CS 590 Reference Sheet

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Asymptotic Notation

- $f(n) = o(g(n)) : 0 \le f(n) < cg(n)$
- $f(n) = O(g(n)): 0 \le f(n) \le cg(n)$ (upper bound)
- $f(n) = \Theta(g(n)) \leftrightarrow \Omega(g(n)) \& O(g(n))$
- $f(n) = \Omega(g(n): 0 \le cg(n) \le f(n)$ (lower bound)
- $f(n) = \omega(g(n): 0 \le cg(n) < f(n)$

Useful Equations

$$\sum_{k=1}^{n} k = \frac{1}{2}n(n+1) = \Theta(n^2)$$

$$\sum_{k=0}^{n} p^k = \frac{1-p^{k+1}}{1-p}$$
Stirling: $n! \approx \sqrt{2\pi n} (\frac{n}{e})^n$

Amortized Analysis

$$\hat{c}_{i} = c_{i} + \Phi(D_{i}) - \Phi(D_{i-1})$$

$$\sum_{i=1}^{n} \hat{c}_{i} = \sum_{i=1}^{n} c_{i} + \Phi(D_{n}) - \Phi(D_{0})$$

Master Method For recurrences of the form T(n) = aT(n/b) + f(n), where a is the number of branches at each level, b is the reduction in size at each level, and f(n) is the initial cost:

Case 1: $f(n) = O(n^{\log_b a - \epsilon}) \to T(n) = \Theta(n \log_b a), \epsilon > 0$: f(n) is polynomially slower than $n^{\log_b a}$, weight increases toward leaves because b < a

Case 2: $f(n) = \Theta(n^{\log_b a}) \to T(n) = \Theta(n^{\log_b a} \log^{k+1} n), k \ge 0$: $f(n) \& n^{\log_b a}$ grow similarly, weight remains constant because b = a

Case 3: $f(n) = \Omega(n^{\log_b a + \epsilon}) \to T(n) = \Theta(f(n))$: weight decreases because b > a

Amortized Analysis The amortized cost (upper bound) \hat{c}_i for the potential function Φ and data D_i is defined as:

$$\hat{c}_{i} = c_{i} + \Phi(D_{i}) - \Phi(D_{i-1})$$

$$\sum_{i=1}^{n} \hat{c}_{i} = \sum_{i=1}^{n} c_{i} + \Phi(D_{n}) - \Phi(D_{0})$$

Hashing A hash function is *perfect* if it causes no collisions, but there is no perfect hash function if |S| > m.

Chaining: put a linked list at each of the slots in the table, cost is proportional to length of the lists. Good hash functions:

$$h(k) = kmod m$$

$$h(k) = (ak + b)mod m$$

$$h(k) = ((ax + b)mod p)mod m$$

A family H of hash functions from U to M is 2-universal if for all $x \neq y$, $Pr(h(x) = h(y)) \leq \frac{1}{m}$. BUT there are u^m functions from U to M, requiring $m \log u$ bits to choose/represent/store. We just pick one at random. For r operations, the expected total work is $r(1 + \frac{s}{m})$.

Good hashing: Let m be a prime number, $(x_0, ..., x_r)$ represent a key $x, \bar{a} = (a_0, ..., a_r)$.

$$h_{\bar{a}}(x) = (\sum_{i=0}^{r} a_i x_i) \mod m$$

$$H = \{h_{\bar{a}}(x) | a_i \in \{0, ..., m-1\}\}$$

Open-addressing: Load factor for number of keys n is $\alpha = \frac{n}{m}$. Search takes $\Theta(1+\alpha)$ expected time. Because elements are stored in the table itself, the load factor cannot exceed 1. The expected number of probes for inserting in a table (or an unsuccessful search) is $\frac{1}{1-\alpha}$. The expected number of probes for a successful search is $\frac{1}{\alpha}ln\frac{1}{1-\alpha}+\frac{1}{\alpha}$.

Complexity Π' is poly-time reducive to Π ($\Pi' \leq \Pi$) iff there is a poly-time function f() that maps inputs x' of Π' into inputs x of Π such that $\Pi'(x') = \Pi(f(x))$.

- 1. If $\Pi \in P$ and $\Pi' \leq \Pi$ then $\Pi' \in P$
- 2. If $\Pi \in NP$ and $\Pi' \leq \Pi$ then $\Pi' \in NP$
- 3. If $\Pi' \leq \Pi$ and $\Pi'' \leq \Pi'$, then $\Pi'' \leq \Pi$

Graph Algorithms A preordering is a list of the vertices in the order that they were first visited by the depth-first search algorithm.

Table 1: Uses of Graph Search

DFS connected components, topological sort

BFS shortest path