

Final Review

Materials

- The final exam is **comprehensive**
- Focus on the materials after midterm

Time and Location

CS4321 R01 Thursday 3pm 325

Topics For Midterm

Topics	Reading
Introduction	1.1-1.3
Induction and Loop invariants	1.4-1.7, GT 1.3
Elementary Algorithmics	Chapter 2
Asymptotic Notation	Chapter 3
Algorithm Analysis <ul style="list-style-type: none">- Analyzing control structures- Worst-case and Average-case- Amortized analysis	4.1-4.6
Solving Recurrences	4.7
Heap and Heap Sort	5.1-5.7
Binomial Heaps	5.8
Binary search tree, Splay Trees	GT 3.1, 3.4
Disjoint Set	5.9
Midterm	

Topics after midterm

Topics	Reading
Greedy Algorithms <ul style="list-style-type: none">- Coin Change- Minimum Spanning Tree- Dijkstra's algorithm- Knapsack- Scheduling	6
Divide-and-Conquer <ul style="list-style-type: none">- Mergesort and quicksort- Median- Strassen's algorithm- Closest pair	7

Topics After Midterm cont.

Topics	Reading
Dynamic Programming <ul style="list-style-type: none">- Binomial coefficient, coin change- 0-1 Knapsack- Floyd's algorithm- Matrix chain	Chapter 8
Exploring Graphs <ul style="list-style-type: none">- Graph Search- Topological sorting- Brand and Bound	Chapter 9
Network Flow	GT 8.1-8.3
Decision Trees and Low Bound Arguments	12.1-12.3
P and NP Problems	12.4-12.6, 13.2

Study Guide

- Study the reviews for Midterm
- Study the questions on quizzes and homework assignments
- Understand how the covered algorithms work

Greedy algorithms

- Know the paradigm
 - Template of a greedy algorithm
 - Be able to design and analyze a greedy algorithm
- Understand the following algorithms
 - Coin change
 - MST (Prim's algorithm and Kruskal's algorithm)
 - Dijkstra's algorithm (single source shortest path)
 - Knapsack
 - Scheduling for shortest total response time

General characteristics of greedy algorithms

```
makeChange(int n)
{
    C = a set of available coins;
    S = ∅; // chosen coins
    s = 0;
    while (C != ∅ && s != n) {
        x = a coin in C with the largest value
            such that s+x ≤ n;
        if no such x exists
            return "no solution found";
        else {
            C = C \ {x};
            S = S ∪ {x};
            s = s+x;
        }
    }
}
```

```
greedy(SET C)
{
    // C is set of candidates
    S = ∅; // S is a partial solution

    while (C != ∅) {
        x = select(C);
        if (feasible(S ∪ {x}))
            S = S ∪ {x};
        if (solution(S))
            return S;
    }
    return "no solutions";
}
```

objective() function is implicit

Kruskal's algorithm -- efficiency

$\Theta(a \log a)$ → sort A by increasing weight;
 $n = \# \text{nodes in } N$;
 $T = \emptyset$;
 make n initial sets, each contains a node in N;

 do { // for all sorted edges
 $e = \langle u, v \rangle$; // shortest edge not yet considered
 $uComponent = \text{find}(u)$;
 $vComponent = \text{find}(v)$;

 if ($uComponent \neq vComponent$) {
 \rightarrow merge($uComponent, vComponent$);
 $T = T \cup \{e\}$;
 }
 } while (!(T contains n-1 edges))
 return T;
 }

called at most a times each

called n-1 times

$\Theta(2a \alpha(2a, n))$

Prim's algorithm

```

Prim(int L[ ][ ])
{
    T = ∅;
    for (i=2; i<=n; i++) {
        nearest[i] = 1;
        mindist[i] = L[i,1];
    }

    for (i=2; i<n; i++) {
        min = ∞;
        for (j=2; j<=n; j++) {
            if (mindist[j] >= 0 && mindist < min) {
                min = mindist[j];
                k = j;
            }
        }
        T = T ∪ <nearest[k], k>;
        mindist[k] = -1;
        for (j=2; j<=n; j++) {
            if (L[j,k] < mindist[j]) {
                mindist[j] = L[j,k];
                nearest[j] = k;
            }
        }
    }
}
  
