# Summary of COMP523 Advanced Algorithm

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# **Symmetry Notation**

#### 1.1 Asymptotic Notation

Asymptotic notation is a way of describing the limiting behavior of a function when the argument tends towards a particular value or infinity. In computer science, asymptotic notation is frequently used to describe the running time or space usage of an algorithm.

- O-notation: f(n) = O(q(n)) if there exist constants c and  $n_0$  such that  $0 \le f(n) \le cq(n)$  for all  $n \ge n_0$ .
- $\Omega$ -notation:  $f(n) = \Omega(g(n))$  if there exist constants c and  $n_0$  such that  $0 \le cg(n) \le f(n)$  for all  $n \ge n_0$ .
- $\Theta$ -notation:  $f(n) = \Theta(g(n))$  if there exist constants  $c_1$ ,  $c_2$  and  $n_0$  such that  $0 \le c_1 g(n) \le f(n) \le c_2 g(n)$  for all  $n \ge n_0$ .
- o-notation: f(n) = o(g(n)) if for any constant c > 0, there exists a constant  $n_0$  such that  $0 \le f(n) < cg(n)$  for all  $n \ge n_0$ .
- $\omega$ -notation:  $f(n) = \omega(g(n))$  if for any constant c > 0, there exists a constant  $n_0$  such that  $0 \le cg(n) < f(n)$  for all  $n \ge n_0$ .

### 1.2 Comparing Functions

#### 1.2.1 Transitivity

- f(n) = O(g(n)) and g(n) = O(h(n)) implies f(n) = O(h(n)).
- $f(n) = \Omega(g(n))$  and  $g(n) = \Omega(h(n))$  implies  $f(n) = \Omega(h(n))$ .
- $f(n) = \Theta(g(n))$  and  $g(n) = \Theta(h(n))$  implies  $f(n) = \Theta(h(n))$ .

For example,  $n^2 = O(n^3)$  and  $n^3 = O(n^4)$  implies  $n^2 = O(n^4)$ .

#### 1.2.2 Reflexivity

- f(n) = O(f(n)).
- $f(n) = \Omega(f(n))$ .
- $f(n) = \Theta(f(n))$ .

For example,  $n^2 = O(n^2)$ .

#### 1.2.3 Symmetry

- f(n) = O(g(n)) implies g(n) = O(f(n)).
- $f(n) = \Omega(g(n))$  implies  $g(n) = \Omega(f(n))$ .
- $f(n) = \Theta(g(n))$  implies  $g(n) = \Theta(f(n))$ .
- f(n) = o(g(n)) implies  $g(n) = \omega(f(n))$ .
- $f(n) = \omega(g(n))$  implies g(n) = o(f(n)).

For example,  $n^2 = O(n^3)$  implies  $n^3 = \Omega(n^2)$ .

#### 1.2.4 Transpose Symmetry

- f(n) = O(g(n)) if and only if  $g(n) = \Omega(f(n))$ .
- $f(n) = \Theta(g(n))$  if and only if  $g(n) = \Theta(f(n))$ .
- f(n) = o(g(n)) if and only if  $g(n) = \omega(f(n))$ .
- $f(n) = \omega(g(n))$  if and only if g(n) = o(f(n)).

For example,  $n^2 = O(n^3)$  if and only if  $n^3 = \Omega(n^2)$ .

#### 1.2.5 sum and maximum

$$f_1(n) + f_2(n) + \dots + f_k(n) = \Theta(\max(f_1(n), f_2(n), \dots, f_k(n)))$$

where k is a constant positive integer.

Let  $f_j(n) = j$ , k = n, then

$$f_1(n) + f_2(n) + \dots + f_k(n) = n(n+1)/2 = \Theta(n^2)$$

#### 1.2.6 Running time hierarchy

- logarithmic:  $O(\log n)$
- linear: O(n)
- $n \log n$ :  $O(n \log n)$
- quadratic:  $O(n^2)$
- polynomial:  $O(n^k)$
- exponential:  $O(c^n)$
- constant: O(1)
- superconstant:  $\omega(1)$
- sublinear: o(n)
- superlinear:  $\omega(n)$
- superpolynomial:  $\omega(n^k)$
- subexponential:  $o(c^n)$

### 1.3 Expect of algorithms

**Correctness**: An algorithm is correct if it halts with the correct output for every input instance.

**Termination**: An algorithm is terminating if it halts for every input instance.

Efficiency: An algorithm is efficient if it halts with the correct output for every input instance and runs in polynomial

time.

