The Need for Data Structures

[A primary concern of this course is efficiency.]

Data structures organize data

⇒ **more efficient programs.** [You might believe that faster computers make it unnecessary to be concerned with efficiency. However...]

- More powerful computers ⇒ more complex applications.
- YET More complex applications demand more calculations.
- Complex computing tasks are unlike our everyday experience. [So we need special training]

Any organization for a collection of records can be searched, processed in any order, or modified. [If you are willing to pay enough in time delay. Ex: Simple unordered array of records.]

- The choice of data structure and algorithm can make the difference between a program running in a few seconds or many days.
Efficiency

A solution is said to be **efficient** if it solves the problem within its **resource constraints**. [Alt: Better than known alternatives (“relatively” efficient)]

- space [These are typical constraints for programs]
- time

[This does not mean always strive for the most efficient program. If the program operates well within resource constraints, there is no benefit to making it faster or smaller.]

The **cost** of a solution is the amount of resources that the solution consumes.
Selecting a Data Structure

Select a data structure as follows:

1. Analyze the problem to determine the resource constraints a solution must meet.
2. Determine the basic operations that must be supported. Quantify the resource constraints for each operation.
3. Select the data structure that best meets these requirements.

[Typically want the “simplest” data structure that will meet requirements.]

Some questions to ask: [These questions often help to narrow the possibilities]

- Are all data inserted into the data structure at the beginning, or are insertions interspersed with other operations?
- Can data be deleted? [If so, a more complex representation is typically required]
- Are all data processed in some well-defined order, or is random access allowed?
Data Structure Philosophy

Each data structure has costs and benefits.

Rarely is one data structure better than another in all situations.

A data structure requires:

- space for each data item it stores, \([\text{Data} + \text{Overhead}]\)
- time to perform each basic operation,
- programming effort. \([\text{Some data structures/algorithms more complicated than others}]\)

Each problem has constraints on available space and time.

Only after a careful analysis of problem characteristics can we know the best data structure for the task.

Bank example:

- Start account: a few minutes
- Transactions: a few seconds
- Close account: overnight
Goals of this Course

1. Reinforce the concept that there are costs and benefits for every data structure. [A worldview to adopt]

2. Learn the commonly used data structures. These form a programmer’s basic data structure “toolkit.” [The “nuts and bolts” of the course]

3. Understand how to measure the effectiveness of a data structure or program.
   - These techniques also allow you to judge the merits of new data structures that you or others might invent. [To prepare you for the future]
Definitions

A **type** is a set of values.

[Ex: Integer, Boolean, Float]

A **data type** is a type and a collection of operations that manipulate the type.

[Ex: Addition]

A **data item** or **element** is a piece of information or a record.

[Physical instantiation]

A data item is said to be a **member** of a data type.

A **simple data item** contains no subparts.

[Ex: Integer]

An **aggregate data item** may contain several pieces of information.

[Ex: Payroll record, city database record]
Abstract Data Types

**Abstract Data Type** (ADT): a definition for a data type solely in terms of a set of values and a set of operations on that data type.

Each ADT operation is defined by its inputs and outputs.

**Encapsulation**: hide implementation details

A **data structure** is the physical implementation of an ADT.

- Each operation associated with the ADT is implemented by one or more subroutines in the implementation.

**Data structure** usually refers to an organization for data in main memory.

**File structure**: an organization for data on peripheral storage, such as a disk drive or tape.

An ADT manages complexity through abstraction: **metaphor**. [Hierarchies of labels]

[Ex: transistors → gates → CPU. In a program, implement an ADT, then think only about the ADT, not its implementation]
Logical vs. Physical Form

Data items have both a **logical** and a **physical** form.

Logical form: definition of the data item within an ADT. [Ex: Integers in mathematical sense: +, −]

Physical form: implementation of the data item within a data structure. [16/32 bit integers: overflow]

Data Type

ADT:
- Type
- Operations

Data Items: Logical Form

Data Structure:
- Storage Space
- Subroutines

Data Items: Physical Form

[In this class, we frequently move above and below “the line” separating logical and physical forms.]
Problems

**Problem**: a task to be performed.

- Best thought of as inputs and matching outputs.
- Problem definition should include constraints on the resources that may be consumed by any acceptable solution. [But NO constraints on HOW the problem is solved]

Problems ⇔ mathematical functions

- A **function** is a matching between inputs (the **domain**) and outputs (the **range**).
- An **input** to a function may be single number, or a collection of information.
- The values making up an input are called the **parameters** of the function.
- A particular input must always result in the same output every time the function is computed.
Algorithms and Programs

**Algorithm**: a method or a process followed to solve a problem. [A recipe]

An algorithm takes the input to a problem (function) and transforms it to the output. [A mapping of input to output]

A problem can have many algorithms.

An algorithm possesses the following properties:

1. It must be **correct**. [Computes proper function]
2. It must be composed of a series of **concrete steps**. [Executable by that machine]
3. There can be **no ambiguity** as to which step will be performed next.
4. It must be composed of a **finite** number of steps.
5. It must **terminate**.

A **computer program** is an instance, or concrete representation, for an algorithm in some programming language.

[We frequently interchange use of “algorithm” and “program” though they are actually different concepts]
Mathematical Background

[Look over Chapter 2, read as needed depending on your familiarity with this material.]

Set concepts and notation [Set has no duplicates, sequence may]
Recursion

Induction proofs

Logarithms [Almost always use log to base 2. That is our default base.]
Summations
Algorithm Efficiency

There are often many approaches (algorithms) to solve a problem. How do we choose between them?

At the heart of computer program design are two (sometimes conflicting) goals:

1. To design an algorithm that is easy to understand, code and debug.
2. To design an algorithm that makes efficient use of the computer’s resources.

Goal (1) is the concern of Software Engineering.

Goal (2) is the concern of data structures and algorithm analysis.

When goal (2) is important, how do we measure an algorithm’s cost?
How to Measure Efficiency?

1. Empirical comparison (run programs).

   [Difficult to do “fairly.” Time consuming.]

2. Asymptotic Algorithm Analysis.

Critical resources:

- Time
- Space (disk, RAM)
- Programmer’s effort
- Ease of use (user’s effort).

Factors affecting running time:

- Machine load
- OS
- Compiler
- Problem size or Specific input values for given problem size

For most algorithms, running time depends on “size” of the input.

Running time is expressed as $T(n)$ for some function $T$ on input size $n$. 
Examples of Growth Rate

Example 1: [As $n$ grows, how does $T(n)$ grow?]

```java
static int largest(int[] array) { // Find largest val
    // all values >=0
    int currLargest = 0; // Store largest val
    for (int i=0; i<array.length; i++) // For each elem
        if (array[i] > currLargest) // if largest
            currLargest = array[i]; // remember it
    return currLargest; // Return largest val
}

[Cost: $T(n) = c_1n + c_2$ steps]

Example 2: Assignment statement [Constant cost]

Example 3:

```
sum = 0;
for (i=1; i<=n; i++)
    for (j=1; j<=n; j++)
        sum++;
```

[Cost: $T(n) = c_1n^2 + c_2$ Roughly $n^2$ steps, with sum being $n^2$ at the end. Ignore various overhead such as loop counter increments.]
Growth Rate Graph

$2^n$ is an exponential algorithm. $10n$ and $20n$ differ only by a constant.
Important facts to remember

• for any integer constants $a, b > 1$ $n^a$ grows faster than $\log^b n$
  
  [any polynomial is worse than any power of any logarithm]
• for any integer constants $a, b > 1$ $n^a$ grows faster than $\log n^b$
  
  [any polynomial is worse than any logarithm of any power]
• for any integer constants $a, b > 1$ $a^n$ grows faster than $n^b$
  
  [any exponential is worse than any polynomial]
Best, Worst and Average Cases

Not all inputs of a given size take the same time.

Sequential search for \( K \) in an array of \( n \) integers:

- Begin at first element in array and look at each element in turn until \( K \) is found.

Best Case: [Find at first position: 1 compare]

Worst Case: [Find at last position: \( n \) compares]

Average Case: \([ (n + 1)/2 \) compares\]

While average time seems to be the fairest measure, it may be difficult to determine.

[Depends on distribution. Assumption for above analysis: Equally likely at any position.]

When is worst case time important?

[algorithms for time-critical systems]
Faster Computer or Algorithm?

What happens when we buy a computer 10 times faster? [How much speedup? 10 times. More important: How much increase in problem size for same time? Depends on growth rate.]

<table>
<thead>
<tr>
<th>$T(n)$</th>
<th>$n$</th>
<th>$n'$</th>
<th>Change</th>
<th>$n'/n$</th>
</tr>
</thead>
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<tr>
<td>$10n$</td>
<td>1,000</td>
<td>10,000</td>
<td>$n' = 10n$</td>
<td>10</td>
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<tr>
<td>$20n$</td>
<td>500</td>
<td>5,000</td>
<td>$n' = 10n$</td>
<td>10</td>
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<td>$5n \log n$</td>
<td>250</td>
<td>1,842</td>
<td>$\sqrt{10n &lt; n'} &lt; 10n$</td>
<td>7.37</td>
</tr>
<tr>
<td>$2n^2$</td>
<td>70</td>
<td>223</td>
<td>$n' = \sqrt{10n}$</td>
<td>3.16</td>
</tr>
<tr>
<td>$2^n$</td>
<td>13</td>
<td>16</td>
<td>$n' = n + 3$</td>
<td>--</td>
</tr>
</tbody>
</table>

[For $n^2$, if $n = 1000$, then $n'$ would be 1003]

$n$: Size of input that can be processed in one hour (10,000 steps).

$n'$: Size of input that can be processed in one hour on the new machine (100,000 steps).

[Compare $T(n) = n^2$ to $T(n) = n \log n$. For $n > 58$, it is faster to have the $\Theta(n \log n)$ algorithm than to have a computer that is 10 times faster.]
Asymptotic Analysis: Big-oh

Definition: For $T(n)$ a non-negatively valued function, $T(n)$ is in the set $O(f(n))$ if there exist two positive constants $c$ and $n_0$ such that $T(n) \leq cf(n)$ for all $n > n_0$.

Usage: The algorithm is in $O(n^2)$ in [best, average, worst] case.

Meaning: For all data sets big enough (i.e., $n > n_0$), the algorithm always executes in less than $cf(n)$ steps [in best, average or worst case].

[Must pick one of these to complete the statement. Big-oh notation applies to some set of inputs.]

Upper Bound.

Example: if $T(n) = 3n^2$ then $T(n)$ is in $O(n^2)$.

Wish tightest upper bound:
While $T(n) = 3n^2$ is in $O(n^3)$, we prefer $O(n^2)$.

[It provides more information to say $O(n^2)$ than $O(n^3)$]
Big-oh Example

Example 1. Finding value $X$ in an array. [Average case]

$T(n) = c_s n/2$. [$c_s$ is a constant. Actual value is irrelevant]
For all values of $n > 1$, $c_s n/2 \leq c_s n$.
Therefore, by the definition, $T(n)$ is in $O(n)$ for $n_0 = 1$ and $c = c_s$.

Example 2. $T(n) = c_1 n^2 + c_2 n$ in average case

$c_1 n^2 + c_2 n \leq c_1 n^2 + c_2 n^2 \leq (c_1 + c_2) n^2$ for all $n > 1$.

$T(n) \leq c n^2$ for $c = c_1 + c_2$ and $n_0 = 1$.
Therefore, $T(n)$ is in $O(n^2)$ by the definition.

Example 3: $T(n) = c$. We say this is in $O(1)$. [Rather than $O(c)$]
Big-Omega

Definition: For $T(n)$ a non-negatively valued function, $T(n)$ is in the set $\Omega(g(n))$ if there exist two positive constants $c$ and $n_0$ such that $T(n) \geq cg(n)$ for all $n > n_0$.

Meaning: For all data sets big enough (i.e., $n > n_0$), the algorithm always executes in more than $cg(n)$ steps.

Lower Bound.

Example: $T(n) = c_1n^2 + c_2n$.

c_1n^2 + c_2n \geq c_1n^2$ for all $n > 1$.

$T(n) \geq cn^2$ for $c = c_1$ and $n_0 = 1$.

Therefore, $T(n)$ is in $\Omega(n^2)$ by the definition.

Want greatest lower bound.
Theta Notation

When big-Oh and Ω meet, we indicate this by using Θ (big-Theta) notation.

Definition: An algorithm is said to be \( \Theta(h(n)) \) if it is in \( O(h(n)) \) and it is in \( \Omega(h(n)) \).

[For polynomial equations on \( T(n) \), we always have \( \Theta \). There is no uncertainty, a “complete” analysis.]

Simplifying Rules:

1. If \( f(n) \) is in \( O(g(n)) \) and \( g(n) \) is in \( O(h(n)) \), then \( f(n) \) is in \( O(h(n)) \).

2. If \( f(n) \) is in \( O(kg(n)) \) for any constant \( k > 0 \), then \( f(n) \) is in \( O(g(n)) \). [No constant]

3. If \( f_1(n) \) is in \( O(g_1(n)) \) and \( f_2(n) \) is in \( O(g_2(n)) \), then \( (f_1 + f_2)(n) \) is in \( O(\max(g_1(n), g_2(n))) \). [Drop low order terms]

4. If \( f_1(n) \) is in \( O(g_1(n)) \) and \( f_2(n) \) is in \( O(g_2(n)) \) then \( f_1(n)f_2(n) \) is in \( O(g_1(n)g_2(n)) \). [Loops]
Running Time of a Program

[Asymptotic analysis is defined for equations. Need to convert program to an equation.]

Example 1: \( a = b; \)

This assignment takes constant time, so it is \( \Theta(1). \) [Not \( \Theta(c) \) – notation by tradition]

Example 2:

\[
\begin{align*}
\text{sum} & = 0; \\
\text{for (i=1; i<=n; i++)} & \\
& \quad \text{sum += n;}
\end{align*}
\]

[\( \Theta(n) \) (even though sum is \( n^2 \))]

Example 3:

\[
\begin{align*}
\text{sum} & = 0; \\
\text{for (j=1; j<=n; j++)} & \quad \text{// First for loop} \\
& \quad \text{for (i=1; i<=j; i++)} \quad \text{// is a double loop} \\
& \quad \quad \text{sum++;} \\
\text{for (k=0; k<n; k++)} & \quad \text{// Second for loop} \\
& \quad \quad A[k] = k;
\end{align*}
\]

[First statement is \( \Theta(1). \) Double for loop is \( \sum i = \Theta(n^2). \) Final for loop is \( \Theta(n). \) Result: \( \Theta(n^2). \)]
More Examples

Example 4.

