure 20-22 shows, the cost of installing each line varies. Because there may be more than one spanning tree, and because the cost of different trees may vary, you need to solve the problem by selecting a spanning tree with the least cost; that is, you must select a spanning tree for which the sum of the edge weights (costs) is minimal. Such a tree is called the minimum spanning tree, and it need not be unique. Although there may be several minimum spanning trees for a particular graph, their costs are equal.

One simple algorithm, called Prim's algorithm, finds a minimum spanning tree that begins at any vertex. Initially, the tree contains only the starting vertex. At each stage, the algorithm selects a leastcost edge from among those that begin with a vertex in the tree and end with a vertex not in the tree. The latter vertex and least-cost edge are then added to the tree. The following pseudocode describes this algorithm:
> // Determines a minimum spanning tree for a weighted,
> // connected, undirected graph whose weights are
> // nonnegative, beginning with any vertex v .
> primsAlgorithm(v: Vertex)
> Mark vertex $v$ as visited and include it in the minimum spanning tree while (there are unvisited vertices)
> \{
> Find the least-cost edge $(v, u)$ from a visited vertex $v$ to some unvisited vertex $u$ Mark u as visited Add the vertex $u$ and the edge $(v, u)$ to the minimum spanning tree \}

Figure 20-23 traces primsA7 gorithm for the graph in Figure 20-22, beginning at vertex $a$. Edges added to the tree appear as solid lines, while edges under consideration appear as dashed lines.

It is not obvious that the spanning tree that primsAlgorithm determines will be minimal. However, the proof that primsAlgorithm is correct is beyond the scope of this book.

FIGURE 20-23 A trace of primsAlgorithm for the graph in Figure 20-22, beginning at vertex a

(a) Mark a, consider edges from a

(c) Mark f, include edge (a, f)

(e) Mark d, include edge (g, d)

(g) Mark c, include edge (d, c)

(b) Mark i, include edge (a, i)

(d) Mark g, include edge ( $\mathrm{f}, \mathrm{g}$ )

(f) Mark h, include edge (d, h)

(h) Mark e, include edge (c, e)

(i) Mark $b$, include edge $(a, b)$

The shortest path between two vertices in a weighted graph has the smallest edge-weight sum

Finding the shortest paths between vertex 0 and all other vertices

### 20.4.4 Shortest Paths

Consider once again a map of airline routes. A weighted directed graph can represent this map: The vertices are cities, and the edges indicate existing flights between cities. The edge weights represent the mileage between cities (vertices); as such, the weights are not negative. For example, you could combine the two graphs in Figure 20-5 to get such a weighted directed graph.

Often for weighted directed graphs you need to know the shortest path between two particular vertices. The shortest path between two given vertices in a weighted graph is the path that has the smallest sum of its edge weights. Although we use the term "shortest," realize that the weights could be a measure other than distance, such as the cost of each flight in dollars or the duration of each flight in hours. The sum of the weights of the edges of a path is called the path's length or weight or cost.

For example, the shortest path from vertex 0 to vertex 1 in the graph in Figure 20-24a is not the edge between 0 and 1 -its cost is 8 -but rather the path from 0 to 4 to 2 to 1 , with a cost of 7 . For convenience, the starting vertex, or origin, is labeled 0 and the other vertices are labeled from 1 to $n-1$. Notice the graph's adjacency matrix in Figure 20-24b.

The following algorithm, which is attributed to E. Dijkstra, actually determines the shortest paths between a given origin and all other vertices. The algorithm uses a set vertexSet of selected vertices and an array weight, where weight $[v]$ is the weight of the shortest (cheapest) path from vertex 0 to vertex $v$ that passes through vertices in vertexSet.

If $v$ is in vertexSet, the shortest path involves only vertices in vertexSet. However, if $v$ is not in vertexSet, then $v$ is the only vertex along the path that is not in vertexSet. That is, the path ends with an edge from a vertex in vertexSet to $v$.

Initially, vertexSet contains only vertex 0 , and weight contains the weights of the single-edge paths from vertex 0 to all other vertices. That is, weight $[\nu]$ equals matrix $[0][v]$ for all $v$, where matrix is the adjacency matrix. Thus, initially weight is the first row of matrix.

After this initialization step, you find a vertex $v$ that is not in vertexSet and that minimizes weight $[v]$. You add $v$ to vertexSet. For all (unselected) vertices $u$ not in vertexSet, you check the values weight $[u]$ to ensure that they are indeed minimums. That is, can you reduce weight $[u]$ - the weight of a path from vertex 0 to vertex $u$-by passing through the newly selected vertex $v$ ?

To make this determination, break the path from 0 to $u$ into two pieces and find their weights as follows:
weight $[v]=$ weight of the shortest path from 0 to $v$ matrix $[v][u]=$ weight of the edge from $v$ to $u$

Then compare weight $[u]$ with weight $[v]+$ matrix $[v][u]$ and let weight $[u]=$ the smaller of the values weight $[u]$ and weight $[v]+$ matrix $[v][u]$

FIGURE 20-24 (a) A weighted directed graph and (b) its adjacency matrix
(a) Origin

(b)

|  | 0 | 1 | 2 | 3 | 4 |
| :--- | :---: | :---: | :---: | :---: | :---: |
| 0 | $\infty$ | 8 | $\infty$ | 9 | 4 |
| 1 | $\infty$ | $\infty$ | 1 | $\infty$ | $\infty$ |
| 2 | $\infty$ | 2 | $\infty$ | 3 | $\infty$ |
| 3 | $\infty$ | $\infty$ | 2 | $\infty$ | 7 |
| 4 | $\infty$ | $\infty$ | 1 | $\infty$ | $\infty$ |

The pseudocode for Dijkstra's shortest-path algorithm is as follows:

```
// Finds the minimum-cost paths between an origin vertex
// (vertex 0) and all other vertices in a weighted directed
// graph theGraph; theGraph's weights are nonnegative.
shortestPath(theGraph: Graph, weight: WeightArray)
// Step 1: initialization
Create a set vertexSet that contains only vertex 0
n = number of vertices in theGraph
for (v = 0 through n - 1)
    weight[v] = matrix[0][v]
// Steps 2 through n
// Invariant: For v not in vertexSet, weight[v] is the
// smallest weight of all paths from 0 to v that pass
// through only vertices in vertexSet before reaching
// v. For v in vertexSet, weight[v] is the smallest
// weight of all paths from 0 to v (including paths
// outside vertexSet), and the shortest path
// from 0 to v lies entirely in vertexSet.
for (step = 2 through n)
{
        Find the smallest weight [v] such that v is not in vertexSet
        Add v to vertexSet
        // Checkweight[u] for all u not in vertexSet
        for (all vertices u not in vertexSet)
            if (weight[u] > weight[v] + matrix[v][u])
                weight[u] = weight[v] + matrix[v][u]
    }
```

The loop invariant states that once a vertex $v$ is placed in vertexSet, weight $[v]$ is the weight of the absolutely shortest path from 0 to $v$ and will not change.

Figure 20-25 traces the algorithm for the graph in Figure 20-24a. The algorithm takes the following steps:

Step 1. vertexSet initially contains vertex 0 , and weight is initially the first row of the graph's adjacency matrix, shown in Figure 20-24b.

Step 2. weight $[4]=4$ is the smallest value in weight, ignoring weight $[0]$ because 0 is in vertexSet. Thus, $v=4$, so add 4 to vertexSet. For vertices not in vertexSet-that is, for $u=1,2$, and 3 -check whether it is shorter to go from 0 to 4 and then along an edge to $u$ instead of directly from 0 to $u$ along an edge. For vertices 1 and 3, it is not shorter to

FIGURE 20-25 A trace of the shortest-path algorithm applied to the graph in Figure 20-24a

| Step | $\underline{\text { v }}$ | vertexSet | [0] | [1] | $\begin{aligned} & \text { eigh } \\ & {[2]} \end{aligned}$ | [3] | [4] |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 1 | - | 0 | 0 | 8 | $\infty$ | 9 | 4 |
| 2 | 4 | 0,4 | 0 | 8 | 5 | 9 | 4 |
| 3 | 2 | 0, 4, 2 | 0 | 7 | 5 | 8 | 4 |
| 4 | 1 | 0, 4, 2, 1 | 0 | 7 | 5 | 8 | 4 |
| 5 | 3 | 0, 4, 2, 1, 3 | 0 | 7 | 5 | 8 | 4 |

FIGURE 20-26 Checking weight $[u]$ by examining the graph: (a) weight $[2]$ in step 2 ;
(b) weight $[1]$ in step 3 ; (c) weight $[3]$ in step 3 ; (d) weight $[3]$ in step 4
(a)


Step 2. The path $0-4-2$ is shorter than 0-2
(b)


Step 3. The path $0-4-2-1$ is shorter than 0-1
(c)


Step 3 continued. The path $0-4-2-3$ is shorter than 0-3
(d)


Step 4. The path $0-4-2-3$ is shorter than $0-4-2-1-3$
include vertex 4 in the path. However, for vertex 2 notice that weight $[2]=\infty>$ weight $[4]$ + matrix $[4][2]=4+1=5$. Therefore, replace weight[2] with 5 . You can also verify this conclusion by examining the graph directly, as Figure 20-26a shows.
Step 3. weight $[2]=5$ is the smallest value in weight, ignoring weight $[0]$ and weight $[4]$ because 0 and 4 are in vertexSet. Thus, $v=2$, so add 2 to vertexSet. For vertices not in vertex-Set-that is, for $u=1$ and 3 -check whether it is shorter to go from 0 to 2 and then along an edge to $u$ instead of directly from 0 to $u$ along an edge. (See Figures 20-26b and 20-26c.)

Notice that
weight $[1]=8>$ weight $[2]+$ matrix $[2][1]=5+2=7$. Therefore, replace weight $[1]$ with 7 .
weight $[3]=9>$ weight $[2]+$ matrix $[2][3]=5+3=8$. Therefore, replace weight $[3]$ with 8 .
Step 4. weight $[1]=7$ is the smallest value in weight, ignoring weight $[0]$, weight $[2]$, and weight [4] because 0,2 , and 4 are in vertexSet. Thus, $v=1$, so add 1 to vertexSet. For vertex 3 , which is the only vertex not in vertexSet, notice that weight $[3]=8<$ weight $[1]$ + matrix $[1][3]=7+\infty$, as Figure $20-26 \mathrm{~d}$ shows. Therefore, leave weight $[3]$ as it is.

Step 5. The only remaining vertex not in vertexSet is 3 , so add it to vertexSet and stop.
The final values in weight are the weights of the shortest paths. These values appear in the last line of Figure 20-25. For example, the shortest path from vertex 0 to vertex 1 has a cost of weight $[1]$, which is 7 . This result agrees with our earlier observation about Figure 20-24. We saw then that the shortest path is from 0 to 4 to 2 to 1 . Also, the shortest path from vertex 0 to vertex 2 has a cost of weight[2], which is 5 . This path is from 0 to 4 to 2 .

The weights in weight are the smallest possible, as long as the algorithm's loop invariant is true. The proof that the loop invariant is true is by induction on step, and is left as a difficult exercise. (See Exercise 20.)

### 20.4.5 Circuits

A circuit is simply another name for a type of cycle that is common in the statement of certain problems. Recall that a cycle in a graph is a path that begins and ends at the same vertex. Typical circuits either visit every vertex once or visit every edge once.

Probably the first application of graphs occurred in the early 1700s when Euler proposed a bridge problem. Two islands in a river are joined to each other and to the river banks by several bridges, as Figure 20-27a illustrates. The bridges correspond to the edges in the multigraph in Figure 20-27b, and the land masses correspond to the vertices. The problem asked whether you can begin at a vertex $v$, pass through every edge exactly once, and terminate at $v$. Euler demonstrated that no solution exists for this particular configuration of edges and vertices.

For simplicity, we will consider an undirected graph rather than a multigraph. A path in an undirected graph that begins at a vertex $v$, passes through every edge in the graph exactly once, and terminates at $v$ is called an Euler circuit. Euler showed that an Euler circuit exists if and only if each vertex touches an even number of edges. Intuitively, if you arrive at a vertex along one edge, you must be able to leave the vertex along another edge. If you cannot, you will not be able to reach all of the vertices.

FIGURE 20-27 (a) Euler's bridge problem and (b) its multigraph representation

An Euler circuit begins at a vertex $v$, passes through every edge exactly once, and terminates at $v$
(a)

(b)


FIGURE 20-28 Pencil and paper drawings
(a)

(b)


FIGURE 20-29 Connected undirected graphs based on the drawings in Figure 20-28

(b)


Finding an Euler circuit is like drawing each of the diagrams in Figure 20-28 without lifting your pencil or redrawing a line, and ending at your starting point. No solution is possible for Figure 20-28a, but you should be able to find one easily for Figure 20-28b. Figure 20-29 contains undirected graphs based on Figure 20-28. In Figure 20-29a, vertices $h$ and $i$ each touch an odd number of edges (three), so no Euler circuit is possible. On the other hand, each vertex in Figure 20-29b touches an even number of edges, making an Euler circuit feasible. Notice also that the graphs are connected. If a graph is not connected, a path through all of the vertices would not be possible.

Let's find an Euler circuit for the graph in Figure 20-29b, starting arbitrarily at vertex $a$. The strategy uses a depth-first search that marks edges instead of vertices as they are traversed. Recall that a depth-first search traverses a path from $a$ as deeply into the graph as possible. By marking edges instead of vertices, you will return to the starting vertex; that is, you will find a cycle. In this example, the cycle is $a, b, e, d, a$ if we visit the vertices in alphabetical order, as Figure 20-30a shows. Clearly this is not the desired circuit, because we have not visited every edge. We are not finished, however.

To continue, find the first vertex along the cycle $a, b, e, d, a$ that touches an unvisited edge. In our example, the desired vertex is $e$. Apply our modified depth-first search, beginning with this vertex. The resulting cycle is $e, f, j, i, e$. Next you join this cycle with the one you found previously. That is, when you reach $e$ in the first cycle, you travel along the second cycle before continuing in the first cycle. The resulting path is $a, b, e, f, j, i, e, d, a$, as Figure 20-30b shows.

The first vertex along our combined cycle that touches an unvisited edge is $i$. Beginning at $i$, our algorithm determines the cycle $i, h, d, c, g, h, k, l, i$. Joining this to our combined cycle results in the Euler circuit $a, b, e, f, j, i, h, d, c, g, h, k, l, i, e, d, a$. (See Figure 20-30c.)

(b)

(c)


### 20.4.6 Some Difficult Problems

The next three applications of graphs have solutions that are beyond the scope of this book.
The traveling salesperson problem. A Hamilton circuit is a path that begins at a vertex $v$, passes through every vertex in the graph exactly once, and terminates at $v$. Determining whether an arbitrary graph contains a Hamilton circuit can be difficult. A well-known variation of this problem-the traveling salesperson problem-involves a weighted graph that represents a road map. Each edge has an associated cost, such as the mileage between cities or the time required to drive from one city to the next. The salesperson must begin at an origin city, visit every other city exactly once, and return to the origin city. However, the circuit traveled must be the least expensive.

A planar graph can be drawn so that no two edges cross

FIGURE 20-31 The three utilities problem


Unfortunately for this traveler, solving the problem is no easy task. Although a solution does exist, it is quite slow, and no better solution is known.

The three utilities problem. Imagine three houses $A, B$, and $C$ and three utilities $X, Y$, and $Z$ (such as telephone, water, and electricity), as Figure 20-31 illustrates. If the houses and the utilities are vertices in a graph, is it possible to connect each house to each utility with edges that do not cross one another? The answer to this question is no.

A graph is planar if you can draw it in a plane in at least one way so that no two edges cross. The generalization of the three utilities problem determines whether a given graph is planar. Making this determination has many important applications. For example, a graph can represent an electronic circuit where the vertices represent components and the edges represent the connections between components. Is it possible to design the circuit so that the connections do not cross? The solutions to these problems are also beyond the scope of this book.

The four-color problem. Given a planar graph, can you color the vertices so that no adjacent vertices have the same color, if you use at most four colors? For example, the graph in Figure 20-11 is planar because none of its edges cross. You can solve the coloring problem for this graph by using only three colors. Color vertices $a, c, g$, and $h$ red, color vertices $b, d, f$, and $i$ blue, and color vertex $e$ green.

The answer to our question is yes, but it is difficult to prove. In fact, this problem was posed more than a century before it was solved in the 1970s with the use of a computer.


Question 1 Describe the graphs in Figure 20-32. For example, are they directed? Connected? Complete? Weighted?

Question 2 Use the depth-first strategy and the breadth-first strategy to traverse the graph in Figure 20-32a, beginning with vertex 0 . List the vertices in the order in which each traversal visits them.

Question 3 Write the adjacency matrix for the graph in Figure 20-32a.
Question 4 Add an edge to the directed graph in Figure 20-14 that runs from vertex $d$ to vertex $b$. Write all possible topological orders for the vertices in this new graph.

Question 5 Is it possible for a connected undirected graph with five vertices and four edges to contain a simple cycle? Explain.

Question 6 Draw the DFS spanning tree whose root is vertex 0 for the graph in Figure 20-33.
Question 7 Draw the minimum spanning tree whose root is vertex 0 for the graph in Figure 20-33.

Question 8 What are the shortest paths from vertex 0 to each vertex of the graph in Figure 2024a? (Note the weights of these paths in Figure 20-25.)

FIGURE 20-32 Graphs for Checkpoint Questions 1, 2, and 3

(a)

(b)

FIGURE 20-33 A graph for Checkpoint Questions 6 and 7 and for Exercises 1 and 4


## Summary

1. The two most common implementations of a graph are the adjacency matrix and the adjacency list. Each has its relative advantages and disadvantages. The choice should depend on the needs of the given application.
2. Graph searching is an important application of stacks and queues. Depth-first search (DFS) is a graph-traversal algorithm that uses a stack to keep track of the sequence of visited vertices. It goes as deep into the graph as it can before backtracking. Breadth-first search (BFS) uses a queue to keep track of the sequence of visited vertices. It visits all possible adjacent vertices before traversing further into the graph.
3. Topological sorting produces a linear order of the vertices in a directed graph without cycles. Vertex $x$ precedes vertex $y$ if there is a directed edge from $x$ to $y$ in the graph.
4. Trees are connected undirected graphs without cycles. A spanning tree of a connected undirected graph is a subgraph that contains all of the graph's vertices and enough of its edges to form a tree. DFS and BFS traversals produce DFS and BFS spanning trees.
5. A minimum spanning tree for a weighted undirected graph is a spanning tree whose edge-weight sum is minimal. Although a particular graph can have several minimum spanning trees, their edge-weight sums will be the same.
6. The shortest path between two vertices in a weighted directed graph is the path that has the smallest sum of its edge weights.
7. An Euler circuit in an undirected graph is a cycle that begins at vertex $v$, passes through every edge in the graph exactly once, and terminates at $v$.
8. A Hamilton circuit in an undirected graph is a cycle that begins at vertex $v$, passes through every vertex in the graph exactly once, and terminates at $v$.

## Exercises

When given a choice of vertices to visit, the traversals in the following exercises should visit vertices in sorted order.

1. Give the adjacency matrix and adjacency list for
a. The weighted graph in Figure 20-33
b. The directed graph in Figure 20-34
2. Show that the adjacency list in Figure $20-8$ b requires less memory than the adjacency matrix in Figure 20-6b.
3. Consider Figure 20-35 and answer the following:
a. Will the adjacency matrix be symmetrical?
b. Provide the adjacency matrix.
c. Provide the adjacency list.
4. Use both the depth-first strategy and the breadth-first strategy to traverse the graph in Figure 20-33, beginning with vertex 0 , and the graph in Figure 20-36, beginning with vertex $a$. List the vertices in the order in which each traversal visits them.
5. By modifying the DFS traversal algorithm, write pseudocode for an algorithm that determines whether a graph contains a cycle.
6. Using the topological sorting algorithm topSort1, as given in this chapter, write the topological order of the vertices for each graph in Figure 20-37.
7. Trace the DFS topological sorting algorithm topSort2, and indicate the resulting topological order of the vertices for each graph in Figure 20-37.
8. Revise the topological sorting algorithm topSort1 by removing predecessors instead of successors. Trace the new algorithm for each graph in Figure 20-37.
9. Trace the DFS and BFS spanning tree algorithms, beginning with vertex $a$ of the graph in Figure 20-11, and show that the spanning trees are the trees in Figures 20-20 and 20-21, respectively.

FIGURE 20-34 A graph for Exercise 1


FIGURE 20-35 A graph for Exercise 3


FIGURE 20-36 A graph for Exercises 4 and 10

10. Draw the DFS and BFS spanning trees rooted at $a$ for the graph in Figure 20-36. Then draw the minimum spanning tree rooted at $a$ for this graph.
11. For the graph in Figure 20-38,
a. Draw all the possible spanning trees.
b. Draw the minimum spanning tree.
12. Write pseudocode for an iterative algorithm that determines a DFS spanning tree for an undirected graph. Base your algorithm on the traversal algorithm dfs .
13. Draw the minimum spanning tree for the graph in Figure 20-22 when you start with
a. Vertex $g$
b. Vertex $c$
*14. Trace the shortest-path algorithm for the graph in Figure 20-39, letting vertex 0 be the origin.
*15. Implement the shortest-path algorithm in $\mathrm{C}++$. How can you modify this algorithm so that any vertex can be the origin?

FIGURE 20-37 Graphs for Exercises 6, 7, and 8
(a)


FIGURE 20-38 A graph for Exercise 11


## FIGURE 20-39 A graph for Exercise 14


*16. Determine an Euler circuit for the graph in Figure 20-40. Why is one possible?
*17. Prove that a connected undirected graph with $n$ vertices and more than $n-1$ edges must contain at least one simple cycle. (See observation 3 in Section 20.4.2.)
*18. Prove that a graph-traversal algorithm visits every vertex in the graph if and only if the graph is connected, regardless of where the traversal starts.
*19. Although the DFS traversal algorithm has a simple recursive form, a recursive BFS traversal algorithm is not straightforward.
a. Explain why this statement is true.
b. Write the pseudocode for a recursive version of the BFS traversal algorithm.
*20.
Prove that the loop invariant of Dijkstra's shortest-path algorithm is true by using a proof by induction on step.

FIGURE 20-40 A graph for Exercise 16


## Programming Problems

1. Write a C++ class derived from GraphInterface, as given in Listing 20-1. Use an adjacency matrix to represent the graph.
2. Repeat the previous programming problem, but represent the graph using an adjacency list instead of an adjacency matrix.
3. Repeat Programming Problems 1 and 2, but allow the graph to be either weighted or unweighted and either directed or undirected.
4. Extend Programming Problem 3 by adding ADT operations such as isConnected and hasCycle. Also, include operations that perform a topological sort for a directed graph without cycles, determine the DFS and BFS spanning trees for a connected graph, and determine a minimum spanning tree for a connected undirected graph.
5. The HPAir problem was the subject of Programming Problems 11 through 14 of Chapter 6. Revise these problems by implementing the ADT flight map as a derived class of the graph class that you wrote for Programming Problem 3.

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## Processing Data in External Storage

## Chapter

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A
11 of the previous dictionary implementations assume that the data items reside in the computer's internal memory. Many real-world applications, however, require a dictionary so large that it greatly exceeds the amount of available internal memory. In such situations, you must store the dictionary on an external storage device such as a disk and perform dictionary operations there.

This chapter considers the problem of data management in an external environment by using a direct access file as a model of external storage. In particular, this chapter discusses how to sort the data in an external file by modifying the merge sort algorithm and how to search an external file by using generalizations of the hashing and searchtree schemes developed previously.

External storage exists after program execution

Generally, there is more external storage than internal memory

Direct access files are essential for external dictionaries

## A file contains

 records that are organized into blocks
### 21.1 A Look at External Storage

You use external storage when your program reads data from and writes data to a $\mathrm{C}++$ file. Also, when you use a word processing program, for example, and choose Save, the program saves your current document in a file. This action enables you to exit the program and then use it later to retrieve your document for revision. This is one of the advantages of external storage: It exists beyond the execution period of a program. In this sense, it is "permanent" instead of volatile like internal memory.

Another advantage of external storage is that, in general, there is far more of it than internal memory. If you have a collection of one million data items, each of which is an object of moderate size, you will probably not be able to store the entire collection in internal memory at one time. On the other hand, this much data can easily reside on an external disk. As a consequence, when dealing with collections of this magnitude, you cannot simply read all of the data into memory when you want to operate on it and then write it back onto the disk when you are finished. Instead, you must devise ways to operate on data-for example, sort it and search it-while it resides externally.

In general, you can create files for either sequential access or direct access. To access the data stored at a given position in a sequential access file, you must advance the file window beyond all the intervening data. In this sense, a sequential access file resembles a linked chain. To access a particular node in the chain, you must traverse the chain from its beginning until you reach the desired node. In contrast, a direct access file allows you to access the data at a given position directly. A direct access file resembles an array in that you can access the element at data[i] without first accessing the elements before data[i].

Without direct access files, it would be impossible to support the dictionary operations efficiently in an external environment. Many programming languages, including C++, support both sequential access and direct access of files. However, to permit a language-independent discussion, we will construct a model of direct access files that illustrates how a programming language that does not support such files might implement them. This model will be a simplification of reality but will include the features necessary for this discussion.

Imagine that a computer's memory is divided into two parts: internal memory and external memory, as Figure 21-1 illustrates. Assume that an executing program, along with its nonfile data, resides in the computer's internal memory; the permanent files of a computer system reside in the external memory. Further assume that the external storage devices have the characteristics of a disk (although some systems use other devices).

A file consists of data records. A data record can be anything from a simple value, such as an integer, to an aggregate structure, such as an employee record. For simplicity, assume that the data records in any one file are all of the same type.

The records of a file are organized into one or more blocks, as Figure 21-2 shows. The size of a block - that is, the number of bits of data it can contain-is determined by both the hardware configuration and the system software of the computer. In general, an individual program has no control over this size. Therefore, the number of records in a block is a function of the size of the records in the file. For example, a file of integer records will have more records per block than a file of employee records.

FIGURE 21-1 Internal and external memory


FIGURE 21-2 A file partitioned into blocks of records


Much as you number the elements of an array, you can number the blocks of a file in a linear sequence. With a direct access file, a program can read a given block from the file by specifying its block number, and similarly, it can write data out to a particular block. In this regard a direct access file resembles an array of arrays, with each block of the file analogous to a single array element, which is itself an array that contains several records.

In this direct access model, all input and output is at the block level rather than at the record level. That is, you can read and write a block of records, but you cannot read or write an individual record. Reading or writing a block is called a block access.

The algorithms in this chapter assume that commands exist for reading and writing blocks. The pseudocode statement

```
buf.readBlock(dataFile, i)
```

will read the $i^{\text {th }}$ block of file dataFile and place it in an object buf. The object must accommodate the many records that each block of file dataFile contains. For example, if each block contains 100 employee records, buf must store at least 100 employee records. The object buf is called a buffer, which is a location that temporarily stores data as it makes its way from one process or location to another.

Once the system has read a block into buf, the program can process-for example, inspect or modify - the records in the block. Also, because the records in the object buf are only copies of the records in the file dataFile, if a program does modify the records in buf, it must write buf back out to dataFile, so that the file also reflects the modifications. We assume that the statement

```
buf.writeBlock(dataFile, i)
```

will write the contents of buf to the $i^{\text {th }}$ block of the file dataFile. If dataFile contains $n$ blocks, the statement

```
buf.writeBlock(dataFile, n + 1)
```

will append a new block to dataFile, and thus the file can grow dynamically, just as a C++ file can.
Again, realize that these input and output commands allow you to read and write only entire blocks. As a consequence, even if you need to operate on only a single record of the file, you must access an entire block. For example, suppose that you want to give employee Smith a $\$ 1,000$ raise. If Smith's record is in block $i$ (how to determine the correct block is discussed later in the chapter), you would perform the following steps:

[^16]Direct access input and output involves blocks instead of records

A buffer stores data temporarily

Updating a portion of a record within a block

Reduce the number of block accesses

File access time is the dominant factor when considering an algorithm's efficiency

```
Find the entry buf.getRecord(j) that contains the
    record whose search key is "Smith"
// Increase the salary portion of Smith's record
(buf.getRecord(j)).setSalary((buf.getRecord(j)).getSalary() + 1000)
// Write changed block back to file dataFile
buf.writeBlock(dataFile, i)
```

The time required to read or write a block of data is typically much longer than the time required to operate on the block's data once it is in the computer's internal memory. ${ }^{1}$ For example, you typically can inspect every record in the buffer buf in less time than that required to read a block into the buffer. As a consequence, you should reduce the number of required block accesses. In the previous pseudocode, for instance, you should process as many records in buf as possible before writing it to the file. You should pay little attention to the time required to operate on a block of data once it has been read into internal memory.

Interestingly, several programming languages, including $\mathrm{C}++$, have commands to make it appear that you can access records one at a time. In general, however, the system actually performs input and output at the block level and perhaps hides this fact from the program. For example, if a programming language includes the statement
rec. readRecord(dataFile, i) // Reads the ith record of file dataFile into rec.
the system probably accesses the entire block that contains the $i^{\text {th }}$ record. Our model of input and output therefore approximates reality reasonably well.

In most external data-management applications, the time required for block accesses typically dominates all other factors. The rest of the chapter discusses how to sort and search externally stored data. The goal will be to reduce the number of required block accesses.

### 21.2 Sorting Data in an External File

This section considers the following problem of sorting data that resides in an external file:
An external file contains 1,600 employee records. You want to sort these records by Social Security number. Each block contains 100 records, and thus the file contains 16 blocks $B_{1}, B_{2}$, and so on to $B_{16}$. Assume that the program can access only enough internal memory to manipulate about 300 records (three blocks' worth) at one time.

Sorting the file might not sound like a difficult task, because you have already seen several sorting algorithms earlier in this book. There is, however, a fundamental difference here in that the file is far too large to fit into internal memory all at once. This restriction presents something of a problem because the sorting algorithms presented earlier assume that all the data to be sorted is available at one time in internal memory (for example, that it is all in an array). Fortunately, however, we can remove this assumption for a modified version of merge sort.

The basis of the merge sort algorithm is that you can easily merge two sorted segments-such as arrays - of data records into a third sorted segment that is the combination of the two. For example, if $S_{1}$ and $S_{2}$ are sorted segments of records, the first step of the merge is to compare the first record of

[^17]each segment and select the record with the smaller sort key. If the record from $S_{1}$ is selected, the next step is to compare the second record of $S_{1}$ to the first record of $S_{2}$. This process is continued until all of the records have been considered. The key observation is that at any step, the merge never needs to look beyond the leading edge of either segment.

This observation makes a merge sort appropriate for the problem of sorting external files, if you modify the algorithm appropriately. Suppose that the 1,600 records to be sorted are in the file $F$ and that you are not permitted to alter this file. You have two work files, $F_{1}$ and $F_{2}$. One of the work files will contain the sorted records when the algorithm terminates. The algorithm has two phases: Phase 1 sorts each block of records, and phase 2 performs a series of merges.

Phase 1. Read a block from $F$ into internal memory, sort its records by using an internal sort, and write the sorted block out to $F_{1}$ before you read the next block from $F$. After you process all 16 blocks of $F, F_{1}$ contains 16 sorted runs $R_{1}, R_{2}$, and so on to $R_{16}$; that is, $F_{1}$ contains 16 blocks of records, with the records within each block sorted among themselves, as Figure 21-3a illustrates. Phase 2. Phase 2 is a sequence of merge steps. Each merge step merges pairs of sorted runs to form larger sorted runs. With each merge step, the number of blocks in each sorted run doubles, and thus the total number of sorted runs is halved. For example, as Figure 21-3b shows, the first

FIGURE 21-3 (a) Sixteen sorted runs, one block each, in file $F_{1}$; (b) Eight sorted runs, two blocks each, in file $F_{2}$; (c) Four sorted runs, four blocks each, in file $F_{1}$; (d) Two sorted runs, eight blocks each, in file $F_{2}$


Merging sorted runs in Phase 2
merge step merges eight pairs of sorted runs from $F_{1}$ ( $R_{1}$ with $R_{2}, R_{3}$ with $R_{4}, \ldots, R_{15}$ with $R_{16}$ ) to form eight sorted runs, each two blocks long, which are written to $F_{2}$. The next merge step merges four pairs of sorted runs from $F_{2}\left(R_{1}\right.$ with $R_{2}, R_{3}$ with $R_{4}, \ldots, R_{7}$ with $R_{8}$ ) to form four sorted runs, each four blocks long, which are written back to $F_{1}$, as Figure 21-3c illustrates. The next step merges the two pairs of sorted runs from $F_{1}$ to form two sorted runs, which are written to $F_{2}$. (See Figure 21-3d.) The final step merges the two sorted runs into one, which is written to $F_{1}$. At this point, $F_{1}$ will contain all of the records of the original file in sorted order.

Given this overall strategy, how can you merge the sorted runs at each step of phase 2? The statement of the problem provides only sufficient internal memory to manipulate at most 300 records at once. However, in the later steps of phase 2, runs contain more than 300 records each, so you must merge the runs a piece at a time. To accomplish this merge, you must divide the program's internal memory into three arrays, in1, in2, and out, each capable of holding 100 records (the block size). You read block-sized pieces of the runs into the two in arrays and merge them into the out array. Whenever an in array is exhausted-that is, when all of its entries have been copied to out-you read the next piece of the run into an in array; whenever the out array becomes full, you write this completed piece of the new sorted run to one of the files.

Consider how you can perform the first merge step. You start this step with the pair of runs $R_{1}$ and $R_{2}$, which are in the first and second blocks, respectively, of the file $F_{1}$. (See Figure 21-3a.) Because at this first merge step each run contains only one block, an entire run can fit into one of the in arrays. You can thus read $R_{1}$ and $R_{2}$ into the arrays $i n 1$ and in 2 , and then merge in 1 and in 2 into out. However, although the result of merging in 1 and in 2 is a sorted run two blocks long ( 200 records), out can hold only one block ( 100 records). Thus, when in the course of the merge out becomes full, you write

FIGURE 21-4 $\quad$ (a) Merging single blocks; (b) merging long runs

its contents to the first block of $F_{2}$, as Figure 21-4a illustrates. The merging of in1 and in2 into out then resumes. The array out will become full for a second time only after all of the records in in1 and in2 are exhausted. At that time, write the contents of out to the second block of $F_{2}$. You merge the remaining seven pairs from $F$ in the same manner and append the resulting runs to $F_{2}$.

This first merge step is conceptually a bit easier than the others, because the initial runs are only one block in size, and thus each can fit entirely into one of the in arrays. What do you do in the later steps when the runs to be merged are larger than a single block? Consider, for example, the merge step in which you must merge runs of four blocks each to form runs of eight blocks each. (See Figure 21-3c.) The first pair of these runs to be merged is in blocks 1 through 4 and 5 through 8 of $F_{1}$.

The algorithm will read the first block of $R_{1}$-which is the first block $B_{1}$ of the file-into in1, and it will read the first block of $R_{2}$ —which is $B_{5}$-into in2, as Figure 21-4b illustrates. Then, as it did earlier, the algorithm merges in 1 and in 2 into out. The complication here is that as soon as you finish moving all of the records from either in1 or in2, you must read the next block from the corresponding run. For example, if you finish in2 first, you must read the next block of $R_{2}$ —which is $B_{6}$-into in2 before the merge can continue. The algorithm thus must detect when the in arrays become exhausted as well as when the out array becomes full.

A high-level description of the algorithm for merging arbitrary-sized sorted runs $R_{\mathrm{i}}$ and $R_{\mathrm{j}}$ from $F_{1}$ into $F_{2}$ is as follows:

```
Read the first block of R into in1
Read the first block of R into in2
while (either in1 or in2 is not exhausted)
{
    Select the smaller "leading" record of in1 and in2 and place it into
    the next position of out (if one of the arrays is exhausted, select the
    leading record from the other)
    if (out is full)
        Write its contents to the next block of F
    if (in1 is exhausted and blocks remain in R 
        Read the next block into in1
    if (in2 is exhausted and blocks remain in R
        Read the next block into in2
}
```

A pseudocode version of the external sorting algorithm follows. Notice that it uses readBlock and writeBlock, as introduced in the previous section, and assumes a function copyFile that copies a file. To avoid further complications, the solution assumes that the number of blocks in the file is a power of 2. This assumption allows the algorithm always to pair off the sorted runs at each step of the merge phase, avoiding special end-of-file testing that would obscure the algorithm. Also note that the algorithm uses two temporary files and copies the final sorted temporary file to the designated output file.
/ / Sorts a file by using an external merge sort.
// Precondition: unsortedFileName is the name of an external
// file to be sorted. sortedFileName is the name that the
// function will give to the resulting sorted file.
// Postcondition: The new file named sortedFileName is sorted.
// The original file is unchanged. Both files are closed.
// Calls: blockSort, mergeFile, and copyFile.