# Recursion and Divide and Conquer techniques

### 2.1 Finding Majority in array

The pesudocode of the algorithm is shown in Algorithm 2.1.

#### Algorithm 1 Finding Majority in array

```
1: procedure MAJORITY(A)
        n \leftarrow \text{length of } A
        if n = 0 then
 3:
 4:
            return -1
        end if
        if n=1 then
            return A[1]
 8:
        if n1 and n is odd then
 9:
10:
        end if
11:
        Array B of size n/2
12:
        set j=0
13:
        for i = 1 to n/2 do
14:
15:
            if A[2i-1] = A[2i] then
                B[j] \leftarrow A[2i-1]
16:
                j \leftarrow j+1
17:
            end if
18:
19:
        end for
20:
        m \leftarrow \mathsf{MAJORITY}(B)
        count \leftarrow 0
21:
        for i=1 to n do
22:
            if A[i] = m then
                count \leftarrow count + 1
24:
            end if
25:
        end for
        if count > n/2 then
27:
            return m
28:
29:
        else
            \mathbf{return} - 1
        end if
32: end procedure
```

#### **Correctness:**

Lemma: If A has a majority element, then the majority element of A is also the majority element of B.

Base case: n = 1, the majority element is A[1].

Induction hypothesis: Assume that the lemma is true for n = k, we will prove that the lemma is true for n = k + 1.

Induction step: If A has a majority element, then the majority element of A is also the majority element of B.

Case 1 (A has a majority element m): Then by the lemma, it is also the majority element of B. Then m appears more than k/2 times in B. Then m appears more than (k+1)/2 times in A.

Case 2 (A has no majority element): Then B has no majority element. Then A has no majority element.

#### **Proof the lemma:**

proof by contradiction. Assume that A has a majority element m and B has a majority element m', but  $m \neq m'$ .

Let x be the numbers of occurrence of m in A.

Let y be the numbers of occurrence of m' in B.

Then 2y times from pairs that are represented in B by a value different from m', and x-2y times, since each occurrence of m in A that is not paired with another occurrence of m in A is paired with an occurrence of m' in B.

In total, this gives 2y + x - 2y = x occurrences of m in A, which is a contradiction.

#### **Running time:**

Recursive formula for the running time:

$$T(n) \le T(n/2) + cn$$

where c is a constant.

The solution to the recurrence is T(n) = O(n).

#### 2.2 Searching in logarithmic time

Searching faster with BinarySearch.

It is a particular case of the divide-and-conquer paradigm.

**Input**: A sorted array A of n elements and a value x.

**Output**: An index i such that A[i] = x or the special value -1 if x does not appear in A.

**Pseudocode** is shown in Algorithm 2.2.

#### Algorithm 2 BinarySearch

```
1: procedure BINARYSEARCH(x, i, j)
       if i = j then
2.
          if A[i] = x then
3:
4:
              return i
          else
5:
              return -1
6:
          end if
7:
       else
8:
          if x = A[|(i+j)/2|] then
9:
              return |(i+j)/2|
10:
          else if x < A[|(i+j)/2|] then
11:
              return BINARYSEARCH(x, i, |(i + j)/2|)
12:
13:
              return BINARYSEARCH(x, |(i+j)/2| + 1, j)
14:
          end if
15:
       end if
17: end procedure
```

#### **Running time**:

The number of comparisons performed by BinarySearch is:

$$T(n) \le T(n/2) + 4$$

Keep calculate:

$$\begin{split} T(n) &\leq T(n/2) + 4 \\ &\leq T(n/4) + 4 + 4 \\ &\leq T(n/8) + 4 + 4 + 4 \\ &\leq T(n/2^k) + 4k \\ &\leq T(n/2^{\log(n-1)}) + 4\log(n-1) \\ &= T(2) + 4(\log n - 1) \\ &\leq 4\log n - 4 \\ &= 4\log n \end{split}$$

proof  $T(n) \leq 4 \log n$ :

Base case: n = 1,  $T(1) = 0 \le 4 \log 1 = 0$ .