```plaintext
sum1 = 0;
for (i=1; i<=n; i++) // First double loop
    for (j=1; j<=n; j++) // do n times
        sum1++;

sum2 = 0;
for (i=1; i<=n; i++) // Second double loop
    for (j=1; j<=i; j++) // do i times
        sum2++;

[First loop, sum is \( n^2 \). Second loop, sum is \( (n + 1)(n)/2 \). Both are \( \Theta(n^2) \).]
```

Example 5.

```plaintext
sum1 = 0;
for (k=1; k<=n; k*=2)
    for (j=1; j<=n; j++)
        sum1++;

sum2 = 0;
for (k=1; k<=n; k*=2)
    for (j=1; j<=k; j++)
        sum2++;

[First is \( \sum_{k=1}^{\log n} n = \Theta(n \log n) \). Second is \( \sum_{k=0}^{\log n-1} 2^k = \Theta(n) \).]
```
Binary Search

static int binary(int K, int[] array, int left, int right) {  
    // Return position in array (if any) with value K  
    int l = left-1;  
    int r = right+1; // l and r are beyond array bounds  
        // to consider all array elements  
    while (l+1 != r) { // Stop when l and r meet  
        int i = (l+r)/2; // Look at middle of subarray  
        if (K < array[i]) r = i; // In left half  
        if (K == array[i]) return i; // Found it  
        if (K > array[i]) l = i; // In right half  
    }  
    return UNSUCCESSFUL; // Search value not in array  
}  

invocation of binary  
int pos = binary(43, ar, 0, 15);  

Analysis: How many elements can be examined in the worst case? \[\Theta(\log n)\]
Other Control Statements

while loop: analyze like a for loop.

if statement: Take greater complexity of then/else clauses.

[If probabilities are independent of $n$.]

switch statement: Take complexity of most expensive case.

[If probabilities are independent of $n$.]

Subroutine call: Complexity of the subroutine.
Analyzing Problems

Use same techniques to analyze problems, i.e. any possible algorithm for a given problem (e.g., sorting)

Upper bound: Upper bound of best known algorithm.

Lower bound: Lower bound for every possible algorithm.

[The examples so far have been easy in that exact equations always yield $\Theta$. Thus, it was hard to distinguish $\Omega$ and $O$. Following example should help to explain the difference – bounds are used to describe our level of uncertainty about an algorithm.]

Example: Sorting
1. Cost of I/O: $\Omega(n)$
2. Bubble or insertion sort: $O(n^2)$
3. A better sort (Quicksort, Mergesort, Heapsort, etc.): $O(n \log n)$
4. We prove later that sorting is $\Omega(n \log n)$
Multiple Parameters

[Ex: 256 colors (8 bits), $1000 \times 1000$ pixels]
Compute the rank ordering for all $C$ (256) pixel values in a picture of $P$ pixels.

```c
for (i=0; i<C; i++) // Initialize count
    count[i] = 0;
for (i=0; i<P; i++) // Look at all of the pixels
    count[value(i)]++; // Increment proper value count
sort(count); // Sort pixel value counts
```

If we use $P$ as the measure, then time is $\Theta(P \log P)$.

But this is wrong because we sort colors
More accurate is $\Theta(P + C \log C)$.
If $C << P$, $P$ could overcome $C \log C$.
Space Bounds

Space bounds can also be analyzed with asymptotic complexity analysis.

Time: Algorithm
Space: Data Structure

Space/Time Tradeoff Principle:
One can often achieve a reduction in time if one is willing to sacrifice space, or vice versa.

- Encoding or packing information
  Boolean flags
- Table lookup
  Factorials

Disk Based Space/Time Tradeoff Principle:
The smaller you can make your disk storage requirements, the faster your program will run.
(because access to disk is typically more costly than "any" computation)
Algorithm Design methods:
Divide et impera

Decompose a problem of size $n$ into (one or more) problems of size $m < n$

Solve subproblems, if reduced size is not "trivial", in the same manner, possibly combining solutions of the subproblems to obtain the solution of the original one ...

... until size becomes "small enough" (typically 1 or 2) to solve the problem directly (without decomposition)

Complexity can be typically analyzed by means of recurrence equations
Recurrence Equations (1)

we have already seen the following

\[ T(n) = a T\left(\frac{n}{b}\right) + cn^k, \text{ for } n > 1 \]
\[ T(1) = d, \]

Solution of the recurrence depends on the ratio \( r = \frac{b^k}{a} \)

\[ T(n) = \Theta(n^{\log_b a}), \text{ if } a > b^k \]
\[ T(n) = \Theta(n^k \log n), \text{ if } a = b^k \]
\[ T(n) = \Theta(n^k), \text{ if } a < b^k \]

Complexity depends on

- relation between \( a \) and \( b \), i.e., whether all subproblems need to be solved or only some do
- value of \( k \), i.e., amount of additional work to be done to partition into subproblems and combine solutions
Recurrence Equations(2)

Examples

• $a = 1, b = 2$ (two halves, solve only one), $k = 0$ (constant partition + combination overhead): e.g., Binary search: $T(n) = \Theta(\log n)$ (extremely efficient!)

• $a = b = 2$ (two halves) and $(k=1)$ (partitioning + combination $\Theta(n)$) $T(n) = \Theta(n \log n)$; e.g., Mergesort;

• $a = b$ (partition data and solve for all partitions) and $k = 0$ (constant partition + combining) $T(n) = \Theta(n \log_b a) = \Theta(n)$, same as linear/sequential processing (E.g., finding the max/min element in an array)

Now we’ll see

1. max/min search as an example of linear complexity

2. other kinds of recurrence equations

• $T(n) = T(n-1) + n$ leads to quadratic complexity: example bubblesort;

• $T(n) = aT(n-1) + k$ leads to exponential complexity: example Towers of Hanoi
MaxMin search(1)

"Obvious" method: sequential search

```java
public class MinMaxPair {
    public int min;
    public int max;
}

public static MinMaxPair minMax (float [] a) {
    //guess a[0] as min and max
    MaxMinPair p = new MaxMinPair(); p.min = p.max = 0;
    // search in the remaining part of the array
    for (int i = 1; i<a.length; i++) {
        if (a[i]<a[p.min]) p.min = i;
        if (a[i]>a[p.max]) p.max = i;
    }
    return p;
}
```

Complexity is $T(n)=2(n-1)=\Theta(n)$

Divide et impera approach: split array in two, find MinMax of each, choose overall min among the two mins and max among the two maxs
public static MinMaxPair minMax (float[] a, int l, int r) {
    MaxMinPair p = new MinMaxPair();
    if(l==r) {p.min = p.max = r; return p;}
    if (l==r-1) {
        if (a[l]<a[r]) {
            p.min=l; p.max=r;
        }
        else {
            p.min=r; p.max=l;
        }
        return p;
    }
    int m = (l+r)/2;
    MinMaxPair p1 = minMax(a, l, m);
    MinMaxPair p2 = minMax(a, m+1, r);
    if (a[p1.min]<a[p2.min]) p.min=p1.min else p.min=p2.min;
    if (a[p1.max]>a[p2.max]) p.max=p1.max else p.max=p2.max;
    return p;
}

Asymptotic complexity analyzable by means of recurrence
T(n) = aT(n/b)+cn^k, for n > 1
T(1) = d,

We have a = b and k = 0 hence T(n) = Θ(n),
apparently no improvement: we need a more precise analysis
Assume for simplicity \( n \) is a power of 2. Here is the tree of recursive calls for \( n = 16 \). There are

- \( n/2 \) leaf nodes, each of which takes 1 comparison
- \( n/2 - 1 \) internal nodes each of which takes 2 comparison
- hence \#comparisons = 
  \[ 2(n/2 - 1) + n/2 = (3/2)n - 2, \] a 25% improvement wrt linear search
bubblesort as a divide et impera algorithm

To sort an array of \( n \) elements, put the smallest element in first position, then sort the remaining part of the array.

Putting the smallest element to first position requires an array traversal (\( \Theta(n) \) complexity)

```java
static void bubsort(Elem[] array) { // Bubble Sort
    for (int i=0; i<array.length-1; i++) // Bubble up
        for (int j=array.length-1; j>i; j--)
            if (array[j].key() < array[j-1].key())
                DSutil.swap(array, j, j-1);
}
```

```
<table>
<thead>
<tr>
<th>i=0</th>
<th>1</th>
<th>2</th>
<th>3</th>
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</tbody>
</table>
```
Towers of Hanoi

Move stack of rings form one pole to another, with following constraints

- move one ring at a time
- never place a ring on top of a smaller one

Divide et impera approach: move stack of \( n - 1 \) smaller rings on third pole as a support, then move largest ring, then move stack of \( n - 1 \) smaller rings from support pole to destination pole using start pole as a support

```java
static void TOH(int n,
    Pole start, Pole goal, Pole temp) {
    if (n==1) System.out.println("move ring from pole " +
            + start + " to pole " + goal);
    else {
        TOH(n-1, start, temp, goal);
        System.out.println("move ring from pole " +
            + start + " to pole " + goal);
        TOH(n-1, temp, goal, start);
    }
}
```

Time complexity as a function of the size \( n \) of the ring stack: \( T(n)=2^{n-1} \)
Exponential complexity
of Towers of Hanoi

Recurrence equation is \( T(n) = 2T(n-1) + 1 \) for \( n > 1 \), and \( T(1) = 1 \).

A special case of the more general recurrence
\( T(n) = aT(n-1) + k \), for \( n > 1 \), and \( T(1) = k \).

It is easy to show that the solution is
\( T(n) = k \sum_{i=0}^{n-1} a^i \) hence \( T(n) = \Theta(a^n) \)


Base: \( T(1) = k = k \sum_{i=0}^{0} a^i \)

Induction:

\( T(n + 1) = aT(n) + k = \)
\( = ak \sum_{i=0}^{n-1} a^i + k = k \sum_{i=1}^{n} a^i + k = k \sum_{i=0}^{n} a^i = \)
\( = k \sum_{i=0}^{(n+1)-1} a^i \)

In the case of Towers of Hanoi \( a = 2, k = 1 \),
hence \( T(n) = \sum_{i=0}^{n-1} 2^i = 2^n - 1 \)
Lists

[Students should already be familiar with lists. Objectives: use alg analysis in familiar context, compare implementations.]

A list is a finite, ordered sequence of data items called elements.

[The positions are ordered, NOT the values.]

Each list element has a data type.

The empty list contains no elements.

The length of the list is the number of elements currently stored.

The beginning of the list is called the head, the end of the list is called the tail.

Sorted lists have their elements positioned in ascending order of value, while unsorted lists have no necessary relationship between element values and positions.

Notation: ( a_0, a_1, ..., a_{n-1} )

What operations should we implement?

[Add/delete elem anywhere, find, next, prev, test for empty.]
List ADT

interface List { // List ADT
public void clear(); // Remove all Objects
public void insert(Object item); // Insert at curr pos
public void append(Object item); // Insert at tail
public Object remove(); // Remove/return curr
public void setFirst(); // Set to first pos
public void next(); // Move to next pos
public void prev(); // Move to prev pos
public int length(); // Return curr length
public void setPos(int pos); // Set curr position
public void setValue(Object val); // Set current value
public Object currValue(); // Return curr value
public boolean isEmpty(); // True if empty list
public boolean isInList(); // True if curr in list
public void print(); // Print all elements
} // interface List

[This is an example of a Java interface. Any Java class using this interface must implement all of these functions. Note that the generic type “Object” is being used for the element type.]
List ADT Examples

List: ( 12, 32, 15 )

MyLst.insert(element);

[The above is an example use of the insert function.
“element” is an object of the list element data type.]
Assume MyLst has 32 as current element:

MyLst.insert(99);

[Put 99 before current element, yielding (12, 99, 32, 15).]
Process an entire list:

for (MyLst.setFirst(); MyLst.isInList(); MyLst.next())
    DoSomething(MyLst.currValue());
Array-Based List Insert

[Push items up/down. Cost: $\Theta(n)$.

Insert 23:

(a) $13\ 12\ 20\ 8\ 3\ \boxed{\ast}\ 
0\ 1\ 2\ 3\ 4\ 5$

(b) $13\ 12\ 20\ 8\ 3\ \boxed{\ast}
0\ 1\ 2\ 3\ 4\ 5$

(c) $23\ 13\ 12\ 20\ 8\ 3\ \boxed{\ast}
0\ 1\ 2\ 3\ 4\ 5$
Array-Based List Class

class AList implements List { // Array-based list

private static final int defaultSize = 10;

private int msize; // Maximum size of list
private int numInList; // Actual list size
private int curr; // Position of curr
private Object[] listArray; // Array holding list

AList() { setup(defaultSize); } // Constructor
AList(int sz) { setup(sz); } // Constructor

private void setup(int sz) { // Do initializations
    msize = sz;
    numInList = curr = 0;
    listArray = new Object[sz]; // Create listArray
}

public void clear() // Remove all Objects from list
{numInList = curr = 0; } // Simply reinitialize values

public void insert(Object it) { // Insert at curr pos
    Assert.notFalse(numInList < msize, "List is full");
    Assert.notFalse((curr >=0) && (curr <= numInList),
        "Bad value for curr");
    for (int i=numInList; i>curr; i--) // Shift up
        listArray[i] = listArray[i-1];
    listArray[curr] = it;
    numInList++;
        // Increment list size
}
public void append(Object it) { // Insert at tail
    Assert.notFalse(numInList < msize, "List is full");
    listArray[numInList++] = it; // Increment list size
}

public Object remove() { // Remove and return Object
    Assert.notFalse(!isEmpty(), "No delete: list empty");
    Assert.notFalse(isInList(), "No current element");
    Object it = listArray[curr]; // Hold removed Object
    for(int i=curr; i<numInList-1; i++) // Shift down
        listArray[i] = listArray[i+1];
    numInList--; // Decrement list size
    return it;
}

public void setFirst() { curr = 0; } // Set to first
public void prev() { curr--; } // Move curr to prev
public void next() { curr++; } // Move curr to next
public int length() { return numInList; }
public void setPos(int pos) { curr = pos; }
public boolean isEmpty() { return numInList == 0; }
public void setValue(Object it) { // Set current value
    Assert.notFalse(isInList(), "No current element");
    listArray[curr] = it;
}

public boolean isInList() // True if curr within list
{ return (curr >= 0) && (curr < numInList); }
} // Array-based list implementation
Link Class

Dynamic allocation of new list elements.

class Link {
    // A singly linked list node
    private Object element; // Object for this node
    private Link next; // Pointer to next node

    Link(Object it, Link nextval) { // Constructor
        element = it; next = nextval;
    }
    Link(Link nextval) { // Constructor
        next = nextval;
    }
    Link next() {
        return next;
    }
    Link setNext(Link nextval) {
        return next = nextval;
    }
    Object element() {
        return element;
    }
    Object setElement(Object it) {
        return element = it;
    }
}
Linked List Position

[Naive approach: Point to current node. Current is 12. Want to insert node with 10. No access available to node with 23. How can we do the insert?]

[Alt implementation: Point to node preceding actual current node. Now we can do the insert. Also note use of header node.]
Linked List Implementation