```

$\Theta(n)$ → for (i=2; i<=n; i++) {
 $\Theta(n^2)$ → for (j=2; j<=n; j++) {

Dijkstra's algorithm

C : candidate set
 S : partial solution set
 $L[i][j]$: weight of edge $\langle i, j \rangle$
 $D[i]$: length of the special path for the source to node i .
 $P[i]$: the previous node of i along its shortest path.

```

Dijkstra(Weight L[ ][ ])
{
    /* initialization */
    C = {i | 2 <= i <= n}; // S = {1}
    for (i=2; i<=n; i++) {
        D[i] = L[1,i];
        P[i] = 1;
    }

    for (i=1; i<=n-2; i++) { // repeat n-2 times
        v = some element of C minimizing D[v];
        C = C - {v}; // S = S ∪ {v}
        for (each w) {
            if (D[v] + L[v,w] < D[w]) {
                D[w] = D[v] + L[v,w];
                P[w] = v;
            }
        }
    }
}
  
```

Analysis of Dijkstra's algorithm: using heap

a : #edges
 n : #nodes

```

Dijkstra(Weight L[ ][ ])
{
    /* initialization */
    C = {i | 2 <= i <= n}; // S = {1}
    for (i=2; i<=n; i++) {
        D[i] = L[1,i];
        P[i] = 1;
    }

    buildHeap(D); // inverted heap

    for (i=1; i<=n-2; i++) { // repeat n-2 times
        v = findMin(D);
        deleteMin(v); // C = C - {v}; S = S ∪ {v}
        for (each w) {
            if (D[v] + L[v,w] < D[w]) {
                D[w] = D[v] + L[v,w];
                percolate(v);
                P[w] = v;
            }
        }
    }
}
  
```

$\Theta(n)$ → for (i=2; i<=n; i++) {
 $\Theta(n)$ → buildHeap(D); // inverted heap
 $\Theta(1)$ → v = findMin(D);
 $O(\log n)$ → deleteMin(v);
 $O(\log n)$ → percolate(v);
 At most a times if use adjacent list

Total: $O((a+n) \log n) = O(a \log n)$

The knapsack problem

- Given
 - n objects numbered from 1 to n. Object i has a positive weight w_i and a positive value v_i
 - a knapsack that can carry a weight not exceeding W
- Problem
 - Fill the knapsack in a way that maximize the value of the included objects, while respecting the capacity constraints
 - In this version, we assume that the objects can be broken into small pieces

A greedy algorithm

```
Knapsack(w[], v[], W)
{
  for (i=1; i<=n; i++)
    x[i] = 0;
  weight = 0;

  while (weight < W) {
    i = select the best remaining object;
    if (weight + w[i] < W)
      x[i] = 1;
    else
      x[i] = (W-weight)/w[i];
  }
  return x;
}
```

The key is
which object
to select

Scheduling

- Minimizing time in the system
 - Know the problem
 - Know the proof
 - Know the algorithm
- Scheduling with deadlines (not required)

Minimizing time in the system

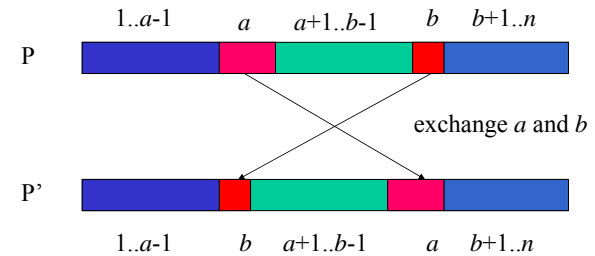
- Assume we submit n jobs into a system at the same time.
 - The service time of job i , t_i , know in advance
- Problem
 - Design an algorithm that minimizing the average response time
 - This is equivalent to
 - Minimizing the total response time
 - $T = \sum_{i=1}^n (\text{response time of job } j)$

A greedy algorithm

- Algorithm
 1. Sort the jobs by their service times
 2. Repeat
 1. Serve the job with minimal service time among the remaining jobs
- Analysis
 - Step 1: $O(n \log n)$
 - Step 2: $\Theta(n)$
 - Total: $O(n \log n)$

Optimality of the greedy scheduling algorithm

- Theorem 6.6.1. The greedy algorithm is optimal



Compares schedules P and P' , job a at P' leaves at the same time as job b in P . Jobs b and $a+1$ to $b-1$ in P' leaves earlier then the corresponding jobs in P .

Optimality of the greedy scheduling algorithm

$$\begin{aligned}
 T(P) &= s_1 + (s_1 + s_2) + \dots (s_1 + s_2 + \dots s_n) \\
 &= ns_1 + (n-1)s_2 + \dots + 1s_n \\
 &= \sum_{k=1}^n (n-k+1)s_k
 \end{aligned}$$

$$T(P) = (n-a+1)s_a + (n-b+1)s_b + \sum_{k=1, k \neq a, b}^n (n-k+1)s_k$$

$$T(P') = (n-a+1)s_b + (n-b+1)s_a + \sum_{k=1, k \neq a, b}^n (n-k+1)s_k$$

$$T(P) - T(P') = (n-a+1)(s_a - s_b) + (n-b+1)(s_b - s_a) = (b-a)(s_a - s_b) > 0$$