Pseudocode to merge sorted runs

A pseudocode mergesort function

```
// Simplifying assumption: The number of blocks in the
// unsorted file is an exact power of 2.
externa7Mergesort(unsortedFileName: string,
    sortedFileName: string)
Associate unsortedFileName with the file variable inFile
    and sortedFi1eName with the file variable outFile
// Phase 1: Sort file block by block and count the blocks
blockSort(inFile, tempFile1, numberOfBlocks)
// Phase 2: Merge runs of size 1, 2, 4, 8,..., number0fB7ocks/2
// (uses two temporary files and a toggle that keeps files for each merge step)
togg7e=1
for(size = 1 through numberOfB7ocks/2 with increments of size)
{
    if (togg7e == l)
        mergeFile(tempFile1, tempFile2, size, number0fB1ocks)
    else
    {
            mergeFile(tempFile2, tempFile1, size, numberOfBlocks)
            togg7e = -togg7e
    }
}
// Copy the current temporary file to outFile
if (toggle== 1)
    copyFile(tempFile1, outFile)
else
    copyFile(tempFile2, outFile)
```

Notice that externalMergesort calls blockSort and mergeFile, which calls mergeRuns. The pseudocode for these functions follows.

```
// Sorts each block of records in a file.
// Precondition: The file variable inFi1e is associated
// with the file to be sorted.
// Postcondition:The file associated with the file variable
// outFile contains the blocks of inFile. Each block is
// sorted; numberOfB1ocks is the number of blocks processed.
// Both files are closed.
// Calls: readBlock and writeBlock to perform direct access
// input and output, and sortBuffer to sort an array.
blockSort(inFile: File, outFile: File, numberOfBlocks: integer)
Prepare inFile for input
Prepare outFile for output
numberOfBTocks = 0
while (more blocks in inFile remain to be read)
{
    numberOfBlocks++
    buffer.readBlock(inFile, numberOfBlocks)
    sortArray(buffer) // Sort with some internal sort
    buffer.writeBlock(outFile, numberOfBlocks)
}
Close inFile and outFile
```

```
// Merges blocks from one file to another.
// Precondition: inFi le is an external file that contains
// numberOfBlocks sorted blocks organized into runs of
// runSize blocks each.
// Postcondition: outFile contains the merged runs of
// inFile. Both files are closed.
// Calls:mergeRuns.
mergeFile(inFile: File, outFile: File,
    runSize: integer, numberOfBlocks: integer)
Prepare inFile for input
Prepare outFile for output
for(next =1 to numberOfBlocks with increments of 2 * runSize)
{
    // Invariant: Runs in outFile are ordered
    mergeRuns(inFile, outFile, next, runSize)
}
Close inFile and outFile
// Merges two consecutive sorted runs in a file.
// Precondition: fromFile is an external file of sorted runs
// open for input. toFiTe is an external file of sorted runs
// open for output. start is the block number of the first
// run on fromFile to be merged; this run contains size
// blocks.
// Run 1: Block start to block start + size - 1
// Run 2: Block start + size tostart + (2 * size) - 1
// Postcondition:The merged runs from fromFi Te are appended
// to toFile. The files remain open.
mergeRuns(fromFile: File, toFile: File,
start: integer, size: integer)
    // Initialize the input buffers for runs 1 and 2
    in1.readBlock(fromFile, first block of Run 1)
    in2.readBlock(fromFile, first block of Run 2)
    // Merge until one of the runs is finished. Whenever an
    // input buffer is exhausted, the next block is read.
    // Whenever the output buffer is full, it is written.
while (neither run is finished)
{
    // Invariant: out and each block in toFile are ordered
    Select the smaller "leading edge" of in1 and in2, and
        place it in the next position of out
    if (out is full)
        out.writeBlock(toFile, next block of toFile)
    if (in1 is exhausted and blocks remain in Run 1)
        in1.readBlock(fromFile, next block of Run 1)
    if (in2 is exhausted and blocks remain in Run 2)
        in2.readBlock(fromFile, next block of Run 2)
}
// Assertion: Exactly one of the runs is complete
// Append the remainder of the unfinished input
// buffer to the output buffer and write it
```

A simple externa dictionary implementation: records stored in search-key order

## Sorted-order

 traversal```
while (in1 is not exhausted)
    // Invariant: out is ordered
    Place next item of in1 into the next position of out
while(in2 is not exhausted)
    // Invariant: out is ordered
    Place next item of in2 into the next position of out
out.writeBlock(toFile, next block of toFile)
// Finish off the remaining complete blocks
while (blocks remain in Run 1)
{
    // Invariant: Each block in toFile is ordered
    in1.readBlock(fromFile, next block of Run l)
    in1.writeBlock(toFile, next block of toFile)
}
while (blocks remain in Run 2)
{
    // Invariant: Each block in toFi1e is ordered
    in2.readB7ock(fromFi1e, next block of Run 2)
    in2.writeBlock(toFi1e, next block of toFi1e)
}
```


### 21.3 External Dictionaries

This section discusses techniques for organizing records in external storage so that you can efficiently perform ADT dictionary operations such as retrieval, insertion, removal, and traversal. Although this discussion will only scratch the surface of this topic, you do have a head start: Two of the most important external dictionary implementations are variations of hashing and the 2-3 tree, which you studied in Chapters 18 and 19.

Suppose that you have a direct access file of records that are to be dictionary items. The file is partitioned into blocks, as described earlier in this chapter. One of the simplest dictionary implementations stores the records in order by their search key, perhaps sorting the file by using the external merge sort algorithm developed in the previous section. Once it is sorted, you can easily traverse the file in sorted order by using the following algorithm:

```
// Traverses the sorted file dataFile in sorted order,
// calling function visit once for each item.
traverse(dataFile: File, numberOfBlocks: integer,
    recordsPerBlock: integer, visit: FunctionType)
// Read each block of file dataFile into an internal buffer buf
for (blockNumber = 1 through numberOfBlocks)
{
    buf.readBlock(dataFile, blockNumber)
    // Visit each record in the block
    for (recordNumber = 1 through recordsPerBlock)
        Visit record buf.getRecord(recordNumber-1)
}
```

To perform the retrieval operation on the sorted file, you can use a binary search algorithm as follows:
// Searches blocks first through last of the file dataFile
// for the record whose search key equals searchkey

```
// Returns the record iffound, else throws NotFoundException.
getItem(dataFile: File, recordsPerBlock: integer,
            first: integer, last: integer,
            searchKey: KeyType): ItemType
if (first > 1ast or nothing is left to read from dataFile)
        throw NotFoundException
    else
    \{
        // Read the middle block of file dataFile into array buf
        mid \(=(\) first +7 last) \(/ 2\)
        buf.readBlock(dataFile, mid)
        if ((searchKey >= (buf.getRecord(0)).getKey ()) \&\&
        (searchKey <= (buf.getRecord(recordsPerBlock-1)).getKey()) )
    \{
        // Desired block is found
        Search buffer buf for record buf.getRecord ( \(j\) ) whose search key equals searchKey
        if (record is found)
            return = buf.getRecord (j)
        else
            throw NotFoundException
    \}
    // Else search appropriate half of the file
    else if (searchKey < (buf.getRecord(0)).getKey())
        return getItem(dataFile, recordsPerBlock, first, mid-1, searchKey)
    e1se
        return getItem(dataFile, recordsPerBlock, mid+1, last, searchKey)
\}
```

The retrieval algorithm recursively splits the file in half and reads the middle block into the internal object buf. Splitting a file segment requires that you know the numbers of the first and last blocks of the segment. You would pass these values as arguments, along with the file variable, to getItem.

Once you have read the middle block of the file segment into buf, you determine whether a record whose search key equals searchKey could be in this block. You can make this determination by comparing searchKey to the smallest search key in buf-which is in buf.getRecord ( 0 ) -and to the largest search key in buf, which is in buf.getRecord(recordsPerBlock-1). If searchKey does not lie between the values of the smallest and largest search keys in buf, you must recursively search one of the halves of the file (which half to search depends on whether searchKey is less than or greater than the search keys in the block you just examined). If, on the other hand, searchKey does lie between the values of the smallest and largest search keys of the block in buf, you must search buf for the record. Because the records within the block buf are sorted, you could use a binary search on the records within this block. However, the number of records in the block buf is typically small, and thus the time required to scan the block sequentially is insignificant compared to the time required to read the block from the file. It is therefore common simply to scan the block sequentially.

This external implementation of the ADT dictionary is not very different from the internal sorted array-based implementation. As such, it has many of the same advantages and disadvantages. Its main advantage is that because the records are sorted sequentially, you can use a binary search to locate the block that contains a given search key. The main disadvantage of the implementation is
add and remove for an external implementation of the ADT dictionary can require many costly file accesses due to shifting records

Advantages of a library catalog

FIGURE 21-5 Shifting across block boundaries

that, as is the case with an array-based implementation, the add and remove operations must shift dictionary items. Shifting records in an external file is, in general, far more costly than shifting array items. A file may contain an enormous number of large records, which are organized as several thousand blocks. As a consequence, the shifting could require a prohibitively large number of block accesses.

Consider, for example, Figure 21-5. If you insert a new record into block $k$, you must shift the records not only in block $k$, but also in every block after it. As a result, you must shift some records across block boundaries. Thus, for each of these blocks, you must read the block into internal memory, shift its records by using a statement such as

```
buf.setRecord(i+1, buf.getRecord(i))
```

and write the block to the file so that the file reflects the change. This large number of block accesses makes the external sorted array-based implementation practical only for dictionaries where insertions and removals are rare.

### 21.3.1 Indexing an External File

Two of the best external dictionary implementations are variations of the internal hashing and searchtree schemes. The biggest difference between the internal and external versions of these implementations is that in the external versions, it is often advantageous to organize an index to the data file rather than to organize the data file itself. An index to a data file is conceptually similar to other indexes with which you are familiar. For example, consider a library catalog. Rather than looking all over the library for a particular title, you can simply search the catalog. The catalog is typically organized alphabetically by title (or by author), so it is a simple matter to locate the appropriate entry. The entry for each book contains an indication (for example, a Library of Congress number) of where on the shelves you can find the book.

Using a catalog to index the books in a library has at least three benefits:

- Because each catalog entry is much smaller than the book it represents, the entire catalog for a large library can fit into a small space. A patron can thus locate a particular book quickly.
- The library can organize the books on the shelves in any way, without regard to how easy it will be for a patron to scan the shelves for a particular book. To locate a particular book, the patron searches the catalog for the appropriate entry.
- The library can have different types of catalogs to facilitate different types of searches. For example, it can have one catalog organized by title and another organized by author.

Now consider how you can use an index to a data file to much the same advantage as the library catalog. As Figure 21-6 illustrates, you can leave the data file in a disorganized state and maintain an organized index to it. When you need to locate a particular record in the data file, you search the index for the corresponding entry, which will tell you where to find the desired record in the data file.

An index to the data file is simply another file, called the index file, that contains an index record for each record in the data file, just as a library catalog contains an entry for each book in the library. An index record has two parts: a key, which contains the same value as the search key of its corresponding record in the data file, and a "pointer," which shows the number of the block in the data file that contains this data record. (Despite its name, an index record's pointer contains an integer, not a C++ pointer.) You thus can determine which block of the data file contains the record whose search key equals searchKey by searching the index file for the index record whose key equals searchKey.

Maintaining an index to a data file has benefits analogous to those provided by the library's catalog:

- In general, an index record will be much smaller than a data record. While the data record may contain many components, an index record contains only two: a key, which is also part of the data record, and a single integer pointer, which is the block number. Thus, just as a library catalog occupies only a small fraction of the space occupied by the books it indexes, an index file is only a fraction of the size of the data file. As you will see, the small size of the index file often allows you to manipulate it with fewer block accesses than you would need to manipulate the data file.
- Because you do not need to maintain the data file in any particular order, you can insert new records in any convenient location, such as at the end of the file. As you will see, this flexibility eliminates the need to shift the data records during insertions and removals.
- You can maintain several indexes simultaneously. Just as a library can have one catalog organized by title and another organized by author, you can have one index file that indexes the data file by one search key (for example, an index file that consists of <name, pointer> records), and a second index file that indexes the data file by another search key (for example, an index file that consists of <socSec, pointer> records). Such multiple indexing is discussed briefly at the end of this chapter.

Although you do not organize the data file, you must organize the index file so that you can search and update it rapidly. Before considering how to organize an index file by using either hashing or search-tree schemes, first consider a less complex organization that illustrates the concepts of

An index to a data file

Advantages of an index file

Organize the index file, but not the data file

FIGURE 21-6 A data file with an index


Data file: blocks of large, unorganized data records

FIGURE 21-7 A data file with a sorted index file

indexing. In particular, let the index file simply store the index records sequentially, sorted by their keys, as shown in Figure 21-7.

To perform the retrieval operation, for example, you can use a binary search on the index file as follows:

```
// Searches the file dataFi le for the record whose search key equals searchKey.
// Returns the record if found, else throws NotFoundException.
getItem(indexFile: File, dataFile: File,
            searchKey: KeyType): ItemType
    if (no blocks are left in indexFile to read)
        throw NotFoundException
    else
    {
    // Read the middle block of indexfi7e into object buf
    mid = number of middle block of indexFile
    buf.readBlock(indexFile, mid)
    if ((searchKey >= (buf.getRecord(0)).getKey()) &&
        (searchKey <= (buf.getRecord(indexrecordsPerBlock-1)).getKey()))
    {
        // Desired block of index file found
        Search buf for index file record whose key value equals searchKey
        if (index record buf.getRecord(j) is found)
        {
            blockNum = number of the data-file block to which buf.getRecord(j) points
            data.readBlock(dataFile, blockNum)
            Find data record whose search key equals searchKey
            return data.getRecord(k)
    }
    e1se
            throw NotFoundException
    }
    e1se if (indexFile is one block in size)
        throw NotFoundException // No more blocks in file
```

```
    // Else search appropriate half of index file
    else if (searchKey < (buf.getRecord(0)).getKey())
    return getItem(first half of indexFile, dataFile,
        searchKey)
    else
        return getItem(second half of indexFile, dataFile,
        searchKey)
```

\}

Because the index records are far smaller than the data records, the index file contains far fewer blocks than the data file. For example, if the index records are one-tenth the size of the data records and the data file contains 1,000 blocks, the index file will require only about 100 blocks. As a result, the use of an index cuts the number of block accesses in getItem down from about $\log _{2} 1000 \approx 10$ to about $1+\log _{2} 100 \approx 8$. (The one additional block access is into the data file once you have located the appropriate index record.)

The reduction in block accesses is far more dramatic for the add and remove operations. In the implementation of an external dictionary discussed earlier in this section, if you insert a record into or remove a record from the first block of data, for example, you have to shift records in every block, requiring that you access all 1,000 blocks of the data file. (See Figure 21-5.)

However, when you perform an insertion or a removal by using the index scheme, you have to shift only index records. When you use an index file, you do not keep the data file in any particular order, so you can insert a new data record into any convenient location in the data file. This flexibility means that you can simply insert a new data record at the end of the file or at a position left vacant by a previous removal (as you will see). As a result, you never need to shift records in the data file. However, you do need to shift records in the index file to create an opening for a corresponding index entry in its proper sorted position. Because the index file contains many fewer blocks than the data file ( 100 versus 1,000 in the previous example), the maximum number of block accesses required is greatly reduced. A secondary benefit of shifting index records rather than data records is a reduction in the time requirement for a single shift. Because the index records themselves are smaller, the time required for the statement buf. setRecord( $\mathrm{i}+1$, buf.getRecord(i)) is decreased.

Removals under the index scheme reap similar benefits. Once you have searched the index file and located the data record to be removed, you can simply leave its location vacant in the data file, and thus you need not shift any data records. You can keep track of the vacant locations in the data file, so that you can insert new data records into the vacancies, as was mentioned earlier. The only shifting required is in the index file to fill the gap created when you remove the index record that corresponds to the deleted data record.

Even though this scheme is an improvement over maintaining a sorted data file, in many applications it is far from satisfactory. The 100 block accesses that could be required to insert or remove an index record often would be prohibitive. Far better implementations are possible when you use either hashing or search trees to organize the index file.

### 21.3.2 External Hashing

The external hashing scheme is quite similar to the internal scheme described in Chapter 18. In the internal hashing scheme, each entry of the array table - the hash table - contains a pointer to the beginning of a chain of items that hash into that location. In the external hashing scheme, each entry of table still contains a pointer to the beginning of a chain, but here each chain consists of blocks of index records. In other words, you hash an index file rather than the data file, as Figure 21-8

An index file reduces the number of required block accesses for dictionary operations

Shift index records instead of data records

An unsorted data file with a sorted index is more efficient than a sorted data file, but other schemes are even better

You hash the index file instead of the data file

FIGURE 21-8 A hashed index file


Data file—each block contains several data records
illustrates. (In many applications the array table is itself so large that you must keep it in external storage-for example, in the first $K$ blocks of the index file. To avoid this extra detail, you can assume here that the array table is an internal array.)

Associated with each entry table[i] is a linked chain of blocks of the index file, as you can see in Figure 21-8. Each block of table[i]'s linked chain contains index records whose keys (and thus whose corresponding data records' search keys) hash into location i. To form the linked chains, you must reserve space in each block for a block pointer-the integer block number of the next block in the chain-as Figure 21-9 illustrates. That is, in this linked chain the pointers are integers, not C++ pointers. A pointer value of -1 is used as a null pointer.

Retrieval under external hashing of an index file. The retrieval operation appears in pseudocode as follows:
// Searches the data file for the record whose search key equals searchKey.
// Returns the record if found, else throws NotFoundException.
getItem(indexFile: File, dataFile: File, searchKey: KeyType): ItemType

FIGURE 21-9 A single block with a pointer

```
// Apply the hash function to the search key
i = h(searchKey)
// Find the first block in the chain of index blocks -
// these blocks contain index records that hash into location i
p = table[i]
// ifp ==-1, no values have hashed into location i
if (p != -1)
    buf.readBlock(indexFile, p)
// Search for the block with the desired index record
while (p != -1 and buf does not contain an index record whose key value equals searchKey)
{
    p = number of next block in chain
    // if p equals -1, you are at the last block in the chain
    if (p != -1)
        buf.readBlock(indexFile, p)
}
```

// Retrieve the data item if present
if ( $p$ ! $=-1$ )
\{
// buf.getRecord(j) is the index record whose key value equals searchKey
blockNum = number of the data-file block to which buf.getRecord ( $j$ ) points
data.readBlock(dataFile, blockNum)
Find data record data.getRecord(k) whose search key equals searchKey
return data.getRecord(k)
\}
else
throw NotFoundException

Insertion under external hashing of an index file. The external hashing versions of the insertion and removal operations are also similar to the internal hashing versions. The major difference is that, in the external environment, you must insert or remove both a data record and the corresponding index record.

To insert a new data record whose search key is searchKey, you take the following steps:

1. Insert the data record into the data file. Because the data file is not ordered, the new record can go anywhere you want. If a previous removal has left a free slot in the middle of the data file, you can insert it there.

If no slots are free, you insert the new data record at the end of the last block, or, if necessary, you append a new block to the end of the data file and store the record there. In either case, let p denote the number of the block that contains this new data record.
2. Insert a corresponding index record into the index file. You need to insert into the index file an index record that has key value searchKey and pointer value $p$. (Recall that $p$ is the number of the block in the data file into which you inserted the new data record.) Because the index file is hashed, you first apply the hash function to searchKey, letting

```
i=h(searchKey)
```

You then insert the index record < searchKey, p> into the chain of blocks that the entry table[i] points to. You can insert this record into any block in the chain that contains a free slot, or, if necessary, you can allocate a new block and link it to the beginning of the chain.

Choose external hashing for basic dictionary operations

## Organize the index

 file as an external 2-3 treeRemoval under external hashing of an index file. To remove the data record whose search key is searchKey, you take the following steps:

1. Search the index file for the corresponding index record. You apply the hash function to searchKey, letting

$$
i=h(\operatorname{searchKey})
$$

You then search the chain of index blocks pointed to by the entry table [i] for an index record whose key value equals searchKey. If you do not find such a record, you can conclude that the data file does not contain a record whose search key equals searchKey. However, if you find an index record <searchKey, p>, you remove it from the index file after noting the block number $p$, which indicates where in the data file you can find the data record to be removed.
2. Remove the data record from the data file. You know that the data record is in block p of the data file. You simply access this block, search the block for the record, remove the record, and write the block back to the file.

Observe that for each of the operations getItem, add, and remove the number of block accesses is very low. You never have to access more than one block of the data file, and at worst you have to access all of the blocks along a single hash chain of the index file. You can take measures to keep the length of each of the chains quite short (for example, one or two blocks long), just as you can with internal hashing. You should make the size of the array table large enough so that the average length of a chain is near one block, and the hash function should scatter the keys evenly. If necessary, you can even structure each chain as an external balanced search tree by using the techniques described in the next section.

The hashing implementation is the one to choose when you need to perform retrieval, insertion, and removal operations on a large external dictionary. As is the case with internal hashing, however, this implementation is not practical for certain other operations, such as sorted traversal, retrieval of the smallest or largest item, and range queries that require ordered data. When these types of operations are added to the basic dictionary operations, you should use a search-tree implementation instead of hashing.

### 21.3.3 B-Trees

Another way to search an external dictionary is to organize it as a balanced search tree. Just as you can apply external hashing to the index file, you can organize the index file, not the data file, as an external search tree. The implementation developed here is a generalization of the 2-3 tree of Chapter 19.

You can organize the blocks of an external file into a tree structure by using block numbers for child pointers. In Figure 21-10a, for example, the blocks are organized into a 2-3 tree. Each block of the file is a node in the tree and contains three child pointers, each of which is the integer block number of the child. A child pointer value of -1 plays the role of a null pointer, and thus, for example, a leaf will contain three child pointers with the value -1 .

If you organized the index file into a 2-3 tree, each node (block of the index file) would contain either one or two index records, each of the form $<$ key, pointer $>$, and three child pointers. The pointer portion of an index record has nothing to do with the tree structure of the index file; pointer indicates the block (in the data file) that contains the data record whose search key equals key. (See Figure 21-10b.) To help avoid confusion, the pointers in the tree structure of the index file will be referred to as child pointers.

FIGURE 21-10 $\quad$ (a) Blocks organized into a 2-3 tree; (b) a single node of the 2-3 tree


You must organize the index records in the tree so that their keys obey the same search-tree ordering property as an internal 2-3 tree. This organization allows you to retrieve the data record with a given value in its search key as follows:

```
// Searches the data file for the record whose search key equals searchKey.
// Returns the record iffound, else throws NotFoundException.
// rootNum is the block number (of the index file) that contains the root of the tree.
getItem(indexFile: File, dataFile: File, rootNum: integer,
            searchKey: KeyType): ItemType
    if (no blocks are left in the index file to read)
    throw NotFoundException
    else
    {
    // Read from index file into internal array buf the block that contains the root of the 2-3 tree
    buf.readBlock(indexFile, rootNum)
    // Search for the index record whose key value equals searchKey
    if (searchKey is in the root)
    {
        blockNum = number of the data-file block that index record specifies
        data.readBlock(dataFile, blockNum)
        Find data record data.getRecord(k) whose search key equals searchKey
        return data.getRecord(k)
    }
```

An external 2-3 tree is adequate, but an improvement is possible

## Keep an external

 search tree short
## Binary search tree:

 the number of records and children per node2-3 tree: the number of records and children per node

## General search tree:

 the number of records and children per node```
// Else search the appropriate subtree
else if (the root is a leaf)
    throw NotFoundException
else
{
    child = block number of root of appropriate subtree
    return getItem(indexFile, dataFile, child, searchKey)
}
}
```

You also can perform insertions and removals in a manner similar to those in the internal version, with the addition that you must insert records into and remove records from both the index file and the data file (as was the case in the external hashing scheme described earlier). In the course of insertions into and removals from the index file, you must split and merge nodes of the tree just as you do for the internal version. You perform insertions into and removals from the data file-which, recall, is not ordered in any way-exactly as described for the external hashing implementation. You thus can support the dictionary operations fairly well by using an external version of the 2-3 tree.

However, you can generalize the 2-3 tree to a structure that is even more suitable for an external environment. Recall the discussion in Chapter 19 about search trees whose nodes can have many children. Adding more children per node reduces the height of the search tree but increases the number of comparisons at each node during the search for a value.

In an external environment, however, the advantage of keeping a search tree short far outweighs the disadvantage of performing extra work at each node. As you traverse the search tree in an external environment, you must perform a block access for each node visited. Because the time required to access a block of an external file is, in general, far greater than the time required to process the data in that block once it has been read in, the overriding concern is to reduce the number of block accesses required. This fact implies that you should attempt to reduce the height of the tree, even at the expense of requiring more comparisons at each node. In an external search tree, you should thus allow each node to have as many children as possible, with only the block size as a limiting factor.

How many children can a block of some fixed size accommodate? If a node is to have $m$ children, clearly you must be able to fit $m$ child pointers in the node. In addition to child pointers, however, the node must also contain index records. Before you can answer the question of how many children a block can accommodate, you must first consider this related question: If a node $N$ in a search tree has $m$ children, how many key values - and thus how many index records-must it contain?

In a binary search tree, if the node $N$ has two children, it must contain one key value, as Figure 21-11a indicates. You can think of the key value in node $N$ as separating the key values in $N$ 's two subtrees - all of the key values in $N$ 's left subtree are less than $N$ 's key value, and all of the key values in $N$ 's right subtree are greater than $N$ 's key value. When you are searching the tree for a given key value, the key value in $N$ tells you which branch to take.

Similarly, if a node $N$ in a 2-3 tree has three children, it must contain two key values. (See Figure 21-11b.) These two values separate the key values in $N$ 's three subtrees-all of the key values in the left subtree are less than $N$ 's smaller key value, all of the key values in $N$ 's middle subtree lie between $N$ 's two key values, and all of the key values in $N$ 's right subtree are greater than $N$ 's larger key value. As is the case with a binary search tree, this requirement allows a search algorithm to know which branch to take at any given node.

In general, if a node $N$ in a search tree is to have $m$ children, it must contain $m-1$ key values to separate the values in its subtrees correctly. (See Figure 21-11c.) Suppose that you denote the subtrees of $N$ as $S_{0}, S_{1}$, and so on to $S_{m-1}$ and denote the key values in $N$ as $K_{1}, K_{2}$, and so on to $K_{m-1}$ (with $K_{1}<K_{2}<\cdots<K_{m-1}$ ). The key values in $N$ must separate the values in its subtrees as follows:

FIGURE 21-11 (a) A node with two children; (b) a node with three children; (c) a node with $m$ children
(a)

(b)

(c)


- All the values in subtree $S_{0}$ must be less than the key value $K_{1}$.
- For all $i, 1 \leq i \geq m-2$, all the values in subtree $S_{i}$ must lie between the key values $K_{i}$ and $K_{i+1}$.
- All the values in subtree $S_{m-1}$ must be greater than the key value $K_{m-1}$.

If every node in the tree obeys this property, you can search the tree by using a generalized version of a search tree's retrieval algorithm as follows:

```
// Searches the data file for the record whose search key equals searchKey.
// Returns the record iffound, else throws NotFoundException.
getItem(indexFile: File, dataFile: File, rootNum: integer,
        searchKey: KeyType): ItemType
    if (no blocks are left in the index file to read)
    throw NotFoundException
    e1se
    { // Read from index file into internal array buf the
        // block that contains the root of the tree
        buf.readBlock(indexFile, rootNum)
        // Search for the index record whose key value equals searchKey
        if (searchKey is one of the Ki in the root)
        {
            blockNum = number of the data-file block that index record specifies
            data.readBlock(dataFile, blockNum)
            Find data record data.getRecord(k) whose search key equals searchKey
            return data.getRecord(k)
        }
        // Else search the appropriate subtree
        else if (the root is a leaf)
            throw NotFoundException
```

Number of children per node

```
    else
    {
        Determine which subtree Si to search
        chi1d = block number of the root of Si
        return getItem(indexFile, dataFile, child, searchKey)
    }
}
```

Now return to the question of how many children the nodes of the search tree can have-that is, how big can $m$ be? If you wish to organize the index file into a search tree, the items that you store in each node will be records of the form <key, pointer>. Thus, if each node in the tree (which, recall, is a block of the index file) is to have $m$ children, it must be large enough to accommodate $m$ child pointers and $m-1$ records of the form <key, pointer>. You should choose $m$ to be the largest integer such that $m$ child pointers (which, recall, are integers) and $m-1<k e y$, pointer> records can fit into a single block of the file. Actually, the algorithms are somewhat simplified if you always choose an odd number for $m$. That is, you should choose $m$ to be the largest odd integer such that $m$ child pointers and $m-1$ index records can fit into a single block.

Ideally, then, you should structure the external search tree so that every internal node has $m$ children, where $m$ is chosen as just described, and all leaves are at the same level, as is the case with full trees and 2-3 trees. For example, Figure 21-12a shows a full tree whose internal nodes each have five children. Although this search tree has the minimum possible height, its balance is too difficult to maintain in the face of insertions and removals. As a consequence, you must make a compromise. You can still insist that all the leaves of the search tree be at the same level-that is, that the tree be bal-anced-but you must allow each internal node to have between $m$ and [ $m / 2$ ] +1 children. (The [ ] notation means greatest integer in. Thus, [5/2] is 2 , for example.)

FIGURE 21-12 (a) A full tree whose internal nodes have five children; (b) the format of a single node

(b)



This type of search tree is known as a B-tree of degree $\boldsymbol{m}$ and has the following characteristics:

- All leaves are at the same level.
- Each node contains between $m-1$ and [ $m / 2$ ] records, and each internal node has one more child than it has records. An exception to this rule is that the root of the tree can contain as few as one record and can have as few as two children. This exception is necessitated by the insertion and removal algorithms described next.

A 2-3 tree is a B-tree of degree 3. Furthermore, the manner in which the B-tree insertion and removal algorithms maintain the structure of the tree is a direct generalization of the 2-3 tree's strategy of splitting and merging nodes. These algorithms are illustrated next by means of an example. Assume that the index file is organized into a B-tree of degree 5-that is, 5 is the maximum and 3 is the minimum number of children that an internal node-other than the root-in the tree can have. (Typically, a B-tree will be of a higher degree, but the diagrams would get out of hand!)

Insertion into a B-tree. To insert a data record with search key 55 into the tree shown in Figure 21-13, you take the following steps:

1. Insert the data record into the data file. First you find block $p$ in the data file into which you can insert the new record. As was true with the external hashing implementation, block $p$ is either any block with a vacant slot or a new block.
2. Insert a corresponding index record into the index file. You now must insert the index record $<55, \mathrm{p}>$ into the index file, which is a B-tree of degree 5 . The first step is to locate the leaf of the tree in which this index record belongs by determining where the search for 55 would terminate.

Suppose that this is the leaf $L$ shown in Figure 21-14a. Conceptually, you insert the new index record into $L$, causing it to contain five records (Figure 21-14b). Since a node can contain only four records, you must split $L$ into $L_{1}$ and $L_{2}$. With an action analogous to the splitting of a node in a 2-3 tree, $L_{1}$ gets the two records with the smallest key values, $L_{2}$ gets the two records with the largest key values, and the record with the middle key value (56) is moved up to the parent $P$. (See Figure 21-14c.)

B-tree of degree $m$

A 2-3 tree is a B-tree of degree 3

FIGURE 21-14 (a through d) The steps for inserting 55 ; (e) splitting the root


In this example, $P$ now has six children and five records, so it must be split into $P_{1}$ and $P_{2}$. The record with the middle key value (56) is moved up to $P$ 's parent, $Q$. Then $P$ 's children must be distributed appropriately, as happens with a 2-3 tree when an internal node is split. (See Figure 21-14d.)

At this point the insertion is complete, as $P$ 's parent $Q$ now contains only three records and has only four children. In general, though, an insertion might cause splitting to propagate all the way up to the root (Figure 21-14e). If the root must be split, the new root will contain only one record and have only two children - the definition of a B-tree allows for this eventuality.

Removal from a B-tree. To remove a data record with a given search key from a B-tree, you take the following steps:

1. Locate the index record in the index file. You use the search algorithm to locate the index record with the desired key value. If this record is not already in a leaf, you swap
it with its inorder successor. Suppose that the leaf $L$ shown in Figure 21-15a contains the index record with the desired key value, 73. After noting the value $p$ of the pointer in this index record (you will need $p$ in step 2 to remove the data record), you remove the index record from $L$ (Figure 21-15b). Because $L$ now contains only one value (recall that a node must contain at least two values), and since $L$ 's siblings cannot spare a value, you merge $L$ with one of the siblings and bring down a record from the parent $P$ (Figure $21-15 \mathrm{c}$ ). Notice that this step is analogous to the merge step for a 2-3 tree. However, $P$ now has only one value and two children, and since its siblings cannot spare a record and child, you must merge $P$ with its sibling $P_{1}$ and bring a record down from $P$ 's parent, $Q$. Because $P$ is an internal node, its children must be adopted by $P_{1}$. (See Figure 21-15d.)

After this merge, $P$ 's parent $Q$ is left with only two children and one record. In this case, however, $Q$ 's sibling $Q_{1}$ can spare a record and a child, so you redistribute children and records among $Q_{1}, Q$ and the parent $S$ to complete the removal. (See Figure 21-15e.) If a removal ever propagates all the way up to the root, leaving it with only one record and only two children, you are finished because the definition of a B-tree allows this situation. If a future removal causes the root to have a single child and no records, you remove the root so that the tree's height decreases by 1 , as Figure 21-15f illustrates. The removal of the index record is complete, and you now must remove the data record.

FIGURE 21-15 (a through e) The steps for removing 73; (f) removing the root

(continues)

Accessing only the search key of each record, but not the data file

FIGURE 21-15 (a through e) The steps for removing 73 ; (f) removing the root (continued)

2. Remove the data record from the data file. Prior to removing the index record, you noted the value $p$ of its pointer. Block $p$ of the data file contains the data record to be removed. Thus, you simply access block p, remove the data record, and write the block back to the file. The high-level pseudocode for the insertion and removal algorithms parallels that of the 2-3 tree and is left as an exercise.

### 21.3.4 Traversals

Now consider the operation traverse in sorted order, which is one of the operations that hashing does not support at all efficiently. Often an application requires only that the traversal display the search keys of the records. If such is the case, the B-tree implementation can efficiently support the operation, because you do not have to access the data file. You can visit the search keys in sorted order by using an inorder traversal of the B-tree, as follows:
// Traverses in sorted order an index file that is organized as a B-tree of degree m.
// blockNum is the block number of the root of the B-tree in the index file.

```
traverse(blockNum: integer, m: integer)
if (blockNum != -1)
{
    // Read the root into internal array buf
    buf.readBlock(indexFile, blockNum)
    // Traverse the children
    // Traverse S0
    Let p be the block number of the Oth child of buf
    traverse(p, m)
    for (i = 1 through m-1)
    {
        Display key Ki of buf
        // Traverse Si
        Let p be the block number of the ith child of buf
        traverse(p, m)
    }
}
```

This traversal accomplishes the task with the minimum possible number of block accesses because each block of the index file is read only once. This algorithm, however, assumes that enough internal memory is available for a recursive stack of $h$ blocks, where $h$ is the height of the tree. In many situations this assumption is reasonable-for example, a 255 -degree B-tree that indexes a file of 16 million data records has a height of no more than 3 . When internal memory cannot accommodate $h$ blocks, you must use a different algorithm. (See Exercise 12.)

If the traversal must display the entire data record (and not just the search key), the B-tree implementation is less attractive. In this case, as you traverse the B-tree, you must access the appropriate

Accessing the entire data record block of the data file. The traversal becomes

```
// Traverses in sorted order a data file that is indexed with a B-tree of degree m.
// blockNum is the block number of the root of the B-tree.
traverse(blockNum: integer, m: integer)
    if (blockNum !=-l)
{
    // Read the root into internal array buf
    buf.readBlock(indexFile, blockNum)
    // Traverse S0
    Let p be the block number of the 0th child of buf
    traverse(p, m)
    for (i = 1 through m-1)
    {
            Let p_i be the pointer in the ith index record of buf
            data.readBlock(dataFile, p_i)
            Extract from data the data record whose search key equals Ki
            Display the data record
            // Traverse Si
            Let p be block number of the ith child of buf
            traverse(p, m)
        }
    }
```

Sorted-order traversal of a data file indexed with a B-tree

Generally, the previous traversal is unacceptable

Multiple index files allow multiple data organizations

This traversal requires you to read a block of the data file before you display each data record; that is, the number of data-file block accesses is equal to the number of data records. In general, such a large number of block accesses would not be acceptable. If you must perform this type of traversal frequently, you probably would modify the B-tree scheme so that the data file itself was kept nearly sorted.

### 21.3.5 Multiple Indexing

Before concluding the discussion of external implementations, let's consider the multiple indexing of a data file. Programming Problem 8 in Chapter 18 asked you to support multiple organizations for data stored in internal memory. Such a problem is also common for data stored externally. For example, suppose that a data file contains a collection of employee records on which you need to perform two types of retrievals:

```
// Retrieves the item whose search key contains the name aName.
retrieveN(aName: NameType): ItemType
// Retrieves the item whose search key contains the
// Social Security Number ssn.
retrieveS(ssn: SSNType): ItemType
```

One solution to this problem is to maintain two independent index files to the data file. For example, you could have one index file that contains index records of the form <name, pointer> and a second index file that contains index records of the form <socSec, pointer>. These index files could both be hashed, could both be B-trees, or could be one of each, as Figure 21-16 indicates. The choice would depend on the operations you wanted to perform with each search key. Similarly, if an application required extremely fast retrievals on socSec and also required operations such as traverse in sorted socSec order and range queries on socSec, it might be reasonable to have two socSec index files-one hashed, the other a B-tree.

FIGURE 21-16 Multiple index files


Although you can perform each retrieval operation by using only one of the indexes (that is, use the name index for retrieveN and the socSec index for retrieveS), insertion and removal operations must update both indexes. For example, the remove-by-name operation removeN(Jones) requires the following steps:

1. Search the name index file for Jones and remove the index record.
2. Remove the appropriate data record from the data file, noting the socSec value ssn of this record.
3. Search the socSec index file for ssn and remove this index record.

In general, the price paid for multiple indexing is more storage space and an additional overhead for updating each index whenever you modify the data file.

This chapter has presented at a very high level the basic principles of managing data in external storage. The details of implementing the algorithms depend heavily on your specific computing system. Particular situations often mandate either variations of the techniques described here or completely different approaches. In future courses and work experience, you will undoubtedly learn much more about these techniques.

A removal by name must update both indexes

Question 1 Consider two files of 1,600 employee records each. The records in each file are organized into sixteen 100 -record blocks. One file is sequential access and the other is direct access. Describe how you would append one record to the end of each file.

Question 2 Trace externalMergesort with an external file of 16 blocks. Assume that the arrays in1, in2, and out are each one block long. List the calls to the various functions in the order in which they occur.
Question 3 Trace the retrieval algorithm for an indexed external file when the search key is less than all keys in the index. Assume that the index file stores the index records sequentially, sorted by their search keys, and contains 20 blocks of 50 records each. Also assume that the data file contains 100 blocks, and that each block contains 10 employee records. List the calls to the various functions in the order in which they occur.
Question 4 Repeat Checkpoint Question 3, but this time assume that the search key equals the key in record 26 of block 12 of the index. Also assume that record 26 of the index points to block 98 of the data file.

## Summary

1. An external file is partitioned into blocks. Each block typically contains many data records, and a block is generally the smallest unit of transfer between internal and external memory. That is, to access a record, you must access the block that contains it.
2. You can access the $i^{\text {th }}$ block of a direct access file without accessing the blocks that precede it. In this sense direct access files resemble arrays.
3. Before you can process (for example, inspect or update) a record, you must read it from an external file into internal memory. Once you modify a record, you must write it back to the file.
4. Block accesses are typically quite slow when compared to other computer operations. Therefore, you must carefully organize a file so that you can perform tasks by using only a few block accesses. Otherwise, response time can be very poor.
5. You can modify the merge sort algorithm, presented in Chapter 11, so that it can sort an external file of records without requiring all of the records to be in internal memory at one time.
6. An index to a data file is a file that contains an index record for each record in the data file. An index record contains both the search key of the corresponding data record and the number of the block in the data file that contains the data record.
7. If a record is inserted into or removed from a data file, you must make the corresponding change to the index file. If a data file has more than one index file, you must update each index file. Thus, multiple indexing has an overhead.
8. You can organize an index file by using either hashing or a B-tree. These schemes allow you to perform the basic dictionary operations by using only a few block accesses.
9. Although external hashing generally permits retrievals, insertions, and removals to be performed more quickly than does a B-tree, it does not support such operations as sorted traversals or range queries. This deficiency is one motivation for multiple indexing.
10. You can have several index files for the same data file. Such multiple indexing allows you to perform different types of operations efficiently, such as retrieval by name and retrieval by Social Security number.

## Exercises

1. Assuming the existence of readBlock and writeBlock functions, write a pseudocode program for shifting data to make a gap at some specified location of a sorted file. Pay particular attention to the details of shifting the last item out of one block and into the first position of the next block. You can assume that the last record of the file is in record lastRec of block lastBlock and that lastBlock is not full. (Note that this assumption permits shifting without allocating a new block to the file.)
2. The problem of managing the blocks of an external data file indexed by either a B-tree or an external hashing scheme is similar to that of managing memory for internal structures. When an external structure such as a data file needs more memory (for example, to insert a new record), it gets a new block from a free list that the system manages. That is, if the file contains $n$ blocks, the system can allocate to it an $(n+1)^{\text {th }}$ block. When the file no longer needs a block, you can deallocate it and return it to the system.

The complication in the management of external storage is that a block allocated to a file may have available space interspersed with data. For example, after you have removed a record from the middle of a data file, the block that contained that record will have space available for at least one record. Therefore, you must be able to keep track of blocks that have space available for one or more records as well as recognize when blocks are completely empty (so that you can return them to the system).

Assuming the existence of allocateBlock and returnBlock functions that get empty blocks from and return empty blocks to the system, write pseudocode implementations of the following external memory-management functions:

```
// Determines the block number (blockNum) and record number (recNum) of an available slot in file
// dataFile. A new block is allocated to the file from the system if necessary.
getSlot(dataFile: File, blockNum: integer,
    recNum: integer)
// Makes record recNum in block blockNum of file dataFile available.
// The block is returned to the system if it becomes empty.
freeSlot(dataFile: File, blockNum: integer,
    recNum: integer)
```

What data structure is appropriate to support these operations? You may assume that you can distinguish slots of a block that do not contain a record from those that do. You can make this distinction either by having a convention for null values within a record or by adding an empty/full flag.
3. Describe pseudocode algorithms for insertion into and removal from a dictionary implemented externally with a hashed index file.
4. Execute the following sequence of operations on an initially empty dictionary book that is implemented as a B-tree of degree 5 . Note that insertion into an empty B-tree will create a single node that contains the inserted item.

```
book.add(10)
book.add(100)
book.add(30)
book.add(80)
book.add(50)
book.remove(10)
book.add(60)
book.add(70)
book.add(40)
book.remove(80)
book.add(90)
book.add(20)
book.remove(30)
book.remove(70)
```

5. Given a B-tree of degree 5 and a height of 3 ,
a. What is the maximum number of nodes (including the root)?
b. What is the maximum number of records that can be stored?
6. Given the B-tree of degree 7 in Figure 21-17, draw the B-tree that results after the insertion of $m, o, y, r, c, i, k$, $w$, and $h$.
7. Given the B-tree of degree 7 in Figure 21-18, draw the B-tree that results after the removal of $s, t, p, m, k$, and $e$.
8. Describe a pseudocode algorithm for finding an item's inorder successor in an external B-tree.
9. Describe pseudocode algorithms for insertion into and removal from an ADT dictionary implemented with an index file organized as a B-tree.
10. Write a rangeQuery function for a B-tree in pseudocode. (See Exercise 3 of Chapter 19.) Assume that only the key values are needed (as opposed to the entire data record).