```
public class LList implements List { // Linked list
private Link head; // Pointer to list header
private Link tail; // Pointer to last Object in list
protected Link curr; // Position of current Object

LList(int sz) { setup(); } // Constructor
LList() { setup(); } // Constructor
private void setup() // allocates leaf node
    { tail = head = curr = new Link(null); }

public void setFirst() { curr = head; }
public void next() 
    { if (curr != null) curr = curr.next(); }

public void prev() { // Move to previous position
    Link temp = head;
    if ((curr == null) || (curr == head)) // No prev
        { curr = null; return; } // so return
    while ((temp != null) && (temp.next() != curr))
        temp = temp.next();
    curr = temp;
}

public Object currValue() { // Return current Object
    if (!isInList() || this.isEmpty() ) return null;
    return curr.next().element();
}

public boolean isEmpty() // True if list is empty
    { return head.next() == null; }
} // Linked list class
```
// Insert Object at current position
public void insert(Object it) {
    Assert.notNull(curr, "No current element");
    curr.setNext(new Link(it, curr.next()));
    if (tail == curr) // Appended new Object
tail = curr.next();
}

// Insert Object at current position
public void insert(Object it) {
    Assert.notNull(curr, "No current element");
    curr.setNext(new Link(it, curr.next()));
    if (tail == curr) // Appended new Object
tail = curr.next();
}
public Object remove() { // Remove/return curr Object
    if (!isInList() || this.isEmpty()) return null;
    Object it = curr.next().element(); // Remember value
    if (tail == curr.next()) tail = curr; // Set tail
    curr.setNext(curr.next().next()); // Cut from list
    return it; // Return value
}
Freelists

System new and garbage collection are slow.

class Link { // Singly linked list node with freelist
    private Object element; // Object for this Link
    private Link next; // Pointer to next Link

    Link(Object it, Link nextval)
    { element = it; next = nextval; }
    Link(Link nextval) { next = nextval; }
    Link next() { return next; }
    Link setNext(Link nextval) { return next = nextval; }
    Object element() { return element; }
    Object setElement(Object it) { return element = it; }

    // Extensions to support freelists
    static Link freelist = null; // Freelist for class

    static Link get(Object it, Link nextval) {
        if (freelist == null) // free list empty: allocate
            return new Link(it, nextval);
        Link temp = freelist; // take from the freelist
        freelist = freelist.next();
        temp.setElement(it);
        temp.setNext(nextval);
        return temp;
    }
    void release() {
        // add current node to freelist
        element = null; next = freelist; freelist = this;
    }
}
Comparison of List Implementations

Array-Based Lists: [Average and worst cases]
- Insertion and deletion are $\Theta(n)$.  
- Array must be allocated in advance.  
- No overhead if all array positions are full.

Linked Lists:
- Insertion and deletion $\Theta(1)$;  
  prev and direct access are $\Theta(n)$.  
- Space grows with number of elements.  
- Every element requires overhead.

Space “break-even” point:

\[
DE = n(P + E); \quad n = \frac{DE}{P + E}
\]

n: elements currently in list  
E: Space for data value  
P: Space for pointer  
D: Number of elements in array (fixed in the implementation)

[arrays more efficient when full, linked lists more efficient with few elements]
Doubly Linked Lists

Simplify insertion and deletion: Add a prev pointer.

class DLink { // A doubly-linked list node
private Object element; // Object for this node
private DLink next; // Pointer to next node
private DLink prev; // Pointer to previous node
DLink(Object it, DLink n, DLink p)
{ element = it; next = n; prev = p; }
DLink(DLink n, DLink p) { next = n; prev = p; }
DLink next() { return next; }
DLink setNext(DLink nextval) { return next=nextval; }
DLink prev() { return prev; }
DLink setPrev(DLink prevval) { return prev=prevval; }
Object element() { return element; }
Object setElement(Object it) { return element = it; }
}
Doubly Linked List Operations

// Insert Object at current position
public void insert(Object it) {
    Assert.notNull(curr, "No current element");
    curr.setNext(new DLink(it, curr.next(), curr));
    if (curr.next().next() != null)
        curr.next().next().setPrev(curr.next());
    if (tail == curr) // Appended new Object
        tail = curr.next();
}

public Object remove() { // Remove/return curr Object
    Assert.notNull(isInList(), "No current element");
    Object it = curr.next().element(); // Remember Object
    if (curr.next().next() != null)
        curr.next().next().setPrev(curr.next());
    else tail = curr; // Removed last Object: set tail
    curr.setNext(curr.next().next()); // Remove from list
    return it; // Return value removed
}
Circularly Linked Lists

- Convenient if there is no last nor first element (there is no total order among elements)
- The "last" element points to the "first", and the first to the last
- tail pointer non longer needed
- Potential danger: infinite loops in list processing
- but head pointer can be used as a marker
Stacks

LIFO: Last In, First Out

Restricted form of list: Insert and remove only at front of list.

Notation:
- Insert: PUSH
- Remove: POP
- The accessible element is called TOP.
Define top as first free position.

class AStack implements Stack{ // Array based stack class
  private static final int defaultSize = 10;
  private int size; // Maximum size of stack
  private int top; // Index for top Object
  private Object [] listarray; // Array holding stack
  AStack() { setup(defaultSize); }
  AStack(int sz) { setup(sz); }

  public void setup(int sz)
  { size = sz; top = 0; listarray = new Object[sz]; }

  public void clear() { top = 0; } // Clear all Objects

  public void push(Object it) // Push onto stack
  { Assert.notFalse(top < size, "Stack overflow");
    listarray[top++] = it; }

  public Object pop() // Pop Object from top
  { Assert.notFalse(!isEmpty(), "Empty stack");
    return listarray[--top]; }

  public Object topValue() // Return top Object
  { Assert.notFalse(!isEmpty(), "Empty stack");
    return listarray[top-1]; }

  public boolean isEmpty() { return top == 0; }
};
public class LStack implements Stack {
    // Linked stack class
    private Link top;       // Pointer to list header

    public LStack() { setup(); }   // Constructor
    public LStack(int sz) { setup(); }    // Constructor

    private void setup() // Initialize stack
    { top = null; }       // Create header node

    public void clear() { top = null; } // Clear stack

    public void push(Object it) // Push Object onto stack
    { top = new Link(it, top); }

    public Object pop() {    // Pop Object from top
        Assert.notFalse(!isEmpty(), "Empty stack");
        Object it = top.element();
        top = top.next();
        return it;
    }

    public Object topValue() // Get value of top Object
    { Assert.notFalse(!isEmpty(), "No top value");
      return top.element(); }

    public boolean isEmpty() // True if stack is empty
    { return top == null; }
} // Linked stack class
Array-based vs linked stacks

- Time: all operations take constant time for both
- Space: linked has overhead but is flexible; array has no overhead but wastes space when not full

Implementation of multiple stacks

- two stacks at opposite ends of an array growing in opposite directions
- works well if their space requirements are inversely correlated
Queues

FIFO: First In, First Out

Restricted form of list:
   Insert at one end, remove from other.

Notation:
   • Insert: Enqueue
   • Delete: Dequeue
   • First element: FRONT
   • Last element: REAR
Array Queue Implementations

Constraint: all elements

1. in consecutive positions
2. in the initial (final) portion of the array

If both (1) and (2) hold: rear element in pos 0, dequeue costs $\Theta(1)$, enqueue costs $\Theta(n)$

Similarly if in final portion of the array and/or in reverse order

If only (1) holds (2 is released)

• both front and rear move to the ”right” (i.e., increase)

• both enqueue and dequeue cost $\Theta(1)$

”Drifting queue” problem: run out of space when at the highest positions

Solution: pretend the array is circular, implemented by the modulus operator, e.g.,
front = (front + 1) % size
A more serious problem: empty queue indistinguishable from full queue

[Application of Pigeonhole Principle: Given a fixed (arbitrary) position for front, there are $n+1$ states (0 through $n$ elements in queue) and only $n$ positions for rear. One must distinguish between two of the states.]

2 solutions to this problem
1. store # elements separately from the queue
2. use a $n+1$ elements array for holding a queue with $n$ elements an most

Both solutions require one additional item of information
Linked Queue: modified linked list.

[Operations are $\Theta(1)$]
A **binary tree** is made up of a finite set of nodes that is either empty (then it is an **empty** tree) or consists of a node called the **root** connected to two binary trees, called the left and right **subtrees**, which are disjoint from each other and from the root.

\[\text{A has depth 0. B and C form level 1. The tree has height 4. Height = max depth + 1.}\]
Notation

(left/right) child of a node: root node of the (left/right) subtree

if there is no left (right) subtree we say that left/(right) subtree is empty

edge: connection between a node and its child (drawn as a line)

parent of a node \( n \): the node of which \( n \) is a child

path from \( n_1 \) to \( n_k \): a sequence \( n_1 \ n_2 \ldots \ n_k \), \( k \geq 1 \), such that, for all \( 1 \leq i < k \), \( n_i \) is parent of \( n_{i+1} \)

length of a path \( n_1 \ n_2 \ldots \ n_k \) is \( k - 1 \) (⇒ length of path \( n_1 \) is 0)

if there is a path from node \( a \) to node \( d \) then
- \( a \) is ancestor of \( d \)
- \( d \) is descendant of \( a \)
hence
- all nodes of a tree (except the root) are descendant of the root
- the root is ancestor of all the other nodes of the tree (except itself)

**depth** of a node: length of a path from the root (⇒ the root has depth 0)

**height** of a tree: 1 + depth of the deepest node (which is a leaf)

**level** \( d \) of a tree: the set of all nodes of depth \( d \) (⇒ root is the only node of level 0)

**leaf** node: has two empty children

**internal** node (non-leaf): has at least one non-empty child
- A: root
- B, C: A’s children
- B, D: A’s subtree
- D, E, F: level 2
- B has only right child (subtree)
- path of length 3 from A to G
- A, B, C, E, F internal nodes
- D, G, H, I leaves
- depth of G is 3, height of tree is 4
Full and Complete Binary Trees

**Full** binary tree: each node either is a leaf or is an internal node with exactly two non-empty children.

**Complete** binary tree: If the height of the tree is $d$, then all levels except possibly level $d - 1$ are completely full. The bottom level has nodes filled in from the left side.

(a) full but not complete
(b) complete but not full
(c) full and complete

[diagram of full and complete binary trees]

[NB these terms can be hard to distinguish]

Question: how many nodes in a complete binary tree?

A complete binary tree is "balanced", i.e., has minimal height given number of nodes

A complete binary tree is full or almost full or "almost full" (at most one node with one son) ]
Making missing children explicit

for a (non-)empty subtree we say the node has a (non-)NULL pointer
Full Binary Tree Theorem

Theorem: The number of leaves in a non-empty full binary tree is one more than the number of internal nodes.

[Relevant since it helps us calculate space requirements.]

Proof (by Mathematical Induction):

- **Base Case**: A full binary tree with 0 internal node has 1 leaf node.
- **Induction Hypothesis**: Assume any full binary tree $T$ containing $n-1$ internal nodes has $n$ leaves.
- **Induction Step**: Given a full tree $T$ with $n-1$ internal nodes ($\Rightarrow n$ leaves), add two leaf nodes as children of one of its leaves ($\Rightarrow$ obtain a tree $T'$ having $n$ internal nodes and $n + 1$ leaves.
Full Binary Tree Theorem Corollary

**Theorem:** The number of empty subtrees in a non-empty binary tree is one more than the number of nodes in the tree.

**Proof:** Replace all empty subtrees with a leaf node. This is a full binary tree, having \#leaves = \#empty subtrees of original tree.

alternative **Proof:**
- by definition, every node has 2 children, whether empty or not
- hence a tree with \( n \) nodes has \( 2n \) children
- every node (except the root) has 1 parent
  \[ \Rightarrow \text{there are } n - 1 \text{ parent nodes (some coincide)} \]
  \[ \Rightarrow \text{there are } n - 1 \text{ non-empty children} \]
- hence \( \#(\text{empty children}) = \#(\text{total children}) - \#(\text{non-empty children}) = 2n - (n - 1) = n + 1. \]
interface BinNode { // ADT for binary tree nodes
    // Return and set the element value
    public Object element();
    public Object setElement(Object v);

    // Return and set the left child
    public BinNode left();
    public BinNode setLeft(BinNode p);

    // Return and set the right child
    public BinNode right();
    public BinNode setRight(BinNode p);

    // Return true if this is a leaf node
    public boolean isLeaf();
} // interface BinNode
Traversals

Any process for visiting the nodes in some order is called a traversal.

Any traversal that lists every node in the tree exactly once is called an enumeration of the tree’s nodes.

Preorder traversal: Visit each node before visiting its children.

Postorder traversal: Visit each node after visiting its children.

Inorder traversal: Visit the left subtree, then the node, then the right subtree.

NB: an empty node (tree) represented by Java’s null (object) value