Divide and Conquer

- Given a problem, know how to design a D&C algorithm
- Know how to analyze a D&C algorithm
 - You need to remember the simple version of the Master Theorem.
- Know the following algorithms
 - Merge sort
 - Quick sort
 - Find median
 - Closest pair

Running-time analysis

- Assume that the l sub-instances have roughly the same size n/b for some constant b
- Let $g(n)$ be the time required by DC on instances of size n , excluding the times need for the recursive calls. We have
 - $t(n) = l \cdot t(n/b) + g(n)$
- If $g(n) \in \Theta(n^k)$ for an integer k , we have

$$t(n) \in \begin{cases} \Theta(n^k) & \text{if } l < b^k \\ \Theta(n^k \log n) & \text{if } l = b^k \\ \Theta(n^{\log_b l}) & \text{if } l > b^k \end{cases}$$

Merge sort

```
mergeSort(int T[n])
{
    if (n is sufficiently small)
        insertionSort(T);
    else {
        int U[n/2], V[(n+1)/2];
        copy T[1..n/2] to U[1..n/2];
        copy T[n/2+1..n] to V[1.., (n+1)/2];
        mergeSort(U[n/2]);
        mergeSort(V[(n+1)/2]);
        merge(U, V, T);
    }
}
```

Merge two sorted arrays

```
Merge(U[m], V[n], T[m+n])
// merge sorted arrays U and V into T
{
    u = 0; // cursor for U
    v = 0; // cursor for V
    U[m] = V[n] = +∞; // sentinels
    for (t=0; t<m+n; t++) { // t is cursor for T
        if (U[u] < V[v]) {
            T[t] = U[u];
            u++;
        } else {
            T[t] = V[v];
            v++;
        }
    }
}
```

What to do if we do not use the two sentinels?

Quick Sort

- Choose an element from the array to be sorted as a pivot
- Partition the array on either side of the pivot such that those no smaller than the pivot are to its right and those no greater are to its left
- Recursive calls on both sides

The algorithm

```
quickSort(T, i, j)
{
    l = pivot(T, i, j); // l is the position of the pivot at end
    if (i < l-1)
        quickSort(T, i, l-1);
    if (l+1 < j)
        quickSort(T, l+1, j);
}
```

Pivot I

```
int pivot(T, i, j)
{
    // choose T[i] as the pivot
    p = T[i];
    l = i; // left cursor
    r = j+1; // right cursor
    do {
        l++;
    } while (T[l] <= p and l < r)

    do {
        r--;
    } while (T[r] > p);
```

```
while (l < r) {
    swap(T[l], T[r]);
    do {
        l++;
    } while (T[l] <= p);
    do {
        r--;
    } while (T[r] > p)
}
swap(T[i], T[r]);
return r;
```

T[i] is at the bound after the algorithm

Analysis

- Worst case: the array is sorted, $\Omega(n^2)$
- Best case
 - $T(n) = 2T(n/2) + \Theta(n)$, $T(n) \in \Theta(n \log n)$
- Average case (not required)

Selection using pseudomedian

```
selection(T[1..n], s)
{
    l = 1; r = n;
    while (true) {
        x = pseudomedian(T[l..r]);
        p = pivot(T[l..r], x);
        if (s < p) r = p-1;
        else if (s > p) l = p+1;
        else return p;
    }
}
```

```
pseudomedian(T[1..n])
{
    if (n <= 5)
        return adhocmedian(T);
    z = ⌊n/5⌋;
    for (i=1; i <= z; i++)
        Z[i] = adhocmedian(T[5i-4..5i]);
    return selection(Z, ⌈z/2⌉);
}
```

We assume the elements are distinct.
You need to know the time complexity of this algorithm

Closest Pair

- Problem
 - Given n points on a two-dimension space, find the closest pair
- A simple algorithm
 - Calculate the distance for all possible pairs, find a smallest one
 - Total $\binom{n}{2}$ pairs
 - Cost: $\Theta(n^2)$
- A better algorithm
 - Divide-and-conquer

Algorithm

```
double closestPair(Points p)
{
    n = p.size();
    mergeSort(p); // by x-coordinate
    return recursiveClosestPair(p, 1, n)
}

double recursiveClosestPair(p, i, j)
{
    if (j-i < 3) {
        return adhocClosest(p, i, j);
        sort p[i..j] by y-coordinate;
    }

    k = (i+j)/2;
    deltaL = recursiveClosestPair(p, i, k);
    deltaR = recursiveClosestPair(p, k+1, j);
    delta = min(deltaL, deltaR);
    return findClosestInStrip(p, i, j, delta);
}
```