Induction hypothesis: Assume that the lemma is true for n = k, we will prove that the lemma is true for n = k + 1. Induction step:  $T(k+1) \le 4\log(k+1)$ .

$$T(k+1) \le T(k/2) + 4$$

$$\le 4 \log(k/2) + 4$$

$$= 4 \log k - 4 + 4$$

$$= 4 \log k$$

$$\le 4 \log(k+1)$$

#### Memory usage:

The memory usage of BinarySearch is:

$$M(n) = O(\log n)$$

#### Comparing BinarySearch and LinearSearch:

$$T_{ ext{BinarySearch}}(n) = O(\log n)$$
 
$$T_{ ext{LinearSearch}}(n) = O(n)$$
 
$$T_{ ext{BinarySearch}}(n) = O(\log n) < O(n) = T_{ ext{LinearSearch}}(n)$$
  $M_{ ext{BinarySearch}}(n) = O(\log n) < O(1) = M_{ ext{LinearSearch}}(n)$ 

### 2.3 Running time of Divide and Conquer algorithms

The Master Theorem:

Suppose that T(n) satisfies the recurrence:

$$T(n) \le aT(n/b) + cn^d$$

where  $a \ge 1$ , b > 1, c > 0 and  $d \ge 0$  are constants.

Then T(n) has the following asymptotic bounds:

$$T(n) = \begin{cases} O(n^d) & \text{if } d > \log_b a \\ O(n^d \log n) & \text{if } d = \log_b a \\ O(n^{\log_b a}) & \text{if } d < \log_b a \end{cases}$$

This theorem is useful for solving recurrences of the form:

$$T(n) = aT(n/b) + f(n)$$

where  $a \ge 1, b > 1$  and f(n) is an asymptotically positive function. **Example**:

```
\begin{split} T(n) &= 8T(n/2) + 100n^2\\ a &= 8, b = 2, f(n) = 100n^2, d = 2, \log_b a = \log_2 8 = 3.\\ d &= 2 < \log_b a = 3, \text{ so } T(n) = O(n^{\log_b a}) = O(n^3). \end{split}
```

#### 2.4 Finding piar of points closest to each other

**Input**: A set P of n points in the plane.

**Output**: The pair of points in P that are closest to each other. **Pseudocode** is shown in Algorithm 2.4. **Running time**:

```
Algorithm 3 ClosestPair
```

```
1: procedure CLOSESTPAIR(P_1, \ldots, P_n)

2: Construct P_x and P_y. P_x is sorted by x-coordinate, P_y is sorted by y-coordinate.

3: return CLOSESTPAIRREC(P_x, P_y)

4: end procedure
```

#### Algorithm 4 ClosestPairRec

```
1: procedure CLOSESTPAIRREC(P_x, P_y)
        if |P_x| = |P_y| \le 3 then
2:
             For each pair of points (P_i, P_j), compute d(P_i, P_j)
 3:
             return the pair of points with the smallest distance
4:
 5:
         end if
        Construct Q_x, Q_y, R_x and R_y.
 6:
        (l_1, l_2) = \text{CLOSESTPAIRREC}(Q_x, Q_y)
 7:
        (r_1, r_2) = \text{CLOSESTPAIRREC}(R_x, R_y)
 8:
        \delta = \min\{d(l_1, l_2), d(r_1, r_2)\}\
9:
10:
        x^* = the largest x-coordinate in Q_x
         L = (x, y) : x = x^*
11:
        S = \{ p \in P : p \in L \text{ and } p \text{ is within } \delta \text{ of } L \}
12:
        Construct S_v
13:
14:
        for p \in S do
             Let q be the point in S_v closest to p
15:
             if d(p,q) < \delta then
16:
                 \delta = d(p,q)
17:
                 (s_1, s_2) = (p, q)
18:
             end if
19:
        end for
20:
        if d(s_1, s_2) < \min\{d(l_1, l_2), d(r_1, r_2)\} then
21:
             return (s_1, s_2)
22:
        end if
23:
        if d(l_1, l_2) < d(r_1, r_2) then
24:
25:
             return (l_1, l_2)
26:
27:
             return (r_1, r_2)
        end if
28:
29: end procedure
```

```
T(n) \le 2T(n/2) + O(n \log n) = O(n \log n) Example:
```

# **Graph Algorithms**

#### 3.1 Graph Definitions

**Graph**: A graph G consists of a set V of vertices and a set E of edges, where each edge is associated with a pair of vertices.