FIGURE 21-17 A B-tree for Exercise 6


## FIGURE 21-18 A B-tree for Exercise 7


11. Integrate calls to the appropriate memory-management functions (see Exercise 2) into the pseudocode for add and remove under both the B-tree and hashing schemes. (See Exercises 3 and 9.)
12. The B-tree traversal algorithm presented in this chapter assumes that internal memory is large enough to accommodate the recursive stack that contains up to $h$ blocks, where $h$ is the height of the B-tree. If you are in an environment where this assumption is not true, modify the traversal algorithm so that the recursive stack contains block numbers rather than the actual blocks. How many block accesses does your algorithm have to perform?
a. Write pseudocode B-tree implementations of traversals and range queries that need to access entire data records, not simply the search keys. How many block accesses do your functions require?
b. To reduce the number of block accesses required by these operations, various modifications of the basic B-tree structure are frequently used. The central idea behind such structures is to keep the data file itself sorted. First, assume that you can keep the data file in sequential sorted order-that is, the records are sorted within each block and the records in $B_{i-1}$ are less than the records in $B_{i}$ for $i=2$, 3, and so on to the number of blocks in the file. Rewrite your implementations of the traversal and range-query operations to take advantage of this fact. How many block accesses do these operations now require?
c. Because it is too inefficient to maintain a sequentially sorted data file in the face of frequent insertions and removals, a compromise scheme is often employed. One such possible compromise is as follows. If a data record belongs in block $B$ and $B$ is full, a new block is allocated and linked to $B$, allowing the new record to be inserted into its proper sorted location. The difficulty is that you must now view each index record in the B-tree as indicating the first of possibly several blocks in a chain of blocks that might contain the corresponding data record. Rewrite the add, remove, getItem, traverse, and rangeQuery operations in terms of this implementation. What is the effect on their efficiency?
14. Write an iterative (nonrecursive) version of the internal merge sort, as given in Chapter 11, that is based on the external version that this chapter describes. That is, merge sorted runs that double in size at each pass of the array.

## Programming Problems

1. a. Implement in $\mathrm{C}++$ the externalMergesort algorithm given in Section 21.2 by using the functions seekg and seekp. Assume that the file to be sorted is a file of type int and that each block contains one integer. Further assume that the file contains $2^{n}$ integers for some integer $n$.
b. Now assume that each block contains many integers. Write C++ functions that simulate readBlock and writeBlock. Implement externa7Mergesort by using these functions.
c. Extend your implementation of externa7Mergesort by removing the restriction that the file contains $2^{n}$ blocks.
2. Implement the ADT dictionary by using a sorted index file, as described in Section 21.3.1.
3. Implement an ADT dictionary that uses a sorted index file using the STL map container.
4. Implement a simple dictionary application, such as the one described in Exercise 4 of Chapter 18, using the external dictionary of Programming Problem 3.
5. Implement the ADT dictionary by using a hashed index file, as described in Section 21.3.2.
6. Implement the ADT dictionary by using a B-tree, as described in Section 21.3.3.
7. Repeat Programming Problem 8 of Chapter 18, using an external dictionary.

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## The Standard Template Library

## C+ + <br> Interlude

## Contents

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This interlude gives an overview of some useful methods available with each of the STL containers and in the STL algorithms library. We do not attempt to describe the entire library.

## C7.1 STL Containers

Throughout this text we have been designing and implementing containers that hold collections of objects. Each ADT was inspired by a problem to solve and was designed to efficiently solve that specific problem. While each ADT can be used in a variety of situations, we need to consider the strengths and weaknesses of each one when choosing which to use.

C++ comes with a library of containers that implement many of the more commonly used ADTs. These classes are defined in the Standard Template Library, or STL. Many of the ADTs that are presented in this text have a corresponding class in the STL. For example, the STL defines a stack class that is similar to the class LinkedStack presented in Chapter 7. Just like the ADTs we designed, each container in the STL has strengths and weaknesses that you need to consider when deciding which one to use.

You may wonder why we spent so much time developing ADTs in this text if they are already provided in the STL. There are many reasons for doing so; here are just a few:

- Developing simple ADTs provides a foundation for learning to develop other ADTs, especially ones that are not in the STL.
- STL containers are not part of a class hierarchy and so cannot take advantage of polymorphism in the same way our ADTs can.
- You might find yourself working in a language that does not provide any predefined ADTs. You need to have the ability to develop ADTs on your own, and hence you must understand the process.
- If the ADTs defined by the language you are using are insufficient, you may need to develop your own or enhance existing ones.

The STL provides support for predefined ADTs through the use of three basic items: containers, iterators, and algorithms. Containers are objects, such as a list, that hold other objects. Iterators provide a way to cycle through the contents of a container and are described in C++ Interlude 6. Algorithms, such as a sorting algorithm, act on containers. STL containers are implemented as templates, similar to the ADTs we implemented in this text.

You already are familiar with one STL class-the class vector-which we used in Chapter 3 to implement the ADT bag method toVector. Moreover, many introductory C++ courses use the STL class vector instead of arrays as a simple way to store multiple items of a similar type

There are three types of containers in the STL:

- Container adapters provide efficient, restricted, position-based access to the collection. They are considered adapters because they use other STL containers to implement their operations.
- Sequence containers provide efficient sequential access to the collection. The containers in this group provide various trade-offs in efficiency for insertions and removals.
- Associative containers provide efficient key-based access to the collection. Each of these containers organizes and accesses the search keys to provide efficient groups of operations.

Each container type implements similar operations, but with a different emphasis in efficiency. You'll find that many of these operations have corresponding operations in the ADTs that we defined in this text, but the STL names might be different.

Two operations that all STL containers define are empty and size. These correspond to isEmpty and getLength, respectively, in this book's ADTs. For sequence and associative containers, the container size is the number of entries it can hold. If the container is empty, not only does it not have any entries, but it cannot hold any. To add a new entry to the container, you must either use one of the methods that increases the container size during the add operation or, in the case of sequence containers, resize the container using the resize method.

## Operations Common to all STL Containers

| Operation | Description | Efficiency |
| :---: | :---: | :---: |
| Constructor | Creates a container; sequence containers allow the user to specify an initial size. | Varies |
| Destructor | Destroys all entries in the container, and then destroys the container. | $\mathrm{O}(n)$ |
| operator= | Assigns entries in the container on the right-hand side to the one on the left-hand side. | $\mathrm{O}(n)$ |
| bool empty () | Returns true if the container is empty (size is 0 ). | $\mathrm{O}(1)$ |
| uint ${ }^{1}$ size() | Returns the number of locations in the container. | $\mathrm{O}(1)$ |

Operations that remove entries from a collection throw an exception when they fail.
To use one of the STL classes, you must include the appropriate header file, which has the same name as the STL class. So to use the class stack, for example, your code must have the following statement:

```
#include <stack>;
```

To instantiate an STL container, you must supply the type of item it will hold, just as we instantiated our ADTs. For example, the following statement creates a vector that can hold 10 strings:

```
vector<string> cities(10); // Create a vector to hold 10 strings
```

All STL containers are part of the std: : namespace. Therefore, if you do not use the statement using namespace std;
in your code, you must precede any reference to the container with std: : as shown here;

```
std::vector<std::string> cities(10); // Create a vector to hold 10 strings
```

In our ADTs, we used ItemType to represent the type of the item stored in the collection. The STL header files use value_type or simply $T$ to represent the item in a collection. If the collection is similar to our ADT dictionary, with key and item types, the STL uses key_type and mapped_type respectively.

## C7.1.1 STL Container Adapters

The STL container adapter classes are stack, queue and priority_queue. These correspond to the ADTs stack, queue, and priority queue and are optimized so that the operations listed below perform at $\mathrm{O}(1)$. These containers grow to accommodate new entries as the entries are added. When entries are removed, the size of the container decreases. Thus, each of these containers has a size exactly equal to the number of entries in it currently.

STL stack Operations

Operation
value_type\& top()
void push(value_type\& item)
void pop()

## DESCRIPTION

Returns a reference to the top entry on the stack.
Pushes item onto the top of the stack.
Removes the top entry from the stack.

## STL queue Operations

Operation
value_type\& front()
value_type\& back()
void push(value_type\& item)
void pop()

## Description

Returns a reference to the front entry in the queue.
Returns a reference to the last (back) entry in the queue.
Adds item onto the back of the queue.
Removes the front entry from the queue.

## STL PRIORITY_QUEUE OPERATIONS

Operation
value_type\& top()
void push(value_type\& item)
void pop()

## DESCRIPTION

Returns a reference to the front entry in the priority queue.
Inserts item into the priority queue.
Removes the front entry in the priority queue.

Listing C7-1 provides an example of how to use the STL stack.

## LISTING C7-1 Example use of the STL stack

```
#include <iostream>
#include <stack>
using namespace std;
int main()
{
    stack<int> aStack;
    // Right now, the stack is empty
    if (aStack.empty())
        cout << "The stack is empty." << end7;
    for (int j = 0; j < 5; j++)
        aStack.push(j); // Places items on top of stack
    while (!aStack.empty())
    {
            cout << aStack.top() << " ";
            aStack.pop();
    } // end while
    return 0;
} // end main
```

Output
The stack is empty.
43210

## C7.1.2 Sequence Containers

The elements in a sequence container are ordered in a linear sequence and can be accessed, inserted, and removed by position. Similar to the arrangement in traditional arrays, the positions in sequence containers are numbered from 0 to size - 1 . The front of these containers is at position 0 , and the back is at position size - 1 .

The STL sequence containers are: array, vector, deque, list, and forward_list. Let's consider each of these containers.

The sequence container array. As part of the recent C11 standard, the STL array provides a fixedsize container that is implemented as a traditional array, so unlike other STL containers, an STL array container cannot grow or shrink in size. You can access individual elements in an STL array
just as you access elements in a traditional array, by using the operator [ ]. An advantage of an STL array over traditional $\mathrm{C}++$ arrays is the method at that allows for range checking. For example, a traditional array declared as

```
int myArray[5];
```

permits the programmer to access elements beyond the memory allocated to the array by using an expression such as myArray [24]. This leads to unpredictable results and is a common source of errors in a program.

An STL array declared as

```
array<int> mySTLArray(5);
```

would have the same issue if the programmer used the expression mySTLArray [24]. In situations where accessing elements outside the bounds of the array is possible, it is better to use an expression such as mySTLArray.at(24). This expression will throw an out_of_bounds exception that you can catch, enabling you to fix the error and allow the program to continue execution safely.

Two other important features of the STL array are its size method-enabling you to determine how many locations are in the array-and the availability of iterators for array traversal, as discussed in C++ Interlude 6.

| STL array Operations |  |  |
| :---: | :---: | :---: |
| Operation | Description | Efficiency |
| value_type\& front() | Returns a reference to the front entry in the array. | $\mathrm{O}(1)$ |
| value_type\& back() | Returns a reference to the last (back) entry in the array. | $\mathrm{O}(1)$ |
| value_type\& at(size_type n) | Behaves the same as [ ] , but performs a bounds check. | $\mathrm{O}(1)$ |
| void fill (const value_type\& val) | Fills array elements with val. | $\mathrm{O}(n)$ |
| iterator begin() | Returns an iterator to the first element in the container. | $\mathrm{O}(1)$ |
| iterator end() | Returns an iterator to the last element in the container. | $\mathrm{O}(1)$ |
| reverse_iterator rbegin() | Returns a reverse iterator to the last element in the container. | $\mathrm{O}(1)$ |
| reverse_iterator rend() | Returns a reverse iterator to the first element in the container. | $\mathrm{O}(1)$ |

Other STL sequence containers have many operations in common but are optimized differently. For example, vector and deque permit efficient access to specific locations using the [ ] operator, but list is better for sorting. Though these containers can be constructed with an initial size, they also grow and shrink as entries are added and removed. If you realize that you need to add a significant number of entries to the container, a resize method lets you adjust the number of elements.

Operations Common to STL Sequence Containers

```
Operation
value_type\& front()
value_type\& back()
```


## DESCRIPTION

Returns a reference to the front entry in the container.
Efficiency

Returns a reference to the last (back) entry in the container.

```
void push_back(value_type& item)
void pop_back(value_type& item)
void resize (uint newSize)
void clear()
voidinsert(uint position,
    value_type& item)
void insert(iterator itPosition,
    value_type& item)
void erase(uint position)
void erase(iterator itPosition)
iterator begin()
iterator end()
reverse_iterator rbegin()
reverse_iterator rend()
```

Returns an iterator to the last element of the container. ..... $\mathrm{O}(1)$
Returns a reverse iterator to the last element of the ..... $\mathrm{O}(1)$container.
Returns a reverse iterator to the first element of the ..... $\mathrm{O}(1)$
Adds item onto the back of the container(grows container).Removes the back entry from the container(shrinks container).
Resizes the container to have newSize elements. ..... $\mathrm{O}(n)$
Destroys all entries in the container, and resizes the ..... $\mathrm{O}(n)$container to a size of 0 .
Inserts item at position and grows the container. ..... Varies
Inserts item at itPosition and grows the container. ..... Varies
Removes element position from the container Varies(shrinks container).
Removes element itPosition from container ..... Varies(shrinks container).
Returns an iterator to the first element of the container. ..... O(1)

The STL class vector. The STL class vector is implemented using dynamic arrays so that it can increase in size as necessary. You access individual elements in a vector just as you do for an STL array by using either [ ] or at. Since vectors are stored as an array in a contiguous sequence of memory locations, vectors are good for accessing elements by either position or iteration, and for adding elements to or removing elements from the back.

## Additional STL vector Operation

## Efficiency

$\mathrm{O}(1)$

The STL container deque. The STL container deque allows access to individual elements using [ ] and enables the efficient addition and removal of entries at both its front and back. Unlike our ADT deque, the STL deque permits insertion (insert) and removal (erase) from the interior of the container.

The STL deque does not use contiguous memory locations, but stores sequences of its entries in multiple blocks of memory that it tracks internally. While this makes the STL deque more complex to implement, it enables it to grow in size more efficiently, since additional memory can be allocated without the need to copy all existing entries into the new elements.

## Additional STL deque Operations

| value_type\& at(size_type n) | Behaves the same as [ ], but performs a bounds check. | $\mathrm{O}(1)$ |
| :--- | :--- | :--- |
| void push_front(value_type\& item) | Adds item onto the front of the deque (grows container). | $\mathrm{O}(1)$ |
| void pop_front (value_type\& item) | Removes the front entry from the deque (shrinks container). | $\mathrm{O}(1)$ |

The STL class 1ist. The STL class list is typically implemented using some form of a doubly linked chain. It is optimized for the insertion and removal of entries but does not provide subscripting, because element access would be inefficient. Bidirectional iterators are provided to access list entries. The STL 7 ist can efficiently insert and remove entries at any location in the container. This performance is improved if an iterator, either forward or reverse, is used to determine position. The nature of the STL list is well suited for sorting and merging operations, and it has methods that implement those algorithms. The STL list does not provide direct access to elements by position. Access to elements other than the front or back must be done using an iterator.

The STL container forward_1 ist. Another C11 STL container is forward_1ist. It is similar in efficiency to the STL list. Since the STL forward_list is implemented as a singly linked chain, it performs better with operations using forward iterators and requires less storage than the STL list. The STL forward_1ist offers the same functionality for sorting, merging, and element access as the STL 1ist.

## Additional STL list and forward_list Operations

| Operation | Description | Efficiency |
| :---: | :---: | :---: |
| void push_front(value_type\& item) | Adds item onto the front of the list (grows container). | $\mathrm{O}(1)$ |
| void pop_front(value_type\& item) | Removes the front entry from the list (shrinks container). | $\mathrm{O}(1)$ |
| void remove(value_type\& val) | Removes all items from the list that are equal to val. | $\mathrm{O}(n)$ |
| void sort() | Sorts the list in ascending order. | $\mathrm{O}(n)$ |
| void merge(list<value_type>\& rhs) | Merges rhs with the current list by placing entries in their correct sorted order, resulting in a sorted list of all entries. | $\mathrm{O}(n)$ |
| voidslice(iterator position, list<value_type>\& rhs) | Inserts entries in rhs into the current list by placing them at position; other forms specify the number of entries. | $\mathrm{O}(n)$ |
| void reverse() | Reverses the order of entries in the list. | $\mathrm{O}(n)$ |

Listing C7-2 contains an example that uses the STL 1ist to maintain a list of groceries. Each entry is inserted in the front of the list, and then the list is sorted.

## LISTING C7-2 Example of using the STL 1ist

```
#include <iostream>
#include <string>
```

```
#include <list>
using namespace std;
int main()
{
    list<string> groceryList; // Create an empty list
    list<string>::iterator myPosition = groceryList.begin();
    groceryList.insert(myPosition, "app1es");
    groceryList.insert(myPosition, "bread");
    groceryList.insert(myPosition, "juice");
    groceryList.insert(myPosition, "carrots");
    cout << "Number of items on my grocery list: "
            << groceryList.size() << endl;
    groceryList.sort();
    cout << "Items are:" << endl;
    myPosition = groceryList.begin();
    while (myPosition != groceryList.end())
    {
        cout << *myPosition << endl;
        ++myPosition;
    } // end while
} // end main
```


## Output

Number of items on my grocery 1ist: 4
Items are:
apples
bread
carrots
juice

## C7.1.3 Associative Containers

STL associative containers store their entries according to search keys, much as our value-based ADT dictionary does. There are two basic associative containers-set and map-with several variations of each. The STL set uses the entry to be stored as the key and uses a binary tree for its structure. The STL map uses a separate entry as the key, similar to the technique used by our ADT dictionary. The STL set and STL map both expect unique keys, and they store entries in ascending order based on these keys.

The STL multiset and STL multimap are variations of the STL set and STL map, respectively, that permit multiple occurrences of the search keys. The STL containers unordered_set, unordered_map, unordered_multiset, and unordered_multimap are hash-based implementations of the corresponding ordered STL containers. ${ }^{2}$

Since they are hash based, the unordered associative containers are faster for direct access of an entry based on its search key. The ordered associative containers are better for iterator-based access. The STL map and STL unordered_map containers permit direct access to entries using the [ ] operator.

[^18]
## Operations Common to the STL set and multiset

void clear()
void insert(value_type\& item)
void erase(value_type\& item)
void erase(iterator\& position)
iterator find(value_type\& item)
uint count(value_type\& item)
iterator lower_bound(value_type\& item)
iterator upper _bound(value_type\& item)
iterator begin()
iterator end()
reverse_iterator rbegin()
reverse_iterator rend()

## Description

Destroys all entries in the container, and resizes Efficiency the container to the size of 0 .

Inserts item and grows the container.
$\mathrm{O}(\log n)$
Removes all entries matching item from the $\mathrm{O}(\log n)$ container (shrinks container).
Removes the entry at position from the container (shrinks container).
Returns the iterator referencing item.
Counts the occurrences of item in the container. For $\mathrm{O}(\log n)$ set this is at most 1 ; it can vary for multiset.
Returns an iterator referencing the first element $\mathrm{O}(\log n)$ not less than item.
Returns an iterator referencing the first element greater than item.
Returns an iterator to the first element of the container.
Returns an iterator to the last element of the container.
Returns a reverse iterator to the last element of the container.

Returns a reverse iterator to the first element of the container.

Generally, the operations of the STL map and its variations have the same behavior as those of the STL set, though the parameters are different since the STL map uses a key for each mapped value stored in the container. For insertion operations, the STL map expects a pair_type that contains both the key and the entry value mapped to that key.

For example, the following code declares a map with a key of type char and a mapped type of int. It then stores several values in the map and uses an iterator to display them:

```
#include <iostream>
#include <map>
using namespace std;
int main ()
{
    map<char, int> myMap;
    myMap.insert(pair<char, int>('A', 100));
    myMap.insert(pair<char, int>('B', 200));
    myMap.insert(pair<char, int>('C', 300));
    map<char, int>::iterator mapIterator;
    for (mapIterator = myMap.begin(); mapIterator != myMap.end(); ++mapIterator)
        cout << "Key: " << mapIterator->first;
        cout << " Value: " << mapIterator->second << endl;
    return 0;
} // end main
```

As an alternative to the insert statement, you can use the [ ] operator:

```
myMap['A'] = 100;
myMap['B'] = 200;
myMap['C'] = 300;
```

Other operations for the STL map use a parameter of key_type or an iterator to identify the entry to act upon.

Operations Common to the STL map and multimap

| Operation | Description | Efficiency |
| :---: | :---: | :---: |
| void clear() | Destroys all entries in the container and changes its size to 0 . | $\mathrm{O}(n)$ |
| void insert(pair_type\& item) | Inserts item (grows container). | $\mathrm{O}(\log n)$ |
| uint erase(key_type\& item) | Removes all entries matching item from the container (shrinks container). | $\mathrm{O}(\log n)$ |
| void erase(iterator\& position) | Removes the entry at position from the container (shrinks container). | $\mathrm{O}(1)$ |
| iterator find(key_type\& item) | Returns an iterator referencing item. | $\mathrm{O}(\log n)$ |
| uint count (key_type\& item) | Counts the occurrences of item in the container. <br> For set, this count is at most 1 ; it can vary for multiset. | $\mathrm{O}(\log n)$ |
| iterator lower_bound (key_type\& item) | Returns an iterator referencing the first entry not less than item. | $\mathrm{O}(\log n)$ |
| iterator upper _bound(key_type\& item) | Returns an iterator referencing the first entry greater than item. | $\mathrm{O}(\log n)$ |
| iterator begin() | Returns an iterator to the first entry of the container. | $\mathrm{O}(1)$ |
| iterator end() | Returns an iterator to the last entry of the container. | $\mathrm{O}(1)$ |
| reverse_iterator rbegin() | Returns a reverse iterator to the last entry of the container. | $\mathrm{O}(1)$ |
| reverse_iterator rend() | Returns a reverse iterator to the first entry of the container. | $\mathrm{O}(1)$ |

As mentioned earlier, the unordered variations of the associative containers are hash-basedthey use a hashing function to determine the position of an entry inserted into the container. When you instantiate an unordered associative container, you can use the default hash function for the type of data you are inserting or you can specify your own. Most often, the default hash function will produce acceptable results, but if you are testing your own hashing algorithm, you have the option to pass it to the constructor of an unordered associative container.

Since hash functions are difficult to write, it is often helpful to take advantage of an STL provided hash. The current C++ standard, C11, does not provide a stand-alone hash function. An unordered associative container, typically an unordered_map, must be instantiated first. Then, the hash function for that container can be used to hash any key of the correct type. Section 18.4.1 in Chapter 18 contains a Programming Tip that demonstrates one way to use this hash within the getHashIndex function.

Let's look at a slightly different way of implementing the getHashIndex function in the HashedDictionary class discussed in Chapter 18. We need to take two steps when using the unorder_map hash function within getHashIndex: Instantiate an unordered_map object, and create a variable to represent the hash function. We can accomplish both of these tasks within the header file of HashedDictionary.

In the private section of our header file, we instantiate our mapper object:

```
std::tr1::unordered_map<KeyType, ItemType> mapper;
```

The sole purpose of the dummy object mapper is to give us access to its hash function.
We then call the method hash_function and store the reference it returns in the variable hashFunction:

```
typename std::tr1::unordered_map<KeyType, ItemType>::hasher hashFunction =
mapper.hash_function();
```

The returned reference is to the hash function used to hash our KeyType. Note that the variable hashFunction is of type hasher. In the C11 standard, hashing functions are functions of this type. They have a single parameter of type KeyType and return an unsigned integer that represents the hash value of the key. The keyword typename is required so that the compiler understands that this statement is based on the template parameters.

Because hashFunction is defined and initialized in the header file, it is available to any method in the HashedDictionary class. When it is used, we need to use the modulo operator, since hashFunction returns an unsigned integer in the range 0 to UINT_MAX-the maximum unsigned integer. After the modulo operator is applied, we can cast the result to be the type of index we need for our dictionary. Thus, we have the following statement:

```
int hashIndex = static_cast<int>(hashFunction(searchKey) % hashTableSize);
```

The code in Listing C7-3 gives an example of how to use a hash function when it is not part of an ADT. This example creates a hash function for string objects. Since the key type of the map determines the hash function, the type of the mapped value is not important; hence we use int.

## LISTING C7-3 Alternative definition of a hashing function

```
#include <iostream>
```

\#include <string>
\#include <tr1/unordered_map> // Note header name since this is in tr1
using namespace std;
// Create a type since this is a long name to use (optional)
typedef std::tr1::unordered_map<string, int> StringKeyMap;
// Create a dummyMap object so we can get its hash function
StringKeyMap dummyMap;
// Capture the hash function for use in program
StringKeyMap::hasher myHashFunction = dummyMap.hash_function();
int main ()
\{
cout << "Hashing a String: " << myHashFunction("Hashing a String:") << end7;
cout << "Smashing a String: " << myHashFunction ("Smashing a String:") << end1;
return 0;
\} // end main

## Output

Hashing a String: 2084157801917477989
Smashing a String: 14048775086903850803

## C7.2 STL Algorithms

One of the strengths of the STL container library is the set of functions defined in the header <al gorithm>. These functions are designed to operate on ranges of elements in any container with iterators. In C++ Interlude 6 you saw a few of these: for_each, find, count, advance, and distance. In most cases, the type of container is not important, but it must support iterators. This makes possible the comparison of differing types of containers. For instance, the entries stored in a vector can be compared to those stored in a map.

The first group of algorithms search and compare items in a range. Though they may change a value, they do not modify the structure of the containers. Several algorithms, such as for_each, find_if, and count_if, apply helper functions to each entry during the traversal in a manner similar to the tree traversals presented in Chapter 15. The table below only shows one form of each function. Most functions have several forms that allow additional customization of the range of container entries to process and the functions to use for comparison. Also, STL map operations, such as upper_ bound and lower_bound can be used as functions and applied to other container types.

STL Search and Compare Algorithms

```
Operation
void for_each(iterator start, iterator end, Function fun)
iterator find(iterator start,
iterator end, value_type\& va1)
iterator find_if(iterator start, iterator end, PredFunction fun)
uint count (iterator start, iterator end, value_type\& val)
uint count_if(iterator start, iterator end, PredFunction fun)
bool equal(iterator start1,
iterator end1,
iterator start2)
value_type\& min(value_type\& item1, value_type\& item2)
value_type\& min_element(iterator start, iterator end)
value_type\& \(\max\) (value_type\& item1, value_type\& item2)
value_type\& max_element(iterator start, iterator end)
```


## Description

Applies fun to the range specified by start and end.

Returns an iterator to val in the range specified by start and end.

Returns an iterator to the entry between start and end that makes the predicate function fun true.

Counts the number of occurrences of val between start and end.

Counts the number of entries between start and end that make the predicate function fun true.

Compares the entries between start1 and end1 to those beginning at start2. Returns true if all items match.

Returns the minimum of item 1 and item2.

Returns the minimum value in the range from start to end.

Returns the maximum of item1 and item2.

Returns the maximum value in the range from start to end.

The sequence-modification algorithms change the entries stored in a container within a range specified by two iterators. The entries can be either

- Replaced by a specific value using fil1
- Copied or moved to new locations using copy or move
- Removed based on specified conditions using remove, or
- Have operations applied to them using transform

The following table lists a few of the more common operations and the syntax used to invoke them.

## STL Sequence Modification Algorithms

```
Operation
iterator copy(iterator start1,
    iterator end1,
    iterator start2)
iterator copy_backward(iterator start1,
                    iterator end1,
                            iterator start2)
void swap(value_type& item1,
    value_type& item2)
iterator transform(iterator start1,
        iterator end1,
        iterator start2,
        UnaryOperator op)
iterator transform(iterator start1,
        iterator end1,
        iterator operand2,
        iterator start2,
        BinaryOperator bop)
void fil1(iterator start1,
    iterator end1,
    value_type& va1)
```


## Description

Copies the entries within the range start1 to end1 to the positions beginning at start2 within the same or another container. Returns an iterator to the last entry copied.
Copies the entries within the range start1 to end1 to the positions beginning at start2 (used if ranges overlap in the same container). Copies entries starting at end1-1. Returns an iterator to the last entry copied.
Swaps the values of two objects, item1 and item2.

Applies the given unary operator op to items within the range start1 to end1 and places the result in elements beginning at start2. Returns an iterator to the last element in the result container.
Uses items within the range start1 to end1 as the left-hand operand for bop and the entries beginning at operand2 as the right-hand operand. Places the result in elements beginning at start2. Returns an iterator to the last element in the result container.
Sets all entries within the range from start1 to end1 to the value val.

The STL algorithm library also provides a number of operations that perform sorting or heap operations on a range of elements in a container. It also includes operations, such as partition and nth_element that can be used to implement user-developed sorting algorithms.

STL Sorting and Heap Algorithms

Operation
void sort(iterator start, iterator end)

Description
Sorts entries within the range start to end into ascending order.

```
void stable_sort(iterator start,
    iterator end)
iterator partition(iterator start,
    iterator end,
    PredFunction fun)
iterator partition_stable(iterator start,
    iterator end,
    PredFunction fun)
void nth_element(iterator start,
    iterator nth,
    iterator end)
void make_heap(iterator start,
    iterator end)
void push_heap(iterator start,
    iterator end)
void pop_heap(iterator start,
    iterator end)
void sort_heap(iterator start,
    iterator end)
```

Sorts entries within the range start to end into ascending order but has stability in ordering - that is, the relative order of equal values is maintained.
Rearranges all entries within the range from start to end so that those that make the predicate function fun true are before those for which the function is false.
Behaves the same as partition, but has stability in ordering - that is, the relative order of equal values is maintained.

Rearranges the entries within the range start to end so that the entry in position nth is in the correct sorted position-that is, it is in the nth position in the sorted sequence.
Rearrange the entries in the range start to end to form a heap.

Assumes that the entries within the range start to end - 1 form a heap and places the entry at position end into the correct location in the heap, extending the heap to position end.
Moves the value at position start to position end - 1 and rebuilds the heap in the range start to end -2 inclusive.
Assumes that the entries within the range start to end - 1 form a heap and turns the heap back into a sorted container.

## Review of C++ Fundamentals

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PrerequisiteKnowledge of a modern programming language

This book assumes that you already know how to write programs in a modern programming language. If that language is $\mathrm{C}++$, you can probably skip this appendix, returning to it for reference as necessary. If instead you know a language such as Java or C, this appendix will introduce you to C++. Also, Appendixes K and L compare C++ to Java and Python, respectively.

It isn't possible to cover all of $\mathrm{C}++$ in these pages. Instead this appendix focuses on the parts of the language used in this book. First we look at the basics of data types, variables, expressions, operators, and simple input/output. We continue with functions, decision constructs, looping constructs, arrays, and strings. Various C++ Interludes, which appear throughout the book as needed, will cover classes, pointers, exceptions, and files.

## A. 1 Language Basics

Let's begin with the elements of the language that allow you to perform simple computations. For example, the $\mathrm{C}++$ program in Figure A-1 computes the volume of a sphere. Running this program produces the following output, where the user's response appears in blue:

Enter the radius of the sphere: 19.1
The volume of a sphere of radius 19.1 is 29186.927734
A typical C++ program consists of several modules, some of which you write and some of which you use from standard libraries. C++ provides a source-inclusion facility, which allows the system to include the contents of a file automatically at a specified point in a program before the program is compiled. For example, our sample program uses a standard library to perform input and output operations. The first line of this program is an include directive that names a standard header iostream, which enables the program to use the input/output module. The second line informs the compiler to use the standard namespace (see Section C1.2 in C++ Interlude 1).

A C++ program is a collection of functions, one of which must be called main. Program execution always begins with the function main. The following paragraphs provide an overview of the basics of $\mathrm{C}++$ and refer to the simple program in Figure A-1 by line number. Note that the only function this simple program contains is main.

## A.1.1 Comments

Each comment line in C++ begins with two slashes // and continues until the end of the line. You can also begin a multiple-line comment with the characters /* and end it with */. However, a

## FIGURE A-1 A simple C++ program

| 1. Enables input and output ------------> | \#include <iostream> |
| :---: | :---: |
| 2. Opens the standard namespace ----> | using namespace std; |
| 3. Begins the function main ------------> | int main() |
|  | // Computes the volume of a sphere of a given radius. |
| 5. Begins body of function -------------> | \{ |
| 6. Defines a constant -------------------> | const doub7e PI = 3.14159; |
| 7. Declares a variable ------------------->> | double radius; |
| 8. Displays a prompt to the user -------> | cout << "Enter the radius of the sphere: "; |
| 9. Reads radius ------------------------> | cin >> radius; |
| 10. Declares and computes volume ---> | double volume $=4$ * PI * radius * radius * radius / 3; |
|  | cout << "The volume of a sphere of radius " |
| 12. Statement continues ------------------> | << radius << " inches is " << volume |
| 13. Statement continues -----------------> | << " cubic inches.\n"; |
| 14. Normal program termination -------> | return 0; |
| 15. Ends body of function --------------->> | \} // end program |

comment that begins with /* and ends with */ cannot contain another comment that begins with /* and ends with */.

Appendix I talks about documentation comments that begin with /** and end with */. We use this style of comment in this book at the beginning of classes, methods, and functions.

## A.1.2 Identifiers and Keywords

A C++ identifier is a sequence of letters, digits, and underscores that must begin with either a letter or an underscore. C++ distinguishes between uppercase and lowercase letters, so be careful when typing identifiers.

You use identifiers to name various parts of the program. Certain identifiers, however, are reserved by $\mathrm{C}++$ as keywords, and you should not use them for other purposes. A list of all $\mathrm{C}++$

C++ is casesensitive reserved keywords appears inside the cover of this book. The keywords that occur within C++ statements in this book appear in color.

## A.1.3 Fundamental Data Types

The fundamental data types in $\mathrm{C}++$ are organized into four categories: boolean, character, integer, and floating point. With the exception of boolean, each category contains several data types. For most applications, you can use

| bool | for boolean values |
| :--- | :--- |
| char | for character values |
| int | for integer values |
| double | for floating-point values |

## FIGURE A-2 Fundamental data types

| Category | Available Data Types by Category |  |  |
| :---: | :---: | :---: | :---: |
| Boolean | bool |  |  |
| Character | char | signed char | unsigned char |
| Signed integer | short | int | long |
| Unsigned integer | unsigned short | unsigned | unsigned long |
| Floating point | float | doub7e | long double |

A boolean value can be either true or false. Characters are represented by their ASCII integer values, which are listed in Appendix J. Integer values are either signed, such as -5 and +98 , or unsigned, such as 5 and 98 . The floating-point types are used for real numbers that have both an integer portion and a fractional portion. Boolean, character, and integer types are called integral types. Integral and floating-point types are called arithmetic types.

Most of the data types are available in several forms and sizes. Although you will probably not need more than the four types given previously, Figure A-2 lists the available fundamental data types for your reference.

The size of a data type affects the range of its values. For example, a long integer can have a larger magnitude than a short integer. The sizes of-and therefore the specific ranges for-a data type depend on the particular computer and version of $\mathrm{C}++$ that you use. $\mathrm{C}++$, however, provides a way to determine these ranges, as you will see later in Section A.1.6.

## A.1.4 Variables

A variable, whose name is a C++ identifier, represents a memory location that contains a value of a particular data type. You declare a variable's data type by preceding the variable name with the data type, as in

```
doub7e radius; // Radius of a sphere
```

Note that you can write a comment on the same line to describe the purpose of the variable.
This declaration is also a definition in that it assigns memory for the variable radius. The memory, however, has no particular initial value and so is said to be uninitialized. The program in Figure A-1 declares radius without an initial value and later reads its value by using cin >> radius.

When possible, avoid uninitialized variables

When possible, you should avoid uninitialized variables. That is, you should initialize a variable when you first declare its data type or, alternatively, declare a variable's data type when you first assign it a value. For example, volume appears for the first time in line 10 of Figure A-1 in the statement

```
double volume = 4 * PI * radius * radius * radius / 3;
```

Because we did not declare volume's data type earlier in the program-thus avoiding an uninitialized value - we declare its data type and assign it a value in the same statement.

## A.1.5 Literal Constants

You use literal constants to indicate particular values within a program. The 4 and 3 in line 10 of Figure A-1 are examples of literal constants that are used within a computation. You can also use a literal constant to initialize the value of a variable. For example, you use true and false as the values of a boolean variable, as mentioned previously.

## FIGURE A-3 Some special character constants

| Constant | Name |
| :---: | :---: |
| \n | New line |
| $\backslash \mathrm{t}$ | Tab |
| \' | Single quote |
| \" | Double quote |
| $\backslash 0$ | Zero |

You write decimal integer constants without commas, decimal points, or leading zeros. The default data type of such a constant is either int, if small enough, or long.

You write floating-point constants, which have a default type of double, with a decimal point. You can specify an optional power-of-10 multiplier by writing e or E followed by the power of 10 . For example, $1.2 \mathrm{e}-3$ means $1.2 \times 10^{-3}$.

Character constants are enclosed in single quotes-for example, 'A' and '2'-and have a default type of char. You write a literal character string as a sequence of characters enclosed in double quotes.

Several characters have names that use a backslash notation, as given in Figure A-3. This notation is useful when you want to embed one of these characters within a literal character string. For example, the program in Figure A-1 uses the new-line character $\backslash n$ in the string "cubic inches. $\backslash n$ " to end the line of output with a carriage return. You will learn about this use of $\backslash n$ in the discussion of output later in this appendix. You also use the backslash notation to specify either a single quote as a character constant ( ' $\backslash$ ' ') or a double quote within a character string.

Programming Tip: Do not begin a decimal integer constant with zero. A constant that begins with zero is either an octal constant or a hexadecimal constant. ${ }^{1}$

## A.1.6 Named Constants

Unlike variables, whose values can change during program execution, named constants have values that do not change. The declaration of a named constant is like that of an initialized variable, but the keyword const precedes the data type. For example, the statement

```
const doub7e PI = 3.14159;
```

declares PI as a named floating constant, as is the case in the sample program in Figure A-1. Once a named constant such as PI is declared, you can use it, but you cannot assign it another value. By using named constants, you make your program both easier to read and easier to modify.

The standard header file climits contains named constants such as INT_MIN and LONG_MAX that specify installation-dependent maximum and minimum values for the integral data types. Likewise, the standard header file cfloat contains named constants that specify installation-dependent maximum and minimum values for the floating data types. You use the include directive to gain access to these header files.

[^19]Do not begin a decimal integer constant with zero

The value of a named constant does not change

Named constants make a program easier to read and modify

## Enumeration

 provides another way to name constantsYou can create an integral data type by naming an enumeration

The typedef statement gives another name to an existing data type, making your program easier to change

## A.1.7 Enumerations

Enumeration provides another way to name integer constants. For example, the statement

```
enum {SUN, MON, TUE, WED, THU, FRI, SAT};
```

is equivalent to the statements

```
const int SUN = 0;
const int MON = 1;
const int SAT = 6;
```

By default, the values assigned to the constants-called enumerators-begin with zero and are consecutive. You can, however, assign explicit values to any or all of the enumerators, as in
enum \{PLUS = '+', MINUS = '-'\};
By naming an enumeration, you create a distinct integral type. For example,
enum Season \{WINTER, SPRING, SUMMER, FALL\};
creates a type Season. The variable whichSeason, declared as
Season whichSeason;
can have values WINTER, SPRING, SUMMER, or FALL. This use of named enumerations instead of int can make your program easier to understand.

## A.1.8 The typedef Statement

You use the typedef statement to give another name to an existing data type. In this way, you can make your program easier to modify and to read. For example, the statement

```
typedef doub7e Rea1;
```

declares Rea 1 as a synonym for doub1e and allows you to use Real and doub7e interchangeably.
Suppose that you revise the program in Figure A-1 by using Real as follows:

```
int main()
{
    typedef doub7e Rea1;
    const Real PI = 3.14159;
    Real radius;
    cout << "Enter the radius of the sphere: ";
    cin >> radius;
    Real volume = 4 * PI * radius * radius * radius / 3;
```

At first glance, this program does not seem to be more advantageous than the original version, but suppose that you decide to increase the precision of your computation by declaring PI, radius, and volume as long double instead of double. In the original version of the program (Figure A-1), you would have to locate and change each occurrence of double to long double. In the revised program, you simply change the typedef statement to

```
typedef long double Rea1;
```

Realize that typedef does not create a new data type, but simply declares a new name for a data type. A new data type requires more than a name; it requires a set of operations. $\mathrm{C}++$, however, does provide a way to create your own data types, as described in $\mathrm{C}++$ Interlude 1.

## A.1.9 Assignments and Expressions

You form an expression by combining variables, constants, operators, and parentheses. The assignment statement

```
volume = 4 * PI * radius * radius * radius / 3;
```

assigns to the previously declared variable volume the value of the arithmetic expression on the righthand side of the assignment operator $=$, assuming that PI and radius have values. The assignment statement

```
double volume = 4 * PI * radius * radius * radius / 3;
```

which appears in line 10 of Figure A-1, also declares volume's data type, since it was not declared previously.