Note: $p[i..j]$ are sorted by x-coordinate before getting in recursiveClosestPair();
sorted by y-coordinate after it returns;

findClosestInStrip

$m = j-i+1$

```
findClosestInStrip(p, i, j, delta)
{
    k = (i+j)/2;    l = p[k].x;
    // p[i..k] sorted by y-coordinate
    // p[k+1..j] sorted by y-coordinate
    merge(p, i, k, j); // p[i..j] sorted by y-coordinate

    t = 0;
    for (k=i; k<=j; k++) {
        if (p[k].x > l - delta
            && p[k].x < l+delta) // in the strip
            v[++t] = p[k];
    }
    for (k=1; k<=t; k++) {
        for (s=k+1; s<=min(t, k+7); s++)
            delta = min(delta, dist(v[k], v[s]));
    }
    return delta;
}
```

$\Theta(m)$

$\Theta(m)$

Cost? $O(m)$

Total: $\Theta(m)$

Dynamic Programming

- Given a problem, know how to design a dynamic programming algorithm
- Know how to analyze a DP algorithm
- Know the following algorithms
 - Calculating Binomial Coefficient
 - Coin Change
 - 0-1 Knapsack
 - Floyd's algorithm
 - Matrix chain

Example: Binomial Coefficient

- We want to calculate $\binom{n}{k}$ which can be defined as follows.

$$\binom{n}{k} = \begin{cases} 1 & \text{if } k=0, n \\ \binom{n-1}{k-1} + \binom{n-1}{k} & \text{if } 0 < k < n \\ 0 & \text{otherwise} \end{cases}$$

A solution using dynamic programming

- We instead calculate bottom-up by filling the following table

n \ k	0	1	2	k-1	k
0	1					
1	1	1				
2	1	2	1			
...						
n-1					$C(n-1, k-1)$	$C(n-1, k)$
n					\rightarrow	$C(n, k)$

Cost: time $\Theta(nk)$ and space $\Theta(k)$

Make change

- The principle of optimality
 - In an optimal sequence of decisions or choices, each subsequences must be also optimal
- Making a change
 - $c[i, j]$ is an optimal solution to for making a change of j using the first i denominations
 - If at lease one coin of denomination i is used, then $c[i, j] = c[i, j - d_i] + 1$
 - Otherwise, $c[i, j] = c[i-1, j]$
 - We obtain $c[i, j] = \min(c[i-1, j], c[i, j - d_i] + 1)$

Example

- Pay 8 units with coins worth 1, 4, and 6 units.
- Construct the following table top-down, left to right.

Amount	0	1	2	3	4	5	6	7	8
d1=1	0	1	2	3	4	5	6	7	8
d2=4	0	1	2	3	1	2	3	4	2
d3=6	0	1	2	3	1	2	1	2	2

$$\begin{aligned} c[3, 8] &= \min(c[2, 8], c[3, 8-6]+1) = \min(2, 3) = 2 \\ c[2, 8] &= \min(c[1, 8], c[2, 4]+1) = \min(8, 2) = 2 \\ c[2, 4] &= \min(c[1, 4], c[2, 0]+1) = \min(4, 1) = 1 \end{aligned}$$

The optimal solution 2. The greedy algorithm returns 3.

0-1 Knapsack

- n objects 1, 2, ..., n. Object i has weight w_i and value v_i
- The knapsack can carry a weight not exceeding W.
- Cannot split an object
- Maximize the total value

$$\bullet \text{ Maximize } \sum_{i=1}^n x_i v_i \quad \text{subject to} \quad \sum_{i=1}^n x_i w_i \leq W,$$

where $v_i, w_i > 0$ and $x_i \in \{0,1\}$ for $1 \leq i \leq n$

By dynamic programming

- Set up a table $C[0..n, 0..W]$ with one row for each available object and one column for each weight from 0 to W. Specifically, $C[0, j] = 0$ for all j.
- $C[i, j]$ is the maximum value if the weight limit is j and only objects 1 to i are available
 - $C[i, j] = \max(C[i-1, j], C[i-1, j-w_i] + v_i)$;
- $C[n, W]$ will be the solution

Algorithm

```
Knapsack0-1(v, w, n, W)
{
  for (w = 0; w <= W; w++) {
    c[0, w] = 0;
  }
  for (i = 1; i <= n; i++) {
    c[i, 0] = 0
    for (w = 1; w <= W; w++) {
      if (w[i] < w) {
        if (c[i-1, w-w[i]] + v[i] > c[i-1, w])
          c[i, w] = c[i-1, w-w[i]] + v[i];
        else c[i, w] = c[i-1, w]
      } else c[i, w] = c[i-1, w]
    } // for w
  } for i
}
```