**Directed Graph**: A directed graph G consists of a set V of vertices and a set E of directed edges, where each directed edge is associated with an ordered pair of vertices.

**Undirected Graph**: An undirected graph G consists of a set V of vertices and a set E of undirected edges, where each undirected edge is associated with an unordered pair of vertices.

**Neighbours of a vertex** v: Set of vertices that are connected to v by an edge.

**Degree of a vertex** v: number of neighbours of v, denoted by deg(v).

Path: A sequence of (non-repeating) nodes with consecutive nodes being connected by an edge.

length = node count - 1 = edge count.

**Distance between two nodes**: The number of edges in the shortest path between the two nodes.

Graph diameter: The maximum distance between any two nodes in the graph.

#### Lines, cycles, trees and cliques:

**Line**: A graph with n vertices and n-1 edges. **Cycle**: A graph with n vertices and n edges.

**cliques**: A graph with n vertices and n(n-1)/2 edges.

**Tree**: A graph with n vertices and n-1 edges.

#### **Graph representations:**

**Adjacency matrix**: A  $n \times n$  matrix A where  $A_{ij} = 1$  if there is an edge between i and j, and  $A_{ij} = 0$  otherwise. examples of adjacency matrices:

Given the following graph:



The adjacency matrix is:

$$\begin{bmatrix} 0 & 1 & 1 & 1 \\ 1 & 0 & 1 & 0 \\ 1 & 1 & 0 & 1 \\ 1 & 0 & 1 & 0 \end{bmatrix}$$

Adjacency matrix for directed graphs: A  $n \times n$  matrix A where  $A_{ij} = 1$  if there is an edge from i to j, and  $A_{ij} = 0$  otherwise.

examples of adjacency matrices for directed graphs: Given the following graph:



The adjacency matrix is:

$$\begin{bmatrix} 0 & 1 & 0 & 0 \\ 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 1 \\ 1 & 0 & 0 & 0 \end{bmatrix}$$

**Adjacency list**: A list of lists, where the ith list contains the neighbours of vertex i. Given the following graph:



The adjacency list is:

$$\begin{bmatrix} 1 & 2 & 3 \\ 0 & 2 \\ 0 & 1 & 3 \\ 0 & 2 \end{bmatrix}$$

**Adjacency list for directed graphs**: A list of lists, where the ith list contains the neighbours of vertex i. Given the following graph:



The adjacency list is:

$$\begin{bmatrix} 1 & 2 \\ 2 \\ 3 \\ 0 \end{bmatrix}$$

#### Adjacency matrix vs adjacency list:

Adjacency matrix	Adjacency list
O(1) to check if there is an edge between $i$ and $j$	O(min(deg(i), deg(j))) to check if there is an edge between $i$ and $j$
O(n) to find the neighbours of $i$	O(deg(j)) to find the neighbours of $i$
$O(n^2)$ space	O(n+m) space

### 3.2 Depth-first search

**Depth-first search**: A graph search algorithm that explores the neighbours of a vertex before exploring the neighbours of its neighbours.

example of depth-first search:



The depth-first search sequence is:

0, 1, 2, 3, 5, 4

Depth-first search algorithm:

#### Algorithm 5 Depth-first search algorithm

```
1: procedure DFS(G, v)
         for e \in V do
2:
             \quad \textbf{if} \ e \ \text{is unexplored then} \\
3:
                  u = \text{head of } e
 4:
                  if u is unexplored then
 5:
                      e is a tree edge
 6:
                      DFS(G, u)
 7:
                  else
 8:
                      e is a back edge
9:
                  end if
10:
             end if
11:
         end for
12:
13: end procedure
```

Running time of depth-first search: O(n+m)