The various kinds of expressions that you can use in an assignment statement are discussed next.
Arithmetic expressions. You can combine variables and constants with arithmetic operators and parentheses to form arithmetic expressions. The arithmetic operators are

+ Binary add or unary plus
- Binary subtract or unary minus
* Multiply
/ Divide
\% Remainder after division
The operators *, /, and \% have the same precedence, ${ }^{2}$ which is higher than that of + and $-;$ unary operators ${ }^{3}$ have a higher precedence than binary operators. The following examples demonstrate operator precedence:

```
a - b / c means a-(b/c) (precedence of / over -)
-5 / a means (-5)/a (precedence of unary operator -)
a / -5 means a/(-5) (precedence of unary operator -)
```

Arithmetic operators and most other operators are left-associative. That is, operators of the same precedence execute from left to right within an expression. Thus,
$\mathrm{a} / \mathrm{b}$ * c
means
(a/b) *c
The assignment operator and all unary operators are right-associative, as you will see later. You can use parentheses to override operator precedence and associativity.
Relational and logical expressions. You can combine variables and constants with parentheses; with the relational, or comparison, operators $<,<=,>=$, and $>$; and with the equality operators $==$ (equal to) and $!=($ not equal to) to form a relational expression. Such an expression is true or false according to the validity of the specified relation. For example, the expression $5=4$ has a value of false because 5 is not equal to 4 . Note that equality operators have a lower precedence than relational operators.

You can combine variables and constants of the arithmetic types, relational expressions, and the logical operators \&\& (and) and \| (or) to form logical expressions, which are either true or false. C++ evaluates logical expressions from left to right and stops as soon as the value of the entire expression

Operators have a set precedence

Operators are either left- or rightassociative

Logical expressions are evaluated from left to right

[^20]Sometimes the value of a logical expression is apparent before it is completely examined

Conversions from one data type to another occur during both assignment and expression evaluation

Use a static cast to convert explicitly from one data type to another
is apparent; that is, $\mathrm{C}++$ uses short-circuit evaluation. For example, $\mathrm{C}++$ determines the value of each of the following expressions without evaluating $(\mathrm{a}<\mathrm{b})$ :

```
(5 == 4) && (a < b) // False since (5 == 4) is false
(5 == 5) || (a < b) // True since (5 == 5) is true
```



Programming Tip: Remember that $=$ is the assignment operator; $==$ is the equality operator.

Conditional expressions. The expression

$$
\text { expression }_{1} ? \text { expression }_{2}: \text { expression }_{3}
$$

has the value of either expression ${ }_{2}$ or expression ${ }_{3}$ according to whether expression ${ }_{1}$ is true or false, respectively. For example, the statement

$$
\text { 1arger }=((a>b) ? a: b) ;
$$

assigns the larger of $a$ and $b$ to larger, because the expression $a>b$ is true if $a$ is larger than $b$ and false if not.

Implicit type conversions. Automatic conversions from one data type to another can occur during assignment and during expression evaluation. For assignments, the data type of the expression on the right-hand side of the assignment operator is converted to the data type of the item on the left-hand side just before the assignment occurs. Floating-point values are truncated-not rounded-when they are converted to integral values.

During the evaluation of an expression, any values of type char or short are converted to int. Similarly, any enumerator value is converted to int if int can represent all the values of that particular enum; otherwise, it is converted to unsigned. These conversions are called integral promotions. After these conversions, if the operands of an operator differ in data type, the data type that is lower in the following hierarchy is converted to one that is higher (int is lowest):

$$
\text { int } \rightarrow \text { unsigned } \rightarrow \text { long } \rightarrow \text { unsigned 1ong } \rightarrow \text { float } \rightarrow \text { double } \rightarrow \text { long double }
$$

For example, if $a$ is long and $b$ is $f 7 o a t, a+b$ is $f 7$ oat. Only a copy of $a$ 's 7 ong value is converted to float prior to the addition, so that the value stored at a is unchanged.
Explicit type conversions. You can explicitly convert from one data type to another using a static cast, with the following notation:

```
static_cast<type> (expression)
```

which converts expression to the data type type. For example, static_cast<int> (14.9) converts the double value 14.9 to the int value 14 . Thus, the sequence

```
doub7e volume = 14.9;
cout << static_cast<int>(volume);
```

displays 14 but does not change the value of volume.
Other assignment operators. In addition to the assignment operator $=, \mathrm{C}++$ provides several twocharacter assignment operators that perform another operation before assignment. For example,

$$
a+=b \text { means } a=a+b
$$

Other operators, such as $-=, *=, /=$, and $\%=$, have analogous meanings.

Two more operators, ++ and --, provide convenient incrementing and decrementing operations:
$\mathrm{a}++$ means $\mathrm{a}+=1$, which means $\mathrm{a}=\mathrm{a}+1$
Similarly,
a-- means a $-=1$, which means $\mathrm{a}=\mathrm{a}-1$
The operators ++ and -- can either follow their operands, as you just saw, or precede them. Although $++a$, for instance, has the same effect as a++, the results differ when the operations are combined with assignment. For example,

$$
\mathrm{b}=++\mathrm{a} \text { means } \mathrm{a}=\mathrm{a}+1 ; \mathrm{b}=\mathrm{a}
$$

Here, the ++ operator acts on a before assigning a's new value to $b$. In contrast,
$\mathrm{b}=\mathrm{a}++$ means $\mathrm{b}=\mathrm{a} ; \mathrm{a}=\mathrm{a}+1$
The assignment operator assigns a's old value to $b$ before the ++ operator acts on $a$. That is, the ++ operator acts on a after the assignment. The operators ++ and -- are often used within loops and with array indices, as you will see later in this appendix. When we use these operators with arithmetic variables, we write the operator after the variable.

In addition to the operators described here, $\mathrm{C}++$ provides several other operators. A summary of all $\mathrm{C}++$ operators and their precedences appears inside the cover of this book.

## A. 2 Input and Output Using iostream

A typical C++ program reads its input from a keyboard and writes its output to a display. Such input and output consist of streams, which are simply sequences of characters that either come from or go to an input or output (I/O) device.

The data type of an input stream is istream, and the data type of an output stream is ostream. The iostream library provides these data types and three default stream variables: cin for the standard input stream, cout for the standard output stream, and cerr for the standard error stream, which also is an output stream. Your program gains access to the iostream library by including the iostream header file. This section provides a brief introduction to simple input and output.

## A.2.1 Input

C++ provides the input operator >> to read integers, floating-point numbers, and characters into variables whose data types are any of the fundamental data types. The input operator has the input stream as its left operand and the variable that will contain the value read as its right operand. Thus,
cin >> x;
reads a value for x from the standard input stream. The $\gg$ operator is left-associative. Thus,

```
cin >> x >> y
```

means

$$
(\operatorname{cin} \gg x) \gg y
$$

That is, both of these expressions read characters for x from the input stream and then read subsequent characters for y .

The operators ++ and -- are useful for incrementing and decrementing a variable

The input operator >> reads from an input stream

The input operator >> skips whitespace

Use get to read whitespace

The output operator << writes to an output stream

The input operator >> skips whitespace, such as blanks, tabs, and new-line characters, that might occur between values in the input data line. For example, after the program segment

```
int ia = 0;
int ib = 0;
doub7e da = 0;
doub7e db = 0;
cin >> ia >> da >> ib;
cin >> db;
```

reads the data line

```
21 -3.45 -6 475.1e-2 <eol>
```

the variable ia contains 21 , da contains -3.45 , ib contains -6 , and db contains 4.751 . A subsequent attempt to read from cin will look beyond the end of the line ( $\langle e o l\rangle$ ) and read from the next data line, if one exists. An error occurs if no data exists for a corresponding variable processed by >> or if the variable's data type does not match the type of the data available. For example, after the previous program segment reads the data line
-1.23 456.1e-2 -7 8 <eol>
the variable ia contains -1 , da contains 0.23 , ib contains 456 , and $d b$ contains 0.001 . The rest of the data line is left for a subsequent read, if any. As another example, if the segment attempts to read a data line that begins with .21 , the read would terminate because ia is int and .21 is not.

An expression such as cin $\gg x$ has a value after the read operation takes place. If the operation is successful, this value is true; otherwise the value is false. You can examine this value by using the selection and iteration statements that are described later in this appendix.

You can also use the >> operator to read individual characters from the input stream into character variables. Again, any whitespace is skipped. For example, after the program segment

```
char ch1 = '';
char ch2 = '';
char ch3 = '';
cin >> ch1 >> ch2 >> ch3;
```

reads the data line
xy z
ch1 contains ' $x$ ', ch2 contains ' $y$ ', and ch3 contains ' $z$ '.
You can read whitespace when reading individual characters into character variables by using the C++ method get. Either of the statements

```
cin.get(ch1);
```

or

```
ch1 = cin.get();
```

reads the next character, even if it is a blank, a tab, or a new-line character, from the input stream into the char variable ch1.

Section A.7, later in this appendix, describes how to read character strings.

## A.2.2 Output

$\mathrm{C}++$ provides the output operator $\ll$ to write character strings and the contents of variables whose data types are any of the fundamental ones. For example, the program segment

```
int count = 5;
doub1e average = 20.3;
cout << "The average of the " << count
    << " distances read is " << average
    << " miles.\n";
```

produces the following output:
The average of the 5 distances read is 20.3 miles.
Like the input operator, the output operator is left-associative. Thus, the previous statements append the string "The average of the" to the output stream, then append the characters that represent the value of count, and so on.

Note the use of the new-line character $\backslash \mathrm{n}$, which you can conveniently embed within a character string. Observe also that the output operator does not automatically introduce whitespace between values that are written; you must do so explicitly. The following statements provide another example of this:

```
int x = 2;
int y = 3;
char ch = 'A';
cout << x << y << ch << "\n"; // Disp7ays 23A
```

Although you can use the output operator to display individual characters, you can also use the put method for this task. Further, you can specify a character either as a char variable or in ASCII. Thus, the statements

```
char ch = 'a';
cout.put(ch); // Displays a
cout.put('b'); // Displays b
cout.put(99); // Displays c, which is 99 in ASCII
cout.put(ch+3); // Displays d
cout.put('\n'); // Carriage return
```

display abcd followed by a carriage return.
Section A.7, later in this appendix, provides further information about writing character strings.

## A.2.3 Manipulators

C++ enables you to gain more control over the format of your output and the treatment of whitespace during input than the previous discussion has indicated. Most of these techniques apply to the format of output.

Suppose, for example, that you have computed your grade point average and you want to display it with one digit to the right of the decimal point. If the floating variable gpa contains 4.0 , the statement

```
cout << "My GPA is " << gpa << "\n";
```

writes 4 without a decimal point. A number of manipulators affect the appearance of your output. You can use these with cout:

```
cout << manipulator;
```

where manipulator has any of the values listed in Figure A-4. A manipulator is a predefined value or function that you use with the input and output operator. For example,

```
cout << showpoint;
```

uses the showpoint manipulator and causes all floating-point output to appear with a decimal point.

You need to explicitly introduce new-line characters and whitespace where desired in a program's output

Use manipulators to specify the appearance of a program's output

Manipulators also affect the appearance of a program's output

FIGURE A-4 Stream manipulators

| Manipulator | Meaning |
| :---: | :---: |
| end7 | Insert new line and flush stream |
| fixed | Use fixed decimal point in floating-point output |
| 1eft | Left-align output |
| right | Right-align output |
| scientific | Use exponential (e) notation in floating-point output |
| setfill(f) | Set fill character to f |
| setprecision(n) | Set floating-point precision to integer n |
| setw( n ) | Set field width to integer n |
| showpoint | Show decimal point in floating-point output |
| showpos | Show + with positive integers |
| ws | Extract whitespace characters |

Even if you use the showpoint manipulator, gpa will likely appear as 4.00000 instead of 4.0. You can specify the number of digits that appear to the right of the decimal point by using the manipulator function setprecision, and you can insert a new-line character and flush the output stream by using the manipulator endl. Thus,

```
cout << showpoint;
cout << setprecision(1) << gpa << endl;
```

displays 4.0 followed by a carriage return.
The effect of setprecision on the output stream remains until another setprecision is encountered. Except for setprecision, however, a manipulator affects the appearance of only the next characters on which << (or >>) operates. For example,
cout $\ll$ right; // Right-align output
cout << "abc" << setw(6) << "def" << "ghi";
displays
abc defghi
Although manipulator values, such as end1, are available when you include iostream in your program, you must also include iomanip to use any of the manipulator functions.

## A. 3 Functions

A C++ program is a collection of functions

A function definition implements a function's task

As was mentioned earlier in this appendix, a C++ program is a collection of functions. Usually, each function should perform one well-defined task. For example, the following function returns the larger of two integers:

```
int computeMax(int x, int y)
{
    if (x > y)
    return x;
    e1se
        return y;
} // end computeMax
```

A function definition, like the one just given, has the following form:

```
type name(parameter-declaration-list)
{
    body
}
```

The portion of the definition before the left brace specifies a return type, the function name, and a list of parameters. The part of the definition that appears between the braces is the function's body.

The return type of a valued function - one that returns a value-is the data type of the value that the function will return. The body of a valued function must contain a statement of the form

```
return expression;
```

where expression has the value to be returned.
Each parameter represents either an input to or an output from the function. You declare a parameter by writing a data type and a parameter name, separating it from other parameter declarations with a comma, as in

```
int x, int y
```

When you call, or invoke, the function computeMax, you pass it arguments that correspond to the parameters with respect to number, order, and data type. For example, the following statements contain two calls to max:

```
int a = 0;
int b = 0;
int c=0;
cin >> a >> b >> c;
int largerAB = computeMax(a, b);
cout << "The largest of " << a << ", " << b << ", "
    << " and " << c << " is " << computeMax(largerAB, c) << ".\n";
```

As written, the definition of computeMax indicates that its arguments are passed by value. That is, the function makes local copies of the values of the arguments-a and $b$, for example-and uses these copies wherever $x$ and $y$ appear in the function definition. Thus, the function cannot alter the arguments that you pass to it. This restriction is desirable in this example because x and y are input parameters, which computeMax does not change.

Alternatively, arguments can be passed by reference. The function does not copy such arguments; rather, it references the argument locations whenever the parameters appear in the function's definition. This allows a function to change the values of the arguments, thus implementing output parameters.

For example, consider the following variation of the function computeMax:

```
void computeMax(int x, int y, int& larger)
{
    larger = ((x > y) ? x : y);
} // end computeMax
```

computeMax is a void function instead of a valued function. That is, its return type is void, and it does not return a value by using a return statement. ${ }^{4}$ Instead, computeMax returns the larger of x and y in

[^21]A valued function must use return to return a value

When you call a function, you pass it arguments that correspond to the parameters in number, order, and data type

An argument passed by value is copied within the function

An argument passed by reference is not copied but is accessed directly within the function

A void function does not use return to return a value

An output argument should be a reference argument

An input argument should be either a value argument or a constant reference argument

An argument that is both an input to and an output from a function is passed by reference
the output parameter 1arger. The \& that follows 1arger's data type int indicates that 1arger is a reference parameter. Thus, computeMax will access and alter the argument that corresponds to larger, whereas the function will make and use copies of the values of the arguments that correspond to the value parameters $x$ and $y$.

The following statements demonstrate how to invoke computeMax:

```
int a = 0;
int b = 0;
int largerAB = 0;
cin >> a >> b;
computeMax(a, b, largerAB);
cout << "The larger of " << a << " and " << b
    << " is " << largerAB << ".\n";
```

If a function's input argument is a large object, like the objects you will encounter in this book, you might not want the function to copy it. Thus, you would not pass the argument by value. Because it is an input argument, however, you do not want the function to be able to alter it. A constant reference parameter is a reference parameter that is tagged as const. The function uses the actual argument that is passed to such a parameter, not a copy of it, yet cannot modify it.

For example, for the function $f$ that begins

```
void f(const int& x, int y, int& z)
```

$x$ is a constant reference parameter, $y$ is a value parameter, and $z$ is a reference parameter. Here $x$ and $y$ are suitable as input parameters because $f$ cannot change them, while $z$ is an output parameter. Note that $z$ can also be an input parameter. That is, the argument corresponding to $z$ can both provide a value to the function and return a value from the function. Such arguments must be passed by reference.

Note: Use reference parameters with caution, as you might inadvertently change an argument. On the other hand, constant reference parameters are safe to use.

If you write another function $f$ that calls computeMax, you must either place the definition of $f$ after the definition of computeMax or precede f's definition with a function declaration for computeMax. For example, you can use either of the following statements to declare the function computeMax:
void computeMax (int $x$, int $y$, int\& max);
or
void computeMax (int, int, int\&);
A function declaration provides the data types of the function's parameters and its return type. Parameter names are optional in a function declaration, although they are helpful stylistically. However, parameter names are required in the function's definition. Although a function declaration ends with a semicolon, a semicolon does not appear in a function definition.

A typical $\mathrm{C}++$ program contains a function declaration for every function used in the program. These declarations appear first in the program, usually with comments that describe each function's purpose, parameters, and assumptions. The program in Listing A-1 demonstrates the placement of a function declaration, function definition, and main function:

```
LISTING A-1 A program that contains a function declaration
#include <iostream>
using namespace std;
/** Returns the larger of two given integers.
@param x An integer.
@param y An integer.
@return The larger of x and y.*/
int computeMax(int x, int y); // A function declaration
int main()
{
    int a = 0;
    int b = 0;
    cout << "Please enter two integers: ";
    cin >> a >> b;
    int largerAB = computeMax (a, b);
    cout << "The larger of " << a << " and " << b
        << " is " << largerAB << ".\n";
} // end main
int computeMax(int x, int y)
{
    return (x > y) ? x : y;
} // end computeMax
```


## A.3.1 Standard Functions

C++ provides many standard functions, such as the square root function sqrt and the input function get. Appendix H provides a summary of the standard functions and indicates which header file you need to include in your program to gain access to them. For example, the standard functions listed in Figure A-5 facilitate character processing and require the header file cctype. Thus, you need to include the statement

```
#include <cctype>
```

in your program when you want to use functions such as isupper and toupper. For the character variable ch, isupper (ch) is true if ch is an uppercase letter, and toupper (ch) returns the uppercase version of the letter ch without actually changing ch.

## A. 4 Selection Statements

Selection statements allow you to choose among several courses of action according to the value of an expression. In this category of statements, C++ provides the if statement and the switch statement.

## A.4.1 The if Statement

You can write an if statement in one of two ways:

```
if (expression)
    statement1
```

Standard functions provide many common operations and require a specific header file

An if statement has two basic forms

Parentheses around the expression in an if statement are required

You can nest if statements

FIGURE A-5 A selection of (a) standard classification functions; and (b) standard conversion functions
(a)

| Function |  | Returns true if ch is |
| :--- | :--- | :--- |
| isalnum (ch) |  | A letter or digit |
| isalpha(ch) |  | A letter |
| isdigit $(\mathrm{ch})$ |  | A digit |
| islower $(\mathrm{ch})$ |  | A lowercase letter |
| isupper $(\mathrm{ch})$ |  | An uppercase letter |

(b)

| Function | Returns |
| :---: | :---: |
| tolower (ch) | Lowercase version of ch |
| toupper (ch) | Uppercase version of ch |
| toascii(ch) | int ASCII code for ch |

or

```
if (expression)
    statement1
else
    statement2
```

where statement 1 and statement 2 represent any $C++$ statement except a declaration. Such statements can be compound; a compound statement, or block, is a sequence of statements enclosed in braces. If the value of expression is true, statementl is executed. Otherwise, the first form of the if statement does nothing, whereas the second form executes statement 2 . Note that the parentheses around expression are required.

For example, the following if statements each compare the values of two integer variables $a$ and $b$ :

```
if (a > b)
    cout << a << " is larger than " << b << ".\n";
cout << "This statement is always executed.\n";
if (a > b)
{
    7argerAB = a;
    cout << a << " is larger than " << b << ".\n";
}
else
{
    1argerAB = b;
    cout << b << " is larger than " << a << ".\n";
} // end if
cout << largerAB << " is the larger value.\n";
```

You can nest if statements in several ways, since either statementl or statement 2 can itself be an if statement. The following example, which determines the largest of three integer variables $a, b$, and c, shows a common way to nest if statements:

```
if ((a >= b) && (a >= c))
    largest = a;
else if (b >= c) // a is not largest at this point
    largest = b;
else
    1argest = c;
```



Note: An arithmetic expression whose value is not zero is treated as true; one having a value of zero is false.

## A.4.2 The switch Statement

When you must choose among more than two courses of action, the if statement can become unwieldy. If your choice is to be made according to the value of an integral expression, you can use a switch statement.

For example, the following statement determines the number of days in a month. The int variable month designates the month as an integer from 1 to 12 , and the boolean variable 1eapYear is true if the year is a leap year.

```
switch (month)
{
    // 30 days hath Sept., Apr., June, and Nov.
    case 9: case 4: case 6: case 11:
        daysInMonth = 30;
        break;
    // Al1 the rest have 31
    case 1: case 3: case 5: case 7:
    case 8: case 10: case 12:
        daysInMonth = 31;
        break;
    // Except February
    case 2: // Assume leapYear is true if a leap year, else is false
        if (leapYear)
            daysInMonth = 29;
        else
            daysInMonth = 28;
        break;
    default:
        cout << "Incorrect value for month.\n";
} // end switch
```

Parentheses must enclose the integral switch expression-month, in this example. The case labels have the form

```
case expression:
```

where expression is a constant integral expression. After the swi tch expression is evaluated, execution continues at the case label whose expression has the same value as the switch expression. Subsequent statements execute until either a break or a return is encountered or the switch statement ends.

Unless you terminate a case with either a break or a return, execution of the switch statement continues. Although this action can be useful, omitting the break statements in the previous example would be incorrect.

## A switch

 statement provides a choice of several actions according to the value of an integral expressionWithout a break statement, execution of a case will continue into the next case

A while statement executes as long as the expression is true

A for statement lists the initialization, testing, and updating steps in one location

If no case label matches the current value of the switch expression, the statements that follow the default label, if one exists, are executed. If no default exists, the switch statement exits.

## A. 5 Iteration Statements

C++ has three statements-the while, for, and do statements-that provide for repetition by iteration, that is, loops. Each statement controls the number of times that another $\mathrm{C}++$ statement-the body-is executed. The body cannot be a declaration and is often a compound statement.

## A.5.1 The while Statement

The general form of the while statement is

```
while (expression)
``` statement

As long as the value of expression is true, statement is executed. Because expression is evaluated before statement is executed, it is possible that statement will not execute at all. Note that the parentheses around expression are required.

Suppose that you wanted to compute the sum of positive integers that you enter at the keyboard. Since the integers are positive, you can use a negative value or zero to indicate the end of the list of integers. The following while statement accomplishes this task:
```

int nextValue = 0;
int sum = 0;
cin >> nextValue;
while (nextValue > 0)
{
sum += nextValue;
cin >> nextValue;
} // end while

```

If 0 was the first value read, the body of the while statement would not execute.
Recall that the expression cin >> nextValue has the value true if the input operation was successful and false otherwise. Thus, you could revise the previous statements as
```

int nextValue = 0;
int sum = 0;
while ( (cin >> nextValue) \&\& (nextValue > 0) )
sum += nextValue;

```

\section*{A.5.2 The for Statement}

The for statement provides for counted loops and has the general form
```

for (initialize; test; update)
statement

```
where initialize, test, and update are expressions. Typically, initialize is an assignment expression that initializes a counter to control the loop. This initialization occurs only once. Then if test, which is usually a logical expression, is true, statement executes. The expression update executes next, usually incrementing or decrementing the counter. This sequence of events repeats, beginning with the evaluation of test, until the value of test is false.

For example, the following for statement displays the integers from 1 to n :
```

for (int counter = 1; counter <= n; counter++)
cout << counter << " ";
cout << endl; // This statement is always executed

```

If n is less than 1 , the for statement does not execute at all. Thus, the previous statements are equivalent to the following while loop:
```

int counter = 1;
while (counter <= n)
{
cout << counter << " ";
counter++;
} // end while
cout << endl; // This statement is always executed

```

In general, the logic of a for statement is equivalent to
```

initialize;

```
while (test)
\{
    statement;
    update;
\}

Note that in a for statement the first expression initialize must have either an arithmetic type or a pointer type. \({ }^{5}\) Note that char in the following example is considered an arithmetic type:
```

for (char ch = 'z'; ch >= 'a'; ch--)
// ch ranges from 'z' to 'a'

```

The initialize and update portions of a for statement each can contain several expressions separated by commas, thus performing more than one action. For example, the following loop raises a floating-point value to an integer power by using multiplication:
```

// Floating-point power equals floating-point x raised to int n;
// assumes int expon
for (power = 1.0, expon = 1; expon <= n; expon++)
power *= x;

```

Both power and expon are assigned values before the body of the loop executes for the first time. The comma here is an example of the comma operator, which evaluates its operand expressions from left to right.

When compared to a while statement, the for statement can make it easier to understand how the loop is controlled because the initialization, testing, and updating steps of the loop are consolidated into one statement. C++ programmers use for statements for loops that process collections or sequences of data.

\section*{A.5.3 The do Statement}

Use the do statement when you want to execute a loop's body at least once. Its general form is
```

do
statement
while (expression);

```

Here, statement executes until the value of expression is false.

For counted loops, a for statement is usually favored over the while statement

A do statement loops at least once

A for statement is equivalent to a while statement

An array is a container of data that has the same type

You can access array elements directly and in any order

Use an index to specify a particular element in an array

An array index has an integer value greater than or equal to 0

You can use an enumerator as an array index

For example, suppose that you execute a sequence of statements and then ask the user whether to execute them again. The do statement is appropriate, because you execute the statements before you decide whether to repeat them:
```

char response;
do
{
// A sequence of statements
cout << "Do it again?";
cin >> response;
} while ((response == 'Y') || (response == 'y') );

```

\section*{A. 6 Arrays and Vectors}

An array is a container of data items, or entries, that have the same data type. An array's memory locations, or elements, have an order: An array has a first element, a second element, and so on, as well as a last element. That is, an array has a finite, limited number of elements. Therefore, you must know the maximum number of elements needed for a particular array when you write your program and before you execute it. Because you can access the array elements directly and in any order, an array is a direct access, or random access, data structure.

\section*{A.6.1 One-Dimensional Arrays}

When you decide to use an array in your program, you must declare it and, in doing so, indicate the data type of its entries as well as its size. The following statements declare a one-dimensional array, maxTemps, which contains the daily maximum temperatures for a given week:
```

const int DAYS_PER_WEEK = 7;
doub7e maxTemps[DAYS_PER_WEEK];

```

The bracket notation [] declares maxTemps as an array. This array can contain at most seven floatingpoint values.

You can refer to any of the floating-point entries in maxTemps directly by using an expression, which is called the index, or subscript, enclosed in square brackets. In C++, array indices must have integer values in the range 0 to size -1 , where size is the number of elements in the array. The indices for maxTemps range from 0 to DAYS_PER_WEEK - 1. For example, the fifth element in this array is maxTemps [4]. If \(k\) is an integer variable whose value is 4 , maxTemps [k] is the fifth element in the array, and maxTemps \([\mathrm{k}+1]\) is the sixth element. Also, maxTemps \([\mathrm{k}++]\) accesses maxTemps \([\mathrm{k}]\) before adding 1 to \(k\). Note that you use one index to refer to an element in a one-dimensional array.

Figure A-6 illustrates the array maxTemps, which at present contains only five temperatures. The last value in the array is in maxTemp [4]; the values of maxTemps [5] and maxTemps [6] are not initialized and therefore are unknown.

You can use enumerators as indices because they have integer values. For example, consider the following definition:
\[
\text { enum Day \{SUN, MON, TUE, WED, THU, FRI, SAT\}; }
\]

Given this definition, maxTemps[THU] has the same meaning as maxTemps [4]. You can also use the enumerators within a loop that processes an array, as in the following for statement:
```

for (Day dayIndex = SUN; dayIndex <= SAT; dayIndex++)
cout << maxTemps[dayIndex] << endl;

```

\section*{FIGURE A-6 A one-dimensional array of at most seven elements}


Clearly, before you access an element of an array, you must assign it a value. You must assign values to array elements one at a time by using the previously described index notation. Note that, if a and \(b\) are arrays of the same type, the assignment \(a=b\) is illegal. \({ }^{6}\)

The data type of maxTemps is a derived type, which is a type that you derive from the fundamental types by using a declaration operator such as []. Naming a derived type by using a typedef is often useful. Thus, you can write
```

const int DAYS_PER_WEEK = 7;
typedef doub7e ArrayType[DAYS_PER_WEEK];
ArrayType maxTemps;

```
and make ArrayType available for use throughout your program.
Initialization. You can initialize the elements of an array when you declare it for the first time. For example,
```

doub7e maxTemps[DAYS_PER_WEEK] = {82.0, 71.5, 61.8, 75.0, 88.3};

```
initializes the first five elements of maxTemps to the values listed and the last two elements to zero.
Passing an array to a function. If you wanted a function that computed the average of the first \(n\) elements in a one-dimensional array, you could declare the function as
```

doub7e getAverageTemp(double temperatures[], int n);

```

Because the compiler does not know the number of items that the array can hold, you must also pass the function either the size of the array or the number of array items to process. Traditionally, the array is listed as the first parameter and the number of items as the second. You can invoke the function by writing, for example,
```

doub7e avg = getAverageTemp(maxTemps, 5);

```
where maxTemps is the previously defined array.
An array is never passed to a function by value, regardless of how you write its parameter. An array is always passed by reference. This restriction avoids the copying of perhaps many array entries. Thus, the function getAverageTemp could modify the elements of maxTemps, even though the array is an input to the function. To prevent such alteration, you can specify the array parameter as a constant reference parameter by preceding its type with const, as follows:
doub7e getAverageTemp(const doub7e temperatures[], int \(n\) );

You can initialize an array when you declare it

Arrays are always passed by reference to a function

\footnotetext{
6 C++ enables you to define your own array data type and array operators so that this assignment would be valid. To do so, you need to use classes (C++ Interlude 1) and overloaded operators (C++ Interlude 5).
}

An array can have more than one dimension

In a two-dimensional array, the first index represents the row, the second index represents the column

\section*{FIGURE A-7 A two-dimensional array}


\section*{A.6.2 Multidimensional Arrays}

You can use a one-dimensional array, which has one index, for a simple collection of data. For example, you can organize 52 temperatures linearly, one after another. A one-dimensional array of these temperatures can represent this organization.

You can also declare multidimensional arrays. You use more than one index to designate an element in a multidimensional array. Suppose that you wanted to represent the minimum temperature for each day during 52 weeks. The following statements declare a two-dimensional array, minTemps:
```

const int DAYS_PER_WEEK = 7;
const int WEEKS_PER_YEAR = 52;
doub7e minTemps[DAYS_PER_WEEK][WEEKS_PER_YEAR];

```

These statements specify the ranges for two indices: The first index can range from 0 to 6 , while the second index can range from 0 to 51 . Most people picture a two-dimensional array as a rectangular arrangement, or matrix, of elements that form rows and columns, as Figure A-7 indicates. The first dimension given in the definition of ArrayType is the number of rows. Thus, minTemps has 7 rows and 52 columns. Each column in this matrix represents the seven daily minimum temperatures for a particular week.

To reference an element in a two-dimensional array, you must indicate both the row and the column that contain the element. You make these indications of row and column by writing two indices, each enclosed in brackets. For example, minTemps [1] [51] is the element in the \(2^{\text {nd }}\) row and the \(52^{\text {nd }}\) column. In the context of the temperature example, this element contains the minimum temperature recorded for the \(2^{\text {nd }}\) day (Monday) of the \(52^{\text {nd }}\) week. The rules for the indices of a one-dimensional array also apply to the indices of multidimensional arrays.

As an example of how to use a two-dimensional array in a program, consider the following program segment, which determines the smallest value in the previously described array minTemps. We use enumerators to reference the days of the week.
```

enum Day {SUN, MON, TUES, WED, THURS, FRI, SAT};
// Initially, assume the lowest temperature is first in the array
doub7e lowestTemp = minTemps[0][0];
Day dayOfWeek = SUN;
int weekOfYear = 1;
// Search array for lowest temperature
for (int weekIndex = 0; weekIndex < WEEKS_PER_YEAR; weekIndex++)
{
for (Day dayIndex = SUN; dayIndex <= SAT; dayIndex++)
{
if (lowestTemp > minTemps[dayIndex][weekIndex])
{
lowestTemp = minTemps[dayIndex][weekIndex];
dayOfWeek = dayIndex;
weekOfYear = weekIndex + 1;
} // end if
} // end for
} // end for
// At this point, lowestTemp is the smallest value in minTemps and
// occurs on the day and week given by dayOfWeek and weekOfYear,
// that is, lowestTemp == minTemps[dayOfWeek][weekOfYear - 1].

```

Although you can declare arrays with more than two dimensions, generally more than three dimensions is unusual. The techniques for working with such arrays, however, are analogous to those for two-dimensional arrays.

Programming Tip: When referencing an element of a multidimensional array, do not use comma-separated indices. For example, myArray \([3,6]\) does not reference the array element myArray [3] [6]. The expression 3,6 is a comma expression whose value is that of the last item listed, namely 6 . Thus, although myArray \([3,6]\) is legal, its meaning is myArray [6], which references the element myArray[0] [6].

Initialization. You can initialize the elements of a two-dimensional array just as you initialize a onedimensional array. You list the initial values row by row. For example, the statement
```

int x[2][3] = { {1, 2, 3},
{4, 5, 6} }; // 2 rows, 3 columns

```
initializes the two-dimensional array x so that it appears as
\(\begin{array}{ll}1 & 2\end{array}\)
456
That is, the statements initialize the elements \(x[0][0], x[0][1], x[0][2], x[1][0], x[1][1]\), and \(x[1][2]\) in that order. In general, when you assign initial values to a multidimensional array, it is the last, or rightmost, index that increases the fastest.

\section*{A.6.3 Vectors}

Another container that can hold data items of the same type is the vector. A vector is similar to a one-dimensional array, but vectors provide additional features for the programmer not found in a simple array. A vector is an object of a standard \(C++\) class named vector. This class is a part of the

\section*{Access vector} entries using [] just as you would an array
size returns the number of elements in the vector
push_back places its argument into a new element at the back of a vector
pop_back removes the last element from a vector

Standard Template Library, or STL. The STL is a collection of data types you can use in your programs. These data types are not part of the official C++ language, but they have been added to the built-in data types. Section A. 9 discusses libraries such as the STL and their usefulness.

To use a vector in your program, you must begin it with the following statements:
\#include <vector>
using namespace std;
You can declare a vector in one of three ways:
- If you know how many elements you want in the vector, you can place the type of data it will hold in angle brackets and the number of elements in parentheses:
```

vector<doub7e> firstVector(10); // Vector to hold 10 doub7es
vector<string> myVector(12); // Vector to hold 12 strings

```

The size you specify when declaring the vector is only its initial size. As you will see, a vector can grow in size when you add entries.
- You can place initial values into a vector when you declare it, by writing a second argument:
```

vector<int> intVector(5, -1); // Vector to hold 5 integers, initialized to -1

```

When the elements in the vector are allocated, they are given the value of the second argument.
- You can also create an empty vector-a vector with no elements-by omitting its size and the parentheses:
```

vector<char> letterVector; // An empty vector of characters

```

You can store or access items in a vector by using the [] operator, just as you would when using an array. As for an array, the subscripts that identify elements in a vector start at 0 and go to \(s-1\), where \(s\) is the current size of the vector. The following statements are examples of accessing an existing value in a vector and changing the value of an existing entry:
```

doub7e x = firstVector[5]; // P7aces sixth entry in x
myVector[3] = "This is a sample string."; // Sets fourth entry's value

```

Using [] stores a value in an existing element. If you are not sure how many elements the vector has, you can call the method size, as in the following example:
cout << intVector.size() << endl; // Displays the number of elements in intVector
By calling the size method, you can determine whether the vector is full. This is an important advantage that a vector has over an array.

If the vector is either full or has no elements-that is, if it was created without elements, as letterVector was previously-you can still add new values by using the method push_back. The push_back method accepts an argument and adds it after the last element of the vector. In other words, it pushes the value onto the back of the vector.

Earlier, we declared a ten-element vector firstVector that could hold data of type double. If that vector was full, and we needed to add the additional values 2.3 and 3.4 , we could use the push_ back method:
firstVector.push_back(2.3); // Grow vector and store value
firstVector.push_back(3.4); // Grow vector and store value
At this point, calling the size method would return 12, since two additional elements have been added to the vector.

You also can reduce the size of a vector by removing either its last element or all of its elements. To remove only the last element, you can use the pop_back method. This method shortens
the vector but does not give you the entry in the removed element. You must save that entry before calling pop_back. For example, the following statements determine the current size of the vector myVector, save the value in the last element, and then remove the last element from the vector:
```

int length = myVector.size(); // Find the current number of elements
string last = myVector[1ength - 1]; // Get the string in the last element
// (subscripts start at 0)
myVector.pop_back(); // Reduce the number of elements

```

A subsequent call to myVector.size() would return length - 1, since the last element was removed.

To remove all elements from a vector and leave the vector empty, you use the method clear:
myVector.clear(); // myVector is now an empty vector
clear removes all elements from a vector so it has a size of 0

After a vector has been cleared, you must use the method push_back to add new entries.
\(\mathrm{C}++\) Interlude 7 provides more information about vector, including these and other methods.

\section*{A. 7 Strings}

Earlier, you saw that \(\mathrm{C}++\) provides literal character strings such as
"This is a string."
You can declare and use variables that contain such strings, and then manipulate the strings as naturally as you manipulate integers by using familiar operators. Our presentation includes only some of the possible operations on strings.

The C++ Standard Library provides the data type string. To use this library, you include the statements
\#include <string>
using namespace std;
in your program.
You can declare a string variable title and initialize it to the empty string by writing
```

string title;

```

You can initialize a string variable to a string literal when you declare it by writing
```

string title = "Wal1s and Mirrors";

```

You can subsequently assign another string to tit 1 e by using an assignment statement such as
```

title = "J Perfect's Diary";

```

In each of the previous examples, title has a length of 17 . You use either of the methods 1 ength or size to determine the current length of a string. Thus, title.length() and title.size() are each 17 .

You can reference the individual characters in a string by using the same index notation that you use for an array. Thus, in the previous example, title[0] contains the character J and title[16] contains the character \(y\).

You can compare strings by using the familiar comparison operators. Not only can you see whether two strings are equal, but you can also discover which of two strings comes before the other.

Examples of true expressions

Use substr to access part of a string

You can use << to display a string

You can use >> to read a string without whitespace

The ordering of two strings is analogous to alphabetic ordering, but you use the ASCII table instead of the alphabet. Thus, the following relationships are all true:
```

"dig" < "dog"
"Star" < "star" (because 'S' < 's')
"start" > "star"
"d" > "abc"

```

You can concatenate two strings to form another string by using the + operator. That is, you place one string after another to form another string. For example, if
```

string str1 = "Com";

```
the statements
```

string str2 = str1 + "puter";
str1 += "puter";

```
assign the string "Computer" to each of str2 and str1. Similarly, you can append a single character to a string, as in
```

str1 += 's';

```

You can manipulate a portion of a string by using the method substr(position, length)

The first argument specifies the position of the beginning of the substring (remember that 0 is the position of the first character in the string). The second argument is the length of the substring. For example,
```

title.substr(2, 7)

```
is the string "Perfect".
To perform input and output with \(\mathrm{C}++\) strings, you must include the library iostream by writing the statement
```

\#include <iostream>

```
in your program. For example, you then can display the contents of a string variable by executing
```

title = "Walls and Mirrors";
cout << title << "\n";

```

The result is Walls and Mirrors. The operator << writes the entire string, including the blanks You can read a string of characters into a string variable. For example, when the statement
```

cin >> title;

```
reads the data line
```

Jamie Perfect's Diary

```
it assigns the string "Jamie" to title. Whitespace in the input line terminates the read operation for a string. To read the entire line of input, including its blank characters, you write
```

getline(cin, title);

```

\section*{A. 8 Classes}

Object-oriented programming, or OOP, views a program not as a sequence of actions but as a collection of components called objects that interact to produce a certain result. A group of objects of the same kind belong to a class, which is a programming construct that defines the object's data type. Chapter 1 talks more about OOP; here we want to discuss how to write a class in C++.

An object contains data and can perform certain operations on or with that data. The class associated with a particular object describes its data and its operations. That is, a class is like a blueprint for creating certain objects. An object's operations, or behaviors, are defined within the class by methods, which are simply functions within a class. These methods, together with the class's data are known as the class's members.

We could use a ball as an example of an object. Because thinking of a basketball, volleyball, tennis ball, or soccer ball probably suggests images of the game rather than the object itself, let's abstract the notion of a ball by picturing a sphere. A sphere of a given radius has attributes such as volume and surface area. A sphere as an object should be able to report its radius, volume, surface area, and so on. That is, the sphere object has methods that return such values.

In C++, a class has the following form:
```

class Sphere
{
private:
// Definitions of private methods
public:
// Definitions of public methods
. . .
};

```
    // Declarations of private data members and methods

By default, all members in a class are private - they are not directly accessible by any program that uses the class-unless you designate them as public. However, explicitly indicating the private and public portions of a class is a good programming practice and one that we will follow in this book. You should always declare a class's data members as private.