The run time performance of this algorithm is $\Theta(nW)$

Finding the objects

```
i = n;
k = W;
while (i > 0 && k > 0) {
  if (C[i, k] < C[i-1, k]) {
    mark the i-th object as in knapsack;
    k = k - w[i];
    i = i - 1;
  } else
    i = i - 1;
}
```

Cost: $O(n+W)$

Floyd's algorithm

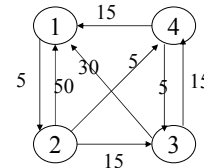
```
Floyd(int L[1..n][1..n])
{
    D = L;
    P = 0;
    for (k=1; k<=n; k++)
        for (i=1; i<=n; i++)
            for (j=1; j<=n; j++)
                if (D[i][k]+D[k][j] < D[i][j]) {
                    D[i][j] = D[i][k]+D[k][j];
                    P[i][j] = k;
                }
}
```

Track the
shortest path

Cost: $\Theta(n^3)$

Why can we use just one array D?

Example



$$D_0 = L = \begin{pmatrix} 0 & 5 & \infty & \infty \\ 50 & 0 & 15 & 5 \\ 30 & \infty & 0 & 15 \\ 15 & \infty & 5 & 0 \end{pmatrix}$$

$$D_1 = \begin{pmatrix} 0 & 5 & \infty & \infty \\ 50 & 0 & 15 & 5 \\ 30 & \underline{35} & 0 & 15 \\ 15 & \underline{20} & 5 & 0 \end{pmatrix}$$

$$D_2 = \begin{pmatrix} 0 & 5 & \underline{20} & \underline{10} \\ 50 & 0 & 15 & 5 \\ 30 & 35 & 0 & 15 \\ 15 & 20 & 5 & 0 \end{pmatrix}$$

$$D_3 = \begin{pmatrix} 0 & 5 & 20 & 10 \\ \underline{45} & 0 & 15 & 5 \\ 30 & 35 & 0 & 15 \\ 15 & 20 & 5 & 0 \end{pmatrix}$$

$$D_4 = \begin{pmatrix} 0 & 5 & \underline{15} & 10 \\ \underline{20} & 0 & \underline{10} & 5 \\ 30 & 35 & 0 & 15 \\ 15 & 20 & 5 & 0 \end{pmatrix}$$

$$P = \begin{pmatrix} 0 & 0 & 4 & 2 \\ 4 & 0 & 4 & 0 \\ 0 & 1 & 0 & 0 \\ 0 & 1 & 0 & 0 \end{pmatrix}$$

Find the shortest path

```
// Print shortest path between
nodes i and j
printShortestPath(int i, int j)
{
    print i;
    printIntermediateNodes(i, j);
    print j;
}
```

```
// Print intermediate nodes in the shortest
path between nodes i and j
printIntermediateNodes(int i, int j)
{
    int k = p[i][j];
    if (k == 0)
        return;
    else {
        printIntermediateNodes(i, k);
        print k;
        printIntermediateNodes(k, j);
    }
}
```

Find the shortest path between nodes 1 and 3 for the example.

Chained Matrix Multiplication

- Problem
 - Find an optimal *parenthesization* of chained matrix multiplication
- Chained matrix multiplication
 - $M = M_1 M_2 \dots M_n$
- Parenthesization
 - A product of matrices is fully parenthesized if it is either a single matrix or the product of two fully parenthesized matrix products, surrounded by parentheses

A recursive equation

- Let $m[i,j]$ be the minimum number of scalar multiplications needed to compute $M_{i..j}$
 - The cost for is $M_{1..n}$ is $m[1,n]$
- Assume that the optimal parenthesization splits the product $M_i M_{i+1} \dots M_j$ between M_k and M_{k+1}
 - Based on the principle of the optimality

$$m[i,j] = \begin{cases} 0, & i = j \\ \min_{i \leq k < j} (m[i,k] + m[k+1,j] + d_{i-1} d_k d_j), & i < j \end{cases}$$

An implementation using dynamic programming

```

matrixChain()
{
    for (i=1; i<=n; i++) // sequence of length 1
        m[i][i]=0;

    for (s=2; s<=n; s++) // s is the length of the sequence
        for (i=1; i<=n-s+1; i++) {
            j = i+s-1;
            m[i][j] = ∞;
            for (k=i; k<=j-1; k++)
                tmp = m[i][k] + m[k+1][j] + d[i-1]d[k]d[j]
                if (tmp < m[i][j]) {
                    m[i][j] = tmp;
                    p[i][j] = k; // split point
                }
            }
        }
}
    
```