```java
void preorder(BinNode rt) // rt is root of subtree {
    if (rt == null) return; // Empty subtree
    visit(rt);
    preorder(rt.left());
    preorder(rt.right());
}
```
Traversals (cont.)

This is a left-to-right preorder: first visit left subtree, then the right one.

Get a right-to-left preorder by switching last two lines

To get inorder or postorder, just rearrange the last three lines.
Binary Tree Implementation

[Leaves are the same as internal nodes. Lots of wasted space.]

[Example of expression tree: \((4x \cdot (2x + a)) - c\). Leaves are different from internal nodes.]
Two implementations of BinNode

class LeafNode implements BinNode { // Leaf node
  private String var; // Operand value

  public LeafNode(String val) { var = val; }
  public Object element() { return var; }
  public Object setElement(Object v)
  { return var = (String)v; }
  public BinNode left() { return null; }
  public BinNode setLeft(BinNode p) { return null; }
  public BinNode right() { return null; }
  public BinNode setRight(BinNode p) { return null; }
  public boolean isLeaf() { return true; }
} // class LeafNode

class IntlNode implements BinNode { // Internal node
  private BinNode left; // Left child
  private BinNode right; // Right child
  private Character opx; // Operator value

  public IntlNode(Character op, BinNode l, BinNode r)
  { opx = op; left = l; right = r; } // Constructor
  public Object element() { return opx; }
  public Object setElement(Object v)
  { return opx = (Character)v; }
  public BinNode left() { return left; }
  public BinNode setLeft(BinNode p) {return left = p;}
  public BinNode right() { return right; }
  public BinNode setRight(BinNode p)
  { return right = p; }
  public boolean isLeaf() { return false; }
} // class IntlNode
Two implementations (cont)

```java
static void traverse(BinNode rt) {  // Preorder
    if (rt == null) return;        // Nothing to visit
    if (rt.isLeaf())              // Do leaf node
        System.out.println("Leaf: " + rt.element());
    else {                        // Do internal node
        System.out.println("Internal: " + rt.element());
        traverse(rt.left());
        traverse(rt.right());
    }
}
```
A note on polymorphism and dynamic binding

The member function `isLeaf()` allows one to distinguish the “type” of a node

- leaf
- internal

without need of knowing its subclass

This is determined dynamically by the JRE (Java Runtime Environment)
Space Overhead

From Full Binary Tree Theorem:
    Half of pointers are NULL.

If leaves only store information, then overhead depends on whether tree is full.

All nodes the same, with two pointers to children:
    Total space required is \((2p + d)n\).
    Overhead: \(2pn\).

If \(p = d\), this means \(2p/(2p + d) = 2/3\) overhead.

[The following is for full binary trees:]
Eliminate pointers from leaf nodes:

\[
\frac{n/2(2p)}{n/2(2p) + dn} = \frac{p}{p + d}
\]

[Half the nodes have 2 pointers, which is overhead.]
This is 1/2 if \(p = d\).

\(2p/(2p + d)\) if data only at leaves ⇒ 2/3 overhead.

Some method is needed to distinguish leaves from internal nodes.  [This adds overhead.]
Array Implementation

[This is a good example of logical representation vs. physical implementation.]

For complete binary trees.

```
Node  0  1  2  3  4  5  6  7  8  9  10  11
```

- Parent\( (r) = \left\lfloor \frac{(r - 1)}{2} \right\rfloor \text{ if } r \neq 0 \text{ and } r < n. \right\rfloor
- Leftchild\( (r) = \left\lfloor 2r + 1 \right\rfloor \text{ if } 2r + 1 < n. \right\rfloor
- Rightchild\( (r) = \left\lfloor 2r + 2 \right\rfloor \text{ if } 2r + 2 < n. \right\rfloor
- Leftsibling\( (r) = \left\lfloor r - 1 \right\rfloor \text{ if } r \text{ is even, } r > 0 \text{ and } r < n. \right\rfloor
- Rightsibling\( (r) = \left\lfloor r + 1 \right\rfloor \text{ if } r \text{ is odd, } r + 1 < n. \right\rfloor

[Since the complete binary tree is so limited in its shape, (only one shape for tree of \( n \) nodes), it is reasonable to expect that space efficiency can be achieved.

NB: left sons’ indices are always odd, right ones’ even, a node with index \( i \) is leaf iff \( i > n.of.nodes/2 \) (Full Binary Tree Theorem)]
Binary Search Trees

Binary Search Tree (BST) Property

All elements stored in the left subtree of a node whose value is \( K \) have values less than \( K \). All elements stored in the right subtree of a node whose value is \( K \) have values greater than or equal to \( K \).

[Problem with lists: either insert/delete or search must be \( \Theta(n) \) time. How can we make both update and search efficient? Answer: Use a new data structure.]
BinNode Class

interface BinNode { // ADT for binary tree nodes
    // Return and set the element value
    public Object element();
    public Object setElement(Object v);

    // Return and set the left child
    public BinNode left();
    public BinNode setLeft(BinNode p);

    // Return and set the right child
    public BinNode right();
    public BinNode setRight(BinNode p);

    // Return true if this is a leaf node
    public boolean isLeaf();
} // interface BinNode

We assume that the datum in the nodes implements interface Elem with a method *key* used for comparisons (in searching and sorting algorithms)

interface Elem {
    public abstract int key();
} // interface Elem
public class BST { // Binary Search Tree implementation
    private BinNode root; // The root of the tree

    public BST() { root = null; } // Initialize root
    public void clear() { root = null; }
    public void insert(Elem val)
        { root = inserthelp(root, val); }
    public void remove(int key)
        { root = removehelp(root, key); }
    public Elem find(int key)
        { return findhelp(root, key); }
    public boolean isEmpty() { return root == null; }

    public void print() {
        if (root == null)
            System.out.println("The BST is empty.");
        else {
            printhelp(root, 0);
            System.out.println();
        }
    }

    private Elem findhelp(BinNode rt, int key) {
        if (rt == null) return null;
        Elem it = (Elem)rt.element();
        if (it.key() > key) return findhelp(rt.left(), key);
        else if (it.key() == key) return it;
        else return findhelp(rt.right(), key);
    }
}
private BinNode inserthelp(BinNode rt, Elem val) {
    if (rt == null) return new BinNode(val);
    Elem it = (Elem) rt.element();
    if (it.key() > val.key())
        rt.setLeft(inserthelp(rt.left(), val));
    else
        rt.setRight(inserthelp(rt.right(), val));
    return rt;
}
private BinNode deletemin(BinNode rt) {
    if (rt.left() == null)
        return rt.right();
    else {
        rt.setLeft(deletemin(rt.left()));
        return rt;
    }
}

private Elem getmin(BinNode rt) {
    if (rt.left() == null)
        return (Elem)rt.element();
    else return getmin(rt.left());
}
private BinNode removehelp(BinNode rt, int key) {
    if (rt == null) return null;
    Elem it = (Elem) rt.element();
    if (key < it.key())
        rt.setLeft(removehelp(rt.left(), key));
    else if (key > it.key())
        rt.setRight(removehelp(rt.right(), key));
    else {
        if (rt.left() == null)
            rt = rt.right();
        else if (rt.right() == null)
            rt = rt.left();
        else {
            Elem temp = getmin(rt.right());
            rt.setElement(temp);
            rt.setRight(deletemin(rt.right()));
        }
    }
    return rt;
}
Cost of BST Operations

Find: the depth of the node being found

Insert: the depth of the node being inserted

Remove: the depth of the node being removed, if it has < 2 children, otherwise depth of node with smallest value in its right subtree

Best case: balanced (complete tree): $\Theta(\log n)$

Worst case (linear tree): $\Theta(n)$

That’s why it is important to have a balanced (complete) BST

Cost of constructing a BST by means of a series of insertions
- if elements inserted in in order of increasing value $\sum_{i=1}^{n} i = \Theta(n^2)$
- if inserted in "random" order almost good enough for balancing the tree, insertion cost is in average $\Theta(\log n)$, for a total $\Theta(n \log n)$
Heaps

Heap: Complete binary tree with the **Heap Property**:

- Min-heap: all values less than child values.
- Max-heap: all values greater than child values.

The values in a heap are **partially ordered**.

Heap representation: normally the array based complete binary tree representation.
Building the Heap

[Max Heap

NB: for a given set of values, the heap is not unique]

(a) requires exchanges (4-2), (4-1), (2-1), (5-2), (5-4), (6-3), (6-5), (7-5), (7-6).

(b) requires exchanges (5-2), (7-3), (7-1), (6-1).

[How to get a good number of exchanges? By induction. Heapify the root's subtrees, then push the root to the correct level.]
The siftdown procedure

To place a generic node in its correct position

Assume subtrees are Heaps

If root is not greater than both children, swap with greater child

Reapply on modified subtree

Shift it down by exchanging it with the greater of the two sons, until it becomes a leaf or it is greater than both sons.
Max Heap Implementation

public class MaxHeap {
    private Elem[] Heap;  // Pointer to the heap array
    private int size;    // Maximum size of the heap
    private int n;       // Number of elements now in heap

    public MaxHeap(Elem[] h, int num, int max) {
        Heap = h; n = num; size = max; buildheap();
    }

    public int heapsize() // Return current size of heap
    { return n; }

    public boolean isLeaf(int pos) // TRUE if pos is leaf
    { return (pos >= n/2) && (pos < n); }

    // Return position for left child of pos
    public int leftchild(int pos) {
        if (pos < n/2)
            return 2*pos + 1;
    }

    // Return position for right child of pos
    public int rightchild(int pos) {
        if (pos < (n-1)/2)
            return 2*pos + 2;
    }

    public int parent(int pos) { // Return pos for parent
        if (pos > 0)
            return (pos-1)/2;
    }
}
Siftdown

For fast heap construction:
- Work from high end of array to low end.
- Call siftdown for each item.
- Don't need to call siftdown on leaf nodes.

```java
public void buildheap() // Heapify contents of Heap
{ for (int i=n/2-1; i>=0; i--) siftdown(i); }

private void siftdown(int pos) { // Put in place
    Assert.notFalse((pos >= 0) && (pos < n),
        "Illegal heap position");
    while (!isLeaf(pos)) {
        int j = leftchild(pos);
        if ((j<(n-1)) && (Heap[j].key() < Heap[j+1].key()))
            j++; // j now index of child with greater value
        if (Heap[pos].key() >= Heap[j].key()) return;
        DSutil.swap(Heap, pos, j);
        pos = j; // Move down
    }
}
```
Cost for heap construction

\[ \log n \sum_{i=1}^{\log n} (i - 1) \frac{n}{2^i} = \Theta(n). \]

\((i - 1)\) is number of steps down, \(n/2^i\) is number of nodes at that level.

cfr. eq(2.7) p.28:

\[ \sum_{i=1}^{n} \frac{i}{2^i} = 2 - \frac{n+2}{2^n} \]

notice that

\[ \sum_{i=1}^{\log n} (i - 1) \frac{n}{2^i} \leq n \sum_{i=1}^{n} \frac{i}{2^i} \]

Cost of removing root is \(\Theta(\log n)\)

Remove element too (root is a special case thereof)
Priority Queues

A priority queue stores objects, and on request releases the object with greatest value.

Example: Scheduling jobs in a multi-tasking operating system.

The **priority** of a job may change, requiring some reordering of the jobs.

Implementation: use a heap to store the priority queue.

To support priority reordering, delete and re-insert. Need to know index for the object.