Most methods are public, but private methods-which only the class can call-can be helpful, as you will see. The definition of a class's method can call any of the class's other methods or use any of its data members, regardless of whether they are private or public.

Classes have special methods, called constructors and destructors, for the creation and destruction of its objects. A constructor creates and initializes new objects, or instances, of a class. A destructor destroys an instance of a class, when the object's lifetime ends. A typical class has several constructors, but only one destructor. For many classes, you can omit the destructor. In such cases, the compiler will generate a destructor for you. For now, the compiler-generated destructor is sufficient. C++ Interlude 2 discusses how and why you would write your own destructor.

In C++, a constructor has the same name as the class. Constructors have no return type-not even void-and cannot use return to return a value. Constructors can have arguments. We discuss constructors in more detail shortly, after we look at an example of a class definition.

\section*{A.8.1 The Header File}

You should place each class definition in its own header file or specification file-whose name by convention ends in .h. The header file Sphere.h shown in Listing A-2 contains a class definition for sphere objects.

A C++ class defines a new data type

An object is an instance of a class

A constructor creates and initializes an object

A destructor destroys an object

Comments in the header file specify the methods

\section*{LISTING A-2 The header file Sphere. \(h\)}
```

/** @file Sphere.h */
const doub7e PI = 3.14159;
/** Definition of a class of Spheres. */
class Sphere
{
private:
doub7e theRadius; // The sphere's radius
public:
/** Defau7t constructor: Creates a sphere and initializes
its radius to a default value.
Precondition: None.
Postcondition: A sphere of radius 1 exists. */
Sphere();
/** Constructor: Creates a sphere and initializes
its radius.
Precondition: initialRadius is the desired radius.
Postcondition: A sphere of radius initialRadius exists. */
Sphere(doub7e initialRadius);
/** Sets (alters) the radius of this sphere.
Precondition: newRadius is the desired radius.
Postcondition: The sphere's radius is newRadius. */
void setRadius(double newRadius);
/** Gets this sphere's radius.
Precondition: None.
Postcondition: Returns the radius. */
doub7e getRadius() const;
/** Gets this sphere's diameter.
Precondition: None.
Postcondition: Returns the diameter. */
doub7e getDiameter() const;
/** Gets this sphere's circumference.
Precondition: PI is a named constant.
Postcondition: Returns the circumference. */
doub7e getCircumference() const;
/** Gets this sphere's surface area.
Precondition: PI is a named constant.
Postcondition: Returns the surface area. */
doub7e getArea() const;
/** Gets this sphere's volume.
Precondition: PI is a named constant.
Postcondition: Returns the volume. */
double getVolume() const;
}; // end Sphere
// End of header file.

```

You should always place a class's data members within its private section. Typically, you provide methods-such as setRadius and getRadius-to access the data members. In this way, you control how and whether the rest of the program can access the data members. This design principle should lead to programs that not only are easier to debug, but also have fewer logical errors from the beginning.

Some method declarations, such as
```

doub7e getRadius() const;

```
are tagged with const. Such methods cannot alter the data members of the class. Making getRadius a const method is a fail-safe technique that ensures that it will only return the current value of the sphere's radius, without changing it.

\section*{A.8.2 The Implementation File}

Let's begin implementing the class Sphere by examining its constructors.
Constructors. A constructor allocates memory for an object and can initialize the object's data to particular values. A class can have more than one constructor, as is the case for the class Sphere.

The first constructor in Sphere is the default constructor:
Sphere() ;
A default constructor by definition has no arguments. Typically, a default constructor initializes data members to values that the class implementation chooses. For example, the implementation

Sphere::Sphere()
\{
theRadius \(=1.0\);
\} // end default constructor
sets theRadius to 1.0. Chapter 3 will show you another way to initialize data members within constructors that is preferable to using assignment statements.

Notice the qualifier Sphere: : that precedes the constructor's name. When you implement any method, you qualify its name with its class type followed by the scope resolution operator : : to distinguish it from other methods that might have the same name.

When you declare an instance of the class, you implicitly invoke a constructor. For example, the statement

Sphere unitSphere;
invokes the default constructor, which creates the object unitSphere and sets its radius to 1.0. Notice that you do not include parentheses after unitSphere.

The next constructor in Sphere is
Sphere(doub7e initialRadius);
It creates a sphere object of radius initialRadius. This constructor needs only to initialize the private data member theRadius to initialRadius. Its implementation is
```

Sphere::Sphere(doub7e initialRadius)
{
theRadius = initialRadius;
} // end constructor

```

You implicitly invoke this constructor by writing a declaration such as
```

Sphere mySphere(5.1);

```

A class's data members should be private
const methods cannot change a class's data members

A default constructor has no arguments

In this case, the object mySphere has a radius of 5.1.
We can make the previous constructor ensure that the given radius is not negative by writing its definition as follows:
```

Sphere::Sphere(doub7e initia1Radius)
{
if (initialRadius > 0)
theRadius = initialRadius;
else
theRadius = 1.0;
} // end constructor

```


Note: If you omit all constructors from your class, the compiler will generate a default constructor-that is, one with no arguments-for you. A compiler-generated default constructor, however, might not initialize data members to values that you will find suitable. If you define a constructor that has arguments but you omit the default constructor, the compiler will not generate one for you. Thus, you will not be able to write statements such as

Sphere defaultSphere;

The implementation file contains the definitions of the class's methods

Typically, you place the implementation of a class's constructors and other methods in an implementation file whose name ends in . cpp. Listing A-3 contains an implementation file for the class Sphere. Notice that within the definition of a method, you can reference the class's data member or invoke its other methods without preceding the member names with Sphere: : In particular, notice how the constructor calls the method setRadius to avoid duplicating the code that ensures a positive radius.
```

LISTING A-3 The implementation file Sphere.cpp
/** @file Sphere.cpp */
\#include "Sphere.h" // Include the header file
using namespace std;
Sphere::Sphere()
{
theRadius = 1.0;
} // end default constructor
Sphere::Sphere(doub7e initialRadius)
{
setRadius(initialRadius); // Sphere:: not needed here
} // end constructor
void Sphere::setRadius(doub7e newRadius)
{
if (newRadius > 0)
theRadius = newRadius;
e7se
theRadius = 1.0;
} // end setRadius

```
```

double Sphere::getRadius() const
{
return theRadius;
} // end getRadius
doub7e Sphere::getDiameter() const
{
return 2.0 * theRadius;
} // end getDiameter
doub7e Sphere::getCircumference() const
{
return PI * getDiameter();
} // end getCircumference
doub7e Sphere::getArea() const
{
return 4.0 * PI * theRadius * theRadius;
} // end getArea
doub7e Sphere::getVolume() const
{
doub7e radiusCubed = theRadius * theRadius * theRadius;
return (4.0 * PI * radiusCubed) / 3.0;
} // end getVolume
// End of implementation file.

```

\section*{Note: Local variables}

You should distinguish between a class's data members and any local variables that the implementation of a method requires. It is inappropriate for such local variables to be data members of the class.

\section*{A.8.3 Using the Class Sphere}

The following simple program demonstrates the use of the class Sphere:
```

\#include <iostream>
\#include "Sphere.h"
using namespace std;
int main()
{
Sphere unitSphere; // Radius is 1.0
Sphere mySphere(5.1); // Radius is 5.1
unitSphere.displayStatistics();
mySphere.setRadius(4.2); // Resets radius to 4.2
cout << mySphere.getDiameter() << endl;
return 0;
} // end main

```

An object such as mySphere can, on request, reset the value of its radius; return its radius; and compute its diameter, surface area, circumference, and volume. These requests to an object are called

A local variable such as radiusCubed should not be a data member

A class derived from the class Sphere
messages and are simply calls to methods. Thus, an object responds to a message by acting on its data. To invoke an object's method, you qualify the method's name-such as setRadius-with the object variable-such as mySphere.

Notice that the previous program included the header file Sphere.h, but did not include the implementation file Sphere. cpp. You compile a class's implementation file separately from the program that uses the class. The way you tell the operating system where to locate the compiled implementation depends on the particular system. Section A. 9 of this appendix and C++ Interlude 1 provide more information about header and implementation files.

The previous program is an example of a client of a class. A client of a particular class is simply a program or module that uses the class. We will reserve the term user for the person who uses a program.

\section*{A.8.4 Inheritance}

A brief discussion of inheritance is provided here, because it is a common way to create new classes in \(\mathrm{C}++\). Further discussions of inheritance occur as needed throughout the book.

Suppose we want to give our spheres a color, knowing that we have already developed the class Sphere. Instead of writing an entirely new class of spheres that have a color, we can reuse the Sphere implementation and add color characteristics and operations by using inheritance. Here is a declaration of the class SphereInColor that uses inheritance:
```

\#include "Sphere.h"
enum Color {RED, BLUE, GREEN, YELLOW};
class SphereInColor : public Sphere
{
private:
Color sphereColor;
public:
SphereInColor(Color initialColor);
SphereInColor(Color initialColor, double initialRadius);
void setColor(Color newColor);
Color getColor() const;
}; // end SphereInColor

```

The class Sphere is called the base class or superclass, and SphereInColor is called the derived class or subclass of the class Sphere.

Any instance of the derived class is also considered to be an instance of the base class and can be used in a program anywhere that an instance of the base class can be used. Also, when the keyword pub7ic precedes the name of the base class in the new class's header, any of the publicly defined methods or data members that can be used with instances of the base class can be used with instances of the derived class. The derived class instances also have the additional methods and data members that are publicly defined in the derived class definition.

The implementation of the methods for the class SphereInColor is as follows:
```

SphereInColor::SphereInColor(Color initialColor): Sphere()
{
sphereColor = initialColor;
} // end constructor
SphereInColor::SphereInColor(Color initialColor, double initialRadius)
{
sphereColor = initialColor;
} // end constructor

```
```

void SphereInColor::setColor(Color newColor)
{
sphereColor = newColor;
} // end setColor
Color SphereInColor::getColor() const
{
return sphereColor;
} // end getColor

```

Notice how the constructors for the class SphereInColor invoke the base-class constructors Sphere() and Sphere(initialRadius). The derived class needs the initialization of the data members in the base class that the base-class constructors can provide. The derived-class constructors then add initializations that are specific to the derived class.

Here is a function that uses the class SphereInColor:
```

void useSphereInColor()
{
SphereInColor bal1(RED);
ba11.setRadius(5.0);
cout << "The ball diameter is " << bal1.getDiameter();
ba11.setColor(BLUE);
} // end useSphereInColor

```

This function uses the constructor and the method setColor from the derived class SphereInColor. It also uses the methods setRadius and getDiameter that are defined in the base class Sphere.

\section*{A. 9 Libraries}

One of the advantages of modular programming is that you can implement modules independently of other modules. You might also find it possible for several different programs to use a particular module. As a result, you can build a library of modules that you can include in future programs.

Any library-a C++ standard library or one that you write-has a corresponding header that provides information about the contents of the library. For standard libraries, the header is simply an abstraction that the compiler may map to a filename or handle in a different manner. Thus, when using the standard libraries, you do not see the. h extension that ends the names of our own header files.

You have already seen some standard libraries, such as the one that provides input and output services. To use the modules contained in a library, you use the include directive with the name of the header associated with the library. For example, you write
```

\#include <iostream>
using namespace std;

```

\section*{Appendix H provides a list of available headers.}

User-defined libraries are typically organized into two files. One file, the header file, contains a definition for each class in the library that is available to your program. This file could also contain, for example, function declarations, constant definitions, typedef statements, enumerations, and other include statements. By convention, the name of a header file associated with a user-defined library ends in .h. The other file-the implementation file-contains definitions of the class methods that the header file declares. Typically, the name of an implementation file ends in .cpp.

The assumption, of course, is that the files are in source form-that is, they need to be compiled. It certainly would be more efficient to compile the method definitions once, independently of any

An instance of a derived class can invoke public methods of the base class
particular program, and then later merge the results of the compilation with any program that you desire. In fact, you should compile the implementation file and then include the header file in source form in your program by using an include directive such as
```

\#include "MyHeader.h"

```

You use double quotes instead of angle brackets to enclose the name of a header file that you have written. The mechanics of incorporating the compiled implementation file into your program are system dependent.

Thus, your program can use previously compiled C++ statements, which are no longer available to you in source form. Maybe you did not even write these statements, just as you did not write the standard \(\mathrm{C}++\) functions such as sqrt. That is, you use a library in the same spirit in which you use standard functions. Because the header file indicates what is available to you, you must think of a library in terms of what it can do for you and not how it is implemented. You should think of all of your modules in this way, even if you eventually implement them yourself.

\section*{Summary}
1. Each comment line in \(\mathrm{C}++\) begins with two slashes // and continues until the end of the line.
2. A C++ identifier is a sequence of letters, digits, and underscores that must begin with either a letter or an underscore.
3. You can use a typedef statement to declare new names for data types. These names are simply synonyms for the data types; they are not new data types.
4. You define named constants by using a statement of the form
const type identifier = value;
5. Enumeration provides another way to name integer constants and to define an integral data type, as in
enum Day \{SUN, MON, TUE, WED, THU, FRI, SAT\};
6. \(\mathrm{C}++\) uses short-circuit evaluation for expressions that contain the logical operators \& (and) and || (or). That is, evaluation proceeds from left to right and stops as soon as the value of the entire expression is apparent.
7. The output operator \(\ll\) places a value into an output stream, and the input operator \(\gg\) extracts a value from an input stream. You can imagine that these operators point in the direction of data flow. Thus, in cout << myVar, the operator points away from the variable myVar-data flows from myVar to the stream-whereas in cin >> myVar, the operator points to the variable myVar-data flows from the stream into myVar.
8. The general form of a function definition is
type name (parameter-declaration-list)
\{
body
\}
A valued function returns a value by using the return statement. Although a void function does not return a value, it can use return to exit.
9. When invoking a function, the actual arguments must correspond to the parameters in number, order, and type.
10. A function makes local copies of the values of any arguments that are passed by value. Thus, the arguments remain unchanged by the function. Such arguments are, therefore, input arguments. A function does not copy arguments that are passed by reference. Rather, it references the actual argument locations whenever the parameters appear in the function's definition. In this way, a function can change the values of the arguments, thus implementing output arguments. However, a function does not copy and cannot change a constant reference argument. If copying an input argument would be expensive, make it a constant reference argument instead of a value argument.
11. The general form of the if statement is
```

if (expression)
statement
else
statement }

```

If expression is true, statement \(_{1}\) executes; otherwise statement \({ }_{2}\) executes.
12. The general form of the switch statement is
```

switch (expression)
{
case constant 1
statement
break;
case constant n
statement
break;
default:
statement
}

```

The appropriate statement executes according to the value of expression. Typically, break (or sometimes return) follows the statement or statements after each case. Omitting break causes execution to continue to the statement(s) after the next case.
13. The general form of the while statement is
```

while (expression)
statement

```

As long as expression is true, statement executes. Thus, it is possible that statement never executes.
14. The general form of the for statement is
for (initialize; test; update)
statement
where initialize, test, and update are expressions. Typically, initialize is an assignment expression that occurs only once. Then if test, which is usually a logical expression, is true, statement executes. The expression update executes next, usually incrementing or decrementing a counter. This sequence of events repeats, beginning with the evaluation of test, until test is false.
15. The general form of the do statement is

> do
statement
while (expression);
Here, statement executes until the value of expression is false. Note that statement always executes at least once. Also note the required semicolon.
16. An array is a container of items that have the same data type. You can refer to these items by using an index that begins with zero. Arrays are always passed to functions by reference.
17. A string is a sequence of characters. You can manipulate the entire string, a substring, or the individual characters.
18. You must be careful that an array index does not exceed the size of the array. \(\mathrm{C}++\) does not check the range of array indices. Similar comments apply to strings.
19. An object encapsulates both data and operations on that data. In \(\mathrm{C}++\), objects are instances of a class, which is a programmer-defined data type.
20. A \(\mathrm{C}++\) class contains at least one constructor, which is an initialization method, and a destructor, which is a cleanup method that destroys an object when its lifetime ends.
21. If you do not define a constructor for a class, the compiler will generate a default constructor- that is, one without arguments-for you. If you do not define a destructor, the compiler will generate one for you. Chapter 4 describes when you need to write your own destructor.
22. Members of a class are private unless you designate them as public. The client of the class - that is, the program that uses the class-cannot use members that are private. However, the implementations of methods can use them. You should make the data members of a class private and provide public methods to access some or all of the data members.
23. Because certain classes have applications in many programs, you should take steps to facilitate their use. You can define and implement a class within header and implementation files, which a program can include when it needs to use the class.
24. A typical \(\mathrm{C}++\) program uses header files that you incorporate by using the include directive. A header file contains class definitions, function declarations, constant definitions, typedef statements, enumerations, and other include statements. The program might also require an implementation file of function definitions that have been compiled previously and placed into a library. The operating system locates the required implementation file and combines it with the program in ways that are system dependent.

\section*{Important Themes in Programming}

\section*{Appendix}

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People read programs, too

Key programming themes

Modularity facilitates programming

Modularity isolates errors

\section*{Prerequisites}

\author{
Appendix A Review of C++ Fundamentals \\ Chapter 1 Data Abstraction: The Walls
}

W
hat are the specific characteristics of good solutions? How can you construct good solutions? This appendix provides some answers to these very difficult questions.

The programming themes that we will discuss should be familiar to you. However, the novice programmer usually does not truly appreciate their importance. After the first course in programming, many students still simply want to "get the thing to run." The discussion that follows should help you realize just how important these themes really are.

One of the most widespread misconceptions held by novice programmers is that a computer program is "read" only by a computer. As a consequence, they tend to consider only whether the computer is able to "understand" the program-that is, does the program compile, execute, and produce the correct output? The truth is, of course, that other people often must read and modify programs. In a typical programming environment, many individuals share a program. One person may write a program, which other people use in conjunction with other programs written by other people, and a year later, a different person may modify the program. It is therefore essential that you take great care to design a program that is easy to read and understand.

You should always keep in mind the following seven themes of programming: modularity, style, modifiability, ease of use, fail-safe programming, debugging, and testing.

\section*{B. 1 Modularity}

As this book continually emphasizes, you should strive for modularity in all phases of the problem-solving process, beginning with the initial design of a solution. Many programming tasks become more difficult as the size and complexity of a program grows. Modularity slows the rate at which the level of difficulty grows. More specifically, modularity has a favorable impact on the following aspects of programming:
- Constructing the program. The primary difference between a small modular program and a large modular program is simply the number of modules each contains. Because the modules are independent, writing one large modular program is not very different from writing many small, independent programs, although the interrelationships among modules make the design much more complicated. On the other hand, working on a large nonmodular program is more like working on many interrelated programs simultaneously. Modularity also permits team programming, in which several programmers work independently on their own modules before combining them into one program.
- Debugging the program. Debugging a large program can be a monstrous task. Imagine that you type a 10,000-line program and eventually get it to compile. Neither of these tasks would be much fun. Now imagine that you execute your program, and after a few hundred lines of output, you notice an incorrect number. You should anticipate spending the next day or so tracing through the intricacies of your program before discovering a problem such as an array index that is too large.

A great advantage of modularity is that the task of debugging a large program is reduced to one of debugging many small programs. When you begin to code a module, you should be almost certain that all other modules coded so far are correct. That is, before you consider a module finished, you should test it extensively, both separately and in context with the other modules, by calling it with actual arguments carefully chosen to induce all possible behaviors of the modules. If this testing is done thoroughly, you can feel fairly sure that any problem is a result of an error in the last module added. Modularity isolates errors.

Formal theoretical techniques are available to verify the correctness of a program. Modular programs are amenable to this verification process.
- Reading the program. A person reading a large program may get lost in its details. Just as a modular design helps the programmer cope with the complexities of solving a problem, so too does a modular program help its reader understand how the program works. A modular program is easy to follow because the reader can get a good idea of what is going on without reading any of the code. A well-written function can be understood fairly well from only its name, initial comments, and the names of the other functions that it calls. Readers of a program need to study actual code only if they require a detailed understanding of how the program operates. Program readability is discussed further in the next section on style.
- Modifying the program. Modifiability is discussed in greater detail in Section B. 3 of this appendix, but as the modularity of a program has a direct bearing on its modifiability, a brief mention is appropriate here. A small change in the requirements of a program should require only a small change in the code. If this is not the case, it is likely that the program is poorly written and, in particular, that it is not modular. To accommodate a small change in the requirements, a modular program usually requires a change in only a few of its modules, particularly when the modules are independent (referred to as being loosely coupled) and each module performs a single well-defined task (referred to as being highly cohesive). This point is of particular importance with iterative development techniques, which produce changes in program requirements at every iteration, as Appendix D will discuss.

When making changes to a program, it is best to make a few at a time. By maintaining a modular design, you can reduce a fairly large modification to a set of small and relatively simple modifications to isolated parts of the program. Modularity isolates modifications.
- Eliminating redundant code. Another advantage of modular design is that you can identify a computation that occurs in many different parts of the program and implement it as a function or class method. Thus, the code for the computation appears only once, resulting in an increase in both readability and modifiability.

\section*{B. 2 Style}

We now consider the following seven issues of style in programming: the use of private data members, the proper use of reference arguments, the proper use of methods, avoidance of global variables in modules, error handling, readability, and documentation. Admittedly, much of the following discussion reflects the personal taste of the authors; certainly other good programming styles are possible. Most organizations that hire programmers publish style guides, so that each programming team can produce code that is easily read and maintained by other teams.

\section*{B.2.1 Use of Private Data Members}

Each object has a set of methods that represents the operations that the object can perform. The object also contains data members for storing information. You should hide the exact representation of these data members from modules that use the object by making all of the data members private. Doing so supports the principle of information hiding. The details of the object's implementation are hidden from view, with methods providing the only mechanism for getting information to and from the object. When the only operations involved with a particular data member are retrieve and modify, the object should provide a simple method-called an accessor-that returns the value of the data member and another method-called a mutator-that sets the value of the data member. For example, a Person object could provide access to the data member theName through the methods getName to return the person's name and setName to change the person's name.

Modular programs are easy to read

Modularity isolates modifications

Modularity eliminates redundancies

Seven issues of style

Data members should always be private

\section*{Reference} arguments return values from a method

When copying the argument is expensive, use pass-by-constantreference instead of pass-by-value

To reduce coupling, restrict what a method can do

Building is coupled to Lamp

\section*{B.2.2 Proper Use of Reference Arguments}

A method interacts, in a controlled fashion, with the rest of the program via its arguments. Value parameters, which are the default in \(\mathrm{C}++\) when you do not write \& after the parameter's data type, pass values into the method. The value of each argument corresponding to a value parameter is copied into the parameter, which is local to the method. Any change that the method makes to these parameters is not reflected in the actual arguments back in the calling program. This communication between the calling program and the method is one-way and supports the notion of an isolated module. Thus, you should use value arguments when possible. However, passing large objects by value is not very efficient because of the copying required.

How and when is it appropriate to use reference parameters? Since a reference parameter becomes an alias (another name) for the corresponding argument, changes to the parameter within the method also occur in the argument in the calling module. The obvious situation requiring reference arguments is when a method needs to return several values to the calling module. Whenever you use a parameter to pass a value out of a method, the parameter must be a reference parameter.

Suppose that you are writing a method that requires access to a large object x , whose value should not be altered. You consider passing \(x\) by value; however, to increase the efficiency of your program, you would like to avoid the computer-time and storage requirements of making a copy of \(x\). So what about passing \(x\) by reference? The problem with passing \(x\) by reference is that it conveys misinformation about the method's relation to the rest of the program. Since a reference parameter is used to communicate a value from the method back to its calling module, the program is more difficult to read if the parameter's value remains unchanged. The program also is more prone to errors if modifications are required. The situation is analogous to using a variable whose value never changes when you really should use a constant. The solution is to precede the parameter's declaration with const, which prevents the method from changing the corresponding argument. Thus, we have the efficiency of pass-by-reference (no copy is made) with the protection of pass-by-value (the argument is input only). This is called pass-by-constant-reference. If objects are large and complex, use pass-by-constant-reference; otherwise pass-by-value works fine.

\section*{B.2.3 Proper Use of Methods}

To reduce coupling among modules, you should restrict the calls that a method can make. A method should call only other methods:
- Defined within the same class
- Of argument objects
- Of objects created within the method
- Of objects contained within the same class as data members

For example, suppose we have the following code for three classes that represent a building, an appliance, and a lamp, respectively:
```

// Original Solution
class Building
{
private:
Appliance appliance;
public:
void turnOnLight()
{
appliance.getLamp().turnOn(); // Improper method use
} // end turnOnLight
}; // end Building

```
```

class Appliance
{
private:
Lamp light;
public:
Lamp getLamp()
{
return light;
} // end getLamp
}; // end Appliance
class Lamp
{
public:
void turnOn();
}; // end Lamp

```

Here, Building's method turnOnLight fails to follow these rules. It calls the method turnOn from the Lamp class, although a Lamp object is not part of the Building class and has not been passed to the turnOnLight method as an argument. This code has coupled the Building class to the Lamp class.

A much better design would decouple the Building and Lamp classes, as follows:
```

// Revised Solution
class Building
{
private:
Appliance appliance;
public:
void turnOnLight()
{
app7iance.turnOnLamp();
} // end turnOnLight
}; // end Building
class Appliance // Improved Appliance class
{
private:
Lamp light;
public:
void turnOnLamp()
{
1ight.turnOn();
} // end turnOnLamp
}; // end Appliance
class Lamp
{
public:
void turnOn();
}; // end Lamp

```

Building and Lamp are not coupled

Now the building asks the appliance to turn on the lamp, and the appliance turns on its light.
Suppose that the lamp's switch is replaced with a dimmer-type switch. We would need to give the turnOn method in the Lamp class an indication of how much to dim the light. Finding the call to turnOn in the original code segment might be difficult, because the Building class does not have explicit access to a Lamp object. You could search the Building class for instances of Lamp, and not find any. This type of dependency makes a solution more difficult to modify.

In the revised code segment, the Appliance class has a Lamp object as a data member. You can easily find the calls to this object, if necessary.

\section*{B.2.4 Avoidance of Global Variables in Modules}

Do not use global variables

In case of an error, methods should return a value or throw an exception, but not display a message

One of the main advantages of the concepts of encapsulation and information hiding is the creation of isolated modules. This isolation is sacrificed when a module accesses a global variable, because the effects of a module's action are no longer self-contained or limited to output arguments. That is, such a module has a side effect. Hence, the isolation of both errors and modifications is greatly compromised when global variables appear in modules.

\section*{B.2.5 Error Handling}

A program should check for errors in both its input and its logic and attempt to behave gracefully and consistently when it encounters them. A method should check for certain types of errors, such as invalid input or argument values. What action should a method take when it encounters an error? Depending on context, the appropriate action in the face of an error can range from ignoring erroneous data and continuing execution to terminating the program. A common technique is for a method to return a boolean value to the calling module to indicate that it has encountered an error. Thus, the method leaves error handling to the calling module. In general, methods should either return a value or throw an exception instead of displaying a message when an error occurs.

\section*{B.2.6 Readability}

For a program to be easy to follow, it should have a good structure and design, a good choice of identifiers, good indentation and use of blank lines, and good documentation. These points are demonstrated in the programs throughout this book. You should avoid clever programming tricks that save a little computer time at the expense of much human time.

Choose identifiers that describe their purpose-that is, that are self-documenting. Distinguish between keywords, such as int, and user-defined identifiers. This book uses the following conventions:
- Keywords are lowercase and appear in color.
- Names of standard functions are lowercase.
- User-defined identifiers use both uppercase and lowercase letters, as follows:
\(\square\) Class names are nouns, with each word in the identifier capitalized.
\(\square\) Method names and function names are verbs or action phrases, with the first letter lowercase and subsequent internal words capitalized.
\(\square\) Variables begin with a lowercase letter, with subsequent words in the identifier capitalized.
\(\square\) Data types declared in a typedef statement and names of enumerations each begin with an uppercase letter.
\(\square\) Named constants and enumerators are entirely uppercase and use underscores to separate words.
- Two other naming conventions are suggested as a learning aid:
\(\square\) Data types declared in a typedef statement end in Type.
- Exception names end in Exception.

Use a good indentation style to enhance the readability of a program. The layout of a program should make it easy for a reader to identify the program's modules. Use a blank line between method or function definitions. Also, within these definitions, you should indent individual blocks of code. These blocks are generally-but are not limited to - the actions performed within a control structure, such as a while loop or an if statement.

You can choose from among several good indentation styles. The four most important general requirements of an indentation style are:
- Blocks should be indented sufficiently so that they stand out clearly.
- Indentation should be consistent: Always indent the same kind of construct in the same manner.
- The indentation style should provide a reasonable way to handle the problem of rightward drift, the problem of nested blocks bumping against the right-hand margin of the page.
- In a compound statement, the open and close braces should line up, and each should appear on its own line:
```

{
statement
statement
.
statement n
}

```
- To prevent future errors, you should make the body of each control structure a compound statement, even if it consists of only one line of code. When you include open and close braces around the single-statement body of a control structure, you enable that control structure to contain multiple statements if needed in the future. A common programmer error is to include a second statement in the single-statement body of a control structure, thinking that it will execute within the body of that control structure. Because of space restrictions, this book will not always follow this style.

Within these guidelines there is room for personal taste. Here is a summary of the style you will see in this book:
```

- A for or while statement is written for a simple action as
while (expression)
statement
and for a compound action as
while (expression)
{
statements
} // end while
- A do statement is written for a simple action as
do
statement
while (expression);
and for a compound action as
do
{
statements
} while (expression);
- An if statement is written for simple actions as

```
```

if (expression)

```
if (expression)
        statement 
        statement 
e1se
e1se
        statement 
```

        statement 
    ```

Guidelines for indentation style

Indentation style in this book
and for compound actions as
```

if (expression)
{
statements
}
else
{
statements
} // end if

```
- One special use of the if statement warrants another style. Nested if statements that choose among three or more different courses of action, such as
```

if (condition }\mp@subsup{)}{1}{
action
else if (condition ()
action2
else if (condition )
action3

```
are written as
if ( condition \(_{1}\) )
        action \(_{1}\)
else if ( \({ }^{1}\) condition \(_{2}\) )
    action \(_{2}\)
else if (condition \({ }_{3}\) )
    action \(_{3}\)

This indentation style better reflects the nature of the construct, which is like a generalized switch statement:
```

case \mp@subsup{condition (: action }{1}{\prime}\mathrm{ ; break;}
case condition}\mp@subsup{n}{2}{}:\mp@subsup{\mathrm{ action }}{2}{}\mathrm{ ; break;
case condition}\mp@subsup{n}{3}{}:\mp@subsup{\mathrm{ action }}{3}{}\mathrm{ ; break;

```
- Braces are used to increase readability, even when they are not a syntactic necessity. For example, in the construct
```

while (expression)
{
if (condition)
statement
else
statement
} // end while

```
the braces are syntactically unnecessary, because an if is a single statement. However, the braces highlight the scope of the while loop.

\section*{B.2.7 Documentation}

A program should be well documented so that others can read, use, and modify it easily. Many acceptable styles for documentation are in use today, and exactly what you should include often depends on

Highly formatted, HTML-based documentation is easy to produce the particular program or your individual circumstances.

This book will use a special form of documentation comment, called a javadoc-style comment because it originated with the \(\mathrm{Java}^{\mathrm{TM}}\) programming language. These comments start with \(/ * *\) and end with */ and contain tags that classify different parts of your documentation. Several utility programs are available to read these comments and generate HTML-based documentation. One of these
programs is called doxygen. When space permits, the source code in the book is commented in this style and can be read by doxygen.

The advantage of using javadoc-style comments-and a documentation system, like doxy-gen-is that you can edit the documentation for your source code as you make changes to that code. The documentation lives directly in the source code-as javadoc-style comments-making it very easy and convenient to keep your documentation up to date.

The following are the essential features of any program's documentation-with associated commenting tags in parentheses:

\section*{Note: Essential features of program documentation}
1. An initial comment at the top of each source code file that includes:
a. File name (@file)
b. Statement of purpose
c. Author (@author)
d. Date (@date)
e. Optional file version number (@version)
2. Initial comments for each class that includes:
a. Name of class and its header file (@class)
b. Statement of purpose
3. Initial comments for each method or function that includes:
a. Statement of purpose
b. Description of each argument in the argument list (@param)
c. Preconditions (@pre)
d. Postconditions (@post)
e. Exceptions thrown (@throw)
f. Return value (@return)
4. Standard C++ comments in the body of each method or function to explain important features or subtle logic

Beginning programmers tend to downplay the importance of documentation because comments do not affect a program's logic. By now, you should realize that people also read programs. Your comments must be clear enough for someone else to either use your module in a program or modify it. Thus, some of your comments are for people who want to use your module, while others are for people who will revise its implementation. You should be conscious of different kinds of comments.

Beginners also have a tendency to document programs as a last step. You should, however, write documentation as you develop the program. Because the task of writing a large program might extend over a period of several weeks, you may find that the module that seemed so obvious when you wrote it last week seems confusing when you try to revise it next week. Why not benefit from your own documentation by writing it now rather than later?

See Appendix I for more details on the use of the javadoc-style comments and the doxygen documentation system.

\section*{B. 3 Modifiability}

After each iteration in the development of a program, the design can change to some degree. This tendency requires that your program be written in a way that makes it easy to modify. This section offers two examples of how you can make a program easy to modify: named constants and typedef statements.

Consider who will read your comments when you write them

You benefit from your own documentation by writing it now instead of later

Named constants make a program easier to modify
typedef
statements make a program easier to modify

\section*{B.3.1 Named Constants}

The use of named constants is a way to enhance the modifiability of a program. For example, the restriction that an array must have a predefined, fixed size causes a bit of difficulty. Suppose that a program uses an array to process the test scores of the computer science majors at your university. When the program was written, there were 202 computer science majors, so the array was declared by
int scores[202];
The program processes the array in several ways. For example, it reads the scores, writes the scores, and averages the scores. The pseudocode for each of these tasks contains a construct such as
for (index \(=0\) through 201)
Process the score
If the number of majors should change, not only do you need to revise the declaration of scores, but you also must change each loop that processes the array to reflect the new array size. In addition, other statements in the program might depend on the size of the array. A 202 here, a 201 there-which to change?

On the other hand, if you use a named constant such as
```

const int NUMBER_OF_MAJORS = 202;

```
you can declare the array by using
```

int scores[NUMBER_OF_MAJORS];

```
and write the pseudocode for the processing loops in this form:
for (index \(=0\) through NUMBER_OF_MAJORS - 1) Process the score

If you write expressions that depend on the size of the array in terms of the constant NUMBER_OF_MAJORS (such as NUMBER_OF_MAJORS - 1), you can change the array size simply by changing the definition of the constant and compiling the program again.

\section*{B.3.2 The typedef Statement}

Suppose that your program performs floating-point computations of type float, but you discover that you need greater precision than float variables provide. To change the relevant float declarations to long double, for example, you would have to locate all such declarations and decide whether to make the change.

You can simplify this change by using a typedef statement, which gives another name to an existing data type. For example, the statement
typedef float RealType;
declares RealType as a synonym for float and allows you to use RealType and float interchangeably. If you declare all the relevant items in the previous program as RealType instead of float, you can make your program easier to modify and to read. To revise the precision of the computations, you would simply change the typedef statement to
typedef long double RealType;

\section*{B. 4 Ease of Use}

Another area in which you need to keep people in mind is the design of the user interface. Humans often process a program's input and output. Here are a few obvious points:
- In an interactive environment, the program should always prompt the user for input in a manner that makes it quite clear what it expects. For example, the prompt "?" is not nearly as enlightening as the prompt "Please enter your account number." You should never assume that the users of your program know what response the program requires.
- A program should always echo its input. Whenever a program reads data, either from a user or from a file, the program should include the values it reads in its output. This inclusion serves two purposes: First, it gives the user a check on the data entered-a safeguard against typos and errors in data transmission. This check is particularly useful in the case of interactive input. Second, the output is more meaningful and self-explanatory when it contains a record of what input generated the output.
- The output should be well labeled and easy to read. An output of
```

1800 6 1
Jones, Q. 223 2234.00 1088.19 N, J Smith, T. 111
110.23 I, Harris, V. }4444000.0022222.2

```
is more prone to misinterpretation than
CUSTOMER ACCOUNTS AS OF 1800 HOURS ON JUNE 1
Account status codes: N=new, J=joint, I=inactive
\begin{tabular}{lrrll} 
NAME & ACC\# & CHECKING & SAVINGS & STATUS \\
Jones, Q. & 223 & \(\$ 2234.00\) & \(\$ 1088.19\) & \(\mathrm{~N}, \mathrm{~J}\) \\
Smith, T. & 111 & \(\$ 110.23\) & - & I \\
Harris, V. & 44 & \(\$ 44000.00\) & \(\$ 22222.22\) & ----
\end{tabular}

These characteristics of a good user interface are only the basics. Several more subtle points separate a program that is merely usable from one that is user-friendly. Students tend to ignore a good user interface, but by investing a little extra time here, you can make a big difference: the difference between a good program and one that only solves the problem. For example, consider a program that requires a user to enter a line of data in some fixed format, with exactly one blank between the items. A free-form input that allows any number of blanks between the items would be much more convenient for the user. It takes so little time to add code that skips blanks, so why require the user to follow an exact format? Once you have made this small additional effort, the code is a permanent part of both your program and your library of techniques, and the user of your program will never have to think about input format.

\section*{B. 5 Fail-Safe Programming}

A fail-safe program is one that will perform reasonably no matter how anyone uses it. Unfortunately, this goal is usually unattainable. A more realistic goal is to anticipate the ways that people might misuse the program and to guard carefully against these abuses.

This discussion considers two types of errors. The first type is an error in input data. For example, suppose that a program expects a nonnegative integer but reads -12 . When a program encounters this type of problem, it should not produce incorrect results or abort with a vague error message. Instead, a fail-safe program provides a message such as

> -12 is not a valid number of children. Please enter this number again.

The second type of error is an error in the program logic. Although a discussion of this type of error belongs in the next section about debugging, detecting errors in program logic is also a characteristic of

Echo the input

Label the output

A good user interface is important

Check for errors in input

Check for errors in logic

This function is not fail-safe
fail-safe programming. A program that appears to have been running correctly may at some point behave unexpectedly, even if the data that it reads is valid. For example, the program may not have accounted for the particular data that elicited the surprise behavior, even though you tried your best to test the program's logic. Or perhaps you modified the program and that modification invalidated an assumption that you made in some other part of the program. Whatever the difficulty, a program should have built-in safeguards against these kinds of errors. It should monitor itself and be able to indicate that something is wrong and that you should not trust the results.

\section*{B.5.1 Guarding Against Errors in Input Data}

Suppose that you are computing statistics about the people in income brackets between \(\$ 10,000\) and \(\$ 100,000\). The brackets are rounded to the nearest thousand dollars: \(\$ 10,000, \$ 11,000\), and so on to \(\$ 100,000\). The raw data is a file of one or more lines of the form
\(G \quad N\)
where \(N\) is the number of people with an income that falls into the \(G\)-thousand-dollar group. If several people have compiled the data, several entries for the same value of \(G\) might occur. As the user enters data, the program must add up and record the number of people for each value of \(G\). From the problem's context, it is clear that \(G\) is an integer in the range 10 to 100 , inclusive, and \(N\) is a nonnegative integer.

As an example of how to guard against errors in input, consider an input function for this problem. The first attempt at writing this function will illustrate several common ways in which a program can fall short of the fail-safe ideal. Eventually you will see an input function that is much closer to this ideal than the original solution.

A first attempt at the function might be
```

const int LOW_END = 10; // 10-thousand-dol7ar income
const int HIGH_END = 100; // 100-thousand-dol7ar income
const int TABLE_SIZE = HIGH_END - LOW_END + 1;
typedef int Tab7eType[TABLE_SIZE];
/** Reads and organizes income statistics.
@param incomeData A TableType of income statistics.
@pre The calling module gives directions to the user.
Input data is error-free, and each input line has the form
G N, where N is the number of people with an income in the
G-thousand-dollar group and LOW_END <= G <= HIGH_END.
An input line with values of zero for both G and N
terminates the input.
@post incomeData[G - LOW_END] is the total number of people
with an income in the G-thousand-dollar group for each
G read. The values read are displayed. */
void readData(Tab1eType incomeData)
{
bool terminateInput = false;
int group, number; // input values
// Clear array
for (group = LOW_END; group <= HIGH_END; group++)
incomeData[group - LOW_END] = 0;
do
{
cout << "Please enter group and number of ";
cout << "people in group, separated by a space: ";
cin >> group >> number;

```
```

            terminateInput = (group == 0) && (number == 0);
            // Group and number are not both 0
            cout << "Income group " << group << " contains "
            << number << " people.\n";
            incomeData[group - LOW_END] += number;
            } while (!terminateInput); // end do-while
    } // end readData

```

This function has some problems. If an input line contains unexpected data, the program will not behave reasonably. Consider two specific possibilities:
- The first integer on the input line, which the function assigns to group, is not in the range LOW_END to HIGH_END. The reference
incomeData[group - LOW_END]
is then incorrect, since it accesses an array element outside the array bounds. This error also will happen when the user tries to end the input by entering zeros for both group and number.
- The second number on the input line, which the function assigns to number, is negative. Although a negative value for number is invalid, because you cannot have a negative number of people in an income group, the function will add number to the group's array entry. Thus, the array incomeData will be incorrect.

After the function reads values for group and number, it must check to see whether group is in the range LOW_END to HIGH_END and whether number is positive. If either value is not in that range, you must handle the input error.

Suppose that instead of checking the value of number before you add it to the array element incomeData[group - LOW_END], you perform the addition and see whether the sum is positive. This approach is insufficient, since you might add a negative value to an entry of incomeData without that entry becoming negative. For example, if number is \(-4,000\) and the corresponding entry in the array incomeData is 10,000 , the sum is 6,000 . Thus, a negative value for number could remain undetected and invalidate the results of the rest of the program.

One possible course of action for the function to take when it detects invalid data is to set an error flag and terminate. Another possibility is for it to set an error flag, ignore the bad input line, and continue. Which action is correct really depends on how the program uses the data once it is read.

The following readData function attempts to be as universally applicable as possible and to make the program that uses it as modifiable as possible. When the function encounters an error in input, it sets a flag, ignores the data line, and continues. By setting a flag, the function leaves it to the calling module to determine the appropriate action-such as abort or continue-when an input error occurs. Thus, you can use the same input function in many contexts and can easily modify the action taken upon encountering an error.
```