#### 3.3 Breadth-first search

**Breadth-first search**: A graph search algorithm that explores the neighbours of a vertex before exploring the neighbours of its neighbours.

exaqmple of breadth-first search:



The breadth-first search sequence starting from vertex 0 is  $0,\,1,\,2,\,3,\,4,\,5.$ 

Breadth-first search algorithm:

#### Algorithm 6 Breadth-first search algorithm

```
1: procedure BFS(G, s)
         initial empty list L
 3:
         L \leftarrow s
 4:
         i \leftarrow 0
         while L[i] \neq \emptyset do
 5:
             L_{i+1} \leftarrow emptylist
 6:
             for v \in L[i] do
 7:
                 for edges (e) incident to v do
 8:
                      if e is unexplored then
 9:
                          w \leftarrow the other end of e
10:
                          if w is unexplored then
11:
                               label e as a tree edge
12:
                               add w to L_{i+1}
13:
                          else
14:
15:
                               label e as a cross edge
16:
                          end if
                      end if
17:
                 end for
18:
             end for
19:
20:
             i \leftarrow i + 1
         end while
21:
22: end procedure
```

Running time of breadth-first search: O(n+m)

### 3.4 Strong Connectivity

Directed graph: A graph where the edges have a direction.

Examples:



#### DFS and BFS on directed graphs:

Very similar to undirected graphs, except that we only consider edges that go out of a vertex.

Running time is O(n+m)

For example graph above the DFS sequence is 0, 1, 2, 3.

The BFS sequence is 0, 1, 2, 3.

#### 3.4.1 Connectivity

Weak connectivity: If we ignore the direction for all edges, there would be a pah from any vertex to any other vertex. Strong Connectivity: For every two nodes u and v, there is a path from u to v and a path from v to u.

#### 3.4.2 Mutual Reachability

Two nodes u and v are mutually reachable if there is a path from u to v and a path from v to u.

**Strong connectivity**: For every pair of nodes u and v, these two nodes are mutually reachable.

**Transitivity:** If u is mutually reachable with v and v is mutually reachable with w, then u is mutually reachable with w.

#### 3.4.3 Testing strong connectivity

```
Algorithm 7 Testing strong connectivity
```

```
1: procedure TESTSTRONGCONNECTIVITY(G)
       define G^R to be the graph with the same vertices as G but with all edges reversed
       Select a node s in G
3:
       BFS(G, s), BFS(G^R, s)
4:
       for each node v do
 5:
          if v is unexplored in either BFS then
              return False
 7:
           end if
8:
       end for
9:
10:
       return True
11: end procedure
```

#### 3.5 Testing bipartiteness

**Bipartite graph**: A graph G = (V, E) is bipartite if any only if the vertices can be partitioned into two sets  $V_1$  and  $V_2$  such that every edge has one end in  $V_1$  and the other end in  $V_2$ .

A Graph G = (V, E) is bipartite if and only if it has no odd cycles.(odd cycle: a cycle with odd number of edges)

#### **Testing bipartiteness:**

```
Given a graph G = (V, E), we want to test if G is bipartite. Given a graph G = (V, E), decide if it is 2-colourable. Given a graph G = (V, E), decide if it has an odd cycle. Colouring the nodes It is quite familiar with BFS:
```

#### Algorithm 8 Colouring the nodes

```
1: procedure Colouring(G, s)
        initial empty list L
 3:
        initial empty list C
 4:
        L \leftarrow s
        C[s] \leftarrow red
 5:
        i \leftarrow 0
 6:
        while L[i] \neq \emptyset do
 7:
 8:
             L_{i+1} \leftarrow emptylist
             for v \in L[i] do
 9:
                 for edges (e) incident to v do
10:
                     if e is unexplored then
11:
                          w \leftarrow the other end of e
12:
                          if w is unexplored then
13:
                              label e as a tree edge
14:
15:
                              add w to L_{i+1}
16:
                              if i+1 is odd then
                                  C[w] \leftarrow green
17:
                              else
18:
                                  C[w] \leftarrow red
19:
20:
                              end if
                          else
21:
                              label e as a cross edge
22:
                              if C[v] = C[w] then
23:
                                   return False
24:
                              end if
25:
                          end if
26:
27:
                     end if
                 end for
28:
             end for
29:
             i \leftarrow i + 1
30:
31:
        end while
        for e(v, w) \in G do
32:
             if C[v] = C[w] then
33:
                 return False
34:
35:
             end if
        end for
36:
        return True
37:
38: end procedure
```

#### Running time of colouring the nodes: O(n+m)

#### **Correctness of colouring the nodes:**

Proof by contradiction.

