# COMP 251: Algorithms & Data Structures

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### 1 Overview of Graph Theory

#### 1.1 Definitions

A graph G=(V,E) is a set V of vertices (a.k.a. nodes) and a set E of edges (denoting vertex pairs). We set n=|V|, and m=|E|. A graph is said to be undirected when for any edge  $(u,v) \in E$  there exists an edge  $(v,u) \in E$  for some nodes u, and v. A graph is said to be directed if it is not undirected. In other words, the edge set of a directed graph consists of ordered pairs where the edge set of an undirected graph consists of unordered pairs.

A walk is a set of vertices  $\{v_0, v_1, \ldots, v_\ell\}$  such that  $(v_i, v_{i+1}) \in E$ ,  $\forall 0 \le i \le \ell$ . A walk where  $v_0 = v_\ell$  is said to be a *circuit* or a *closed walk*. A circuit where every edge in the graph is used exactly once is known as an *Eulerian circuit*. A *cycle* is a walk  $\{v_0, v_1, \ldots, v_\ell\}$  such that every vertex is distinct except  $v_0 = v_\ell$ . A cycle where every vertex of the graph is used exactly once is known as a *Hamiltonian cycle*. A walk where every vertex is distinct is said to be a *path*.

A graph is said to be *connected* if for each  $u, v \in V$  there exists a walk from u to v. A graph is said to be *disconnected* id it is not connected. Each connected subgraph of a graph is called a *component*. A connected graph therefore has exactly one component.

A connected component with no cycles is called a *tree*. A graph whose components are all trees is said to be a *forrest*. A tree is said to be *spanning* if it contains every vertex in the graph. A vertex in a tree with at most one neighbour is called a *leaf*.

A matching is a set of vertex-disjoint edges i.e. each edge is incident to at most one other edge in a matching. A matching is said to be *perfect* if every vertex is incident to exactly one edge in the matching.

A *clique* is a set of pairwise adjacent vertices. In *independent set* (a.k.a. a *stable set*) is set of pairwise non-adjacent vertices.

A bipartite graph is a graph such that the vertex set V can be partitioned as  $V = X \cup Y$  where each edge has one node in X and the other node in Y. Note that X and Y are necessarily independent sets.

### 1.2 Some Theorems for Undirected Graphs

**Theorem:** (Handshaking Lemma) Let G = (V, E) be an undirected graph, let  $\Gamma(v) := \{u : (u, v) \in E\}$  be the set of neighbours of a node v, and let the degree  $\deg(v)$  of a vertex v equal the cardinality of  $\Gamma(v)$ . Then there are an even number of vertices with odd degree.

**Proof:** First note that since we're double-counting the number of pairs where

(v,e) is an edge incident to v

$$2 \cdot |E| = \sum_{v \in V} \deg(v)$$

Since the degree of a vertex is either even or odd, we can partition V into a set of odd-degree vertices  $\mathcal{O}$ , and a set of even-degree vertices  $\mathcal{E}$ . This gives us

$$\sum_{v \in V} \deg(v) = \sum_{v \in \mathcal{O}} \deg(v) + \sum_{v \in \mathcal{E}} \deg(v)$$

which implies

$$\sum_{v \in \mathcal{O}} \deg(v) = 2 \cdot |E| - \sum_{v \in \mathcal{E}} \deg(v)$$

since both the  $2\cdot |E|$  term is even (obvious) and the  $\sum_{v\in\mathcal{E}}\deg(v)$  term is even

(sum of even numbers) then the  $\sum_{v \in \mathcal{O}} \deg(v)$  term must also be even.

**Theorem:** (Euler's Theorem) If G is an undirected graph then G contains an Eulerian circuit if and only if every vertex has even degree.

**Proof:** Easy proof by induction

**Lemma:** A tree T with  $n \geq 2$  vertices has at least one leaf vertex.

**Proof:** Trees are connected so there exists no vertices with degree 0 when  $n \geq 2$ . Suppose each vertex has degree of at least 2. Then consider the longest path  $P \subseteq T$ ,  $P = \{v_1, v_2, \dots, v_{\ell-1}, v_\ell\}$ . Since  $\deg(v_\ell) \geq 2$ ,  $\exists$  a neighbour (of  $v_\ell$ )  $x \in P$  with  $x \neq v_{\ell-1}$ . If  $x = v_{\ell+1}$  then P is not the longest path, a contradiction. Therefore, for P to be the longest path, x must be somewhere else in P, but this creates a cycle, another contradiction. Thus there must exist at least one node v such that  $0 < \deg(v) < 2$  – a leaf.

**Theorem:** A tree with n vertices has exactly n-1 edges.

**Proof:** Simple proof by induction.

Base case: A tree with one vertex trivially has 0 edges.