/** Reads and organizes income statistics.
@param incomeData A TableType of income statistics.
@pre The calling program gives directions to the user.
Each input Tine contains exactly two integers in the form G N,
where N is the number of people with an income in the G-thousand-
dollar group and LOW_END <= G <= HIGH_END. An input line with
values of zero for both G and N terminates the input.
@post incomeData[G - LOW_END] is the total number of people with
an income in the G-thousand-dollar group. The values read are
displayed. If either G or N is erroneous (either G < LOW_END,
G > HICH_END, or N < 0), the function ignores the data line and
continues execution.
@return False if either G or N are erroneous (either G < LOW_END,

```

Test for invalid input data

A function that includes fail-safe programming
```

    G > HIGH_END, or N < 0) for any data line read. In this case,
    the calling program should take action. The return value is
    true if the data is error-free. */
    bool readData(Tab1eType incomeData)
{
int group = -1;
int number = -1; // Input values
boo1 dataCorrect = true; // No data error found as yet
bool userFinished = false;
for (group = LOW_END; group <= HIGH_END; group++)
incomeData[group - LOW_END] = 0;
do
{
cout << "Please enter group and number of ";
cout << "peop1e in group, separated by a space: ";
cin >> group >> number;
cout << "Input line specifies that income group "
<< group << "\ncontains " << number
<< " people.\n";
if ((group >= LOW_END) \&\& (group <= HIGH_END) \&\&
(number >= 0))
// Input data is valid - add it to tally
incomeData[group - LOW_END] += number;
else if ((group == 0) \&\& (number == 0))
userFinished = true; // User signaled input is finished
else
{
// Error in input data:
// Set error flag and ignore input line
dataCorrect = false;
} // end if
} while( (dataCorrect \&\& !userFinished ); // end do-while
return dataCorrect;
} // end readData

```

Although this input function will behave gracefully in the face of most common input errors, it is not completely fail-safe. What happens if an input line contains only one integer? What happens if an input line contains a noninteger? The function would be more fail-safe if it read its input character by character, converted the characters to an integer, and checked for the end of the input line. In most contexts, this processing would be a bit extreme. However, if the people who enter the data frequently err by typing nonintegers, you could alter the input function easily because the function is an isolated module. In any case, the function's initial comments should include any assumptions it makes about the data and an indication of what might make the program abort abnormally.

\section*{B.5.2 Guarding Against Errors in Program Logic}

Now consider the second type of error that a program should guard against: errors in its own logic. These are errors that you may not have caught when you debugged the program or that you may have introduced through program modification.

Unfortunately, a program cannot reliably let you know when something is wrong with it. (Could you rely on a program to tell you that something is wrong with its mechanism for telling you that something is wrong?) You can, however, build into a program checks that ensure that certain conditions always hold when the program is correctly implementing its algorithm. For example, all integers in the array incomeData of the previous example must be greater than or equal to zero. Although we
argued that the function readData should not check the validity of the entries of incomeData instead of checking number, it could do so in addition to checking number. For example, if the function finds that an entry in the array incomeData is outside some range of believability, it can signal a potential problem to its users.

Another general way in which you should make a program fail-safe is to make each function check its precondition. For example, consider the following function, factorial, which returns the factorial of an integer:
```

```
/** Computes the factorial of an integer.
```

```
/** Computes the factorial of an integer.
@param n The given integer.
@param n The given integer.
@pre n >= 0.
@pre n >= 0.
@post None.
@post None.
@return n* (n - 1)* ...*1, if n>0, or 1, if n == 0.*/
@return n* (n - 1)* ...*1, if n>0, or 1, if n == 0.*/
int factorial (int n)
int factorial (int n)
{
{
    int fact = 1;
    int fact = 1;
    for (int i = n; i > 1; i-)
    for (int i = n; i > 1; i-)
        fact *= i;
        fact *= i;
    return fact;
    return fact;
} // end factorial
```

} // end factorial

```
```

        fact (= i; il>1, i-)
    ```
```

        fact (= i; il>1, i-)
    ```

The initial comments in this function contain a precondition-information about what assump-
are made-as should always be the case. The value that this function returns is valid only if the
In the context of the program for which this function was written, it may be reasonable to make
assumption that n will never be negative. That is, if the rest of the program is working correctly, it
call factoria only with correct values of \(n\). Ironically, this last observation gives you a good
on for factorial to check the value of \(n\) : If \(n\) is less than zero, the warning that results from the
The initial comments in this function contain a precondition-information about what assump-
tions are made-as should always be the case. The value that this function returns is valid only if the
precondition is met. If \(n\) is less than zero, the function will return the incorrect value of 1 .
In the context of the program for which this function was written, it may be reasonable to make
the assumption that \(n\) will never be negative. That is, if the rest of the program is working correctly, it
will call factorial only with correct values of \(n\). Ironically, this last observation gives you a good
reason for factorial to check the value of \(n\) : If \(n\) is less than zero, the warning that results from the
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will call factorial only with correct values of n . Ironically, this last observation gives you a good
reason for factorial to check the value of n : If n is less than zero, the warning that results from the check indicates that something may be wrong elsewhere in the program.

Another reason the function factorial should check whether \(n\) is less than zero is that the function should be correct outside the context of its program. That is, if you borrow the function for use in another program, the function should warn you if you use it incorrectly by passing it an \(n\) that is negative. A stronger check than simply the statement of the precondition in a comment is desirable.


Note: A function should state its assumptions and, when possible, check whether its arguments conform to these assumptions.

In this example, factorial could check the value of n and, if it is negative, return zero, because factorials are never zero. The program that uses factorial could then check for this unusual value. Alternatively, factorial could abort execution if its argument is negative. Many programming languages, including \(\mathrm{C}++\), support a mechanism for error handling called an exception. A module indicates that an error has occurred by throwing an exception. A module reacts to an exception that another module throws by catching the exception and executing code to deal with the error condition. \(\mathrm{C}++\) Interlude 3 provides more information about exceptions.
\(\mathrm{C}++\) also provides a convenient function assert (expression) that both displays an informative message and aborts a program if expression is zero. You can use assert to check for both error conditions and the validity of preconditions within your program. C++ Interlude 3 provides more information about assert.

Functions should enforce their preconditions

Functions should check the values of their arguments

Use breakpoints, single-stepping, watches, and temporary cout statements to find logic errors

\section*{Systematically} check a program's logic to determine where an error occurs

\section*{B. 6 Debugging}

No matter how much care you take in writing a program, it will contain errors that you need to track down. Fortunately, programs that are modular, clear, and well documented are generally amenable to debugging. Fail-safe techniques, which guard against certain errors and report them when they are encountered, are also a great aid in debugging.

Without a systematic approach, finding a small mistake in a large program can indeed be a difficult task. Many people have difficulty in debugging a program, perhaps because they believe that their program is really doing what it is supposed to do. For example, on receiving an execution-time error message at line 1098, a person might say, "That's impossible. The statement at line 1098 was not even executed, because it is in the e1se clause, and I am positive that it was not executed." However, the proper approach is either to trace the program's execution by using available debugging facilities or to add cout statements that show which part of the if statement was executed. By doing so, you verify the value of the expression in the if statement. If the expression is 0 , for example, when you expect it to be 1 , the next step is to determine how it became 0 .

How can you find the point in a program where something becomes other than what it should be? A typical Integrated Development Environment, or IDE, allows you to trace a program's execution either by single-stepping through the statements in the program or by setting breakpoints at which execution will halt. You also can examine the contents of particular variables by either establishing watches or inserting temporary cout statements. The key to debugging is simply to use these techniques to tell you what is going on. This may sound pretty mundane, but the real trick is to use these debugging aids in an effective manner. After all, you do not simply put breakpoints, watches, and cout statements at random points in the program and have them report random information.

The main idea is systematically to locate the points of the program that cause the problem. A program's logic implies that certain conditions should be true at various points in the program. If the program's results differ from your expectations, an error occurs. To correct the error, you must find the first point in the program at which this difference is evident. By inserting either breakpoints and watches or cout statements at strategic locations of a program-such as at the entry and departure points of loops and functions-you can systematically isolate the error.

These diagnostic techniques should inform you whether things start going wrong before or after a given point in the program. Thus, after you run the program with an initial set of diagnostics, you should be able to trap the error between two points. For example, suppose that things are fine before you call method \(M_{1}\), but that something is wrong by the time you call \(M_{2}\). This kind of information allows you to focus your attention between these two points. You continue the process until eventually the search is limited to only a few statements. There is really no place in a program for an error to hide.

Programming Tip: A statement of truth about some aspect of a program's logic is known as an assertion. You can express an assertion either as a comment or by using the assert macro. By including assertions in your program, you facilitate the debugging process.

The ability to place breakpoints, watches, and cout statements in appropriate locations and to have them report appropriate information comes in part from thinking logically about the problem and in part from experience. Here are a few general guidelines.

\section*{B.6.1 Debugging Functions and Methods}

You should examine the values of the arguments passed to a function or method as at its beginning and end by using either watches or cout statements. Ideally, you should debug each major function or method separately before using it in your program.

\section*{B.6.2 Debugging Loops}

You should examine the values of key variables at the beginnings and ends of loops, as the comments in this example indicate:
```

// Check values of start and stop before entering loop
for (index = start; index <= stop; index++)
{
// Check values of index and key variables
// at the beginning of iteration
// Check values of index and key variables
// at the end of iteration
} // end for
// Check values of start and stop after exiting loop

```

\section*{B.6.3 Debugging if Statements}

Just before an if statement, you should examine the values of the variables within its expression. You can use either breakpoints or cout statements to determine which branch the if statement takes, as this example indicates:
```

// Check variab7es within expression before executing if
if (expression)
{
cout << "Condition is true (value of expression is 1).";
}
else
{
cout << "Condition is false (value of expression is 0).";
...
} // end if

```

\section*{B.6.4 Using cout Statements}

Sometimes cout statements can be more convenient than watches. Such cout statements should report both the values of key variables and the location in the program at which the variables have those values. You can use a comment to label the location, as follows:
```

// This is point A
cout << "At point A in the method computeResults:\n"
<< "x = " << x << ", y = " << y << endl;

```

Remember to either disable or remove these statements when your program finally works.

\section*{B.6.5 Using Special Dump Functions}

Often the variables whose values you wish to examine are arrays or other, more complex data structures. If so, you should write dump functions to display the data structures in a highly readable manner. You can easily move the single statement that calls each dump function from one point in the

You can never test a program too much

Unit testing checks individual modules

Integration testing checks many parts of a solution together

System testing checks the integration of the solution with other systems
program to another as you track down an error. The time you spend on these functions often proves to be worthwhile, as you can call them repeatedly while debugging different parts of the program.

Hopefully, this discussion has conveyed the importance of the effective use of diagnostic aids in debugging. Even the best programmers have to spend some time debugging. Thus, to be a truly good programmer, you must be a good debugger.

\section*{B. 7 Testing}

Programs need to be tested to make sure that they correctly solve their current requirements. Good testing cannot happen without a plan. Software is complex, and determining whether it works correctly is not easy. Software testing is so important that some iterative development processes argue that tests should be developed before writing any code.

\section*{B.7.1 Levels of Testing}

Several levels of testing should take place. The first level of testing is called unit testing, and it happens on the individual modules. You should test individual methods first, and then test the classes. The second level of testing is called integration testing, and it tests the interactions among modules. The next level of testing is called system testing, in which the entire program is tested. Finally, a special type of testing called acceptance testing shows whether the system as a whole complies with its requirements.

Unit testing is something that you have probably done before. It is closely related to some of the ideas talked about in the previous debugging section. A function is tested to make sure that it conforms to its requirements. Often with small programs, testing does not go beyond this level.

One way to perform integration tests is to make sure that the various objects within a program correctly implement the interactions in your design. Integration testing can be challenging, because there are many ways in which the objects in a program can interact. It is difficult to make sure they do what they are designed to do and nothing more.

System testing entails running the program in the environment in which it was designed to work. Here the program is tested to make sure that it interacts with external systems correctly. The external systems might be other programs-such as a government tax collection system-or hardware devices.

\section*{B.7.2 Kinds of Testing}

Two kinds of testing can be applied to a system:
- Open-box testing. With open-box, white-box, or glass-box testing, you design tests knowing how the modules work. You carefully choose a method's actual arguments in a way that exercises all lines of code in the method. For example, you must make sure that all branches of if and switch statements are executed.
- Closed-box testing. With closed-box, black-box, or functional testing, you know the name of the method being tested, as well as its argument list, its return type, and its operation contract. Using only this information, you develop tests across a range of valid input, and then check the output against hand-calculated results. Choosing some actual arguments outside of their valid range tests how the module handles incorrect input.

\section*{B.7.3 Developing Test Data}

The test data that you use must test a wide range of conditions. It should include valid, typical input and values at the extreme limits of valid data-called boundary values-as well as values that
should never occur in normal execution. For example, suppose that you are testing a simple linear search algorithm:
```

for (int i = 0; i < size; i++)
{
if (array[i] == target)
return i;
} // end for

```

You should use all of the following values of target:
- Just less than index[0] (boundary at the beginning of the array)
- Equal to index[0] (boundary at the beginning of the array)
- Equal to some entry in the middle of the array
- Within the range of values stored in the array, but unequal to any
- Equal to index[size - 1] (boundary at the end of the array)
- Just greater than index[size - 1] (boundary at the end of the array)

You should also test this algorithm with
- An empty array
- An array with one entry in it
- An array with many entries

For each of these tests, you should manually trace the data so that the results are known ahead of time. In this way, test data can be the input to a module, and you can compare the output with known results.

\section*{B.7.4 Testing Techniques}

As indicated in the debugging section, assert statements can be useful for checking assertions. However, the assert statement is not powerful enough for the full-scale testing of a system. Other techniques need to be utilized.

A common technique is to write a block of code that tests a module. When you have finished testing the system, you can simply comment it out-rather than deleting it-as more testing will be required after you make modifications during the next iteration. Alternatively, you can turn these testing blocks of code on and off globally by using one of the techniques that we discuss next.

The first technique uses a global boolean variable to grant access to testing code contained within the body of if statements. Setting the variable to false turns testing code "off," and setting it to true enables testing. For example, the following test is disabled:
```

const boo1 TESTING = false; // Disables testing code
if (TESTING)
{
... // Testing code
} // end if

```

You could define several different boolean variables to turn different tests "on" and "off." All that is required is to change the value of the boolean variable and to recompile.

A second, more efficient, technique uses the preprocessor to the compiler. You define a macro that guards the testing code. Then you place the test code within preprocessor directives that test whether that macro has been defined. For example:
```

\#define TESTING
\#ifdef TESTING
...// Testing code in here.
\#endif

```

Always carefully test boundary values

Values for testing a search

Disable, but do not remove, testing code after use

A stub is an incomplete method that acknowledges that it has been called
sortData is a stub

A driver is a module that tests another module

To turn off testing, you comment out the definition of the macro (// \#define TESTING) and recompile. Because the macro is no longer defined, the test code is not included in the executable. This has the advantage that the test code is completely removed from the executable, making it smaller. Again, you can define several different macros to enable you to turn different blocks of testing code "on" and "off."

\section*{B.7.5 Use of Stubs}

Testing a class before implementing all of its methods is often useful. Instead of simply omitting some methods, you partially implement them so that each one acknowledges that it has been called. Such methods are called stubs. Stubs allow you to see that the method has been called in the correct place in the overall flow of a program.

For example, suppose that a class has an array data of integers as a data member and a method sortData. You could write the following stub for sortData:
```

void sortData() // Stub
{
cout << "sortData has been called\n"
data[0] = 1;
data[1] = 2;
data[2] = 3;
} // end sortData

```

The stub sortData not only indicates that it has been called, but also changes the data into a sorted form. Thus, this stub could be used for simple integration testing as well. Later, you could implement sortData by using an appropriate sorting algorithm.

\section*{B.7.6 Use of Drivers}

Given the previous class-let's name it Example-the following main function creates an Example object and then calls its methods:
```

int main()
{
Example t;
t.dump();
t.sortData();
t.dump();
return 0;
} // end main

```

Notice the use of Example's method dump, whose purpose is to display the data in an Example object. First, main dumps the object \(t\), then it sorts \(t\) 's data, and finally it dumps \(t\) again. This main function is known as a driver. It does not do anything algorithmically interesting; it simply tests an Example object.

As is evident from the preceding discussion, testing takes a lot of planning and is not easy to do. Like many things in programming, testing well takes some time and experience.

\title{
The Unified Modeling Language
}

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\section*{Prerequisite}

C++ Interlude 4 Class Relationships and Reuse

The Unified Modeling Language visually represents object-oriented designs

A use case is a textual story; a scenario describes the system's behavior under certain circumstances from the perspective of the user

An example of a main success scenario

Thhe Unified Modeling Language, or UML, is a modeling language used to express object-oriented designs. The UML provides specifications for many types of diagrams and text-based descriptions of various aspects of a design. The diagrams are particularly useful in showing specific, interesting aspects of the elements in the domain of the problem.

\section*{C. 1 The Use Case}

A use case is a set of textual scenarios or stories that describe the proposed solution. A use case is not object-oriented, but it is an important part of the analysis of the problem. A single use case usually has several scenarios. The main success scenario (or "happy path") describes how the system satisfies the goals of the user when everything goes smoothly. Other alternate scenarios describe the interaction between the user and the system when specific things do not go well or when exceptional conditions apply.

For example, suppose that a customer wants to withdraw money from his or her checking account. A main success scenario might look like this:

Customer makes a request to withdraw money from a bank account. Bank identifies and authenticates the customer. Bank gets from the customer the type of account (savings or checking), the account number, and the amount of the withdrawal. Bank verifies that the account balance is greater than the amount of the withdrawal. Bank generates a receipt for the transaction. Bank counts out the correct amount of money for the customer. Bank gives the customer the receipt and money. Customer leaves the bank.

An example of an alternate scenario

A scenario is written from the perspective of the user of the system. In this way, the analysis will focus on the responsibilities of the system to meet a user's goals. That is, the analysis will focus on what the system needs to accomplish to meet the goals of its users.

Notice that the scenario does not describe how the system works. For example, "Bank identifies and authenticates the customer" does not indicate how this will occur. The focus should be on the interactions between the user and the system but not on the user interface to the system, which will be explored during design.

Alternate scenarios should be written only for stories that are interesting or tricky. For example, one such scenario describes what happens when the customer fails authentication:

Customer makes request to withdraw money from a bank account. Bank identifies but fails to authenticate the customer. Bank refuses to process the customer request. Customer leaves the bank.

Other alternate scenarios could include the denial of a withdrawal request because insufficient funds are in the account or because the bank does not have sufficient cash on hand.

From the scenarios in the use case, you generate a list of objects by listing the nouns in the use case. (This is where the analysis becomes object oriented.) For example, bank, money, customer, account, account type, account number, transaction amount, and receipt are all found in the main success scenario. The final solution might not use all of the objects discovered.

\section*{Note: The use case during object-oriented analysis}
1. Describes the problem domain in terms of scenarios involving the solution that satisfies user goals
2. Discovers noteworthy objects, attributes, and associations within the scenarios

\section*{C. 2 UML Sequence Diagrams}

After you create a use case and list potential objects, you need to explore how these objects will interact with one another. You accomplish this by using UML sequence diagrams that model your use case scenarios. A UML sequence diagram, or interaction diagram, shows how two or more objects will interact with one another over time within a single scenario. This diagram allows you to visualize the messages sent among the objects in a scenario and the order in which those messages happen. In addition to giving a sense of the flow of the logic, the diagram is important when defining the responsibilities of the objects: What must the object "remember," or keep track of, and what must the object do for other objects? Many of the responsibilities are described in the use case. Figure C-1 shows a UML sequence diagram for the main success scenario given previously.

Although the development of the use case is part of object-oriented analysis (OOA), the creation of sequence diagrams is part of object-oriented design (OOD). The sequence diagram shows a scenario's objects from left to right horizontally and their actions in order of occurrence vertically. The UML represents an object as a square-cornered box containing the name of the object-if it has one-followed by a colon and the type of the object, all underlined. In Figure C-1, the object of type Bank is named bank, and the objects of type Customer and Account are nameless.

Each object in a sequence diagram has a lifeline denoted by a vertical dashed line below the object box extending forward in time. Each lifeline can have one or more activation bars, each represented by an open box on the lifeline. An activation bar indicates when an object is activeand represents a responsibility of the object's class. The class needs a method to handle that responsibility. To represent the sending of a message between objects, you draw a solid arrow from the activation bar of the calling object to the top of a new activation bar on the lifeline of the called

FIGURE C-1 Sequence diagram for the main success scenario

object. Such arrows are labeled with a message expression indicating what task the object needs to perform. The message expression corresponds to the name of a method for handling that task. The message expression may also include parameters necessary for performing the task. Ultimately, these parameters will correspond to method arguments. When an object finishes a task, the activation bar associated with that task ends. If the task returns something of interest to the calling object, a dashed arrow may be shown pointing back to the calling object's activation bar. You can label this arrow with the name of the variable receiving the returned value.

In Figure C-1, the customer asks the bank for authorization by giving its name and identifierpassed as arguments - to the authorize message. The bank then processes this message-as indicated by the activation bar associated with the bank object's lifeline-and then signals whether or not the customer has been authenticated by returning a value to the variable 0 K within the Customer object. Note that you should write the messages in a general, language-independent fashion. A Bank class could declare the authorize message as follows:

An activation bar on an object's lifeline represents a responsibility of the object's class
```

class Bank

```
class Bank
{
{
public:
public:
    bool authorize(string name, string identifier);
    bool authorize(string name, string identifier);
}; // end Bank
```

}; // end Bank

```

\section*{A stereotype} identifies a special characteristic of an element

FIGURE C-2 Sequence diagram showing the creation of a new object


An object may also send messages to itself, as is done in the generateReceipt and countCash messages in Figure C-1. Notice the piggybacked activation bars, which indicate that the Bank class is calling its own member methods.

A sequence diagram can also show the creation of a new object. In Figure C-2, a new account is created. This diagram can represent a part of an alternate scenario in a use case, in which a customer comes into the bank but does not yet have an account. In the UML, the notation <<create>> is what is known as a stereotype. Stereotypes use left and right guillemets ( \(\ll\) and \(\gg\) ), but when drawing a sequence diagram by hand, you can simply label the message arrow with create. A stereotype is used to identify a unique characteristic of an element in any UML diagram.

The UML sequence diagram is a design document that gives guidance on several levels to the programmer implementing a solution. It gives a view of how the objects will dynamically interact (collaborate) to solve a use case scenario. In particular, a sequence diagram shows
- Which objects collaborate to solve a particular use-case scenario
- The synchronous nature of the messages that are passed between the objects in the scenario, providing the order in which tasks are accomplished
- Which object is responsible for what task or tasks by handling a particular message


\section*{Note: UML sequence diagrams}
- Provide a visual representation of the flow of messages sent between objects during a single scenario
- Indicate the responsibilities of an object
- Indicate the collaborations among objects

\section*{C. 3 UML Class Diagrams}

The last design element we will introduce is the UML class diagram, or static diagram. The class diagram shows the attributes and operations of an individual class and how multiple classes are related to one another. These design elements do not change after they have been created within a solution - that is, they are static, or unchanging.

The class diagram shown in Figure C-3 represents a class of banks. Each diagram consists of a square-cornered box containing the name of the class centered within its borders. Figure \(\mathrm{C}-3 \mathrm{a}\) shows the simplest of class diagrams: a class with a name but no attributes or operations. Figure C-3b shows the bank's attributes in the middle section and its operations in the bottom section. If a class has no attributes, as in Figure C-3c, an empty middle section must be present.

\section*{FIGURE C-3 Three possible class diagrams for a class of banks}
(a)
Bank
(b)

(c)
\begin{tabular}{|c|}
\hline Bank \\
\hline \\
\hline createAccount() \\
\hline
\end{tabular}

The class diagram represents a conceptual model of a class of objects in a language-independent way. That is, the diagram does not dictate how the class is implemented. It just gives you a design to work from. However, you can specify many qualities associated with the attributes and operations of the class as an option. These qualities are easily implemented in \(\mathrm{C}++\), and they increase the amount of information given in a design.

The UML syntax for the attributes of a class is
[visibility] name [:type][=defaultValue] [\{property\}]

UML syntax for the attributes of a class
where
- All elements within square brackets are optional.
- visibility can be + for public accessibility, - for private accessibility, or \# for protected accessibility. (C++ Interlude 4 covers protected accessibility.) If omitted, the visibility defaults to private accessibility.
- name is the name of the attribute.
- type is the data type of the attribute.
- defaultValue is the default value of the attribute.
- property indicates some property that applies to this attribute.

At a minimum, the name of an attribute should be given. The defaultValue is given only in situations where the design dictates that a default value is appropriate for that attribute. In certain cases it might be appropriate to omit the type of an attribute, leaving this detail until the implementation phase.

The UML specifies that programming-language-independent names be used for the type of an attribute. This text will follow the UML recommendations by using integer for integral storage, float for floating-point storage, bool ean for boolean storage, and string for storing strings.

The property of an attribute can have one of the following values:
- changeable indicates a normal, modifiable attribute and is usually omitted.
- frozen indicates a constant or write-once attribute.

The attributes for the Bank class in Figure C-3b could now be written as
```

-name: string
-routingNum: integer

```

The visibility of these attributes is private and adheres to the object-oriented programming principle of information hiding.

The UML syntax for a class's operations is more involved:

UML syntax for the operations of a class
where
- All elements within square brackets are optional.
- visibility is the same as specified for attributes, except if omitted operations default to public accessibility.
- name is the name of the operation.
- type is the data type of the result returned from the operation. If the operation returns nothing, type can be either omitted or specified as void. The same language-independent types apply here as for attribute types.
- property indicates some property that applies to this operation. The only property relevant to us is query, which indicates that the operation will not modify any of the data members in the calling object.
- parameterList is either empty-in which case the parentheses are still required-or contains a comma-delimited list of parameters to the operation. The syntax for each parameter looks like the following:
[direction] name:type \([=\) defaultValue \(]\)
where
- All elements within square brackets are optional.
- direction shows whether the parameter is used for input (in), output (out), or both input and output (inout).
- name is the name of the parameter.
- type is the data type of the parameter. The same language-independent types apply here as for attribute types.
- defaultValue is the value of this parameter if no corresponding actual argument is provided when this operation is called.

In Figure C-3b, the Bank class operations could be written as
```

+authorize(in name: string, in identifier: string): void {query}
-createAccount()

```

The authorize operation has public accessibility, and the createAccount operation is private. Because the design of the authorize operation has the query property specified, this operation will not change any of the data in the calling object.

The UML provides for several types of relationships among the classes in a class diagram. For example, Figure C-4 shows a static (class) diagram for a banking system. Because the bank has both checking accounts and savings accounts for its customers, the diagram shows five classes of objects: a Bank class, a Customer class, an Account class, a Checking class, and a Savings class. These classes collaborate by sending each other messages. Through these collaborations, the classes are related to one another in various ways.

For example, the Bank and Customer classes have an association relationship indicated by the solid line with no arrowheads. The numbers at the ends of the association-and other relationshipsare called multiplicities and are optional. Each bank object is associated with zero or more customers, but each customer is associated with one bank. Each customer can have multiple accounts of any type, but an account can belong to only one customer.

The Bank and Account classes are in one kind of part/whole relationship called an aggregation, or containment, relationship that is denoted by the open diamond arrowhead near the containing class. In an aggregation relationship, the lifetime of the containing object and the object contained are not necessarily the same. Banks "live" longer than the accounts that they contain. Another part/whole relationship, called composition, uses a filled-in diamond arrowhead near the containing class. This

Aggregation indicates that one class contains an instance of another class

An association indicates that the classes know about each other

FIGURE C-4 A UML class diagram of a banking system

relationship denotes a stronger containment, where the lifetimes of the container and the contained objects are the same. For example, a ballpoint pen is a pen that has a ball at the tip. When the pen "dies," the ball "dies" with it. Thus, the pen and its ball are in a composition relationship. By also looking at the multiplicities, we can say that a bank stores-that is, contains-zero or more accounts. The Customer and Account classes also have an aggregation relationship.

Each of the bank's accounts stores its account number and current balance and allows a customer to ask for the current balance, withdraw funds, and deposit funds. The Account class specifies these attributes and operations common to all accounts. The Account class is related to the Checking and Savings classes through a generalization relationship, indicated by the open triangular arrowhead pointing to the general (or parent) class. The generalization relationship denotes inheritance. The attributes and operations of the ancestor class, Account, are inherited by the descendant classes, Checking and Savings.

Note that the Account class name is in italic in Figure C-4. This denotes an abstract base class. The getBalance operation in the Account class is also in italic. This indicates that getBalance is not implemented in the Account class but must be implemented within the descendant class. In fact, the Checking and Savings classes each implement the inherited getBalance operation. This operation will take on polymorphic behavior when called from a parent class object. Many of these points, both in design and implementation, will be covered in detail in C++ Interlude 4.


\section*{Note: UML Class (Static) Diagrams}
- Show the attributes and operations of individual classes
- Show the unchanging relationships among the classes in the solution

Composition is a stronger form of aggregation; the lifetimes of the container and the contained objects are the same

UML diagrams are not documentation

The UML is a tool for exploration and communication

\section*{C. 4 Keeping The UML In Perspective}

Although the UML can be thought of as a tool for documenting a solution to a problem, this is not how it should be used. Rather, you should use the UML to quickly draw various aspects of the current problem to better understand it during OOA and to create a set of objects during OOD that can collaborate to solve this problem. Often small teams of programmers will draw several UML diagrams on a whiteboard and, after working out the kinks in the diagrams, take digital pictures of them for future reference. These design diagrams are considered an artifact, or document, associated with the analysis and design. However, the ultimate implementation-that is, the solution-might not exactly match the original design. Thus, the design diagrams might not document the solution.

The UML is a useful tool for creating a design for a solution. It enables programmers to explore and understand the problem domain and to develop a design that is close to the required solution. Although the UML will not necessarily make better designs than other techniques, it can conceptually model a problem domain in terms of software objects, independently of a programming language. Because the UML gives a visual representation of the proposed system, humans-as visual creatures - can benefit from using it. After all, a picture is worth a thousand words. The UML enables the members of a programming team to communicate visually with one another, thereby gaining a common understanding of the system that is being built.

\section*{The Software Life Cycle}

\section*{Appendix} D

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PrerequisitesA general knowledge of software

The software life cycle describes the phases of software development

Iterative development forms a solution after many iterations; each iteration builds on the previous iteration until a complete solution is achieved

Thhis appendix examines what a major piece of software "experiences" during its lifetime. The life cycle of software describes the phases through which software will progress from conception to replacement to deletion from a hard disk forever. We will look at the development phases, from project conception to deployment to end users. Beyond this development process, software will need maintenance to correct errors and extend it to do new tasks. Eventually the software will be retired.

The development of good software is a complex affair. Over the last several decades many techniques have been developed to try to control that complexity. Some of these techniques have been more successful than others.

\section*{D. 1 Iterative and Evolutionary Development}

An iterative development of a solution to a problem progresses through many short, fixed-length iterations, where each iteration cycles through the analysis, design, implementation, testing, and integration of a small portion of the problem domain. The early iterations build the core of the system. Subsequent iterations build on that core. After many iterations, the entire solution is developed.

To manage the time spent on an iteration, each one is assigned a predetermined duration - called a timebox - at the beginning of the project. Typical timeboxes are two

A critique of each iteration influences the next iteration

\section*{Iterative}
development enables you to adapt a solution to a changing environment

The RUP gives structure to the software development process
to four weeks in length. The portion of the system that can be developed in less than two weeks is typically too small to be meaningful. In more than four weeks, the partial system becomes too large. At the end of each iteration, the partial system should be functional and completely tested. Each iteration makes relatively few changes to the partial system produced by the previous iteration, so it can occur quickly. Thus, you do not expend too much effort on something that you may have misunderstood about the system requirements.

This process creates the solution incrementally. Thus, there are many places for end users and developers to check that it solves the correct problem. Because each iteration produces a functional but partial system, developers can engage end users to generate feedback. This feedback influences the next iteration and can change the direction of development so a more correct solution is reached. Figure D-1 illustrates the progress of an iterative development. We see that early iterations may be way off base, reflecting the fact that the requirements of the solution are not clearly understood. This is not anyone's fault; even the people that have requested the software to be built, the domain experts, may not be able to see what is required in the final system. Later iterations will probably not vary so much, as the requirements of the final system will have become more stable and in line with the correct solution. In an incremental fashion, the correct solution is created.

FIGURE D-1 Iterative development's progress toward the correct system


Iterative, incremental development gives a programmer the ability to let the requirements analysis and design evolve over time to meet a changing environment. Thus, the process is sometimes called evolutionary development.


Note: Iterative and Evolutionary Development
- Determines timebox lengths at the start of a project
- Keeps timebox lengths short (two to four weeks)
- Gets end-user and domain-expert feedback from iteration \(n\) to influence the direction of iteration \(n+1\)

\section*{D. 2 Rational Unified Process Development}

This section introduces the widely used Rational Unified Process (RUP) as a basis for examining the phases of the development process. The RUP will give structure to the use of the analysis and design tools you learned earlier. As the RUP is a large and complex process, most of its workings are beyond the scope of this text. Thus, we will simply introduce the process.

The RUP organizes development into four phases:
- Inception-feasibility study, project vision, rough estimates of time and cost
- Elaboration-refined project vision, iterative development of core system, development of system requirements, more accurate time and cost estimates
- Construction-iterative development of remaining system
- Transition-testing and deployment of the system

We will look at the characteristics of each phase. Figure D-2 shows how the RUP development phases relate to iterations.

\section*{FIGURE D-2 RUP development phases}


\section*{D.2.1 Inception Phase}

During the inception phase, we define the scope of the project and decide whether the project is feasible. If a team has already implemented solutions for similar problems, and the solution is clearly feasible, the inception phase may be very brief.

At a minimum, the inception phase must define an initial set of system requirements and choose the length of the iteration timebox. While defining the initial requirements, you generate a core set of use case scenarios and determine which elements of the solution involve the greatest risk. This set of scenarios and the highest-risk aspects of the solution are the focus of the first iterations in the elaboration phase. Note that the inception phase is not an iteration. You should not try to develop all-or even most-of the use case scenarios during inception. Usually only about 10 percent of the use case scenarios are written in any detail during the inception phase.

\section*{D.2.2 Elaboration Phase}

During the elaboration phase, the system's core architecture is iteratively developed, high-risk elements of the system are addressed, and most of the system requirements are defined. The core architecture is composed of those software components that are central to the system. For example, the objects represented in the Unified Modeling Language (UML) class diagram in Figure C-4 of Appendix C would make up the core architecture of that banking system.

The elaboration phase extends over two or more iterations. If the elaboration phase has only one iteration, no feedback can come into play. Multiple iterations are necessary when developing the correct solution incrementally and adaptively.

The usual reason for a solution to fail is that high-risk elements were not handled adequately. Early in the elaboration phase, high-risk elements need to be addressed and resolved. If these

The inception phase defines a project's scope and feasibility

The elaboration phase develops a core system and addresses high-risk elements

The elaboration phase refines a system's requirements and cost

The construction phase develops the remaining system

During the transition phase, beta testing and system development occur
elements are irresolvable, either the scope of the project must be changed to remove them or the project must be abandoned. In either case, little time, money, and effort will have been wasted. Leaving the high-risk elements until later in the development of a project can lead to time and cost overruns-and poor time and cost estimates-because the resolution of the high-risk issues will consume considerable resources. Moreover, if these issues cause the project to fail, much is lost. Therefore, you should address high-risk issues early.

Each iteration starts with an object-oriented analysis of the system requirements and input from the testing of the previous iteration. This analysis leads to a set of use case scenarios to analyze, design, implement, and test. Those scenarios are examined for objects and attributes. UML sequence and class diagrams are drawn to clearly show the objects' collaborations, responsibilities to the system, and relationships to one another. This design is then coded and tested. If this iteration is not one of the first, the code may need to be integrated into the existing subsystem. After final testing, end users and domain experts use the system to produce feedback for subsequent iterations.

\section*{D.2.3 Construction Phase}

The construction phase begins once most of the system requirements have been formalized. Each iteration in this phase still consists of analysis, design, implementation, and testing. At this point in the development of the system, however, the analysis and design components require less attention, and increased effort is focused on implementation and testing.

\section*{D.2.4 Transition Phase}

The transition phase begins as the system is put into the environment in which it was designed to work. This may involve beta testing with advanced end users or one of several other techniques to move the system into a production environment.

Figure D-3 shows an example of the relative levels of work done during each of the phases in analysis, design, implementation, and testing. This example simply suggests the work levels; do not take them literally.

FIGURE D-3 Relative amounts of work done in each development phase


\section*{D. 3 What About the Waterfall Method of Development?}

The RUP is probably new to you, as is another development process-the waterfall method-in which the solution is developed by sequentially moving through phases such as requirements analysis, design, implementation, testing, and deployment. We think the waterfall method is inferior to the RUP. However, you doubtless will meet programmers during your career who know only the waterfall method. For this reason, you should know what it is.

The waterfall method often results in the following actions:
Systems analysts produce a set of documents specifying the requirements of the proposed system. The requirements documents are passed to a design team that produces a set of design documents, which are passed to programmers who code the design. The code is passed to a testing team, which verifies that the code correctly implements the requirements of the system, as specified by the systems analysts at the beginning of the process.

Notice that the phases occur one after the other. If the systems analysts do not understand everything about all aspects of the system and its environment, it is likely that they will specify the wrong system. Such aspects include the system's end users and their needs, other computer systems that must be used, the global business environment, and so on. Incorrect requirements analysis will create a design that, when implemented, leads to a correct system for the wrong problem. In other words, the system solves the problem defined by the requirements analysis, but the requirements analysis describes the wrong problem.

The difficulty with trying to specify the requirements of the proposed system up front is that system requirements are not predictable. Most things in the world today, especially in the high-tech and business worlds, are highly speculative. Change is constant. This implies that the requirements of any system might change at any time. The development process must allow the system under development to evolve and adapt to the changing environment. The waterfall method requires analysts to be omnipotent. For a system to succeed, the waterfall method should not be used for its development. Using iterative, incremental, evolutionary development processes, however, allows feedback to guide the adaptation of the system to a correct solution.

Be careful not to impose the waterfall method on the RUP development phases. For example, the RUP's inception phase is not the same as the waterfall requirements phase. During the inception phase, some, but not all, system requirements are specified, and programmers also make sure the system is feasible, estimate how much the system will cost to develop, and identify the significant highrisk issues that might doom the project. Trying to develop all use case scenarios during the inception phase makes it the same as the waterfall method.

Further, the RUP's elaboration phase is not the same as the waterfall design phase. During the elaboration phase, some, but not all, design is done, and programmers also do analysis, coding, and testing. If the elaboration phase degrades to having only one iteration, the development process effectively becomes the waterfall method. Similarly, the RUP's construction phase is not only about implementation, and it is not the same as the waterfall implementation phase.

The waterfall method is outdated and should not be used

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\section*{Mathematical Induction}

\section*{Appendix}

Many proofs of theorems or invariants in computer science use a technique called mathematical induction, or simply induction. Induction is a principle of mathematics that is like a row of dominoes standing on end. If you push the first domino, all the dominoes will fall one after another. What is it about the dominoes that allows us to draw this conclusion? If you know that when one domino falls the next domino will fall, then you know that pushing the first domino will cause them all to fall in succession. More formally, you can show that all the dominoes will fall if you can show that the following two facts are true:
- The first domino falls.
- For any \(k \geq 1\), if the \(k^{\text {th }}\) domino falls, the \((k+1)^{\text {th }}\) domino will fall.

The principle of mathematical induction is an axiom that is stated as follows:
Axiom E-1. The principle of mathematical induction. A property \(P(n)\) that involves an integer \(n\) is true for all \(n \geq 0\) if the following are true:
1. \(P(0)\) is true.
2. If \(P(k)\) is true for any \(k \geq 0\), then \(P(k+1)\) is true.

A proof by induction on \(\boldsymbol{n}\) is one that uses the principle of mathematical induction. Such a proof consists of the two steps given in Axiom E-1. The first step is called the basis, or base case. The second step is the inductive step. We usually break the inductive step into two parts: the inductive hypothesis ("if \(P(k)\) is true for any \(k \geq 0\) ") and the inductive conclusion ("then \(P(k+1)\) is true").

\section*{Example 1}

The following recursive function, which is given here in pseudocode, computes \(x^{n}\) :
```