Suppose that G is not bipartite.

Then G has an odd cycle.

Suppose to the contrary that the algorithm return True.

That means that the algorithm did not detect the odd cycle.

### 3.6 DAGs and Topological Ordering

**DAG**: A directed acyclic graph (DAG) is a directed graph with no directed cycles. examples of DAGs:



**Topological ordering**: Given a graph G = (V, E), a topological ordering of G is an ordering of the nodes  $u_1, u_2, \ldots, u_n$  such that for every edge  $(u_i, u_j)$ , we have i < j.

Intutively, a topological ordering is an ordering of the nodes such that every edge goes from left to right. example of topological ordering based on given graph above:

#### Topological ordering implies DAG:

- If G has a topological ordering, then G is a DAG.
- Suppose by contradiction that G has a topological ordering  $u_1, u_2, \ldots, u_n$  but G also has a cycle C.
- Let  $u_i$  be the smallest element of C in the topological ordering.
- Let  $u_i$  be its predecessor in C.
- $u_i$  must appear before  $u_i$  in the topological ordering.
- This contradicts the fact that  $u_j$  is the smallest element of C in the topological ordering.

#### DAG implies topological ordering:

Proof by induction: Base case: If G has one or two nodes, then G has a topological ordering.

Induction steps: Assume that a DAG up to k nodes has a topological ordering(induction hypothesis). we will prove that a DAG with k+1 nodes has a topological ordering.

- By our lemma, there is at least one source node in G, and let u be the node.
- ullet Put u at the beginning of the topological ordering.
- ullet Consider the graph G', obtained by G by removing u and its incident edges.
- G' is a DAG with k nodes.
- It has a topological ordering  $u_1, u_2, \dots, u_k$  by the induction hypothesis.
- Append this ordering to u to get a topological ordering of G.

Here is the algorithm:

#### Algorithm 9 Topological Sorting

- 1: **procedure** TopologicalSorting(G)
- 2: find a source vertex u
- 3: set u as the first element of the topological ordering
- 4:  $G' \leftarrow G$  with u and its incident edges removed
- 5:  $L \leftarrow \text{TopologicalSorting}(G')$
- 6: append L to u
- 7: end procedure

Running time of the algorithm is  $O(n^2)$ 

#### **Modified Topological Sorting:**

Running time of the algorithm is O(n+m)

#### Algorithm 10 Modified Topological Sorting

```
1: procedure ModifiedTopologicalSorting(G)
        L \leftarrow emptylist
3:
        S \leftarrow \text{set of all source vertices}
       while S \neq \emptyset do
4:
 5:
           remove a vertex u from S
           append u to L
6:
           for each edge (u, v) do
7:
                remove edge (u, v) from G
8:
                if v is a source vertex then
9.
                    add v to S
10:
                end if
11:
            end for
12:
       end while
13:
       if G has edges then
14:
15:
            return G has a cycle
16:
            return L
17:
       end if
18:
19: end procedure
```

#### 3.7 Finding strongly connected components

**connected components**: A connected component of an undirected graph is subgraph of the graph where any two nodes are connected by a path.

**strongly connected components**: A strongly connected component of a directed graph is a subgraph of the graph where any two nodes are mutually reachable. (mutually reachable: there is a path from u to v and a path from v to u)

Finding strongly connected components:

Kosaraju's algorithm:

#### Algorithm 11 Kosaraju's algorithm

```
1: procedure KOSARAJU(G)
2: Initialise stack S
3: Select a arbitrary node s
4: DFS_tree=DFS(G, s)
5: S \leftarrow nodes in DFS_tree
6: G^R \leftarrow nodes in order of S
7: DFS(G^R, s)
8: return the nodes in the DFS tree
9: end procedure
```