Cost:

$$\begin{aligned}
 & n + \sum_{s=2}^n \sum_{i=1}^{n-s+1} (s-1) \\
 &= n + \sum_{s=2}^n (n-s+1)(s-1) \\
 &= n + n \sum_{l=1}^{n-1} l - \sum_{l=1}^{n-1} l^2 \\
 &= n + \frac{n^2(n-1)}{2} - \frac{n(n-1)(2n-1)}{6} \\
 &= (n^3 + 5n)/6 \\
 &\in \Theta(n^3)
 \end{aligned}$$

Example

M1	13×5
M2	5×89
M3	89×3
M4	3×34

i \ j	1	2	3	4
1	0	5785	1530	2856
2		0	1335	1845
3			0	9078
4				0

$m[1][3] = \min(m[1][1] + m[2][3] + 13 \cdot 5 \cdot 3, m[1][2] + m[3][3] + 13 \cdot 89 \cdot 3) = \min(1530, 9256) = 1530$
 $m[2][4] = \min(m[2][2] + m[3][4] + 5 \cdot 89 \cdot 34, m[2][3] + m[4][4] + 5 \cdot 3 \cdot 34) = \min(24208, 1845) = 1845$
 $m[1][4] = \min(m[1][1] + m[2][4] + 13 \cdot 5 \cdot 34, m[1][2] + m[3][4] + 13 \cdot 89 \cdot 34, m[1][3] + m[4][4] + 13 \cdot 3 \cdot 34)$
 $= \min(4055, 54201, 2856) = 2856$

Graph Traversal

- Know the depth search, breadth search, and topological sort algorithms
- Know how to use graph search to solve problems
- Topics
 - Tree traversal
 - Preconditioning
 - find ancestor in a tree
 - Graph search
 - Breadth first
 - Depth first
 - Articulation points (not required)
 - Topological sort

Graph search: some concepts

- To keep track of progress, graph search colors each node *white*, *gray*, or *black*
 - All nodes start with *white*
 - A node is *discovered* at the first time it is encountered during the search, at which time it becomes *non-white*
 - Different search distinguishes itself by a different way to *blacken* or *gray* nodes

Breadth-first search

- Given a graph $G = \langle N, E \rangle$, and a source node, s , start breadth-first search from s .
- Expands the frontier between discovered and undiscovered nodes uniformly across the breadth of the frontier
 - Discovers all nodes at distance k from s before discovering any nodes at distance $k+1$.
- Coloring: if $(u, v) \in E$ and vertex u is black, then node v is either black or gray
 - Black node: discovered and the node itself is finished
 - Gray node: discovered but not finished

Breadth-first search algorithm

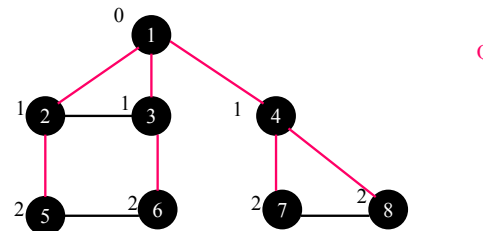
```
BFS(G,s)
{
  for each node  $u \in N - \{s\}$  {
    color[u] = WHITE;
    d[u] =  $\infty$ ;
     $\pi[u]$  = null;
  }

  color[s] = GRAY;
  d[s] = 0;
  enqueue(Q, s);

  while (!empty(Q)) {
    u = dequeue(Q);
    for each v adjacent to u {
      if (color[v] == WHITE) {
        color[v] = GRAY;
        d[v] = d[u] + 1;
         $\pi[v]$  = u;
        enqueue(Q, v);
      }
    }
    color[u] = BLACK;
  }
}
```

$d[\cdot]$: tracks shortest distance, assuming each edge's weight is 1
 $\pi[\cdot]$: tracks the parent-child relationship in the breadth-first tree

Breadth-first Example



Depth-first search

- Search deeper in the graph whenever possible
 - Edges are explored out of the most recently discovered node v that still has undiscovered edges leaving it
 - When all of v 's edges have been explored, the search “backtracks” to explore edges leaving the vertex from which v was discovered
 - This process finishes until all nodes reachable from the original source are discovered
 - Select one undiscovered node as the new source and continue the process

Depth-first search coloring and time stamps

- Coloring
 - Each nodes is initially *white*
 - A node is *grayed* if it is discovered during the search and *blackened* if it is finished, that is, when its adjacency list has been examined completely
- Timestamps
 - Each node v has two timestamps
 - $d[v]$ records when v is discovered (grayed)
 - $f[v]$ records when v is finished (blackened)

Depth-first search algorithm

```
DFS(G)
{
  for each node  $u \in N$  {
    color[u] = WHITE;
     $\pi[u]$  = null;
  }

  time = 0;

  for each node  $u \in N$  {
    if (color[u] == WHITE)
      DFS-Visit(u);
  }
}
```

```
DFS-Visit(u)
{
  color[u] = GRAY;
  d[u] = ++time;

  for each  $v$  adjacent to  $u$  {
    if (color[v] == WHITE) {
       $\pi[v]$  =  $u$ ;
      DFS-Visit(v);
    }
  }

  color[u] = BLACK;
  f[u] = ++time;
}
```