```java
// Remove value at specified position
public Elem remove(int pos) {
    Assert.notFalse((pos >= 0) && (pos < n),
        "Illegal heap position");
    DSutil.swap(Heap, pos, --n); // Swap with last value
    while (Heap[pos].key() > Heap[parent(pos)].key())
        DSutil.swap(Heap, pos, parent(pos)); // push up
    if (n != 0) siftdown(pos); // push down
    return Heap[n];
}
```
A **tree** $T$ is a finite set of nodes such that it is empty or there is one designated node $r$ called the root of $T$, and the remaining nodes in $(T - \{r\})$ are partitioned into $n \geq 0$ disjoint subsets $T_1, T_2, \ldots, T_k$, each of which is a tree.

[Note: disjoint because a node cannot have two parents.]
General Tree ADT

[There is no concept of "left" or "right" child. But, we can impose a concept of "first" (leftmost) and "next" (right).]

```java
public interface GTNode {
    public Object value();
    public boolean isLeaf();
    public GTNode parent();
    public GTNode leftmost_child();
    public GTNode right_sibling();
    public void setValue(Object value);
    public void setParent(GTNode par);
    public void insert_first(GTNode n);
    public void insert_next(GTNode n);
    public void remove_first(); // remove first child
    public void remove_next(); // remove right sibling
}

public interface GenTree {
    public void clear();
    public GTNode root();
    public void newroot(Object value, GTNode first,
                        GTNode sib);
}
```
General Tree Traversal

[preorder traversal]

static void print(GTNode rt) { // Preorder traversal
    if (rt.isLeaf()) System.out.print("Leaf: ");
    else System.out.print("Internal: ");
    System.out.println(rt.value());
    GTNode temp = rt.leftmost_child();
    while (temp != null) {
        print(temp);
        temp = temp.right_sibling();
    }
}

[RACDEBF]
General Tree Implementations

Lists of Children

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<th>Par</th>
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</tbody>
</table>

[Hard to find right sibling.]
Leftmost Child/Right Sibling

[Note: Two trees share same array.]
Linked Implementations

[Allocate child pointer space when node is created.]
Sequential Implementations

List node values in the order they would be visited by a preorder traversal.

Saves space, but allows only sequential access.

Need to retain tree structure for reconstruction.

For binary trees: Use symbol to mark NULL links.

```
AB/D/CEG/FH/I/
```
Sequential Implementations (cont.)

Full binary trees: Mark leaf or internal.

[Need NULL mark since this tree is not full.]

\[A'B'/DC'E'G/F'H]I\]

General trees: Mark end of each subtree.

\[RAC\)D\)E\)\)BF\)\)]
Convert to Binary Tree

Left Child/Right Sibling representation essentially stores a binary tree.

Use this process to convert any general tree to a binary tree.

A **forest** is a collection of one or more general trees.

[Dynamiic implementation of “Left child/right sibling.”]
K-ary Trees

Every node has a fixed maximum number of children

fixed $\# \text{ children} \Rightarrow$ easy to implement, also in array

$K$ high $\Rightarrow$ potentially many empty subtrees $\Rightarrow$ different implementation for leaves becomes convenient

Full and complete K-ary trees similar to binary trees

Theorems on $\# \text{ empty subtrees}$ and on relation between $\# \text{ internal nodes}$ and $\# \text{ leaves}$ similar to binary trees
Graphs

**Graph** $G = (V, E)$: a set of **vertices** $V$, and a set of **edges** $E$; each edge in $E$ is a connection between a pair of vertices in $V$, which are called **adjacent** vertices.

# vertices written $|V|$; # edges written $|E|$. $0 \leq |E| \leq |V|^2$.

A graph is
- **sparse** if it has "few" edges
- **dense** if it has "many" edges
- **complete** all possible edges
- **undirected** as in figure (a)
- **directed** as in figure (b)
- **labeled** (figure (c)) if it has labels on vertices
- **weighted** (figure (c)) if it has (numeric) labels on edges
Graph Definitions (Cont)

A sequence of vertices \( v_1, v_2, ..., v_n \) forms a **path** of length \( n - 1 \) (\( \Rightarrow \) length = \# edges) if there exist edges from \( v_i \) to \( v_{i+1} \) for \( 1 \leq i < n \).

A path is **simple** if all vertices on the path are distinct.

In a **directed** graph

- a path \( v_1, v_2, ..., v_n \) forms a **cycle** if \( n > 1 \) and \( v_1 = v_n \). The cycle is **simple** if, in addition, \( v_2, ..., v_n \) are distinct

- a cycle \( v, v \) is a **self-loop**

- a directed graph with no self-loops is **simple**

In an **undirected** graph

- a path \( v_1, v_2, ..., v_n \) forms a (simple) **cycle** if \( n > 3 \) and \( v_1 = v_n \) (and, in addition, \( v_2, ..., v_n \) are distinct)

  - hence the path ABA is *not* a cycle, while ABCA *is* a cycle
Graph Definitions (Cont)

Subgraph \( S = (V_S, E_S) \) of a graph \( G = (V, E) \):
\[ V_S \subset V \text{ and } E_S \subset E \text{ and both vertices of any edge in } E_S \text{ are in } V_S \]

An undirected graph is connected if there is at least one path from any vertex to any other.

The maximal connected subgraphs of an undirected graph are called connected components.

A graph without cycles is acyclic.

A directed graph without cycles is a directed acyclic graph or DAG.

A free tree is a connected, undirected graph with no cycles. Equivalently, a free tree is connected and has \(|V - 1|\) edges.
Connected Components

A graph with (composed of) 3 connected components
Graph Representations

**Adjacency Matrix**: space required $\Theta(|V|^2)$.

**Adjacency List**: space required $\Theta(|V| + |E|)$.

[Instead of bits, the graph could store edge, weights.]
Graph Representations (cont)

Adjacency list efficient for sparse graphs (only existing edges coded)

Matrix efficient for dense graphs (no pointer overload)

Algorithms visiting each neighbor of each vertex more efficient on adjacency lists, especially for sparse graphs
interface Graph { // Graph class ADT
    public int n(); // Number of vertices
    public int e(); // Number of edges
    // Get first edge having v as vertex v1
    public Edge first(int v);
    // Get next edge having w.v1 as the first edge
    public Edge next(Edge w);
    public boolean isEdge(Edge w); // True if edge
    public boolean isEdge(int i, int j); // True if edge
    public int v1(Edge w); // Where from
    public int v2(Edge w); // Where to
    public void setEdge(int i, int j, int weight);
    public void setEdge(Edge w, int weight);
    public void delEdge(Edge w); // Delete edge w
    public void delEdge(int i, int j); // Delete (i, j)
    public int weight(int i, int j); // Return weight
    public int weight(Edge w); // Return weight

    // Set Mark of vertex v
    public void setMark(int v, int val);

    // Get Mark of vertex v
    public int getMark(int v);
} // interface Graph

Edges have a double nature:
seen as pairs of vertices or as aggregate objects.

Vertices identified by an integer \( i, 0 \leq i \leq |V| \)
interface Edge { // Interface for graph edges
    public int v1(); // Return the vertex it comes from
    public int v2(); // Return the vertex it goes to
} // interface Edge

// Edge class for Adjacency Matrix graph representation
class Edgem implements Edge {
    private int vert1, vert2; // The vertex indices

    public Edgem(int vt1, int vt2) //the constructor
    { vert1 = vt1; vert2 = vt2; }
    public int v1() { return vert1; }
    public int v2() { return vert2; }
} // class Edgem
Implementation: Adjacency Matrix

class Graphm implements Graph { // Adjacency matrix
    private int[][] matrix; // The edge matrix
    private int numEdge; // Number of edges
    public int[] Mark; // The mark array, initially all 0

    public Graphm(int n) { // Constructor
        Mark = new int[n];
        matrix = new int[n][n];
        numEdge = 0;
    }

    public int n() { return Mark.length; }
    public int e() { return numEdge; }

    public Edge first(int v) { // Get first edge
        for (int i=0; i<Mark.length; i++)
            if (matrix[v][i] != 0)
                return new Edgem(v, i);
        return null; // No edge for this vertex
    }

    public Edge next(Edge w) { // Get next edge
        if (w == null) return null;
        for (int i=w.v2()+1; i<Mark.length; i++)
            if (matrix[w.v1()][i] != 0)
                return new Edgem(w.v1(), i);
        return null; // No next edge;
    }

    Class Graphm implements interface Graph
    Class Edgem implements interface Edge
public boolean isEdge(Edge w) { // True if an edge
    if (w == null) return false;
    else return matrix[w.v1()][w.v2()] != 0;
}

public boolean isEdge(int i, int j) // True if edge
    { return matrix[i][j] != 0; }

public int v1(Edge w) {return w.v1();} // Where from
public int v2(Edge w) {return w.v2();} // Where to

public void setEdge(int i, int j, int wt) {
    Assert.notFalse(wt!=0, "Cannot set weight to 0");
    if (matrix[i][j] == 0) numEdge++;
    matrix[i][j] = wt;
}

public void setEdge(Edge w, int weight) // Set weight
    { if (w != null) setEdge(w.v1(), w.v2(), weight); }

public void delEdge(Edge w) { // Delete edge w
    if (w != null)
        if (matrix[w.v1()][w.v2()] != 0)
            { matrix[w.v1()][w.v2()] = 0; numEdge--; }
}

public void delEdge(int i, int j) { // Delete (i, j)
    if (matrix[i][j] != 0)
        { matrix[i][j] = 0; numEdge--; }
}

NB: matrix[i][j]==0 iff there is no edge (i,j)
If there is no edge (i,j) then
weight(i,j)==Integer.MAX_VALUE (INFINITY)
public int weight(int i, int j) { // Return weight
    if (matrix[i][j] == 0) return Integer.MAX_VALUE;
    else return matrix[i][j];
}

public int weight(Edge w) { // Return edge weight
    Assert.notNull(w,"Can’t take weight of null edge");
    if (matrix[w.v1()][w.v2()] == 0)
        return Integer.MAX_VALUE;
    else return matrix[w.v1()][w.v2()];
}

public void setMark(int v, int val)
    { Mark[v] = val; }

public int getMark(int v) { return Mark[v]; }
} // class Graphm
Graph Traversals

Some applications require visiting every vertex in the graph exactly once.

Application may require that vertices be visited in some special order based on graph topology.

Example: Artificial Intelligence
- Problem domain consists of many “states.”
- Need to get from Start State to Goal State.
- Start and Goal are typically not directly connected.

To insure visiting all vertices:

void graphTraverse(Graph G) {
    for (v=0; v<G.n(); v++)
        G.setMark(v, UNVISITED); // Initialize mark bits
    //next for needed to cover all the graph in case
    //of graph composed of several connected components
    for (v=0; v<G.n(); v++)
        if (G.getMark(v) == UNVISITED)
            doTraverse(G, v);
}

[Two traversals we will talk about: DFS, BFS.]
Depth First Search

static void DFS(Graph G, int v) { // Depth first search
    PreVisit(G, v); // Take appropriate action
    G.setMark(v, VISITED);
    for (Edge w = G.first(v); G.isEdge(w); w = G.next(w))
        if (G.getMark(G.v2(w)) == UNVISITED)
            DFS(G, G.v2(w));
    PostVisit(G, v); // Take appropriate action
}

Cost: \( \Theta(|V| + |E|) \).

[The directions are imposed by the traversal. This is the Depth First Search Tree.]

If PreVisit simply prints and PostVisit does nothing then DFS prints
A C B F D E
Breadth First Search

Like DFS, but replace stack with a queue. Visit the vertex's neighbors before continuing deeper in the tree.