Induction Hypothesis: Assume any tree with n-1 vertices has n-2 edges. Inductive Step: Take a tree with  $n \geq 2$  vertices. By the previous lemma this tree contains a leaf vertex v. This implies that  $T \setminus \{v\}$  is a tree with n-1 vertices and by the induction hypothesis  $T \setminus \{v\}$  is a tree with n-2 edges, which implies that T is a tree with n-1 edges.

**Theorem:** (Hall's Theorem) Let  $G = (X \cup Y, E)$  with |X| = |Y| be a bipartite graph. G contains a perfect matching if and only if  $\forall B \subseteq X$ ,  $|\Gamma(B)| \ge |B|$  (Hall's condition).

**Proof:** Firstly, the  $(\Rightarrow)$  direction is fairly obvious. If  $B \subseteq X$  with  $\Gamma(B) < |B|$  then the graph can't have a perfect matching. The  $(\Leftarrow)$  direction is a bit trickier. Suppose Hall's condition is satisfied. Then, take the maximum cardinality

matching M is the graph. If M is perfect then we are done. Otherwise there must exist an unmached vertex  $b_0$ .

- Since Hall's condition holds, we have  $|\Gamma(\{b_0\})| \ge |\{b_0\}| = 1$  so  $b_0$  must have at least one neighbour  $s_0$ .
- Suppose  $s_0$  is matched in M to  $b_1$ .
- Since Hall's condition holds, we have  $|\Gamma(\{b_0, b_1\})| \ge |\{b_0, b_1\}| = 2$  so  $\{b_0, b_1\}$  must have at least one neighbour  $s_1 \ne s_0$ .
- Suppose  $s_1$  is matched in M to  $b_2$ .
- Since Hall's condition holds, we have  $|\Gamma(\{b_0, b_1, b_2\})| \ge |\{b_0, b_1, b_2\}| = 3$  so  $\{b_0, b_1, b_2\}$  must have at least one neighbour  $s_2 \notin \{s_0, s_1\}$ .

• ...

we repeat this argument as long as we can. Since the graph contains a finite number of vertices this process must terminate, but it can only terminate when we reach an unmatched node  $s_k$ . Using the edges we've formed in M we can create a path P from  $b_0$  to  $s_k$  that alternates between using non-matching edges and using matching edges. Swapping the matching edges with the non-matching edges gives us one more matching edge (as we have an odd number of edges.) This is still a valid matching as the internal nodes of P are still incident to exactly one matching edge. Also, the end nodes,  $b_0$  and  $s_k$  were previously unmatched but are now incident to exactly one edge in the new matching. Thus M isn't the maximum capacity matching – a contradiction.

### 1.3 """Data Structures""" for Representing Graphs

#### 1.3.1 Adjacency Matrices

For a graph, an adjacency matrix M is a matrix suxh that

- 1. There is a row for each vertex
- 2. There is a column for each vertex

3. The 
$$ij-th$$
 entry is defined as  $M_{ij} = \begin{cases} 1, (i,j) \in E \\ 0, (i,j) \notin E \end{cases}$ 

Note that in an undirected graph the matric is symmetric around the diagonal because  $(i, j) \sim (j, i)$ . Of course this is not necessarily true of directed graphs.

#### 1.3.2 Adjacency Lists

An adjacency list of an undirected graph is such that for each vertex v of V we store a list of its neighbours. For a directed graph we have two lists: one in which we store the in-neighbours of v and one in which we store the out-neighbours of v.

#### 1.3.3 Adjacency Matrices vs. Adjacency Lists

The main difference between the two is the amount of storage required to implement them.

- An adjacency matrix requires we store  $\Theta(n^2)$  numbers
- An adjacency list requires we store  $\Theta(m)$  numbers

In any graph  $m = O(n^2)$ . This means that for a sparse graph adjacency lists are highly favourable in terms of space complexity.

Verifying whether an edge exists, however, is much faster in an adjacency matrix – when using the array representation of a matrix it takes O(1) time, where verifying the existance of an edge takes  $O(\log n)$  time for an ordered adjacency list (using binary search), and O(n) time if the adjacency list is not ordered (using sequential search).

## 2 Divide & Conquer

A  $divide\ and\ conquer\ algorithm$  ideally breaks up a problem of size n into smaller sub-problems such that:

- ullet There are exactly a sub-problems
- Each sub-problem has a size of at most  $\frac{1}{b} \cdot n$
- Once solved, the solutions to the sub-problems must be combined in  $O(n^d)$  time to produce a solution to the original problem

Therefore the time-complexity of a divide and conquer algorithm satisfies a recurrence relation given by

$$T(n) = a \cdot T\left(\frac{n}{b}\right) + O(n^d)$$

#### 2.0.1 The Master Theorem

Lemma 1:  $\sum_{k=0}^{\ell} \tau^k = \frac{1-\tau^{\ell+1}}{1-\tau}$ , for any  $\tau \neq 1$ .