Classification of graph edges

- After depth-first search of a directed graph, we can classify the graph edges into four categories
 - **Tree edge**
 - An edge in the search tree
 - **Back edge**
 - An edge (u, v) not in search tree and v is an ancestor of u
 - Indicates a loop
 - **Forward edge**
 - An edge (u, v) not in search tree and u is an ancestor of v
 - **Cross edge**
 - An edge (u, v) not in search tree and v is neither an ancestor nor a descendant of u

Example: depth-first search directed graph

The graph consists of 8 nodes labeled 1 through 8. The edges are as follows:

- Tree edges (pink):** 1 → 2 (5), 1 → 3 (4), 1 → 4 (6), 4 → 8 (7), 8 → 7 (9).
- Forward edges (green):** 1 → 8 (12).
- Back edges (blue):** 3 → 2 (3), 6 → 5 (14), 7 → 4 (8).
- Cross edges (black):** 2 → 5 (13), 5 → 6 (16), 6 → 3 (15).

Legend:

- Tree edge: pink arrow
- Forward edge: green arrow
- Back edge: blue arrow
- Cross edge: black arrow

Topological Sort

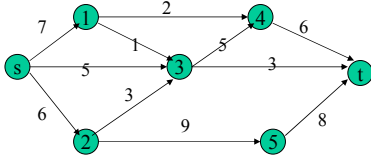
- Given an acyclic directed graph, topological sort finds a topological ordering of the nodes such that if there exists an edge (u, v) , then node u precedes node v in the ordering list.
- The finished time numbering gives us a reverse topological ordering
 - A node is finished after all the nodes it reaches have finished

Branch-and-Bound

- Not required

Maximum Flow Problem

- Given a weighted directed graph
 - Each edge is a pipe whose weight denotes its capacity: the maximum amount it can transport
 - Use $c(e)$ for the capacity of edge e
 - Given a source, s , and a sink, t , find the maximum amount (flow) can transfer from s to t

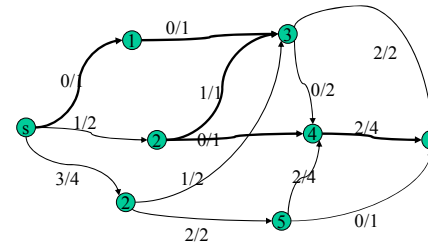


```
graph LR; s((s)) -- 7 --> 1((1)); s -- 5 --> 3((3)); s -- 6 --> 2((2)); 1 -- 2 --> 4((4)); 1 -- 1 --> 3; 2 -- 3 --> 3; 2 -- 9 --> 5((5)); 3 -- 5 --> 4; 3 -- 3 --> t((t)); 4 -- 6 --> t; 5 -- 8 --> t
```

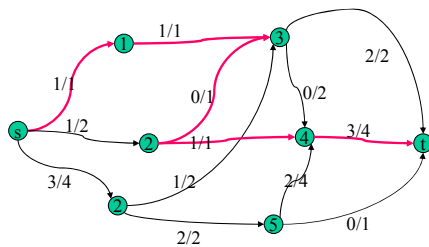
Concepts

- You need to know the follow concepts
 - Flow of a network
 - Capacity and flow of an edge
 - Cut
 - Capacity and flow of a cut
 - Residual capacity of an edge and a path
 - Augmenting Path
 - Residual capacity

Example of the Ford-Fulkerson Algorithm



Example of the Ford-Fulkerson Algorithm



The Ford-Fulkerson Algorithm

```

maxFlowFordFulkerson(N)
// N=(G, c, s, t)
{
  for each edge e in N do {
    f(e) = 0;
    stop = false;
  }
  while (!stop) {
    traverse G starting at s to find an augmenting path for f;
    if an augmenting  $\pi$  path exists {
       $\Delta$  = minimum  $\Delta_f(e)$  along  $\pi$ ;
      for each edge e in  $\pi$  {
        if (e is an forward edge) f(e) +=  $\Delta$ ; else f(e) -=  $\Delta$ ;
      }
    } else
      stop = true;
  }
}
  
```


The Edmonds-Karp Algorithm

- Try to find a “good” augmenting path each time
 - Choose an augmenting path with the smallest number of edges
 - Can be implemented using BFS traversal