```java
static void BFS(Graph G, int start) {
    Queue Q = new AQueue(G.n()); // Use a Queue
    Q.enqueue(new Integer(start));
    G.setMark(start, VISITED);
    while (!Q.isEmpty()) { // Process each vertex on Q
        int v = ((Integer)Q.dequeue()).intValue();
        PreVisit(G, v); // Take appropriate action
        for (Edge w=G.first(v); G.isEdge(w); w=G.next(w))
            if (G.getMark(G.v2(w)) == UNVISITED) {
                G.setMark(G.v2(w), VISITED);
                Q.enqueue(new Integer(G.v2(w)));
            }
        PostVisit(G, v); // Take appropriate action
    }
}
```

If PreVisit simply prints and PostVisit does nothing then BFS prints A C E B D F

![Diagram](image1)

![Diagram](image2)

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Topological Sort

Problem: Given a set of jobs, courses, etc. with prerequisite constraints, output the jobs in an order that does not violate any of the prerequisites. (NB: the graph must be a DAG)

static void topsort(Graph G) { // Topo sort: recursive
    for (int i=0; i<G.n(); i++) // Initialize Mark array
        G.setMark(i, UNVISITED);
    for (int i=0; i<G.n(); i++) // Process all vertices
        if (G.getMark(i) == UNVISITED)
            tophelp(G, i); // Call helper function
}

static void tophelp(Graph G, int v) { // Topsort helper
    G.setMark(v, VISITED);
    for (Edge w = G.first(v); G.isEdge(w); w = G.next(w))
        if (G.getMark(G.v2(w)) == UNVISITED)
            tophelp(G, G.v2(w));
    printout(v); // PostVisit for Vertex v
}

[Prints in reverse order: J7, J5, J4, J6, J2, J3, J1]

It is a DFS with a PreVisit that does nothing]
Queue-based Topological Sort

static void topsort(Graph G) { // Topo sort: Queue
    Queue Q = new AQueue(G.n());
    int[] Count = new int[G.n()];
    int v;
    for (v=0; v<G.n(); v++) Count[v] = 0; // Initialize
    for (v=0; v<G.n(); v++) { // Process every edge
        for (Edge w=G.first(v); G.isEdge(w); w=G.next(w))
            Count[G.v2(w)]++; // Add to v2’s count
    }
    for (v=0; v<G.n(); v++) // Initialize Queue
        if (Count[v] == 0) // Vertex has no prereqs
            Q.enqueue(new Integer(v));
    while (!Q.isEmpty()) { // Process the vertices
        v = ((Integer)Q.dequeue()).intValue(); // PreVisit for Vertex V
        printout(v);
        for (Edge w=G.first(v); G.isEdge(w); w=G.next(w)) {
            Count[G.v2(w)]--; // One less prerequisite
            if (Count[G.v2(w)] == 0) // This vertex now free
                Q.enqueue(new Integer(G.v2(w)));
        }
    }
}
Sorting

Each record is stored in an array and contains a field called the **key**.

Linear (i.e., total) order: comparison.

\[ a < b \text{ and } b < c \Rightarrow a < c. \]

**The Sorting Problem**

Given a sequence of records \( R_1, R_2, ..., R_n \) with key values \( k_1, k_2, ..., k_n \), respectively, arrange the records into any order \( s \) such that records \( R_{s_1}, R_{s_2}, ..., R_{s_n} \) have keys obeying the property \( k_{s_1} \leq k_{s_2} \leq ... \leq k_{s_n} \).

[Put keys in ascending order. ]

NB: there can be records with the same key

A sorting algorithm is **stable** if after sorting records with the same key have the same relative position as before

Measures of cost:

- Comparisons
- Swaps (when records are large)
Sorting (cont)

Assumptions: for every record type there are functions

- R.key returns the key value for record R
- DSutil.swap(array, i, j) swaps records in positions i and j of the array

Measure of the "degree of disorder" of an array in the number of INVERSIONS

\[ \forall el = a[i], \]
\[ \#inversions = \#elements > el \text{ which are in a position } j < i \]

\[ \#inversions \text{ for the entire array } = \]
\[ = \sum \#inversions \text{ of each array element} \]

For a sorted array \( \#inversions = 0 \)

For an array with elements in decreasing order \( \#inversions = \Theta(n^2) \)
Insertion Sort

static void inssort(Elem[] array) { // Insertion Sort
    for (int i=1; i<array.length; i++) // Insert record
        for (int j=i; (j>0) &&
             (array[j].key()<array[j-1].key()); j--)
             DSutil.swap(array, j, j-1);
}

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<th></th>
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<th>3</th>
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<td>23</td>
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<td>28</td>
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<td>15</td>
<td>15</td>
<td>15</td>
<td>15</td>
<td>42</td>
</tr>
</tbody>
</table>

Best Case: [0 swaps, \( n - 1 \) comparisons]

Worst Case: [\( n^2/2 \) swaps and compares]

Average Case: [\( n^2/4 \) swaps and compares: \# inner loop iterations for an element in position \( n = \#\text{inversione} = n/2 \) in the average]

[At each iteration takes one element to its place and does only that; it works only on the sorted portion of the array]

[Nearly best performance when input "nearly sorted" ⇒ used in conjunction with mergesort and quicksort small array segments]
Bubble Sort

```java
static void bubblesort(Elem[] array) { // Bubble Sort
    for (int i=0; i<array.length-1; i++) // Bubble up
        for (int j=array.length-1; j>i; j--)
            if (array[j].key() < array[j-1].key())
                DSutil.swap(array, j, j-1);
}
```

[Using test “j > i” saves a factor of 2 over “j > 0”.]

```
<table>
<thead>
<tr>
<th></th>
<th>0</th>
<th>1</th>
<th>2</th>
<th>3</th>
<th>4</th>
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<td>23</td>
<td>28</td>
<td>28</td>
<td>28</td>
<td>42</td>
<td>42</td>
</tr>
</tbody>
</table>
```

Best Case: \([n^2/2\text{ compares, 0 swaps}]\)

Worst Case: \([n^2/2\text{ compares, }n^2/2\text{ swaps}]\)

Average Case: \([n^2/2\text{ compares, }n^2/4\text{ swaps}]\)

[At each iteration takes the smallest to its place, but it moves also other ones;]

NB: it works also on the unsorted part of the array;

No redeeming features to this sort.]
Selection Sort

static void selsort(Elem[] array) { // Selection Sort
    for (int i=0; i<array.length-1; i++) { // Select i’th
        int lowindex = i; // Remember its index
        for (int j=array.length-1; j>i; j--) // Find least
            if (array[j].key() < array[lowindex].key())
                lowindex = j; // Put it in place
        DSutil.swap(array, i, lowindex);
    }
}

[Select the value to go in the &ith position.]

<table>
<thead>
<tr>
<th>i=0</th>
<th>1</th>
<th>2</th>
<th>3</th>
<th>4</th>
<th>5</th>
<th>6</th>
</tr>
</thead>
<tbody>
<tr>
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<td>13</td>
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<td>13</td>
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<td>17</td>
<td>42</td>
<td>42</td>
<td>42</td>
<td>42</td>
</tr>
</tbody>
</table>

Best Case: [0 swaps (n – 1 as written), n^2/2 compares.]

Worst Case: [n – 1 swaps, n^2/2 compares]

Average Case: [O(n) swaps, n^2/2 compares]

[It minimizes # swaps]
Pointer Swapping

[For large records.]
This is what done in Java, when records are objects
## Exchange Sorting

### Summary

<table>
<thead>
<tr>
<th></th>
<th>Insertion</th>
<th>Bubble</th>
<th>Selection</th>
</tr>
</thead>
<tbody>
<tr>
<td><strong>Comparisons:</strong></td>
<td></td>
<td></td>
<td></td>
</tr>
<tr>
<td>Best Case</td>
<td>$\Theta(n)$</td>
<td>$\Theta(n^2)$</td>
<td>$\Theta(n^2)$</td>
</tr>
<tr>
<td>Average Case</td>
<td>$\Theta(n^2)$</td>
<td>$\Theta(n^2)$</td>
<td>$\Theta(n^2)$</td>
</tr>
<tr>
<td>Worst Case</td>
<td>$\Theta(n^2)$</td>
<td>$\Theta(n^2)$</td>
<td>$\Theta(n^2)$</td>
</tr>
<tr>
<td><strong>Swaps:</strong></td>
<td></td>
<td></td>
<td></td>
</tr>
<tr>
<td>Best Case</td>
<td>0</td>
<td>0</td>
<td>$\Theta(n)$</td>
</tr>
<tr>
<td>Average Case</td>
<td>$\Theta(n^2)$</td>
<td>$\Theta(n^2)$</td>
<td>$\Theta(n)$</td>
</tr>
<tr>
<td>Worst Case</td>
<td>$\Theta(n^2)$</td>
<td>$\Theta(n^2)$</td>
<td>$\Theta(n)$</td>
</tr>
</tbody>
</table>
Mergesort

List mergesort(List inlist) {
    if (inlist.length() <= 1) return inlist;;
    List l1 = half of the items from inlist;
    List l2 = other half of the items from inlist;
    return merge(mergesort(l1), mergesort(l2));
}

Analyze first the algorithm for merging sorted sublists

- examine first element of each sublist
- pick the smaller element (it is the smallest overall)
- remove it from its sublist and put it in the output list
- when one sublist is exhausted, pick from the other

Complexity of merging two sorted sublist: \( \Theta(n) \)
Mergesort Implementation

Mergesort is tricky to implement. Main question: how to represent lists?

Linked lists

- merging does not require direct access, but...
- splitting requires a list traversal ($\Theta(n)$), whether list size is known (take as two sublists the first and second halves) or unknown (assign elements alternating between the two lists)

Lists represented by arrays

- splitting very easy ($\Theta(1)$) if array bounds are known
- merging easy ($\Theta(n)$) only if sub-arrays merged into a second array (hence double the space requirement!)
- avoid the need for a distinct additional array for each recursive call by first copying sub-arrays into auxiliary array and then merging them back to the original array (hence can use only one array for the overall process)
Mergesort Implementation (2)

```java
static void mergesort(Elem[] array, Elem[] temp, int l, int r) {
    if (l == r) return; // One element list
    int mid = (l+r)/2; // Select midpoint
    mergesort(array, temp, l, mid); // Sort first half
    mergesort(array, temp, mid+1, r); // Sort second half
    merge(array, temp, l, mid, mid+1, r);
}

static void merge(Elem[] array, Elem[] temp, int l1, int r1, int l2, int r2) {
    for (int i=l1; i<=r2; i++) // Copy subarrays
        temp[i] = array[i];
    // Do the merge operation back to array
    int i1 = l1; int i2 = l2;
    for (int curr=l1; curr<=r2; curr++) {
        if (i1 > r1) // Left sublist exhausted
            array[curr] = temp[i2++];
        else if (i2 > r2) // Right sublist exhausted
            array[curr] = temp[i1++];
        else if (temp[i1].key() < temp[i2].key())
            array[curr] = temp[i1++]; // Get smaller val
        else array[curr] = temp[i2++];
    }
}
```
Complexity of Mergesort

Ad hoc analysis

- depth of recursion is \( \log n \)
- at each recursion depth \( i \)
  - \( 2^i \) recursive calls
  - each recursive call has array length \( n/2^i \), hence...
  - total length of merged arrays is \( n \) at every depth
- therefore total cost is \( T(n) = \Theta(n \log n) \)

Alternative analysis: use recurrence equation

\[
T(n) = aT(n/b) + cn^k = 2T(n/2) + cn, \quad T(1) = d
\]

We have \( a = b = 2, \ k = 1 \) and therefore \( a = b^k \), hence

\[
T(n) = \Theta(n \log n)
\]
Heapsort uses a max-heap.