**Proof:** 

$$(1 - \tau) \sum_{k=0}^{\ell} \tau^k = \sum_{k=0}^{\ell} \tau^k - \sum_{k=1}^{\ell+1} \tau^k$$
$$= \tau^0 - \tau^{\ell+1}$$
$$= 1 - \tau^{\ell+1}$$
$$\sum_{k=0}^{\ell} \tau^k = \frac{1 - \tau^{\ell+1}}{1 - \tau}$$

**Lemma 2:**  $x^{\log_b y} = y^{\log_b x}$  for any base  $b \in \mathbb{R}$ .

**Proof:** From the power rule of logarithms we have

$$\log_b x \cdot \log_b y = \log_b \left( y^{\log_b x} \right)$$

similarily, we have

$$\log_b x \cdot \log_b y = \log_b \left( x^{\log_b y} \right)$$

therefore

$$\log_b\left(x^{\log_b y}\right) = \log_b\left(y^{\log_b x}\right)$$

thus

$$x^{\log_b y} = y^{\log_b x}$$

Theorem: (The Master Theorem) If a recurrence relation is of the form  $T(n) = a \cdot T\left(\frac{n}{b}\right) + O(n^d)$ , for constants a > 0, b > 1, and  $d \ge 0$ , then

$$T(n) = \begin{cases} O(n^d) & \text{, if } a < b^d \text{ [Case I]} \\ O(n^d \cdot \log n) & \text{, if } a = b^d \text{ [Case II]} \\ O(n^{\log_b a}) & \text{, if } a > b^d \text{ [Case III]} \end{cases}$$

**Proof:** By adding dummy numbers, we may assume that n is a power of b, i.e.  $n=b^{\ell}$ , for some  $\ell \in \mathbb{N}_0$ . Then

$$T(n) = n^d + a\left(\frac{n}{b}\right)^d + a^2\left(\frac{n}{b^2}\right)^d + \dots + a^\ell\left(\frac{n}{b^\ell}\right)^d$$

simlifying, we get

$$T(n) = n^d \left( 1 + \frac{a}{b^d} + \left( \frac{a}{b^d} \right)^2 + \dots + \left( \frac{a}{b^d} \right)^\ell \right)$$

we now have three cases:

Case I:  $\frac{a}{b^d} < 1$ Set  $\tau := \frac{a}{b^d}$  then we have

$$T(n) = n^d \sum_{k=0}^{\ell} \tau^k$$

applying lemma 1 we get

$$T(n) = n^d \left(\frac{1 - \tau^{\ell+1}}{1 - \tau}\right) \le n^d \left(\frac{1}{1 - \tau}\right)$$

but  $\frac{1}{1-\tau}$  is a constant so  $T(n) \le n^d$  and thus

$$T(n) = O(n^d)$$

Case II:  $\frac{a}{b^d} = 1$ Then we have that

$$T(n) = n^d (1 + 1 + 1^2 + \dots + 1^\ell) = n^d (\ell + 1)$$

but  $n = b^{\ell} \Rightarrow \ell = \log_b n$ , thus

$$T(n) = O(n^d \cdot \log n)$$

Case III:  $\frac{a}{b^d} > 1$ Set  $\tau := \frac{a}{b^d}$  then we have

$$T(n) = n^d \sum_{k=0}^{\ell} \tau^k$$

applying lemma 1 we get

$$T(n) = n^d \left(\frac{\tau^{\ell+1} - 1}{\tau - 1}\right) \le n^d \left(\frac{\tau^{\ell+1}}{\tau - 1}\right)$$

but since  $\tau - 1$  is a constant we get

$$\begin{split} T(n) &= n^d O(\tau^{\ell+1}) \\ &= O(n^d \tau^{\ell+1}) \\ &= O(n^d \tau^\ell) \\ &= O\left(\left(\frac{a}{b^d}\right)^\ell n^d\right) \\ &= O\left(\left(\frac{n}{b^\ell}\right)^d a^\ell\right) \end{split}$$

but  $n = b^{\ell}$  so

$$T(n) = O(a^\ell)$$

and  $\ell = \log_b a$  so

$$T(n) = O(a^{\log_b n})$$

and applying lemma 2 gives

$$T(n) = O(n^{\log_b a})$$

This completes the proof.

3 Greedy Algorithms

4 Dynamic Programming

## 5 Network Flows

## 6 Data Structures