Running time of Kosaraju's algorithm: O(n+m)Correctness of Kosaraju's algorithm:

- Define a meta-graph of G, called  $G^{SCC} = (V^{SCC}, E^{SCC})$ .
- Supposed that G has strongly connected components (SCCs)  $C_1, C_2, \ldots, C_k$ , for some k.
- $V^{SCC} = \{C_1, C_2, \dots, C_k\}$  contains some of the SCCs of G.
- There is an edge  $(C_i, C_j)$  in  $E^{SCC}$  if G contains a directed edge (x, y) such that  $x \in C_i$  and  $y \in C_j$ , crossing different components.

Examples:



The SCCs are  $\{0,1,2,3\}$  and  $\{4,5\}.$  The meta-graph is:



# **Greedy Algorithms**

#### The greedy approach:

- The goal is to find a global solution to a problem.
- The solution will be built up in small consecutive steps.
- For each step, we choose the best option available to us at that moment.

### 4.1 Interval Scheduling

#### **Interval Scheduling:**

A set of requests  $R = \{1, 2, \dots, n\}$ .

- Each request i has a start time  $s_i$  and a finish time  $f_i$ .
- Alternative view: every request is an interval  $[s_i, f_i]$ .

Two requests i and j are compatible if  $[s_i, f_i]$  and  $[s_j, f_j]$  do not overlap.

**Goal**: Find a maximum-size subset of compatible requests.

Example:

#### Interval scheduling.

- Job j starts at s<sub>j</sub> and finishes at f<sub>j</sub>.
- Two jobs compatible if they don't overlap.
- Goal: find maximum subset of mutually compatible jobs.



Figure 4.1: Interval Scheduling

#### **Interval Scheduling Algorithm:**

#### Algorithm 12 Interval Scheduling Algorithm

```
1: procedure IntervalScheduling([s_1, f_1], [s_2, f_2], \dots, [s_n, f_n])
        R is the set of requests
2:
        A \leftarrow \emptyset
3.
        while R \neq \emptyset do
4:
            select a request i in R with the smallest finishing time
 5:
6:
 7:
            remove all requests from R that are incompatible with i
        end while
8:
9:
        return A
10: end procedure
```

Running time of Interval Scheduling Algorithm:  $O(n \log n)$ 

**Correctness of Interval Scheduling Algorithm**: Since the algorithm always selects the request with the smallest finishing time, it is clear that the algorithm will always select a compatible request. **Arguing optimality**:

### 4.2 Minimum Spanning Trees

Consider a connected graph G = (V, E), such that each edge e = (v, w) of E, there is an associated cost  $c_e$ . **Goal**: Find a spanning tree T of E so that the graph G' = (V, T) has minimum cost. **Example**:



#### Greedy approach 1:

- Start with an empty set of edges T.
- ullet Repeat until T forms a spanning tree:
  - Select an edge e of minimum cost.
  - If  $T \cup \{e\}$  does not contain a cycle, then add e to T.

#### krukals algorithm:

#### Algorithm 13 Krukals algorithm

```
1: procedure KRUKALS(G)
         T \leftarrow \emptyset
2:
3:
         while T is not a spanning tree do
             select an edge e of minimum cost
4:
             if T \cup \{e\} does not contain a cycle then
5:
                  \operatorname{add} e \operatorname{to} T
 6:
             end if
 7:
         end while
 8.
9.
         return T
10: end procedure
```

#### Running time of Krukals algorithm: $O(m \log n)$ Greedy approach 2:

- Start with an empty set of edges T.
- Start with a node s.
  - Add an edge e = (s, v) of minimum cost to T.
- Repeat until T forms a spanning tree:

#### Prims algorithm:

#### Algorithm 14 Prims algorithm

```
1: procedure PRIMS(G)
2: T \leftarrow \emptyset
3: s \leftarrow an arbitrary node
4: while T is not a spanning tree do
5: add an edge e = (s, v) of minimum cost to T
6: s \leftarrow v
7: end while
8: return T
9: end procedure
```

# Running time of Prims algorithm: $O(m \log n)$ minimum spanning tree of example graph:

the minimum spanning tree sequence is d, a, c, b, e, f, h, g.