pow2 (x: integer, $n$ : integer)
if ( $\mathrm{n}==0$ )
return 1
else
return $x$ * $\operatorname{pow2}(x, n-1)$

```

You can prove that pow2 returns \(x^{n}\) for all \(n \geq 0\) by using the following proof by induction on \(n\).

Basis. Show that the property is true when \(\boldsymbol{n}=\mathbf{0}\). That is, you must show that pow \(2(\mathrm{x}, 0)\) returns \(x^{0}\), which is 1 . As you can see from the definition of pow2, pow \(2(x, 0)\) is 1 , which establishes your basis.

Now you must establish the inductive step. By assuming that the property is true when \(n=k\) (the inductive hypothesis), you must show that the property is true when \(n=k+1\) (the inductive conclusion).
Inductive Hypothesis. Assume that the property is true when \(\boldsymbol{n}=\boldsymbol{k}\). That is, assume that
\[
\operatorname{pow} 2(\mathrm{x}, \mathrm{k})=x^{k}
\]

Inductive Conclusion. Show that the property is true when \(\boldsymbol{n}=\boldsymbol{k}+\mathbf{1}\). That is, you must show that pow \(2(\mathrm{x}, \mathrm{k}+1)\) returns the value \(x^{k+1}\). By definition of the function pow2,
\[
\operatorname{pow} 2(\mathrm{x}, \mathrm{k}+1)=x * \operatorname{pow} 2(\mathrm{x}, \mathrm{k})
\]

By the inductive hypothesis, pow2 ( \(\mathrm{x}, \mathrm{k}\) ) returns the value \(x^{k}\), so
\[
\begin{aligned}
\operatorname{pow} 2(\mathrm{x}, \mathrm{k}+1) & =x * x^{k} \\
& =x^{k+1}
\end{aligned}
\]
which is what you needed to show to establish the inductive step.
The inductive proof is thus complete. We demonstrated that the two steps in Axiom E-1 are true, so the principle of mathematical induction guarantees that pow2 returns \(x^{n}\) for all \(n \geq 0\). (End of proof.)

\section*{Example 2}

Prove that
\[
1+2+\cdots+n=\frac{n(n+1)}{2} \text { when } n \geq 1
\]

It will be helpful to let \(S_{n}\) represent the sum \(1+2+\cdots+n\).
Basis. Sometimes the property to be proven is trivial when \(n=0\), as is the case here. You can use \(n=1\) as the basis instead. (Actually, you can use any value of \(n \geq 0\) as the basis, but a value of 0 or 1 is typical.)

You need to show that the sum \(S_{1}\), which is simply 1 , is equal to \(1(1+1) / 2\). This fact is obvious.
Inductive Hypothesis. Assume that the formula is true when \(n=k\); that is, assume that \(S_{k}=k(k+1) / 2\).
Inductive Conclusion. Show that the formula is true when \(n=k+1\). To do so, you can proceed as follows:
\[
\begin{aligned}
S_{k+1} & =(1+2+\cdots+k)+(k+1) & & \left(\text { definition of } S_{k+1}\right) \\
& =S_{k}+(k+1) & & \text { (definition of } \left.S_{k}\right) \\
& =k(k+1) / 2+(k+1) & & \text { (inductive hypothesis) } \\
& =(k(k+1)+2(k+1)) / 2 & & \text { (common denominator) } \\
& =(k+1)(k+2) / 2 & & \text { (factorization) }
\end{aligned}
\]

The last expression is \(n(n+1) / 2\) when \(n\) is \(k+1\). Thus, if the formula for \(S_{k}\) is true, the formula for \(S_{k+1}\) is true. Therefore, by the principle of mathematical induction, the formula is true when \(n \geq 1\). (End of proof.)

\section*{Example 3}

Prove that \(2^{n}>n^{2}\) when \(n \geq 5\).
Basis. Here is an example in which the base case is not \(n=0\) or 1 , but instead is \(n=5\). It is obvious that the relationship is true when \(n=5\), because
\[
2^{5}=32>5^{2}=25
\]

Inductive Hypothesis. Assume that the relationship is true when \(n=k \geq 5\)-that is, assume that \(2^{k}>k^{2}\) when \(k \geq 5\).

Inductive Conclusion. Show that the relationship is true when \(n=k+1\)-that is, show that \(2^{k+1}\) \(>(k+1)^{2}\) when \(k \geq 5\). To do so, you can proceed as follows:
\[
\begin{aligned}
(k+1)^{2} & =k^{2}+(2 k+1) & & \text { (square } k+1) \\
& <k^{2}+k^{2} \text { when } k \geq 5 & & \left(2 k+1<k^{2}\right) \\
& <2^{k}+2^{k} \text { when } k \geq 5 & & \text { (inductive hypothesis) } \\
& =2 k+1 & &
\end{aligned}
\]

Therefore, by the principle of mathematical induction, \(2 n>n^{2}\) when \(n \geq 5\). (End of proof.)
Sometimes, the inductive hypothesis in Axiom E-1 is not sufficient. That is, you may need to assume more than \(P(k)\). The following axiom is a stronger form of the principle of mathematical induction:

Axiom E-2. The principle of mathematical induction (strong form). A property \(P(n)\) that involves an integer \(n\) is true for all \(n \geq 0\) if the following are true:
1. \(P(0)\) is true.
2. If \(P(0), P(1), \ldots, P(k)\) are true for any \(k \geq 0\), then \(P(k+1)\) is true.

Notice that the inductive hypothesis of Axiom E-2 ("If \(P(0), P(1), \ldots, P(k)\) are true for any \(k \geq 0\) ") includes the inductive hypothesis of Axiom E-1 ("If \(P(k)\) is true for any \(k \geq 0\) ").

\section*{Example 4}

Prove that every integer greater than 1 can be written as a product of prime integers.
Recall that a prime number is one that is divisible only by 1 and itself. The inductive proof is as follows:

Basis. The statement that you must prove involves integers greater than 1. Thus, the base case is \(n=2\). However, 2 is a prime number, and therefore it trivially is a product of prime numbers.
Inductive Hypothesis. Assume that the property is true for each of the integers \(2,3, \ldots, k\), where \(k \geq 2\).
Inductive Conclusion. Show that the property is true when \(n=k+1\); that is, show that \(k+1\) can be written as a product of prime numbers.

If \(k+1\) is a prime number, then there is nothing more to show. However, if \(k+1\) is not a prime number, it must be divisible by an integer \(x\) such that \(1<x<k+1\). Thus,
\[
k+1=x \times y
\]
where \(1<y<k+1\). Notice that \(x\) and \(y\) are each less than or equal to \(k\), so the inductive hypothesis applies. That is, \(x\) and \(y\) can each be written as a product of prime numbers. Clearly, the product \(x \times y\), which is equal to \(k+1\), must be a product of prime numbers. Because the formula holds for \(n=k+1\), it holds for all \(n \geq 2\) by the principle of mathematical induction. (End of proof.)

\section*{Example 5}

Chapter 2 discusses the following recursive definition:
```

$\operatorname{rabbit}(1)=1$
$\operatorname{rabbit}(2)=2$
$\operatorname{rabbit}(n)=\operatorname{rabbit}(n-1)+\operatorname{rabbit}(n-2)$ when $n>2$

```

Prove that
\[
\operatorname{rabbit}(n)=\left(a^{n}-b^{n}\right) / \sqrt{5}
\]
where \(a=(1+\sqrt{5}) / 2\) and \(b=(1-\sqrt{5}) / 2=1-a\).
Basis. Because \(\operatorname{rabbit}(0)\) is undefined, begin at \(n=1\). Some algebra shows that \(\operatorname{rabbit}(1)=\) \(\left(a^{1}-b^{1}\right) / \sqrt{5}=1\). However, notice that \(\operatorname{rabbit}(2)\) is also a special case. That is, you cannot compute \(\operatorname{rabbit}(2)\) from rabbit(1) by using the recurrence relationship given here. Therefore, the basis in this inductive proof must include \(n=2\).

When \(n=2\), some more algebra will show that \(\operatorname{rabbit}(2)=\left(a^{2}-b^{2}\right) / \sqrt{5}=1\). Thus, the formula is true when \(n\) is either 1 or 2 .

Inductive Hypothesis. Assume that the formula is true for all \(n\) such that \(1 \leq n \leq k\), where \(k\) is at least 2 .
Inductive Conclusion. Show that the formula is true for \(n=k+1\). To do so, you can proceed as follows:
\[
\begin{aligned}
\operatorname{rabbit}(k+1) & =\operatorname{rabbit}(k)+\operatorname{rabbit}(k-1) & & \text { (recurrence relation) } \\
& =\left[\left(a^{k}-b^{k}\right)+\left(a^{k-1}-b^{k-1}\right)\right] / \sqrt{5} & & \text { (inductive hypothesis) } \\
& =\left[a^{k-1}(a+1)-b^{k-1}(b+1)\right] / \sqrt{5} & & \text { (factorization) } \\
& =\left[\left(a^{k-1}\left(a^{2}\right)-b^{k-1}\right)\left(b^{2}\right) /\right] \sqrt{5} & & \left(a+1=a^{2} ; b+1=b^{2}\right) \\
& =\left(a^{k+1}-b^{k+1}\right) / \sqrt{5} & &
\end{aligned}
\]

Because the formula holds for \(n=k+1\), it holds for all \(n>2\) by the principle of mathematical induction. (End of proof.)

Note that the previous proof requires that you show that \(a+1=a^{2}\) and \(b+1=b^{2}\). Although simple algebra will demonstrate the validity of these equalities, exactly how did we discover them after the factorization step? Some experience with inductive proofs will give you the confidence to determine and verify the auxiliary relationships-such as \(a+1=a^{2}\)-that are necessary in a proof. Here, after we introduced the factors \((a+1)\) and \((b+1)\), we observed that if these factors were equal to \(a^{2}\) and \(b^{2}\), respectively, we could finish the proof. Thus, we tried to show that \(a+1=a^{2}\) and \(b+1=\) \(b^{2}\); indeed, we were successful. Inductive proofs often require adventurous algebraic manipulations!

\section*{Algorithm Verification} F

\section*{Prerequisite}

Appendix E Mathematical Induction

You can prove the correctness of some algorithms

Formal, theoretical techniques are available for proving that an algorithm is correct. Although research in this area is incomplete, it is useful to mention some aspects of the verification process.

An assertion is a statement about a particular condition at a certain point in an algorithm. Preconditions and postconditions are simply assertions about conditions at the beginning and end of methods and functions. An invariant is a condition that is always true at a particular point in an algorithm. A loop invariant is a condition that is true before and after each execution of an algorithm's loop. As you will see, loop invariants can help you write correct loops. By using invariants, you can detect errors before you begin coding and thereby reduce your debugging and testing time. Overall, invariants can save you time.

Proving that an algorithm is correct is like proving a theorem in geometry. For example, to prove that a method or function is correct, you would start with its precon-ditions-which are analogous to the axioms and assumptions in geometry-and demonstrate that the steps of the algorithm lead to the postconditions. To do so, you would consider each step in the algorithm and show that an assertion before the step leads to a particular assertion after the step.

By proving the validity of individual statements, you can prove that sequences of statements, and then methods and functions, classes, and finally the program are correct. For example, suppose you show that if assertion \(A_{1}\) is true and statement \(S_{1}\) executes, then assertion \(A_{2}\) is true. Also suppose you have shown that assertion \(A_{2}\) and statement \(S_{2}\) lead to assertion \(A_{3}\). You can then conclude that if assertion \(A_{1}\) is true, executing the sequence of statements \(S_{1}\) and \(S_{2}\) will lead to assertion \(A_{3}\). By continuing in this manner, you eventually will be able to show that the program is correct.

Clearly, if you discovered an error during the verification process, you would correct your algorithm and possibly modify the analysis and/or design. Thus, when you use invariants, your algorithm will likely contain fewer errors before you begin coding. As a result, you will spend less time debugging your program.

You can formally prove that particular constructs such as if statements, loops, and assignments are correct. An important technique uses loop invariants to demonstrate

Loop invariant

Steps to establish the correctness of an algorithm
the correctness of iterative algorithms. For example, we will prove that the following simple loop computes the sum of the first \(n\) elements in the array item:
```

// Computes the sum of item[0], item[1], . . ., item[n-1]
// for any n >= 1
int sum = 0;
int j = 0;
while (j < n)
{
sum += item[j];
j++;
} // end while

```

Before this loop begins execution, sum is 0 and j is 0 . After the loop executes once, sum is item[0] and j is 1 . In general,
sum is the sum of the elements item [0] through item[j-1]
This statement is the invariant for this loop.
The invariant for a correct loop is true at the following points:
- Initially, after any initialization steps, but before the loop begins execution
- Before every iteration of the loop
- After every iteration of the loop
- After the loop terminates

For the previous loop example, these points are as follows:
```

int sum = 0;
int j = 0;
while (j < n)
{

```

```

                            \longleftarrow The invariant is true here
    sum += item[j];
        j++;
    } // end while
\longleftarrow< The invariant is true here
$\longleftarrow$ The invariant is true here

```

You can use these observations to prove the correctness of an iterative (loop-controlled) algorithm. For the previous example, you must show that each of the following four points is true:
1. The invariant must be true initially, before the loop begins execution for the first time. In the previous example, sum is 0 and j is 0 initially. In this case, the invariant states that sum contains the sum of the elements item[0] through item[-1]; the invariant is true because there are no elements in this range.
2. An execution of the loop must preserve the invariant. That is, if the invariant is true before any given iteration of the loop, you must show that it is true after the iteration. In the example, the loop adds item[j] to sum and then increments \(j\) by 1 . Thus, after an execution of the loop, the most recent element added to sum is item \([j-1]\); that is, the invariant is true after the iteration.
3. The invariant must capture the correctness of the algorithm. That is, you must show that if the invariant is true when the loop terminates, the algorithm is correct. When the loop in the previous example terminates, \(j\) contains \(n\), and the invariant is true: sum contains the sum of the elements item[0] through item[n-1], which is the sum that you intended to compute.
4. The loop must terminate. That is, you must show that the loop will terminate after a finite number of iterations. In the example, \(j\) begins at 0 and then increases by 1 at each execution of the loop. Thus, j eventually will equal n for any \(\mathrm{n} \geq 1\). This fact and the nature of the while statement guarantee that the loop will terminate.

Not only can you use invariants to show that your loop is correct, but you can also use them to show that your loop is wrong. For example, suppose that the expression in the previous while statement was \(\mathrm{j}<=\mathrm{n}\) instead of \(\mathrm{j}<\mathrm{n}\). Steps 1 and 2 of the previous demonstration would be the same, but step 3 would differ: When the loop terminated, j would contain \(\mathrm{n}+1\) and, because the invariant would be true, sum would contain the sum of the elements item [0] through item[n]. Because this is not the desired sum, you know that something is wrong with your loop.

Notice the clear connection between steps 1 through 4 and mathematical induction. \({ }^{1}\) Showing the invariant to be true initially, which establishes the base case, is analogous to establishing that a property of the natural numbers is true for 0 . Showing that each iteration of the loop preserves the invariant is the inductive step. This step is analogous to showing that if a property is true for an arbitrary natural number \(k\), then the property is true for the natural number \(k+1\). After performing the four steps just described, you can conclude that the invariant is true after every iteration of the loopjust as mathematical induction allows you to conclude that a property is true for every natural number.

Identifying loop invariants will help you write correct loops. You should state the invariant as a comment that either precedes or begins each loop, as appropriate. For example, in the previous example, you might write the following:
```

// Invariant: 0 <= j <= n and
// sum = item[0] + ... + item[j-1]
while (j < n)

```

You should confirm that the invariants for the following unrelated loops are correct. Remember that each invariant must be true both before the loop begins and after each iteration of the loop, including the final one. Also, you might find it easier to understand the invariant for a for loop if you temporarily convert it to an equivalent while loop:
```

// Compute n! for an integer n >= 0
int factorial = 1;
// Invariant: factorial == (j-1)!
for (int j = 1; j <= n; j++)
factorial *= j;
// Compute an approximation of ex for a real x
doub7e term = 1.0;
doub7e series = 1.0;
int k = 1;
// Invariant: term == xk-1/(k-1)! and
// series == 1+x+x2/2!+...+x-1/(k-1)!
while (k <= n)
{
term *= x / k;
series += term;
k++;
} // end while

```

State loop invariants in your programs

Examples of loop invariants

1 A review of mathematical induction appears in Appendix E.

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\section*{C++ File Fundamentals}

\section*{Appendix}

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\section*{Prerequisite}

Appendix A Review of C++ Fundamentals

\section*{G. 1 File Input and Output}

You have used files ever since you wrote your first program. In fact, your C++ source program is in a file that you probably created by using a text editor. You can create and access such files outside of, and independently of, any particular program. Files can also contain data that is either read or written by your program. It is this type of file that concerns us here.

A file is a sequence of bytes that resides in auxiliary storage, often a disk. Files are useful because they can be large and can exist after program execution terminates. In contrast, variables, for example, represent memory that is accessible only within the program that creates them. When program execution terminates, the operating system reuses this memory and changes its contents.

Because files can exist after program execution, they not only provide a permanent record for human users, they also allow communication between programs. Program A can write its output into a file that program B can use later for input. However, files that you discard after program execution are not unusual. Chapter 16 discusses one such file that can help balance a binary search tree. You can also use such a file as a scratch pad during program execution when you have too much data to retain conveniently in memory all at once.

Files are classified as follows. A text file is a file of characters that are organized logically into lines. The files that you create-by using an editor-to contain your C++ programs are text files. Because text files consist of characters, and accessing characters by position number is usually not convenient, you typically process a text file sequentially. A file that is not a text file is called a binary file or, sometimes, a general file or nontext file.

A text file contains lines of characters

Files end with a special end-of-file symbol

Use a stream variable to access a file

\section*{G. 2 Text Files}

Text files are designed for easy communication with people. As such, these files are flexible and easy to use, but they are not as efficient with respect to computer time and storage as binary files.

One special aspect of text files is that they appear to be divided into lines. This illusion is often the source of much confusion. In reality, a text file-like any other file-is a sequence of bytes. That is, a text file is a sequence of characters. A special end-of-line character creates the illusion that a text file contains lines by making the file behave as if it were divided into lines.

When you create a text file by typing data at your keyboard, each time you press the Enter, or Return, key, you insert one end-of-line character into the file. When an output device, such as a printer or monitor, encounters an end-of-line character in a text file, the device moves to the beginning of the next line. In \(\mathrm{C}++\), this end-of-line symbol is the character \(\backslash \mathrm{n}\).

In addition, you can think of a special end-of-file character that follows the last component in a file. Such a symbol may or may not actually exist in the file, but C++ behaves as if one did. The predefined constant EOF, for "end of file," represents this symbol within your program. This book assumes that all text files-including the empty file-end with both an end-of-line symbol and an end-of-file symbol. Figure G-1 depicts a text file with these special symbols.

FIGURE G-1 A text file with end-of-line and end-of-file symbols


Any program that uses files must access the standard C++ file stream library. You enable this access by including the following statements in your program:
```

\#include <fstream>
using namespace std;

```

The C++ file stream library provides three stream types: ifstream for input file streams, ofstream for output file streams, and fstream for file streams that are for both input and output. You use a stream variable of one of these types to access a file. Files are sometimes referred to as streams since you can think of a file as a stream of characters flowing into or out of your program.

To use a file in your program, there are three simple steps:
1. Open the file.
2. Read data from the file or write data to the file.
3. Close the file.

\section*{G.2.1 Opening and Closing Text Files}

You must initialize, or open, a file before you can use it

Opening a file. Before you can read from or write to a file, you need to open it. That is, you need to initialize the file and associate its name with a stream variable. One way to open a file is to provide the file's name when you declare the stream variable. For example,
ifstream inFile("Ages.DAT"); // Input file
declares an input stream variable \(\mathrm{inFile}^{2}\) and associates it with the file named Ages.DAT. The file name can be a literal constant, as it is here, or a string variable. Once the file is open, you no longer need the file's name; it is used only to link the file on the disk to the stream variable that you use in your program.

Alternatively, you can declare an input stream variable by writing
ifstream inFile;
and then later use the open method to associate it with the file's name:
```

inFile.open("Ages.DAT");

```

Regardless of how you open a file, you can check the stream variable to see if the process was successful by writing
```

if (!inFile)
processError(); // Deal with failure to open file

```

Closing a file. You close a particular file-that is, disassociate it from a stream variable and free system resources-by using the close method:
```

myFile.close();

```

Once a file is closed, it is no longer available for input or output until a program opens it again.

\section*{G.2.2 Reading and Writing Text Files}

Associated with each file in a program is a file window, which marks the current position within the file. Opening a file positions the file window over the first byte in the file, as Figure G-2 illustrates. Because each byte in a text file is a character, the file window for a text file moves from character to character. The following sections describe the behavior of the file window.

FIGURE G-2 The file window after an existing text file is opened for input


Character input. Suppose that you have declared an input stream variable inFile and associated it with the name of the text file by writing
```

ifstream inFile(fileName);

```
where fileName is a string variable. When you use a file for input, the file window is over the byte that you will read next. Thus, after you open a file, you are ready to read the first byte, as you saw in Figure G-2. As you read the characters from a text file, the file window advances sequentially from one character to another. After reading several characters, you will see the file window shown in Figure G-3.

The input operator >> and methods such as get, whose use with cin was described in Appendix A, are also used with files. For the character variable ch and the stream variable inFile, either of the statements
```

inFile >> ch; // Preferred technique.

```

You can check whether a file was opened successfully

A file window marks the current position within a file

Opening a text file for input positions the file window over the first character in the file

FIGURE G-3 A file window over the component to be read next

or
inFile.get(ch);
means
\(\mathrm{ch}=\) the value at the file window
Advance the file window to the next component
as Figure G-4 illustrates.

FIGURE G-4 The effect of inFile >> ch on a text file inFile

Before inFile >> ch;


After inFile >> ch;


You can detect when the file window has reached either the end of a line or the end of the file by using the peek method. This method returns the character in the file window but does not advance the window to the next character. For example, the loop
```

while (inFile.peek() != '\n') // Loop until end of Tine.
cout << (inFile.get()); // Get character from inFile and display it.

```
displays a line of a text file, and the loop
```

while (inFile.peek() != EOF) // Loop until end of entire file.
cout << (inFile.get()); // Get character from inFile and display it.

```
displays the contents of an entire text file.

To summarize, consider the text file inFile that appears in Figure G-4. If ch is a character variable, the statements in the following sequence assign values to ch as indicated:
```

ifstream.inFile(fileName);
inFile >> ch; // ch = 'a'.
inFile.get(ch); // ch = 'b' - alternate technique 1.
ch = inFile.get(); // ch = 'c' - alternate technique 2.
inFile.ignore(10, '\n'); // Skip 'd' and end-of-file symbol.
inFile >> ch; // ch = 'e'.
ch = inFile.peek(); // ch = 'f'.
inFile >> ch; // ch = 'f'.
inFile >> ch; // ch = '\n'.
inFile.ignore(1); // Skip end-of-file symbol.
inFile >> ch; // Error: attempted read beyond end of file.

```

You also can read the characters in the file as strings by using getline, get, and read in the same way that you used them with cin.

Character output. Suppose that you have declared an output stream variable outFile and associated it with the name of the text file, as in
```

ofstream outFile(fileName);

```

If you are creating a new file, the file window will be positioned at the beginning (and the end) of the new file, which is empty. If the file already exists, opening it erases the data in the file and positions the window at the beginning-and the end-of the now empty file.

The output operator \(\ll\) and methods such as put, whose use with cout was described in Appendix A, are also used with files. For the character variable ch and the stream variable outFile, either of the statements
```

outFile << ch; // Preferred technique.

```
or
```

outFile.put(ch);

```
means
Write the value of ch at the file-window position Advance the file window

Figure G-5 illustrates these steps when ch contains the character \(X\). Note that if ch contains \(\backslash n\), either of the previous statements writes the end-of-line symbol to the file.

FIGURE G-5 The effect of outFile << ch on a text file outFile when ch contains the character \(X\)
\begin{tabular}{|c|c|c|c|c|c|c|c|c|c|c|}
\hline \multicolumn{11}{|r|}{Before outFile << ch; After outFile << ch;} \\
\hline & & & & & & & & & & T \\
\hline R & Z & W & Y & -•• & R & Z & W & Y & X & -• \\
\hline \multicolumn{11}{|c|}{\(\square\)} \\
\hline
\end{tabular}

You can also write strings to a text file by using either \(\ll\) or the method write, as described in Appendix A.

Numeric data within text files. As you know, you can read integer and floating-point values from the standard input stream into variables with arithmetic data types. You also know that the standard input stream is a sequence of characters. Likewise, a text file is a sequence of characters, so it should not surprise you that integer and floating-point values can be read from and written to a text file. Although this presentation uses int values to illustrate the concepts, the other arithmetic data types follow by analogy.

When your program reads from a text file into an int variable, the system expects a sequence of characters that it can convert into an integer. For example, if the text file contains the character sequence \(2,3,4\), and you read from the text file into the int variable \(x\), the system will convert these three characters into the computer's internal representation for the integer 234 and assign this value to x. More precisely, the text file contains the ASCII codes for the characters 2, 3, and 4-which are, respectively, the decimal values 50,51 , and 52 . However, these codes appear in the file in binary, as Figure G-6a indicates. If you read those characters into the integer variable \(\mathrm{x}, \mathrm{x}\) will contain the computer's internal representation for the integer 234, which appears in binary as shown in Figure G-6b. Thus, the representation of digits in a text file differs from the representation in memory of the number that those digits symbolize.

FIGURE G-6 (a) The ASCII characters 2, 3, and 4 represented in binary in a text file;
(b) the internal binary representation of the integer 234
(a) Text file

(b) Memory for the variable x


To summarize, if inFile is an input stream variable that is associated with a text file of valid integers, and \(x\) is an int variable, the statement
inFile >> \(x\);
has the following effect:

> Skip to the first nonblank character
> Convert into an integer the sequence of characters that begins at the current position of inFile's window and ends just before the next character c that is not a digit
> Assign this integer value to x
> Advance the file window so that it is over the character c

Figure G-7 illustrates these steps. Observe that if the sequence begins with a character other than,+- , or 0 through 9 , reading will terminate. For example, the system cannot convert the sequence w123 into an integer. It will, however, read the integer 123 from the sequence \(123 w r t\).

When your program writes an integer value such as 234 to a text file, the system first converts the integer from the computer's internal binary representation \((0000000011101010)\) to the character

FIGURE G-7 Reading an integer from a text file
Before inFile >> x;


After inFile >> x ;

123

sequence 2, 3, 4 and then writes these characters to the file. For the output stream variable outFile and the integer variable \(x\), the statement
```

outFile << x;

```
has the following effect:
Convert the value of x into a sequence of characters
Append this sequence of characters to the file
Position the file window just past the last character written
The following function reads and displays the contents of an entire text file of integers:
```

void echoFile(string fileName)
{
ifstream inFile(fileName);
int x = 0;
while (inFile >> x) // Read fails at end of file.
cout << x << " ";
cout << end7;
inFile.close();
} // end echoFile

```

The function ignores the end-of-line symbols in the file and displays the integers all on one line.
Suppose that you want to display each line in the file regardless of the number of integers per line. You can use the peek and ignore methods to accomplish this. Note that you always pass a stream variable to a function as a reference argument.
```

// Skips blanks in a text file.
void skipBlanks(ifstream\& inFile)
{
while (inFile.peek() == ' ') // While the current character is a blank:
inFile.ignore(1); // Ignore character, advance file window.
} // end skipBTanks
// Displays one line of a text file.
void echoLine(ifstream\& inFile)

```
```

{
int x = 0;
while (inFile.peek() != '\n') // While the current character is not eol:
{
inFile >> x; // Read the next integer into x.
cout << x << " "; // Display the number to user.
skipBlanks(inFile); // Skip blanks following number.
} // end while
inFile.ignore(1); // Advance beyond \n.
cout << "\n"; // For user display.
} // end echoLine
// Displays the contents of a text file.
void echoFile(string fileName)
{
ifstream inFile(fileName);
skipB1anks(inFile);
while (inFile.peek() != EOF)
echoLine(inFile);
inFile.close();
} // end echoFile

```

\section*{G.2.3 Manipulating Text Files}

Copying a text file. Suppose that you wanted to make a copy of the text file associated with the stream variable originalFile. Copying a text file requires some work and provides a good example of the statements you have just studied. The approach taken by the following function copies the file one character at a time, taking into account both the end-of-line symbols and the end-of-file symbol:
```

/*** Makes a duplicate copy of a text file.
@pre The name of an existing external text file and the name of the
text file to be created are given as strings.
@post The original text file is duplicated.
@param originalFileName Name of file to copy.
@param copyFileName Name of final copy. */
void copyTextFile(string originalFileName, string copyFileName)
{
ifstream originalFile(originalFileName); // Input file
ofstream copyFile(copyFileName); // Output file
char ch = '';
// Copy characters one at a time from given file to new file.
while (origina1File.get(ch)) // Fails at end of file.
copyFile << ch; // Write character to new file.
// Close the files.
originalFile.close();
copyFile.close();
} // end copyTextFile

```

Notice that this function copies each end-of-line symbol just as it copies any other character. To do so, the expression
```

origina1File.get(ch)

```
is necessary because
```

origina1File >> ch

```
skips whitespace, including the end-of-line symbol.

Adding to a text file. When you open a file, you can specify a second argument in addition to the file's name. This second argument has the form
```

ios::mode

```
where mode has values such as in, out, or app. Until now, this argument has been omitted because ifstream files are by default opened for input, whereas ofstream files are opened for output. You can append a component to an ofstream file outFile by first opening it in append mode. Either the statement
```

ofstream outFile("Sample.DAT", ios::app);

```
or the statements
ofstream outFile;
outFile.open("Samp1e.DAT", ios::app);
prepare the file for output and position the file window after the file's last component. Thus, the old contents of the file are retained, and you can write additional components.

Searching a text file sequentially. Suppose that you have a text file of data about a company's employees. For simplicity, assume that this file contains two consecutive lines for each employee. The first line contains the employee's name, and the next line contains data such as salary.

Given the name of an employee, you can search the file for that name and then determine the other information about this person. A sequential search examines the names in the order in which they appear in the file until the desired name is located. The following function performs such a sequential search. If the person is found, it returns that worker's salary; otherwise, it returns -1.0 .
```

doub7e getSalary(string fileToSearch, string desiredName)
{
bool found = false;
string nextName;
double nextSalary = -1.0;
ifstream inFile(fileToSearch);
while ( !found \&\& getline(inFile, nextName) )
{
inFile >> nextSalary; // Salary for nextName.
inFile.ignore(); // Ignores remaining whitespace in 7ine.
if (nextName == desiredName)
found = true;
} // end while
inFile.close(); // Always close file before return statement.
return nextSalary;
} // end getSalary

```

This function needs to look at all the names in the file before determining that a particular name does not occur. If the names were in alphabetical order, you could determine when the search passed the place in the file that should have contained the desired name, if it existed. In this way, you could terminate the search before you needlessly searched the rest of the file.

Accessing a text file directly. Although you usually process a text file sequentially, you can access the character stored at a given position directly without first reading the preceding characters.

The characters in a text file are numbered sequentially in order of appearance in the file, beginning with zero. The seekg method provides access to any character in the file, given the character's number. For example,
```

myFile.seekg(15)

```
advances the file window to the character numbered 15 , which is actually the \(16^{\text {th }}\) character in the file. Immediately following this operation, you can read the character.

You can also locate a character relative to either the beginning of the file, the current position in the file, or the end of the file by specifying-as the second argument to seekg-
ios::mode
where mode is one of beg, cur, or end. Thus,
```

myFile.seekg(2, ios::cur)

```
seeks the second character after the one at the present location of the file window.

\section*{G. 3 Binary Files}

Files that are not text files are called binary (or general or nontext) files. Like a text file, a binary file is a sequence of bytes grouped together into file components. It is important to emphasize that each file component is an indivisible entity. For example, each component of a binary file of integers is an integer in the computer's internal representation. If you write the integer value 234 to a binary file, the system would write the computer's internal representation of 234 , which is 0000000011101010 in binary, to the file, rather than the three ASCII characters 2, 3, 4, which are, respectively, 00110010 , 00110011 , and 00110100 in binary. Similar comments are true for a binary file of floating-point numbers. If you could use a text editor to look at the file, you would see gibberish. You create a binary file not by using an editor - as you can for a text file-but rather by running a program.

The statement
```

ofstream outFile(myFileName, ios::binary);

```
associates the stream variable outFile with the external binary file whose name is in the string variable myFileName. A binary file has an end-of-file symbol at its end, just as a text file does. However, the notion of lines does not exist for a binary file, although a binary file might contain data that coincidentally looks like an end-of-line symbol. Except for the differences noted here, \(\mathrm{C}++\) treats binary files in the same way that it treats text files.

\section*{C++ Header Files and Standard Functions}

\section*{Appendix}

Here is a list of commonly used \(\mathrm{C}++\) header files. Other header files are introduced in the chapters when needed.

\section*{cassert}

This library contains only the function assert. You use
assert(assertion);
to test the validity of an assertion. If assertion is false, assert writes an error message and terminates program execution. You can disable all occurrences of assert in your program by placing the directive \#define NDEBUG before the include directive.

\section*{cctype}

Most functions in this library classify a given ASCII character as a letter, a digit, and so on. Two other functions convert letters between uppercase and lowercase.

The classification functions return a true value if ch belongs to the specified group; otherwise they return false.
\begin{tabular}{ll} 
isalnum(ch) & Returns true if ch is either a letter or a decimal digit \\
isalpha(ch) & Returns true if ch is a letter \\
iscntrl (ch) & Returns true if ch is a control character (ASCII 127 or 0 to 31) \\
isdigit(ch) & Returns true if ch is a decimal digit \\
isgraph(ch) & Returns true if ch is printable and nonblank \\
islower(ch) & Returns true if ch is a lowercase letter \\
isprint(ch) & Returns true if ch is printable (including blank) \\
ispunct(ch) & Returns true if ch is a punctuation character \\
isspace(ch) & \begin{tabular}{l} 
Returns true if ch is a whitespace character: space, tab, carriage return, new line, \\
or form feed
\end{tabular} \\
isupper(ch) & Returns true if ch is an uppercase letter \\
isdigit(ch) & \begin{tabular}{l} 
Returns true if ch is a hexadecimal digit
\end{tabular} \\
toascii (ch) & Returns the ASCII code for ch
\end{tabular}
```

tolower(ch) Returns the lowercase version of ch if ch is an uppercase letter; otherwise
returns ch
toupper(ch) Returns the uppercase version of ch if ch is a lowercase letter; otherwise
returns ch

```

\section*{cmath}

The C++ functions in this library compute certain standard mathematical functions. These functions are overloaded to accommodate float, double, and long double. Unless otherwise indicated, each function has one argument, with the return type being the same as the argument type (either float, doub7e, or long double).
\begin{tabular}{|c|c|}
\hline \(\operatorname{acos}(x)\) & Returns the arc cosine \\
\hline \(\operatorname{asin}(x)\) & Returns the arc sine \\
\hline \(\operatorname{atan}(x)\) & Returns the arc tangent \\
\hline \(\operatorname{atan} 2(x, y)\) & Returns the arc tangent of \(x / y\) \\
\hline ceil (x) & Rounds up \\
\hline \(\cos (x)\) & Returns the cosine \\
\hline \(\cosh (x)\) & Returns the hyperbolic cosine \\
\hline \(\exp (x)\) & Returns \(e\) raised to the power \(x\) \\
\hline fabs (x) & Returns the absolute value \\
\hline floor (x) & Rounds down \\
\hline \(f m o d(x, y)\) & Returns x modulo y \\
\hline \(\log (\mathrm{x})\) & Returns the natural log \\
\hline \(\log 10(x)\) & Returns the log base 10 \\
\hline \(\operatorname{modf}(x, i p t r)\) & For arguments x and iptr, returns the fractional part of x and sets \(i p t r\) to point to the integer part of x \\
\hline \(\operatorname{pow}(x, y)\) & Returns \(x\) raised to the power \(y\) \\
\hline \(\sin (x)\) & Returns the sine \\
\hline \(\sinh (x)\) & Returns the hyperbolic sine \\
\hline sqrt (x) & Returns the square root \\
\hline \(\tan (\mathrm{x})\) & Returns the tangent \\
\hline \(\tanh (x)\) & Returns the hyperbolic tangent \\
\hline
\end{tabular}

\section*{cstd1ib}
\begin{tabular}{ll} 
abort() & Terminates program execution abnormally \\
abs \((x)\) & Returns the absolute value of an integer \\
atof(s) & Converts a string argument to floating point \\
atoi \((s)\) & Converts a string argument to an integer
\end{tabular}
\begin{tabular}{ll} 
exit( \(x)\) & Terminates program execution, returns \(x\) to the operating system \\
\(\operatorname{rand}()\) & Returns a pseudorandom integer \\
srand ( \(x\) ) & Initializes the pseudorandom number generator to \(x\) \\
srand() & Initializes the pseudorandom number generator to 1
\end{tabular}

\section*{cstring}

This library enables you to manipulate C strings that end in \(\backslash 0\). Unless noted otherwise, these functions return a pointer to the resulting string in addition to modifying an appropriate argument. The argument ch is a character, n is an integer, and the other arguments are strings.
\begin{tabular}{|c|c|}
\hline strncat(toS, fromS, n) & Copies at most n characters of froms to the end of toS and appends ' \(\backslash 0\) ' \\
\hline strcmp(str1, str2) & Returns an integer that is negative if \(\operatorname{str} 1<\mathrm{str} 2\), zero if \(\operatorname{str} 1==\) str2, and positive if str1> str2 \\
\hline stricmp(str1, str2) & Behaves like strcmp, but ignores case \\
\hline strncmp(str1, str2, n) & Behaves like strcmp, but compares the first n characters of each string \\
\hline strncpy (toS, fromS, n) & Copies n characters of fromS to toS, truncating or padding with ' \(\backslash 0\) ' as necessary \\
\hline strspn(str1, str2) & Returns the number of initial consecutive characters of str1 that are not in str2 \\
\hline strcspn(str1, str2) & Returns the number of initial consecutive characters of str1 that are in str2 \\
\hline strlen(str) & Returns the length of str, excluding ' \(\backslash 0\) ' \\
\hline strlwr(str) & Converts any uppercase letters in str to lowercase without altering other characters \\
\hline strupr(str) & Converts any lowercase letters in str to uppercase without altering other characters \\
\hline strchr (str, ch) & Returns a pointer to the first occurrence of ch in str; otherwise returns nullptr \\
\hline strrchr (str, ch) & Returns a pointer to the last occurrence of ch in str; otherwise returns nullptr \\
\hline strpbrk(str1, str2) & Returns a pointer to the first character in str1 that also appears in str2; othewise returns nul1ptr \\
\hline strstr (str1, str2) & Returns a pointer to the first occurrence of str2 in str1; otherwise returns nullptr \\
\hline strtok(str1, str2) & Finds the next token in str1 that is followed by str2, returns a pointer to the token, and writes nu11ptr immediately after the token in str1 \\
\hline
\end{tabular}

\section*{fstream}

Defines the \(\mathrm{C}++\) classes that support file I/O.

\section*{iomanip}

The manipulators in this library affect the format of stream operations. Note that iostream contains additional manipulators.
```

setbase(b) Sets number base to b=8,10, or 16
setfill(f) Sets fill character to f
setprecision(n) Sets floating-point precision to integer n
setw(n) Sets field width to integer n

```

\section*{iostream}

The manipulators in this library affect the format of stream operations. Note that iomanip contains additional manipulators.
dec Tells the subsequent operation to use decimal representation
end 1 Inserts the new-line character \(\backslash n\) and flushes the output stream
ends Inserts the null character \(\backslash 0\) into an output stream
flush Flushes an output stream
hex Tells the subsequent I/O operation to use hexadecimal representation
oct Tells the subsequent I/O operation to use octal representation
ws Extracts whitespace characters from the input stream

\section*{stdexcept}

Defines several exception classes that can be thrown or extended for specific exceptional conditions.

\section*{string}

This library enables you to manipulate \(\mathrm{C}++\) strings. Described here is a selection of the methods that this library provides. In addition, you can use the following operators with \(\mathrm{C}++\) strings: \(=,+,==,!=,<\), \(<=,>,>=, \ll\), and \(\gg\). Note that positions within a string begin at 0 .
\begin{tabular}{ll} 
erase() & Makes the string empty \\
erase(pos, len) & \begin{tabular}{l} 
Removes the substring that begins at position pos and contains len \\
characters
\end{tabular} \\
find (subString) & \begin{tabular}{l} 
Returns the position of the substring subString within the string \\
length() \\
replace(pos, 1en, str)
\end{tabular} \\
\begin{tabular}{l} 
Returns the number of characters in the string (same as size) \\
Replaces the substring that begins at position pos and contains 1en \\
characters with the string str
\end{tabular} \\
size() & \begin{tabular}{l} 
Returns the number of characters in the string (same as length) \\
substr(pos, len)
\end{tabular} \\
\begin{tabular}{l} 
Returns the substring that begins at position pos and contains len \\
characters
\end{tabular}
\end{tabular}

\section*{C++} Appendix

\section*{Documentation Systems} I

Documenting programs has never been a very enjoyable task. With the advent of the javadoc documentation system, commenting Java programs is at least less tedious. Inspired by javadoc, others have produced utility software to process documentation for other programming languages, including C++. Documentation tools like doxygen (doxygen.org) and DOC++ (docpp.sourceforge.net) read C++ source code and generate HTML-based documentation from javadoc-style comments embedded within the code. The comments in this book assume that doxygen will prepare the documentation.

The doxygen tool extracts the header for your class, the headers for all public methods, and comments that are written in a certain form. No method bodies and no private items are extracted.

For a comment to be extracted, the comment must satisfy two conditions:
- The comment must occur immediately before a public class definition or the header of a public method.
- The comment must begin with /** and end with */ and is a special form of the \(\mathrm{C}++\) multiline comment /* ... */.

Inside the comment, doxygen tags identify different aspects of the documentation, such as the programmer's name and a method's parameters and return value. Tags begin with the symbol @. We will describe only some of the available tags in this appendix. Note that some doxygen tags differ from the tags used by javadoc.
- @author lists the name of a class's programmer.
- @file identifies the name of the file containing a class.
- @param identifies a method's parameter by name and includes its description.
- @post identifies a method's postcondition.
- @pre identifies a method's precondition.
- @return describes a method's return value.
- @throw lists an exception that a method can throw.

You write a @param tag for each parameter in a method. You should list these tags in the order in which the parameters appear in the method's header. After the @param tag, you give the name and description of the parameter. For example, for the parameter customer, you might write
@param customer The string that names the customer.

You write a @return tag for every method that returns a value, even if you have already described the value in the method's description. Try to say something more specific about this value here. This tag must come after any @param tags in the comment. Do not use this tag for void methods and constructors.

Next, if a method can throw an exception, you name it by using a @throw tag. You list multiple exceptions alphabetically by name.

Here is a sample comment for a method. We usually begin such comments with a brief description of the method's purpose. This is our convention; javadoc has no tag for it.
```