```java
static void heapsort(Elem[] array) { // Heapsort
    MaxHeap H = new MaxHeap(array, array.length,
                            array.length);
    for (int i=0; i<array.length; i++) // Now sort
        H.removemax(); // Put max value at end of heap
}
```

Cost of Heapsort: $[\Theta(n \log n)]$

Cost of finding $k$ largest elements: $[\Theta(k \log n + n)$.

Time to build heap: $\Theta(n)$.

Time to remove least element: $\Theta(\log n)$.

[Compare to sorting with BST: this is expensive in space (overhead), potential bad balance, BST does not take advantage of having all records available in advance.]

[Heap is space efficient, balanced, and building initial heap is efficient.]
Heapsort Example

Original Numbers

| 73 | 6  | 57 | 88 | 60 | 42 | 83 | 72 | 48 | 85 |

Build Heap

| 88 | 85 | 83 | 72 | 73 | 42 | 57 | 6  | 48 | 60 |

Remove 88

| 85 | 73 | 83 | 72 | 60 | 42 | 57 | 6  | 48 | 88 |

Remove 85

| 83 | 73 | 57 | 72 | 60 | 42 | 48 | 6  | 85 | 88 |

Remove 83

| 73 | 72 | 57 | 6  | 60 | 42 | 48 | 83 | 85 | 88 |
## Empirical Comparison

### [MS Windows – CISC]

<table>
<thead>
<tr>
<th>Algorithm</th>
<th>10</th>
<th>100</th>
<th>1000</th>
<th>10,000</th>
</tr>
</thead>
<tbody>
<tr>
<td>Insert. Sort</td>
<td>0.10</td>
<td>9.5</td>
<td>957.9</td>
<td>98,086</td>
</tr>
<tr>
<td>Bubble Sort</td>
<td>0.13</td>
<td>14.3</td>
<td>1470.3</td>
<td>157,230</td>
</tr>
<tr>
<td>Select. Sort</td>
<td>0.11</td>
<td>9.9</td>
<td>1018.9</td>
<td>104,897</td>
</tr>
<tr>
<td>Shellsort</td>
<td>0.09</td>
<td>2.5</td>
<td>45.6</td>
<td>829</td>
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<tr>
<td>Quicksort</td>
<td>0.15</td>
<td>1.8</td>
<td>23.6</td>
<td>291</td>
</tr>
<tr>
<td>Quicksort/O</td>
<td>0.10</td>
<td>1.6</td>
<td>20.9</td>
<td>274</td>
</tr>
<tr>
<td>Mergesort</td>
<td>0.12</td>
<td>2.4</td>
<td>36.8</td>
<td>505</td>
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<tr>
<td>Mergesort/O</td>
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<td>1.8</td>
<td>28.0</td>
<td>390</td>
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<tr>
<td>Heapsort</td>
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<td>50.0</td>
<td>60.0</td>
<td>880</td>
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<tr>
<td>Radix Sort/1</td>
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<td>8.6</td>
<td>89.5</td>
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<td>0.19</td>
<td>1.2</td>
<td>11.5</td>
<td>115</td>
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</table>

### [UNIX – RISC]

<table>
<thead>
<tr>
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<th>10</th>
<th>100</th>
<th>1000</th>
<th>10,000</th>
</tr>
</thead>
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<tr>
<td>Insert. Sort</td>
<td>0.66</td>
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<td>Select. Sort</td>
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<td>7,973</td>
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<tr>
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<td>679</td>
<td>6,895</td>
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<tr>
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<td>18.7</td>
<td>160</td>
<td>1,678</td>
</tr>
<tr>
<td>Radix Sort/8</td>
<td>4.10</td>
<td>11.5</td>
<td>97</td>
<td>808</td>
</tr>
</tbody>
</table>

[Clearly, $n \log n$ superior to $n^2$. Note relative differences on different machines.]
Upperbound and Lowerbound for a Problem

**Upperbound**: asymptotic cost of the fastest *known* algorithm

**lowerbound** best possible efficiency of *any* possible (known or unknown) algorithm

**open problem**: upperbound different from (greater than) lowerbound

**closed problem**: upperbound equal to lowerbound
Sorting Lower Bound

Want to prove a lower bound sorting problem based on key comparison.

Sorting I/O takes $\Omega(n)$ time. (no algorithm can take less than I/O time)

Sorting is $O(n \log n)$.

Will now prove $\Omega(n \log n)$ lower bound.

Form of proof:
- Comparison based sorting can be modeled by a binary tree.
- The tree must have $\Omega(n!)$ leaves (because there are $n!$ permutations of $n$ elements).
- The tree cannot be less than $\Omega(n \log n)$ levels deep (a tree with $k$ nodes has at least $\log k$ levels).

this comes from the fact that
$$
\log n! = \Theta(n \log n)
$$

which is due to Stirling’s approximation of $n!:
$$
n! \approx \sqrt{2\pi n} \left(\frac{n}{e}\right)^n
$$

from which $\log n! \approx n \log n$
Decision Trees

[Illustration of Insertion sort. Lower part of table shows possible output (sorted version of input array) after each check]

There are \( n! \) permutations, and at least 1 node for each permutation.

Where is the worst case in the decision tree?
Primary vs. Secondary Storage

review following sections of textbook

9.1 on Primary vs secondary storage

9.2 on Disk & Tape drives

9.3 on Buffers and Buffer Pools
Buffer Pools

A series of buffers used by an application to cache disk data is called a buffer pool.

Virtual memory uses a buffer pool to imitate greater RAM memory by actually storing information on disk and “swapping” between disk and RAM.

Caching Same technique to imitate greater CACHE memory by storing info on RAM and swapping between RAM and CACHE.

Organization for buffer pools: which one to use next?

- First-in, First-out: Use the first one on the queue.
- Least Frequently Used (LFU): Count buffer accesses, pick the least used.
- Least Recently Used (LRU): Keep buffers on linked list. When a buffer is accessed, bring to front. Reuse the one at the end.
Programmer’s View of Files

Logical view of files:

• An array of bytes.

• A file pointer marks the current position.

Three fundamental operations:

• Read bytes from current position (move file pointer).

• Write bytes to current position (move file pointer).

• Set file pointer to specified byte position.
Java File Functions

RandomAccessFile(String name, String mode)

close()

read(byte[] b)

write(byte[] b)

seek(long pos)
External Sorting

Problem: Sorting data sets too large to fit in main memory.

- Assume data stored on disk drive.

To sort, portions of the data must be brought into main memory, processed, and returned to disk.

An external sort should minimize disk accesses.
Model of External Computation

Secondary memory is divided into equal-sized **blocks** (512, 2048, 4096 or 8192 bytes are typical sizes).

The basic I/O operation transfers the contents of one disk block to/from main memory.

Under certain circumstances, reading blocks of a file in sequential order is more efficient. (When?) [1) Adjacent logical blocks of file are physically adjacent on disk. 2) No competition for I/O head.]

Typically, the time to perform a single block I/O operation is sufficient to Quicksort the contents of the block.

Thus, our primary goal is to minimize the number of block I/O operations.

Most workstations today must do all sorting on a single disk drive.
Key Sorting

Often records are large while keys are small.

- Ex: Payroll entries keyed on ID number.

Approach 1: Read in entire records, sort them, then write them out again.

Approach 2: Read only the key values, store with each key the location on disk of its associated record.

If necessary, after the keys are sorted the records can be read and re-written in sorted order.

[But, this is not usually done. (1) It is expensive (random access to all records). (2) If there are multiple keys, there is no “correct” order.]
External Sort: Simple Mergesort

Quicksort requires random access to the entire set of records.

Better: Modified Mergesort algorithm

- Process \( n \) elements in \( \Theta(\log n) \) passes.

1. Split the file into two files.
2. Read in a block from each file.
3. Take first record from each block, output them in sorted order.
4. Take next record from each block, output them to a second file in sorted order.
5. Repeat until finished, alternating between output files. Read new input blocks as needed.
6. Repeat steps 2-5, except this time the input files have groups of two sorted records that are merged together.
7. Each pass through the files provides larger and larger groups of sorted records.

A group of sorted records is called a **run**.
Problems with Simple Mergesort

Is each pass through input and output files sequential?  [yes]

What happens if all work is done on a single disk drive?  [Competition for I/O head eliminates advantage of sequential processing.]

How can we reduce the number of Mergesort passes?  [Read in a block (or several blocks) and do an in-memory sort to generate large initial runs.]

In general, external sorting consists of two phases:

1. Break the file into initial runs.
2. Merge the runs together into a single sorted run.
Breaking a file into runs

General approach:
• Read as much of the file into memory as possible.
• Perform and in-memory sort.
• Output this group of records as a single run.
General Principals of External Sorting

In summary, a good external sorting algorithm will seek to do the following:

- Make the initial runs as long as possible.
- At all stages, overlap input, processing and output as much as possible.
- Use as much working memory as possible. Applying more memory usually speeds processing.
- If possible, use additional disk drives for more overlapping of processing with I/O, and allow for more sequential file processing.
Search

Given: Distinct keys $k_1, k_2, \ldots, k_n$ and collection $T$ of $n$ records of the form

$$(k_1, I_1), (k_2, I_2), \ldots, (k_n, I_n)$$

where $I_j$ is information associated with key $k_j$ for $1 \leq j \leq n$.

**Search Problem:** For key value $K$, locate the record $(k_j, I_j)$ in $T$ such that $k_j = K$.

**Exact match query:** search records with a specified key value.

**Range query:** search records with key in a specified range.

**Searching** is a systematic method for locating the record (or records) with key value $k_j = K$.

A **successful** search is one in which a record with key $k_j = K$ is found.

An **unsuccessful** search is one in which no record with $k_j = K$ is found (and presumably no such record exists).
Approaches to Search

1. Sequential and list methods (lists, tables, arrays).
2. Direct access by key value (hashing).
3. Tree indexing methods.

[recall: sequences: duplicate key values allowed; sets no key duplication]
Searching Ordered Arrays

Sequential Search

Binary Search

static int binary(int K, int[] array, int left, int right) {
    // Return position of element (if any) with value K
    int l = left-1;
    int r = right+1; // l and r are beyond array bounds
    while (l+1 != r) { // Stop when l and r meet
        int i = (l+r)/2; // Look at middle of subarray
        if (K < array[i]) r = i; // In left half
        if (K == array[i]) return i; // Found it
        if (K > array[i]) l = i; // In right half
    }
    return UNSUCCESSFUL; // Search value not in array
}

<table>
<thead>
<tr>
<th>Position</th>
<th>0</th>
<th>1</th>
<th>2</th>
<th>3</th>
<th>4</th>
<th>5</th>
<th>6</th>
<th>7</th>
<th>8</th>
<th>9</th>
<th>10</th>
<th>11</th>
<th>12</th>
<th>13</th>
<th>14</th>
<th>15</th>
</tr>
</thead>
<tbody>
<tr>
<td>Key</td>
<td>11</td>
<td>13</td>
<td>21</td>
<td>26</td>
<td>29</td>
<td>36</td>
<td>40</td>
<td>41</td>
<td>45</td>
<td>51</td>
<td>54</td>
<td>56</td>
<td>65</td>
<td>72</td>
<td>77</td>
<td>83</td>
</tr>
</tbody>
</table>

Improvement: **Dictionary Search**, expected record position computed from key value; value of key found there used as in binary search
Lists Ordered by Frequency

Order lists by (expected) frequency of occurrence ⇒ Perform sequential search.

Cost to access first record: 1; second record: 2

Expected (i.e., average) search cost:

\[ \overline{C}_n = 1p_1 + 2p_2 + \ldots + np_n \]

\([p_i \text{ is probability of } i\text{th record being accessed.}]\]

Example: all records have equal frequency

\[ \overline{C}_n = \sum_{i=1}^{n} \frac{i}{n} = \frac{(n + 1)}{2}. \]

Example: Exponential frequency

\[ p_i = \begin{cases} \frac{1}{2^i} & \text{if } 1 \leq i \leq n - 1 \\ \frac{1}{2^{n-1}} & \text{if } i = n \end{cases} \]

\([\text{Second line is to make probabilities sum to 1.}]\)

\[ \overline{C}_n \approx \sum_{i=1}^{n} \left( \frac{i}{2^i} \right) \approx 2. \]

\([\text{very good performance, because assumption (exp. freq.) is strong}]\)
Zipf Distributions

Applications:

- Distribution for frequency of word usage in natural languages.
- Distribution for populations of cities, etc.

Definition: Zipf frequency for item $i$ in the distribution for $n$ records as $1/i\mathcal{H}_n$.

$[\mathcal{H}_n = \sum_{i=1}^{n} \frac{1}{i} \approx \log_e n.]$

$$\overline{C}_n = \sum_{i=1}^{n} \frac{i}{i\mathcal{H}_n} = \frac{n}{\mathcal{H}_n} \approx \frac{n}{\log_e n}$$

80/20 rule: 80% of the accesses are to 20% of the records.

For distributions following the 80/20 rule,

$$\overline{C}_n \approx 0.122n.$$
Self-Organizing Lists

Self-organizing lists modify the order of records within the list based on the actual pattern of record access.

Based on assumption that past searches provide good indication of future ones

This is a **heuristic** similar to those for managing buffer pools.

- **Order by actual historical frequency of access.** (Similar to LFU buffer pool replacement strategy.)  [COUNT method: slow reaction to change]

- **Move-to-Front**: When a record is found, move it to the front of the list.  [Not worse than twice “best arrangement”; easy to implement with linked lists, not arrays]

- **Transpose**: When a record is found, swap it with the record ahead of it.  [A pathological, though unusual case: keep swapping last two elements.]
Advantages of self-organizing lists

• do not require sorting
• cost of insertion and deletion low
• no additional space
• simple (hence easy to implement)
Example of Self-Organizing Tables

Application: Text compression.

Keep a table of words already seen, organized via Move-to-Front Heuristic.

If a word not yet seen, send the word. Otherwise, send the (current) index in the table.

[NB: sender and receiver maintain identical lists, so they agree on indices]

The car on the left hit the car I left.

The car on 3 left hit 3 5 I 5.
Hashing

**Hashing**: The process of mapping a key value to a position in a table.

A **hash function** maps key values to positions. It is denoted by \( h \).

A **hash table** is an array that holds the records. It is denoted by \( T \).

[NB: records not necessarily ordered by key value or frequency]

The hash table has \( M \) slots, indexed from 0 to \( M - 1 \).

For any value \( K \) in the key range and some hash function \( h \),

\[
h(K) = i, \quad 0 \leq i < M, \quad \text{such that } T[i].\text{key}() = K.
\]
Hashing (continued)

Hashing is appropriate only for sets (no duplicates).

Good for both in-memory and disk based applications.

[Very good for organizing large databases on disk]

Answers the question “What record, if any, has key value $K$?”

[Not good for range queries.]

Trivial Example: Store the $n$ records with keys in range 0 to $n - 1$.

- Store the record with key $i$ in slot $i$.
- Use hash function $h(K) = K$.

Typically, there are however many more values in the key range than slots in the hash table
Collisions

More reasonable example:

- Store about 1000 records with keys in range 0 to 16,383.
- Impractical to keep a hash table with 16,384 slots.
- We must devise a hash function to map the key range to a smaller table.

Given: hash function $h$ and keys $k_1$ and $k_2$. $\beta$ is a slot in the hash table.
If $h(k_1) = \beta = h(k_2)$, then $k_1$ and $k_2$ have a collision at $\beta$ under $h$.

**Perfect Hashing**: hash function devised so that there are no collisions

Often impractical, sometimes expensive but worthwhile

It works when the set is **very stable** (e.g., a database on a CD)
Collisions (cont)

Search for the record with key $K$:
1. Compute the table location $h(K)$.
2. Starting with slot $h(K)$, locate the record containing key $K$ using (if necessary) a collision resolution policy.

Collisions are inevitable in most applications.
- Example: 23 people are likely to share a birthday ($p = \frac{1}{2}$).

Example: store 200 students, in a table $T$ with 365 records, using hash function $h$: birthday
A hash function MUST return a value within the hash table range.

To be practical, a hash function SHOULD evenly distribute the records stored among the hash table slots.

Ideally, the hash function should distribute records with equal probability to all hash table slots. In practice, success depends on the distribution of the actual records stored.

If we know nothing about the incoming key distribution, evenly distribute the key range over the hash table slots.

If we have knowledge of the incoming distribution, use a distribution-dependant hash function.
Hash Functions (cont.)

Reasons why data values are poorly distributed

- Natural distributions are exponential (e.g., populations of cities)
- collected (e.g., measured) values are often somehow skewed (e.g., rounding when measuring)
- coding and alphabets introduce uneven distributions (e.g., words in natural language have first letter poorly distributed)
Example Hash Functions

static int h(int x) {
    return(x % 16);
}

This function is entirely dependant on the lower 4 bits of the key, likely to be poorly distributed.

**Mid-square method**: square the key value, take the middle $r$ bits from the result for a hash table of $2^r$ slots.

[Works well because all bits contribute to the result.]

Sum the ASCII values of the letters and take results modulo $M$.

static int h(String x, int M) {
    int i, sum;
    for (sum=0, i=0; i<x.length(); i++)
        sum += (int)x.charAt(i);
    return(sum % M);
}

[Only good if the sum is large compared to the size of the table $M$.]

[This is an example of a folding method]

[NB: order of characters in the string is immaterial]
Open Hashing

What to do when collisions occur? **Open hashing** treats each hash table slot as a bin.

Open: collisions result in storing values outside the table

Each slot is the head of a linked list
Open Hashing Performance

Factors influencing performance
- how records are ordered in a slot's list (e.g., by key value or frequency of access)
- ration $N/M$ (records/slots)
- distribution of record key values

NB: Open Hash table must be kept in main memory (storing on disk would defeat the purpose of hashing)
Bucket Hashing

Divide the hash table slots into buckets.
- Example: 8 slots/bucket.

Include an overflow bucket.

Records hash to the first slot of the bucket, and fill bucket. Go to overflow if necessary.

When searching, first check the proper bucket. Then check the overflow.
Closed Hashing

Closed hashing stores all records directly in the hash table.

Each record $i$ has a **home position** $h(k_i)$.

If $i$ is to be inserted and another record already occupies $i$’s home position, then another slot must be found to store $i$.

The new slot is found by a **collision resolution policy**.

Search must follow the same policy to find records not in their home slots.
Collision Resolution

During insertion, the goal of collision resolution is to find a free slot in the table.

**Probe Sequence**: the series of slots visited during insert/search by following a collision resolution policy.

Let $\beta_0 = h(K)$. Let $(\beta_0, \beta_1, ...)$ be the series of slots making up the probe sequence.