#### Greedy approach 3:

- Start with the full graph G = (V, E).
- Delete an edge from G
  - the edge of maximum cost
- Repeat until G forms a spanning tree:

#### Reverse-delete algorithm:

#### Algorithm 15 Reverse-delete algorithm

```
1: procedure REVERSEDELETE(G)
2: T \leftarrow G
3: while T is not a spanning tree do
4: delete an edge e of maximum cost from T
5: end while
6: return T
7: end procedure
```

For when two edges have the same cost, use distinct labels to distinguish them.

#### **Optimal with Priorty Queue:**

Add PQ to Prim's algorithm.

#### Algorithm 16 Optimal with Priorty Queue

```
1: procedure OPTIMAL(G)
        T \leftarrow \emptyset
 2:
        s \leftarrow \text{an arbitrary node}
 3:
 4:
        PQ \leftarrow empty priority queue
 5:
        for each node v do
            add v to PQ with key \infty
 6:
        end for
 7:
        decrease key of s to 0
 8:
        while PQ is not empty do
9:
            v \leftarrow \text{node} with minimum key in PQ
10:
            add an edge e = (s, v) of minimum cost to T
11:
12:
            for each edge e = (v, w) incident to v do
13:
                if w is in PQ then
14:
                    decrease key of w to c_e
15:
16:
                end if
            end for
17:
        end while
18:
19:
        return T
20: end procedure
```

Running time of Optimal with Priorty Queue:  $O(m \log n)$ 

#### 4.3 Clustering

- a collection of n objects
- they have different degrees of similarity
- we want to organise them into coherent groups
- there is a notion of distance between objects

#### **Definition**:

- Given a set U of n elements, a k-clustering of U is a partition of U into non-empty subsets  $C_1, C_2, \ldots, C_k$ .
- The spacing of a k-clustering is the minimum distance between any pair of points in different clusters.

**Goal**:Among all possible k-clusterings, find one with minimum spacing. **Example**:



#### Greedy approach:

- Pick two objects  $p_i$  and  $p_j$  with minimum distance  $d(p_i, p_j)$ .
- Connect them with an edge  $e = (p_i, p_j)$ .

- ullet Continue like this until we have k clusters.
- If the edge e under consideration connects two object  $p_i$  and  $p_j$  already in the same cluster, then discard e.

#### kruskals algorithm:

```
Algorithm 17 kruskals algorithm for clustering
```

```
Require: A graph G = (V, E)
Ensure: A minimum spanning tree of G with k clusters
 1: procedure KRUSKAL(G, k)
        T \leftarrow \emptyset
 2:
        C \leftarrow \{\{v\} \mid v \in V\}
                                                                                                                 ▶ Initial clusters
 3:
        Sort edges in E in increasing order of weight
 4:
        for \{u,v\} \in E do
 5:
            if C contains k clusters then
 6:
                break
 7:
            end if
 8:
            if clusters containing u and v are different in C then
 9:
                 T \leftarrow T \cup \{\{u,v\}\}
10:
                merge clusters containing u and v in C
11:
            end if
12:
13:
        end for
        \mathbf{return}\ T
14:
15: end procedure
```

For Given example, the result of divide them into 3 clusters is:

$$\{a,b,c,d\},\{e,f,h\},\{g\}$$

# **Dynamic Programming**

The paradigm of dynamic programming: Given a problem P, define a sequence of subproblems, with the following properties:

- The subproblems are ordered from the simplest to the largest
- ullet The largest problem is our original problem P
- The optimal solution of a subproblem can be structured from the optimal solutions of smaller subproblems.

Solve the subproblems from the smallest to the largest. When you solve a subproblem, store the solution and use it to solve larger subproblems.

#### 5.1 Weighted Interval Scheduling

- A set of requests  $R = \{1, 2, ..., n\}$ .
  - Request i has a start time  $s_i$  and a finish time  $f_i$ , and a value  $v_i$ .
  - Alternative view: every request is an interval  $[s_i, f_i]$  associated with a value  $v_i$ .
- Two requests i and j are compatible if  $[s_i, f_i]$  and  $[s_j, f_j]$  do not overlap.

Why it not working with greedy approach?