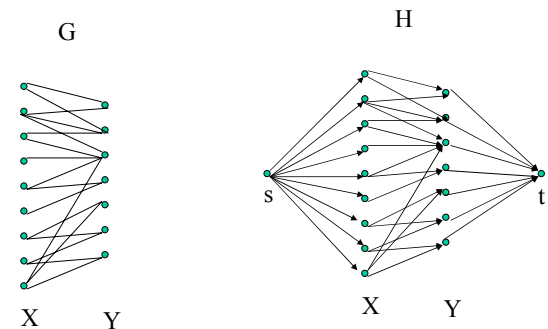
Maximum Bipartite Matching

- Bipartite graph
 - a graph with vertices partitioned into two sets X and Y, such that every edge has one endpoint in X and the other in Y
- Matching in a bipartite graph
 - A set of edges that has no end points in common
- Maximum bipartite matching
 - The matching with the greatest number of edges

Reduction to the Maximum Flow Problem

- Let G be a partite graph whose vertices are partitioned into sets X and Y. Create a flow network H as follows
 - Add each vertex of G into H plus a source vertex s and a sink vertex t.
 - Add edges of G into H and make each edge orient from an endpoint in X to an endpoint in Y
 - Insert a directed edge from s to each vertex in X
 - Insert a directed edge from each vertex in Y to t
 - Assign each edge in H a capacity of 1

An example of reduction



All edges with capacity 1

Reduction to the Maximum Flow

Problem

- Given the maximum flow f of H , define M as a set of edges such that $e \in M$ iff $f(e) = 1$
 - M is a matching
 - M is a maximum matching
- Reverse transformation: given a matching M in H , define a flow f
 - For each edge e of H that is also in G , $f(e) = 1$ if $e \in M$ and $f(e) = 0$ otherwise.
 - For each edge of H incident to s or t and v be the other endpoint, $f(e) = 1$ if v is an endpoint of some edge of M and $f(e) = 0$ otherwise

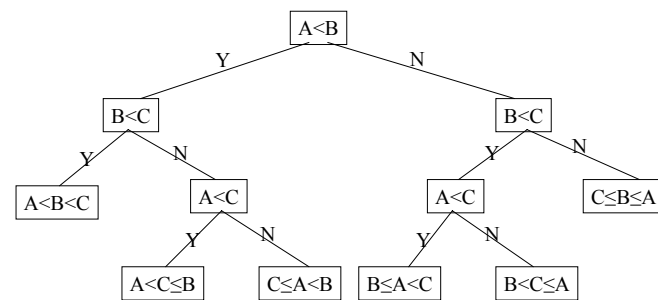
Complexity

- Arguing a lower bound
 - Information-theoretic arguments
- Given a comparison based problem
 - Know how to construct a decision tree
 - Know how to argue the lower bound using decision tree
- Adversary arguments
- Proof equivalence of complexity or compare complexity
 - Linear reductions
 - NP-completeness

Information-theoretic arguments

- Particularly applies to those problem involving comparisons
- Uses a decision tree to represent the working process of an algorithm on all possible data of a given size
 - A decision tree is a binary tree where
 - Each internal node contains a test on the data
 - Each leaf contains an output, called *verdict*
 - A *trip* through the tree starts from the root and recursively goes to the left subtree or the right subtree depending on whether the answer to the root is “yes” or “no”
 - The trip ends when it reaches a leaf (verdict)

A decision tree for sorting



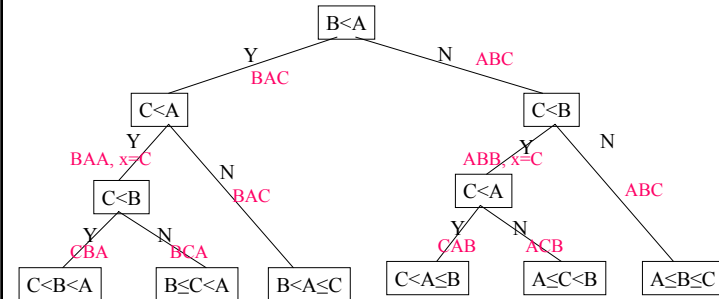
Note that the number of leaves = The number of possible outputs

Insertion Sort

```
void insertionSort(int A[], int n)
{ // array index starts from 1 here
  int i, j, x;

  for (i=2; i<=n; i++) {
    x = A[i];
    j = i-1;
    while (j>0 && x<A[j]) {
      A[j+1] = A[j];
      j--;
    }
    A[j+1] = x;
  }
}
```

A three-item insertion sort decision tree



Observations

- The number possible outputs = The number of leaves (verdicts)
- The worst-case time is the height of the tree
- The average time is the average depth of leaves assuming equal distribution

Theorem

- Any binary tree with k leaves has an average height of at least $\lg k$
- Any comparison-based algorithm takes a worst case time and average case time $\Omega(n \log n)$