/** Adds a new entry to a roster.
@param newEntry The object to be added to the roster.
@param newPosition The position of newEntry within the roster.
@pre None.
@post If the addition is successful, the roster contains the new entry;
otherwise, the roster is unchanged.
@return True if the addition is successful, or false otherwise. */

```

Here is the beginning of a sample class definition that lives in the file someClass.h:
```

/*% An example of a class.
@file someClass.h
@author Ima Student */

```

\section*{Programming Tip: Comments}
- Use /** ... */ comments before the headers of classes and methods to document their use and specifications. Use tags within a comment to identify aspects of the documentation.
- Use // for comments within the body of a class declaration or method definition that describe the details of the implementation.
- Use /* ... */ during debugging to temporarily disable a portion of your code.

\section*{Appendix}

\section*{ASCII}

\section*{Character Codes}
\begin{tabular}{|c|c|c|c|c|c|c|c|}
\hline Code & Character & Code & Character & Code & Character & Code & Character \\
\hline 0 & NUL & 32 & (blank) & 64 & @ & 96 & '(reverse quote) \\
\hline 1 & STX & 33 & ! & 65 & A & 97 & a \\
\hline 2 & SOT & 34 & " & 66 & B & 98 & b \\
\hline 3 & ETX & 35 & \# & 67 & C & 99 & c \\
\hline 4 & EOT & 36 & \$ & 68 & D & 100 & d \\
\hline 5 & ENQ & 37 & \% & 69 & E & 101 & e \\
\hline 6 & ACK & 38 & \& & 70 & F & 102 & f \\
\hline 7 & BEL & 39 & ' (apostrophe) & 71 & G & 103 & g \\
\hline 8 & BS & 40 & ( & 72 & H & 104 & h \\
\hline 9 & HT & 41 & ) & 73 & I & 105 & i \\
\hline 10 & LF & 42 & * & 74 & J & 106 & j \\
\hline 11 & VT & 43 & + & 75 & K & 107 & k \\
\hline 12 & FF & 44 & , (comma) & 76 & L & 108 & 1 \\
\hline 13 & CR & 45 & - & 77 & M & 109 & m \\
\hline 14 & SO & 46 & . & 78 & N & 110 & n \\
\hline 15 & SI & 47 & 1 & 79 & O & 111 & o \\
\hline 16 & DLE & 48 & 0 & 80 & P & 112 & p \\
\hline 17 & DC1 & 49 & 1 & 81 & Q & 113 & q \\
\hline 18 & DC2 & 50 & 2 & 82 & R & 114 & r \\
\hline 19 & DC3 & 51 & 3 & 83 & S & 115 & s \\
\hline 20 & DC4 & 52 & 4 & 84 & T & 116 & t \\
\hline 21 & NAK & 53 & 5 & 85 & U & 117 & u \\
\hline 22 & SYN & 54 & 6 & 86 & V & 118 & v \\
\hline 23 & ETB & 55 & 7 & 87 & W & 119 & w \\
\hline 24 & CAN & 56 & 8 & 88 & X & 120 & x \\
\hline 25 & EM & 57 & 9 & 89 & Y & 121 & y \\
\hline 26 & SUB & 58 & : & 90 & Z & 122 & z \\
\hline 27 & ESC & 59 & ; & 91 & [ & 123 & \{ \\
\hline 28 & FS & 60 & \(<\) & 92 & 1 & 124 & | \\
\hline 29 & GS & 61 & \(=\) & 93 & ] & 125 & \} \\
\hline 30 & RS & 62 & > & 94 & \(\wedge\) & 126 & \(\sim\) \\
\hline 31 & US & 63 & ? & 95 & _ (underscore) & 127 & DEL \\
\hline
\end{tabular}

Note: The codes 0 through 31 and 127 are for control characters that do not print.

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\section*{C++ for Java Programmers}

\section*{Appendix} \(T\)

\section*{Contents}
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This appendix covers some of the differences between Java and C++ that a beginning programmer may encounter. While there are many additional differences, they either are not applicable to the code and examples used in this text or are explained when needed. This appendix should be used in conjunction with Appendix A to fully understand the features provided in \(\mathrm{C}++\).

\section*{K. 1 General Concepts}

Transitioning from Java to C++ is not too difficult, since most of the syntax and many of the concepts are the same. C++ gives you much greater control of how and where you create variables and objects. Thus, a C++ programmer can create programs that are flexible and efficient in speed and use of storage. This flexibility and greater control require you to take great care in managing data and memory, because C++ does not provide garbage collection.

Both Java and C++ are compiled languages. While Java programs are compiled into byte-code for the Java Virtual Machine (JVM), C++ programs are compiled into object code that is specific to an operating system. This means that you need to recompile your \(\mathrm{C}++\) code for each operating system that you want your program to run on. The code in this book follows the recent ANSII C11 Standard, so you should be able to compile and run your programs on whichever operating system you choose if your compiler supports that standard.

Java classes and interfaces can be grouped into packages that are stored in the same directory hierarchy. When your Java code is compiled, the compiler searches this hierarchy for classes that you use. Any class or method in the package is automatically
found by the compiler and later the JVM during program execution. If you need a class outside of the package, you must import it using the Java import statement.

The \(\mathrm{C}++\) compiler is stricter and requires the programmer to declare all functions, classes, and variables in a file before they are used. Because classes and constants are often used in multiple programs and multiple locations in a program, \(\mathrm{C}++\) provides a mechanism called the pre-processor that lets you specify classes and data in a separate header file and then include that file in your program. Commands, or directives, to the pre-processor are lines that begin with the character \(\#\), such as
\#include <iostream> // Include the declaration of the iostream classes \#include "MyClass.h" // Include the declaration of MyClass

Notice that system files use < >, while user-defined files use " ", to delimit the file name. Section A.8.1 of Appendix A, C++ Interlude 1, and Section K. 4 of this appendix show how to develop C++ class declarations.

Although both Java and C++ are object-oriented languages, C++ has many elements that reflect its roots in the imperative language C. An imperative programming language does not use classes to group data and operations on that data. Instead, it is a sequence of statements and functions-methods that are not part of a class. As such, variables, constants, and functions in a C++ program can be declared anywhere in a file; you are not forced to declare them within a class, as you are in Java. The only restriction to this capability is that a function cannot declare them inside another function.
Comments. Comments are an important part of any program, and \(\mathrm{C}++\) supports the same commenting standards as Java:
```

// This is a comment that appears on its own Tine.
int sum = 0; // This is a comment that appears within a C++ statement
// The following multiline comments end in */.
/** This is a special multiline comment used for documentation systems
such as doxygen. It begins with /**. See Appendix I for details. */
/* This is a multiline comment that begins with /*.
This sty7e is often used to temporarily disable a group of
statements during debugging. */

```

\section*{K. 2 Data Types, Variables, Constants, and Arrays}

C++ uses primitive data types and variable declarations that are similar to those in Java. The Java types char, int, float, and double are used in C++ and behave the same as in Java. In C++, int can be modified with short, long, signed, and unsigned. All Java variables are initialized to either zero, if a primitive data type, or nu11, if an object reference. In contrast, you must initialize all C++ variables. If you do not, and an uninitialized variable is used, the result of the equation or statement is unpredictable.

Boolean types. One very important difference between Java and C++ data types occurs with boolean variables. In Java, boolean is a distinct type that has the value true or false. In C++, the analogous data type is bool. It also has a value of true or fa1se, but in C++, true and false are special integers. That is, variables of type bool are actually integers. This affects the type of statement that is used as the predicate or test clause in an if statement, for loop, or while loop, as you will see in the next section. The C++ value of false is 0 , and an integer value of 0 can be considered false in a boolean test. The value of true is 1 , and any non-zero integer is considered true.
String types. Java has a special class String to represent character strings. C++ also provides a special class string (note the lower case s) that implements many of the same features. To use this class, you must write
as one of the first lines in your program. More information on the \(\mathrm{C}++\) class string is in Section A. 7 of Appendix A.

Constants. Constants in Java are declared within a class or method using the syntax
```

public static final int MAX_SIZE = 20;

```
which creates a publicly accessible constant MAX_SIZE with the value 20. A similar statement in C++ uses the keyword const to indicate a constant value:
const int MAX_SIZE = 20;
Constants in \(\mathrm{C}++\) are available anywhere within the scope of their declaration. If a constant is declared outside of a class or file, that constant is available to any function or class defined after it in the file and is said to be global, even though it is restricted to that file.

Java has a constant nu 17 that is used to indicate that an object has not been assigned to a variable. \(\mathrm{C}++\) uses nul1ptr for this purpose.

Arrays. Although you declare and reference arrays in Java and C++ using similar statements, their implementations are very different. Arrays in Java are objects and have a public data field 1ength that contains the number of locations in the array. If a program attempts to access a position beyond the end of the array, an exception is thrown.
\(\mathrm{C}++\) arrays are not objects, but are sequences of memory locations of the data type declared for the array. Entries in a C++ array are accessed using [ ], just as they are in Java. C++ arrays have no data field that tracks the number of positions in the array, and most importantly, \(\mathrm{C}++\) permits the programmer to access positions in the array that do not exist. This is a very common source of errors in a \(\mathrm{C}++\) program and is one of the areas where you must take great care. The STL class vector, which is discussed in Section A.6.3 of Appendix A, is an alternative to C++ arrays that provides some of the features of the Java array.

The following statement declares an array of 20 integers in Java:
```

int[] myArray = new int[MAX_SIZE];

```

This statement does the same thing in \(\mathrm{C}++\) :
```

int myArray[MAX_SIZE];

```

You can also declare a C++ array by using syntax similar to that of Java:
```

int myArray[] = new int[MAX_SIZE];

```

One difference is the location of the square brackets, but you must also take responsibility for deleting the array when you are finished using it. You do so with the de1ete statement:
```

delete[] myArray;

```

Notice the [ ] following the de1ete operator to indicate that myArray is an array. In C++ Interlude 2, we discuss the use of the operators new and de1ete in \(\mathrm{C}++\).

\section*{K. 3 Flow Control and Functions}

Flow control. Since the syntax of Java is derived from C++, the structure of flow-control statements in C++ is the same as those you are familiar with in Java. This is true for the selection statements if, if-e1se, and switch. It is also true for loop statements while, do-while, and for. The significant differences in \(\mathrm{C}++\) are related to the test expression that controls the selection or repetition.

Consider the following Java if-el se statement that compares the integer \(x\) to 6 and increments \(y\) by 10 if \(x\) is 6 , and by 20 otherwise:
```

if (x == 6)
y += 10;
else
y += 20;

```

Because the syntax for an if-else statement is the same in \(\mathrm{C}++\), we could place this statement in a \(\mathrm{C}++\) program and it would behave in the same way.

The test \(x==6\) is a predicate expression that returns true if the variable \(x\) is equal to 6 and false otherwise. In Java, a predicate expression such as this returns a value of type boolean. A Java flowcontrol statement requires its test expression to have a boolean value, so you must use a predicate expression as the test expression.

A C++ flow-control statement does not restrict the data type of its test expression. Flow-control decisions are made based on whether the expression returns zero or nonzero. If the test expression returns zero, the test is false. Any nonzero value is considered true. The equality operator in the example above returns true, which is 1 , if \(x\) has the value of 6 . Otherwise, it returns false, which is 0 . The if-e 1 se statement determines its course of action based on these values.

This behavior of the \(\mathrm{C}++\mathrm{if}-\mathrm{e} 7\) se statement gives you great flexibility in constructing tests for flow-control statements. It also can create confusion and subtle logic errors. For example, consider the following if-e 1 se statement:
```

if (x = 6)
y += 10;
else
y += 20;

```

At first glance, this may appear to be the same as the earlier example, but upon closer inspection you can see that the test expression is the assignment statement \(x=6\). The Java compiler will reject this as an error, since it is not a predicate expression. The \(\mathrm{C}++\) compiler will accept this statement, and when the program runs, the if-el se statement will be evaluated as follows:
```

if (x = 6) // The variable x is assigned the value of 6.
y += 10; // 6 is non-zero, so it is "true"; increment y by 10.

```

Every time this code executes, \(x\) will be assigned the value 6 , and \(y\) will be incremented by 10 . The original value of \(x\) is overwritten by the if clause's test expression and is never used! Tracking down subtle errors, such as this one, has caused this author many sleepless nights. When performing equality tests in C++ flow-control statements, always check that the test uses \(==\) and not \(=\).

Functions. As mentioned earlier, \(\mathrm{C}++\) functions do not need to be part of a class. Thus, you do not need to precede C++ functions with access modifiers, as you would for Java methods. Otherwise, the syntax for the declaration and invocation of a C++ function is similar to that of a Java method. Functions can be defined anywhere inside a \(\mathrm{C}++\) file, except inside another function.

Here is an example of a \(\mathrm{C}++\) function, realCube, that has a parameter of type double and returns the cube of a number:
```

doub7e rea1Cube(doub7e x)
{
return x * x * x;
} // end realCube

```

Any function defined after rea1Cube in the file can invoke rea1Cube. But what if a function is defined before realCube and needs to call it? In this case, the compiler would issue an error stating that
realCube was not defined. C++ requires that each function, constant, variable, or class be defined before it can be called.

In a simple program with a few functions, it is easy to define the functions in the reverse of the order in which they are needed so the compiler sees each function definition before it is needed. In more complex programs, this order is more difficult or even impossible to achieve.

The part of a function definition that the compiler needs is the header of the function and C++ provides a mechanism that allows the programmer to provide the compiler a function's header separate from the actual definition of the function. This mechanism is the function prototype and consists of the function's header followed by a semicolon. For example, the function prototype of the function realCube is
doub7e rea1Cube(doub7e x);
To eliminate the chance that a function will be needed prior to its definition within a \(\mathrm{C}++\) file, a common practice is to list the functions' prototypes at the beginning of the file. In larger programs, function prototypes are placed in header files, as described in Section K.1. You then include the header files in your program. Prototypes are also important in the creation of classes, as you will see in the next section.

Note that you can pass a function as an argument to another function or method. This is discussed in Section 15.2.1 of Chapter 15.

\section*{K. 4 Classes}

Many object-oriented concepts implemented in Java classes also apply to C++ classes. This appendix covers only some of the syntax differences that you may encounter in a simple class. Additional information about C++ classes appears in Section A. 8 of Appendix A and in C++ Interludes 1 and 2.

Consider the following Java class:
```

public class PlainBox
{
private String item;
public PlainBox()
{
} // end defau7t constructor
public PlainBox(String theItem)
{
item = theItem;
} // end constructor
public void setItem(String theItem)
{
item = theItem;
} // end setItem
public String getItem()
{
return item;
} // end getItem
} // end PlainBox class

```

Though is it possible to define a class in \(\mathrm{C}++\) in a single file, we will use two files, since this is a more common and flexible approach. The first file, is the class header file; it contains the data fields
for the class and the prototypes of all methods defined in the class. It is a description of the class and does not normally contain executable code. For example, here is the header file for the C++ class PlainBox:
```

/** @file PTainBox.h */
\#include <string> // Needed because we use string objects.
using namespace std; // Strings are part of the system std 7ibrary.
class PlainBox // This is a class declaration for PlainBox.
{
private: // Items defined below have private access.
string item;
pub1ic: // Items defined below have public access.
// Prototypes for constructors and methods.
PlainBox();
PlainBox(string theItem);
void setItem(string theItem);
string getItem();
}; // end P7ainBox - Note semicolon after declaration.

```

Notice that each data member and method does not need to have an access modifier. Private data members and methods are grouped together, and public methods are grouped with each other. A colon follows each access modifier. It is possible to have more than one private, protected, or public section, but it is not common. The class declaration must end with a semicolon after the closing brace, as shown above.

The implementation file, or source file, for the class typically has the same name as the header file, but ends in .cpp to indicate that it is a \(\mathrm{C}++\) file. The implementation file contains the definitions for the methods declared in the accompanying header file. Here is the implementation file for the \(\mathrm{C}++\) class PlainBox:
```

/** @file PlainBox.cpp */
\#include "PlainBox.h" // We need to tell the compiler about our data
// fields and methods.
PlainBox::PlainBox()
{
} // end default constructor
PlainBox::PlainBox(string theItem)
{
item = theItem;
} // end constructor
void PlainBox::setItem(string theItem)
{
item = theItem;
} // end setItem
string PlainBox::getItem()
{
return item;
} // end getItem

```

This implementation file begins by including the header file, so that the compiler has the class declaration and knows the data fields and methods of this class. We then define each method. In the implementation file there is no "master" set of braces enclosing all of the methods. Each method is
individually defined. To let the compiler know that a method belongs to the P1ainBox class, we added the prefix PlainBox: : to the method name. Observe that the prefix comes after the method's return type. The prefix is an extension of the method name only.

To create a P1ainBox object in Java, you could use the following statement:
PlainBox myBox = new PlainBox("Jewe1ry");
\(\mathrm{C}++\) provides several forms for creating objects also:
PlainBox myBox = PlainBox("Jewelry");
PlainBox myBox("Jewe1ry");
These statements are equivalent in \(\mathrm{C}++\) and create a PlainBox object that is local to the function or method in which it is created. In these examples the \(\mathrm{C}++\) keyword new is not used.

The following example does use the keyword new to create a PlainBox object that exists until the programmer specifically deletes it:
```

P1ainBox* someBoxPtr = new PlainBox("Jewe1ry");

```

The syntax is very similar to Java's, except that PlainBox is followed by the * character. The * indicates that someBoxPtr is a pointer to a PlainBox object and is not the object itself. Pointers are discussed in more detail in C++ Interlude 2.

If an object in \(\mathrm{C}++\) is instantiated using the new operator, as in the example above, it is the responsibility of the programmer to delete that object from memory when it is no longer needed, because \(\mathrm{C}++\) does not provide garbage collection. To delete the object pointed to by someBoxPtr, use the statements
```

de1ete someBox; // Frees memory used by the object.
someBox = nullptr; // A safety precaution.

```

Each time you use the new operator to create an object, you must write a corresponding delete operator to free the object's memory. Otherwise, your application will have a memory leak.

\section*{K. 5 Basic Input and Output}

Although at the system level, Java and C++ treat keyboard input and console-display output similarly, the high-level constructs used to accomplish these tasks differ greatly.

For a Java application to accept input from the keyboard, you create a Scanner object with System. in and use it to accept input from the user. Output to the display is accomplished by calling either the print or println method of the System. out object. The following code creates a Scanner object, displays a prompt for the user, accepts a response from the user, and echoes the response back to the user:
```

Scanner keyboard = new Scanner(System.in);
System.out.print("How many apples are in a box? ");
int applesPerBox = keyboard.nextInt();
System.out.print("You entered " + applesPerBox + " apples per box. ");

```

C++ does not have a System. in or System. out object; instead you use cin and cout objects for keyboard input and console-display output, respectively. To use cin and cout, you include the following statements at the start of your program:
```

\#include <iostream>
using namespace std;

```

The following \(\mathrm{C}++\) statements are equivalent to the previous Java statements:
```

cout << "How many apples are in a box? ";
int applesPerBox = 0;
cin >> applesPerBox;
cout << "You entered " << applesPerBox << " apples per box. " << endl;

```

Notice the use of the output stream operator \(\ll\) to send string literals and variables to the output object cout. Similarly, the input stream operator \(\gg\) sends input from the input object cin to a variable. It is easy to remember which operator to use if you imagine them as arrows pointing in the direction of the data flow.

The constant end 1 is defined in iostream and represents the new-line character. Additional information about the iostream library, and about basic input and output in \(\mathrm{C}++\), is in Section A. 2 of Appendix A.

\section*{C++ for Python Programmers}

\section*{Appendix}


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This appendix covers some of the differences between Python and C++ that a beginning programmer may encounter. While there are many additional differences, they either are not applicable to the code and examples used in this text or are explained when needed. This appendix should be used in conjunction with Appendix A to fully understand the features provided in \(\mathrm{C}++\).

\section*{L. 1 General Concepts}

Transitioning from the Python programming language to \(\mathrm{C}++\) is similar to graduating from high school and going through basic training in the military. Python is a weakly typed, interpreted language with a readable syntax. C++, on the other hand, is a strongly typed, compiled language that follows strict punctuation rules for its syntax. The good news is that after completing boot camp, you have the chance to become an Army Ranger or Navy Seal. This text can help you be all you can as a computer scientist and C++ programmer.

Python programs are interpreted and can execute on any computer that has a Python interpreter. Python interpreters have even been written in JavaScript, so that you can run Python code inside a window in your web browser! C++ programs, on the other hand, are compiled into object code that is specific to an operating system. This means that you need to recompile your \(\mathrm{C}++\) code for each operating system that you want your program to run on. The code in this book follows the recent ANSII C11 Standard, so you should be able to compile and run your programs on whichever operating system you choose if your compiler supports that standard.

\section*{L. 2 Comparison of Language Elements}

When it comes to writing a C++ program, there are several important differences you need to keep in mind.

Comments. Comments are an important part of any program, and C++ supports single and multiline comments with different delimiters:
```

// This is a comment that appears on its own line.
int sum = 0; // This is a comment that appears within a C++ statement.
// The following comment begins with /** and ends with */.
/** This is a special multiline comment used for documentation systems
such as doxygen. See Appendix I for details. */
// The following comment begins with /* and ends with */.
/* This is a multiline comment that is often used to temporarily
disable a group of statements during debugging. */

```

Statements. White space and indentation in C++ have no significance. C++ statements end with a semicolon. Statements in a block, known as a compound statement, are analogous to a suite in Python and are grouped together by braces \{ \}.

Variables. In Python, a variable can be used as soon as it is assigned a value. Variables in C++ must be declared and given a type before they can be used. Though the C++ compiler does not require it, variables should also be given an initial value before being used. Once a variable has been declared with a given a type, it maintains that type throughout the scope in which it was declared.

Relational operators. Variables in C++ can be used only with one relational operator at a time. Python expressions such as \(5<x<10\) must be written as \((5<x) \& \&(x<10)\) in \(C++\). The operator \&\& is used for a logical AND, and \| is used for a logical OR.

Predicate expressions. Predicate expressions in conditional and iteration flow-control statements must be in parentheses.

Assignment statements. Multiple assignment is not allowed in \(\mathrm{C}++\). Python statements such as \(\mathrm{x}, \mathrm{y}\) \(=5,6\) cannot be used.

Iteration statements. \(\mathrm{C}++\) has a while statement that behaves like the Python while statement. The C++ for statement is a special case of the while statement. To use a C++ for statement as you would use a Python form statement, you must use iterators, which are explained in C++ Interlude 6. C++ has a do-whi le statement, which is a post-test loop. Python has no equivalent statement.

Basic input and output. In a Python application, you use the raw_input function to accept input from the keyboard. This function returns a string. If you need an integer or a value of some other data type, you must use another function to transform the string into the desired input. Output to the display is accomplished by calling the print function. For example, the following Python code displays a prompt for the user, accepts a response as an integer, and echoes the response back to the user:
```

userNumber = int (raw_input ("How many apples are in a box? "))
print "You entered %d apples per box."%( userNumber)

```

C++ uses two objects, cin and cout, for keyboard input and console display output, respectively. To use cin and cout, you should include the following statements at the start of your program:
```

\#include <iostream>
using namespace std;

```

The following C++ statements that are equivalent to the previous Python statements:
```

cout << "How many apples are in a box? ";
int applesPerBox = 0;
cin >> applesPerBox;
cout << "You entered" << applesPerBox << " apples per box." << endl;

```

Notice the use of the output stream operator \(\ll\) to send string literals and variables to the output object cout. Similarly, the input stream operator >> sends input from the input object cin to a variable. It is easy to remember which operator to use if you imagine them as arrows pointing in the direction of the data flow.

The constant endl is defined in iostream and represents the new-line character. Additional information about the iostream library and basic input and output in C++ is in Section A. 2 of Appendix A.

Example. Listing L-1 gives a short interactive Python program using the statements and expressions just discussed. Listing L-2 shows the equivalent program in C++. C++ programs begin execution at the main function. You can see some of the differences in user input and output in this example also.

\section*{LISTING L-1 Short Python program to test and modify a user-entered value}
```

sma11, large = 1, 1000
print "Your number was between %d and %d large." %(sma11, large)
userNumber = int (raw_input ("Enter something: "))
if small < userNumber < large :
print "Your number was between %d and %d large."%(sma11, large)
elif userNumber < smal1 :
while userNumber <= small :
userNumber = userNumber + 1

```
else :
    while True :
        userNumber = userNumber - 1
        if userNumber < 1000 :
            break
        print "I fixed your number to be between \%d and \%d!" \%(sma11, 1arge)

\section*{LISTING L-2 Short C++ program equivalent to the Python program in Listing L-1.}
```

\#include <iostream>
using namespace std;
int main()
{
int sma11 = 1;
int large = 1000;
int userNumber = 0;
cout << "Enter something: ";
cin >> userNumber;

```
```

    if ((sma11 < userNumber) && (userNumber < 1arge) )
    {
        cout << "Your number was between " << smal1 << " and ";
        cout << large << end1;
    }
    else if (userNumber < 1)
    {
    while (userNumber <= 1)
    {
        userNumber++;
    }
    }
e1se
{
do
{
userNumber--;
} while (userNumber >= 1000);
cout << "I fixed your number to be between " << smal1 << " and ";
cout << large << end7;
}
}

```

\section*{L. 3 Functions}

As with its variables, \(\mathrm{C}++\) functions are strongly typed. When a \(\mathrm{C}++\) function is defined, the parameters are given a type. If the function returns a value, the function must indicate the type of the return value. This information is given in the header, that is, the first line, of the function. If a function does not return a value, its return type is void. The definition of the function follows the header and is enclosed in \{ \}, as in other C++ blocks.

Here is a short Python function that computes the cube of its single parameter:
```

def realCube(x) :
return x * x * x

```

Here is the equivalent \(\mathrm{C}++\) function:
```

doub7e rea1Cube(doub7e x)
{
return x * x * x;
}

```

Notice that the function and its parameter are each given a type. C++ functions are discussed in greater detail in Section A. 2 of Appendix A.

Any function defined after realCube in the file can invoke realCube. But what if a function is defined before realCube and needs to call it? In this case, the compiler would issue an error stating that realCube was not defined. C++ requires that each function, constant, variable, or class be defined before it can be called.

In a simple program with a few functions, it is easy to order the function definitions so that the compiler sees each one before it encounters the function call. In more complex programs, this is more difficult, or even impossible, to achieve. However, the compiler needs only the header of a function before processing a call to it, and C++ provides a way for you to satisfy the compiler. You simply write
the function's header followed by a semicolon-that is, its prototype-at the beginning of the program that calls it. For example, the prototype of the function realCube is
```

doub7e rea1Cube(doub7e x);

```

Then you can place the function's definition anywhere within your program.
In larger programs, function prototypes are placed into one or more header files, as described in Section K. 1 of Appendix K. You then include the header files in your program. Prototypes are also important in the creation of classes, as you will see in the next section. Section A.8.1 of Appendix A also discusses header files.

Note that you can pass a function as an argument to another function or method. This is discussed in Section 15.2.1 of Chapter 15.

\section*{L. 4 Classes}

Python classes permit clients of the classes to access all data fields and methods. In \(\mathrm{C}++\), this is called public access and is recommended only for methods. In Python, when you create a class member that should be used only by class methods, you typically begin the member's name with two underscores, _. This convention is only a signal to the programmer; the Python interpreter does not enforce it. C++ permits the class designer to restrict access to certain data fields and methods by labeling them as private, and the compiler enforces this restriction by permitting only class methods to access private members.

Classes in C++ have constructors that are similar to the __init__ method in a Python class. When writing a method in a Python class, you precede the names of data fields by self to indicate that they are instance variables of the class. Instance variables in a \(\mathrm{C}++\) class do not need a special prefix, if they are referenced from within a method of the class.

Here is a simple Python class, P 7 ainBox:
```

class PlainBox
def __init__(self, theItem):
self.item = theItem
def setItem(self, theItem):
self.item = theItem
def getItem(self) :
return self.item

```

You also can define the same class in a form that is closer to that of a \(\mathrm{C}++\) class:
```

class PlainBox
__item = " "
def ___init__(self, theItem):
self.__item = theItem
def setItem(se1f, theItem):
self.__item = theItem
def getItem(se1f) :
return self. item

```

Though is it possible to define a class in \(\mathrm{C}++\) in a single file, we will use two files, since this is a more common and flexible approach. The first file is the class header file; it contains the data fields for the class and prototypes for all methods defined in the class. It is a description of the class and does
not normally contain executable code. For example, here is the header file for the \(\mathrm{C}++\) version of the previous Python class:
```

/** @ file PTainBox.h */
\#include <string> // Needed because we use string objects.
using namespace std; // Strings are part of the system std Tibrary.
class PlainBox // This is a class declaration for PlainBox.
{
private: // Items defined below have private access.
string item;
public: // Items defined below have public access.
// Prototypes for constructors and methods:
PlainBox();
PlainBox(string theItem);
void setItem(string theItem);
string getItem();
}; // end PTainBox - Note semi-colon after declaration.

```

Notice that each data member and method does not need to have an access modifier. Private data members and methods are grouped together, and public methods are grouped with each other. A colon follows each access modifier. It is possible to have more than one private, protected, or public section, but it is not common. The class declaration must end with a semi-colon after the closing brace, as shown above.

The implementation file, or source file, for the class typically has the same name as the header file, but ends in . cpp to indicate that it is a \(\mathrm{C}++\) file. The implementation file contains the definition for the methods declared in the accompanying header file. Here is the implementation file for the \(\mathrm{C}++\) class PlainBox:
```

/** @ file PlainBox.cpp */
\#include "PlainBox.h" // We need to te7l the compiler about our data
// fields and methods.
P1ainBox::P1ainBox()
{
} // end default constructor
PlainBox::PlainBox(string theItem)
{
item = theItem;
} // end constructor
void PlainBox::setItem(string theItem)
{
item = theItem;
} // end setItem
string PlainBox::getItem()
{
return item;
} // end getItem

```

This implementation file begins by including the header file, so that the compiler has the class declaration and knows the data fields and methods of this class. We then define each method. In the implementation file there is no "master" set of braces enclosing all of the methods. Each method is individually defined. To let the compiler know that a method belongs to the PlainBox class, we add
the prefix PlainBox: : to the method name. Observe that the prefix comes after the method's return type. The prefix is an extension of the method name only.

To create a PlainBox object in Python, you could use the following statement:
```

myBox = P1ainBox("Jewe1ry");

```
\(\mathrm{C}++\) provides several forms for creating objects:
PlainBox myBox = PlainBox("Jewelry");
PlainBox myBox("Jewe1ry");
These statements are equivalent in \(\mathrm{C}++\) and create a PlainBox object that is local to the function or method in which it is created. In these examples, the \(\mathrm{C}++\) keyword new is not used.

The following example uses the keyword new to create a PlainBox object that exists until the programmer specifically deletes it:

PlainBox* someBoxPtr = new PlainBox("Jewe1ry");
Notice that PlainBox is followed by the * character to indicate that someBoxPtr is a pointer to a PlainBox object and is not the object itself. Pointers are discussed in more detail in C++ Interlude 2 .

If an object in \(\mathrm{C}++\) is instantiated using the new operator, as in the example above, it is the responsibility of the programmer to delete that object from memory when it is no longer needed. In Python, garbage collection deletes unused objects and frees the memory that was used by those objects. C++ does not provide garbage collection, so you must free unused memory. To delete the object pointed to by someBoxPtr, use the statements
```

delete someBox; // Frees memory used by the object.
someBox = nul1ptr ; // A safety precaution.

```

Each time you use the new operator to create an object, you must write a corresponding delete statement to free the object's memory. Otherwise, your application will have a memory leak.

Additional information about C++ classes appears in Section A. 8 of Appendix A, Chapter 1, and in C++ Interludes 1 and 4.

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\section*{C++ Operators}

Operators within the same box have the same precedence, which is higher than operators in lower boxes.
\begin{tabular}{|llll|}
\multicolumn{1}{l}{ Operator } & Meaning & Associativity & Usage \\
\hline\(::\) & global & right & \(::\) name \\
\(::\) & scope resolution & left & class_name: : member_name \\
\hline
\end{tabular}
\begin{tabular}{|c|c|c|c|}
\hline -> & member selection & left & pointer->member \\
\hline & member selection & left & object.member \\
\hline [] & array index & left & array_name [expr] \\
\hline () & function call & left & function_name(expr_list) \\
\hline () & type construction & left & type (expr_list) \\
\hline ++ & post-increment & right & lvalue++ \\
\hline -- & post-decrement & right & lvalue-- \\
\hline typeid & type identification & right & \begin{tabular}{l}
typeid (type) \\
typeid (expr)
\end{tabular} \\
\hline dynamic_cast & checked conversion & right & dynamic_cast<type> (expr) \\
\hline static_cast & checked conversion & right & static_cast<type> (expr) \\
\hline reinterpret_cast & unchecked conversion & right & reinterpret_cast<type> (expr) \\
\hline
\end{tabular}
\begin{tabular}{|c|c|c|c|}
\hline sizeof & size of type & right & sizeof (type) \\
\hline sizeof & size of object & right & sizeof expr \\
\hline ++ & pre-increment & right & ++lvalue \\
\hline -- & pre-decrement & right & --lvalue \\
\hline \(\sim\) & bitwise complement & right & \(\sim\) expr \\
\hline ! & logical NOT & right & !expr \\
\hline + & unary plus & right & +expr \\
\hline - & unary minus & right & -expr \\
\hline * & dereference & right & *expr \\
\hline \& & address of & right & \&lvalue \\
\hline () & cast & right & (type) expr \\
\hline \multirow[t]{4}{*}{new} & allocate & right & new type \\
\hline & & & new type (expr_list) \\
\hline & & & new (expr_list) type \\
\hline & & & new (expr_list) type(expr_list) \\
\hline de7ete & deallocate & right & delete pointer delete[] pointer \\
\hline
\end{tabular}```


[^0]:    1 Pseudocode in this book appears in blue.

[^1]:    2 If you are familiar with Java, the interfaces for our ADTs correspond to Java interfaces. Both provide a public interface consisting of methods that the programmer of our ADT must implement.

[^2]:    1 You will see this notation in the rest of the book to represent a portion of an array.

[^3]:    2 Chapters 5 and 11 present other practical, efficient applications of recursion.

[^4]:    3 The binary search algorithm also has an iterative formulation.

[^5]:    2 Each time you encounter an open brace, you increment the count; each time you encounter a close brace, you decrement the count. If this count ever falls below zero, or if it is greater than zero when the end of the string is reached, the string is unbalanced.

[^6]:    3 A statement such as this one is called a loop invariant. It is a condition that is true before and after each execution of an algorithm's loop. Invariants are discussed further in Appendix F.

[^7]:    1 Recall that the return statement returns a copy of its argument.

[^8]:    1 Although omitting multiplication operators is common in algebra, we indicate them explicitly here to facilitate counting them.

[^9]:    2 The graph of $f(n)=1$ is omitted because the scale of the figure makes it difficult to draw. It would, however, be a straight line parallel to the $x$ axis through $y=1$.

[^10]:    3 CLOCKS_PER_SECOND is a C++ predefined constant.

[^11]:    1 To allow for duplicate data items, ascending is used here to mean nondecreasing and descending to mean nonincreasing.

[^12]:    // Copy the item from the last node and place it into the root
    items[0] = items[itemCount - 1]
    // Remove the last node
    itemCount--

[^13]:    1 Section 4.5 of Chapter 4 discussed how the expected and maximum number of items in an ADT affect an array-based implementation.

[^14]:    3 D. E. Knuth, Searching and Sorting, vol. 3 of The Art of Computer Programming (Menlo Park, CA: Addison-Wesley, 1973).

[^15]:    Transform each 2-node into a 3-node or a 4-node

[^16]:    // Read block i from file dataFile into buffer buf
    buf.readBlock(dataFile, i)

[^17]:    1 Data enters or leaves a buffer at a rate that differs from the record-processing rate. (Hence, a buffer between two processes compensates for the difference in the rates at which they operate on data.)

[^18]:    2 The unordered hash-based containers are part of C11-TR1 (Technical Report 1) and should be preceded by the tr1: : namespace indicator, even if you include the statement using namespace std; in your program.

[^19]:    1 Octal and hexadecimal constants are also available, but they are not used in this book. An octal constant begins with 0 , a hex

[^20]:    2 A list of all C++ operators and their precedences appears inside the cover of this book.
    3 A unary operator requires only one operand; for example, the - in -5. A binary operator requires two operands; for example, the + in $2+3$.

[^21]:    4 Whereas valued functions must contain a statement of the form return expression, void functions cannot contain such a statement. A void function can, however, contain return without an expression. Such a statement causes the function to return to the statement that follows its call. This book does not use return with void functions.