```cpp
void hashInsert(Elem R) { // Insert R into hash table T
    int home; // Home position for R
    int pos = home = h(R.key()); // Initial pos on sequence
    for (int i=1; T[pos] != null; i++) {
        // Find next slot: p() is the probe function
        pos = (home + p(R.key()), i)) % M;
        Assert.notFalse(T[pos].key() != R.key(),
            "Duplicates not allowed");
    }
    T[pos] = R; // Insert R
}
```
Collision Resolution (cont.)

// p(K, i) probe function returns offset from home position
// for ith slot of probe sequence of K
ELEM hashSearch(int K) { // Search for record w/ key K
    int home; // Home position for K
    int pos = home = h(K); // Initial pos on sequence
    for (int i = 1; (T[pos] != null) &&
        T[pos].key() != K; i++)
        pos = (home + p(K, i)) % M; // Next pos on sequence
    if (T[pos] == null) return null; // K not in hash table
    else return T[pos]; // Found it}
Linear Probing

Use the probe function

\[
\text{int } p(\text{int } K, \text{ int } i) \{ \text{return } i; \}
\]

This is called **linear probing**.

Linear probing simply goes to the next slot in the table.

If the bottom is reached, wrap around to the top.

To avoid an infinite loop, one slot in the table must always be empty.
**Linear Probing Example**

Assuming hash function $h(x) = x \mod 11$

![Hash Table Example](image)

**Primary Clustering**: Records tend to cluster in the table under linear probing since the probabilities for which slot to use next are not the same.

[notation: prob($x$) is the probability that next element goes to position $x$]

[For (a): prob(3) = 4/11, prob(4) = 1/11, prob(5) = 1/11, prob(6) = 1/11, prob(10) = 4/11.]

[For (b): prob(3) = 8/11, prob(4,5,6) = 1/11 each.]

[small clusters tend to merge ⇒ long probe sequences]
Improved Linear Probing

Instead of going to the next slot, skip by some constant \( c \).

Warning: Pick \( M \) and \( c \) carefully.

The probe sequence SHOULD cycle through all slots of the table.

[If \( M = 10 \) with \( c = 2 \), then we effectively have created 2 hash tables (evens vs. odds).]

Pick \( c \) to be relatively prime to \( M \).

There is still some clustering.

- Example: \( c = 2 \). \( h(k_1) = 3 \). \( h(k_2) = 5 \).
- The probe sequences for \( k_1 \) and \( k_2 \) are linked together.
Pseudo Random Probing

The ideal probe function would select the next slot on the probe sequence at random.

An actual probe function cannot operate randomly. (Why?)

[Execution of random procedure cannot be duplicated when searching]

**Pseudo random probing:**

- Select a (random) permutation of the numbers from 1 to $M - 1$:

  $$r_1, r_2, \ldots, r_{M-1}$$

- All insertions and searches use the same permutation.

Example: Hash table of size $M = 101$

- $r_1 = 2, r_2 = 5, r_3 = 32$.
- $h(k_1) = 30, h(k_2) = 28$.
- Probe sequence for $k_1$ is: [30, 32, 35, 62]
- Probe sequence for $k_2$ is: [28, 30, 33, 60]

[The two probe sequences diverge immediately]
Quadratic Probing

Set the \( i \)'th value in the probe sequence as

\[
(h(K) + i^2) \mod M.
\]

Example: \( M = 101 \).

• \( h(k_1) = 30, \ h(k_2) = 29 \).
• Probe sequence for \( k_1 \) is:

\[
\begin{align*}
[30, 31, 34, 39] & = \\
[30, 30+1^2, 30 + 2^2, 30 + 3^2]
\end{align*}
\]

• Probe sequence for \( k_2 \) is:

\[
\begin{align*}
[29, 30, 33, 38] & = \\
[29, 29+1^2, 29 + 2^2, 29 + 3^2]
\end{align*}
\]

Problem: not all slots in the hash table are necessarily in the probe sequence
Double Hashing

Pseudo random probing eliminates primary clustering.

If two keys hash to same slot, they follow the same probe sequence. This is called **secondary clustering**.

To avoid secondary clustering, need a probe sequence to be a function of the original key value, not just the home position.

**Double hashing:**

\[ p(K, i) = i \times h_2(K) \text{ for } 0 \leq i \leq M - 1. \]

Be sure that all probe sequence constants are relatively prime to \( M \) \[\text{just like in improved linear probing}\].

Example: Hash table of size \( M = 101 \)

- \( h(k_1) = 30, h(k_2) = 28, h(k_3) = 30 \).
- \( h_2(k_1) = 2, h_2(k_2) = 5, h_2(k_3) = 5 \).
- Probe sequence for \( k_1 \) is: \([30, 32, 34, 36]\)
- Probe sequence for \( k_2 \) is: \([28, 33, 38, 43]\)
- Probe sequence for \( k_3 \) is: \([30, 35, 40, 45]\)
Analysis of Closed Hashing

The expected cost of hashing is a function of how full the table is.

The **load factor** is $\alpha = N/M$ where $N$ is the number of records currently in the table.

Expected # accesses (NB: accesses are due to collisions) vs $\alpha$
- solid lines: random probing
- dashed lines: linear probing
Deletion

1. Deleting a record must not hinder later searches.
2. We do not want to make positions in the hash table unusable because of deletion.

Both of these problems can be resolved by placing a special mark in place of the deleted record, called a **tombstone**.

A tombstone will not stop a search, but that slot can be used for future insertions.

Unfortunately, tombstones do add to the average path length.

Solutions:

1. Local reorganizations to try to shorten the average path length.
2. Periodically rehash the table (by order of most frequently accessed record).
Indexing

Goals:
- Store large files.
- Support multiple search keys.
- Support efficient insert, delete and range queries.

**Entry sequenced** file: Order records by time of insertion.  [Not practical as a database organization.]

Use sequential search.

**Index file**: Organized, stores pointers to actual records.  [Could be a tree or other data structure.]

**Primary key**: A unique identifier for records. May be inconvenient for search.

**Secondary key**: an alternate search key, often not unique for each record. Often used for search key.
Linear Indexing

**Linear Index**: an index file organized as a simple sequence of key/record pointer pairs where the key values are in sorted order.

Features:

- If the index is too large to fit in main memory, a second level index may be used.
- Linear indexing is good for searching variable length records.
- Linear indexing is poor for insert/delete.
**Tree Indexing**

Linear index is poor for insertion/deletion.

Tree index can efficiently support all desired operations (typical of a database):
- Insert/delete
- Multiple search keys
- Key range search

Storing a (BST) tree index on disk causes additional problems:
1. Tree *must* be balanced. [Minimize disk accesses.]
2. Each path from root to a leaf should cover few disk pages.

Use buffer pool to store recently accessed pages; exploit locality of reference

But only mitigates the problem
Tree indexing (cont.)

Rebalance a BST after insertion/deletion can require much rearranging

Example of insert(1)

```
(3)
  5
 /    \
/      /\n3      7
 /   /   /   \
2   4   6
```

```
(4)
  2
 /  \
1   3
 /   /   \
5   6
```

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2-3 Tree

A 2-3 Tree has the following shape properties:

1. A node contains one or two keys.
2. Every internal node has either two children (if it contains one key) or three children (if it contains two keys).
3. All leaves are at the same level in the tree, so the tree is always height balanced.

The 2-3 Tree also has search tree properties analogous to BST

1. values in left subtree < first node value
2. values in center subtree ≥ first node value
3. values in center subtree < second node value (if existing)
4. (if both existing) values in right subtree ≥ first node value
The advantage of the 2-3 Tree over the BST is that it can be updated at low cost.
- always insert at leaf node
- search position for key to be inserted
- if there is room (1 free slot) then finished
- otherwise must add a node (split operation)
- from 1 node with 2 keys get 2 nodes with 1 key and promote middle valued key
- recursively, insert promoted key into parent node
- if splitting repeated until root of the tree then its depth increases (but tree remains balanced)
2-3 Tree Insertion

[Insert 14]

[Insert 55. Always insert at leaf node.]
2-3 Tree Splitting

[Insert 19 into node 20-21 ⇒ split and promote 20 into node 23-30 ⇒ split and promote 23, this becomes new root, tree is 1 level deeper]

[NB: All operations are local to original search path.]
B-Trees

The B-Tree is a generalization of the 2-3 Tree.

The B-Tree is now the standard file organization for applications requiring insertion, deletion and key range searches.

1. B-Trees are always balanced.
2. B-Trees keep related records on a disk page, which takes advantage of locality of reference.
3. B-Trees guarantee that every node in the tree will be full at least to a certain minimum percentage. This improves space efficiency while reducing the typical number of disk fetches necessary during a search or update operation.
B-Trees (Continued)

A B-Tree of order $m$ has the following properties.

- The root is either a leaf or has at least two children.
- Each node, except for the root and the leaves, has between $\lceil m/2 \rceil$ and $m$ children.
- All leaves are at the same level in the tree, so the tree is always height balanced.

NB: A 2-3 Tree is a B-Tree of order 3

A B-Tree node is usually selected to match the size of a disk block.

A B-Tree node could have hundreds of children $\Rightarrow$ depth is $\approx \log_{100} n$.

A block implemented in a disk block

A pointer implemented by a disk block reference
B-Tree Example

Search in a B-Tree is a generalization of search in a 2-3 Tree.

1. Perform a binary search on the keys in the current node. If the search key is found, then return the record. If the current node is a leaf node and the key is not found, then report an unsuccessful search.

2. Otherwise, follow the proper branch and repeat the process.

A B-Tree of order 4

Example: search for record with key 47
B-Tree Insertion

Obvious extension of 2-3 Tree insertion

NB: split and promote process ensures all nodes are half full

Example: node with 4 keys + add one key

SPLIT ⇒ promote middle key + 2 nodes with 2 keys each
**B⁺-Trees**

The most commonly implemented form of the B-Tree is the B⁺-Tree.

Internal nodes of the B⁺-Tree do not store records – only key values to guide the search.

Leaf nodes store records or pointers to records.

A leaf node may store more or less records than an internal node stores keys.

Requirement: leaf nodes always half full.

Leaf nodes doubly linked in a list ⇒ can traverse it in any order ⇒ very good for range queries.

Search: similar to B-Tree search: must always go to the leaf (internal nodes do not store records)
B⁺-Tree Example: search

[Assume leaves can store 5 values, internal notes 3 (4 children).]

[Example: search key 33]
**B⁺-Tree Insertion**

Insertion similar to B-Tree insertion:
- find leaf that should contain inserted key
- if not full, insert and finish
- else split and promote a copy of least valued key of the newly formed right node
B+ - Tree Example: Insertion

[Note special rule for root: May have only two children.]

((b) Add 50.)

Add 45, 52, 47 (split), 18, 15, 31 (split), 21, 20.

[Add 30 (split).]
**B⁺-Tree Deletion**

- locate level $N$ containing key to be deleted
- if more than half full, remove and finish
- else (underflow) must restructure the tree
  - if possible get spare values from adjacent siblings ($\Rightarrow$ possibly keys in parent node must be updated)
  - if siblings cannot give values (they are only half full)
    - $N$ gives its values to them and is removed (possible because its siblings are only half full and $N$ is underflowing)
    - this can cause underflow in the parent node ($\Rightarrow$ propagate upwards, possibly eventually causing two children or root to merge and tree to lose one level)
**B⁺-Tree Example: Deletion**

[Simple delete – delete 18 from original example.]

[NB: do not need to delete 18 from internal node: it is a placeholder, can still be used to guide the search]

[Delete of 12 from original example: Borrow from sibling.]
B-Tree Space Analysis

B⁺-Tree nodes are always at least half full.

The B*-Tree splits two pages for three, and combines three pages into two. In this way, nodes are always 2/3 full.

Improves performance, makes implementation very complex

Tradeoff between space utilization and efficiency and complexity of implementation

Asymptotic cost of search, insertion and deletion of records from B-Trees, B⁺-Trees and B*-Trees is $\Theta(\log n)$. (The base of the log is the (average) branching factor of the tree.)

Ways to reduce the number of disk fetches:
- Keep the upper levels in main memory.
- Manage B⁺-Tree pages with a buffer pool.
B-Tree Space Analysis: Examples

Example: Consider a B$^+$-Tree of order 100 with leaf nodes containing 100 records.

1 level B$^+$-Tree: [Max: 100]

2 level B$^+$-Tree: [Min: 2 leaves of 50 for 100 records. Max: 100 leaves with 100 for 10,000 records.]

3 level B$^+$-Tree: [Min: $2 \times 50$ nodes of leaves for 5000 records. Max: $100^3 = 1,000,000$ records.]

4 level B$^+$-Tree: [Min: 250,000 records ($2 \times 50 \times 50 \times 50$). Max: 100 million records ($100 \times 100 \times 100 \times 100$).]