

i^{th} order statistics algorithm runs in $O(n)$. That is the selection of the i^{th} item can be done in $O(n)$ running time.

Dynamic programming: Prove both *Optimal substructure* – (if an optimal solution to the problem contains within it optimal solutions to subproblems) and *overlapping subproblems*– (subproblems must be solved several time throughout solving the original problem.)

Optimal Substructure:

1. Show that the problem consists of making a choice that leaves one or more subproblems to solve.
2. Suppose that for a given problem you are given the choice that leads to the optimal solution.
3. Determine which subproblems ensue and how to best characterize the resulting space of subproblems
4. Show that the solutions to the subproblems used within the optimal solution to the problem must themselves be optimal by cut-n-paste contradiction method.

Transforming a recursive solution into a bottom-up dynamic programming solution: Example

MCM: Matrix A_i has dimensions $p_{i-1} \times p_i$. $m[i,j]$ is the minimum number of scalar multiplications needed to compute the matrix $A_{i,j}$. So

$$m[i, j] = \begin{cases} 0 & \text{if } i = j \\ \min\{m[i, k] + m[k+1, j] + p_{i-1}p_kp_j\} & \text{if } i < j \end{cases}$$

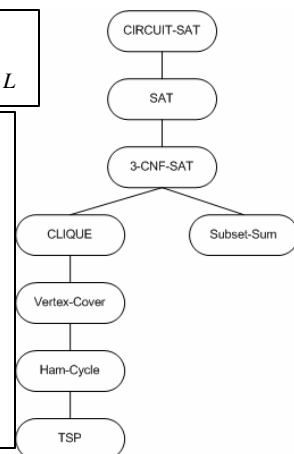
$$c[i, j] = \begin{cases} 0 & \text{if } i = 0 \text{ or } j = 0 \\ c[i-1, j-1] + 1 & \text{if } i, j > 0 \text{ and } x_i = y_i \\ \max\{c[i, j-1], c[i-1, j]\} & \text{if } i, j > 0 \text{ and } x_i \neq y_i \end{cases}$$

LCS: Let $X = \langle x_1, x_2, \dots, x_m \rangle$, we define the i^{th} prefix of X , for $i=0, \dots, m$ as $X_i = \langle x_1, \dots, x_i \rangle$. Let $Y = \langle y_1, \dots, y_n \rangle$ and $Z = \langle z_1, \dots, z_k \rangle$ be any LCS of X and Y . Then Optimal substructure argument

1. If $x_m = y_n$, then $z_k = x_m = y_n$ and $Z - z_k$ is an LCS of X_{m-1} and Y_{n-1}
2. If $x_m \neq y_n$, then $z_k \neq x_m \rightarrow Z$ is an LCS of X_{m-1} and Y .
3. If $x_m \neq y_n$, then $z_k \neq y_n \rightarrow Z$ is an LCS of X and Y_{n-1} .

Recursive solution at right.

Ogden's lemma: $s = uv^mxy^nz$; vy contains at least one distinguished position; wxy contains no more than n distinguished positions; x contains at least 1 (or 2) distinguished position(s) " $m = 0, uv^mxy^nz^l$ "



<p>Matrix -Chain -Order(p)</p> <pre>n = length[p] - 1 for i=n do {m[i,i]=0} for L=2 to n for i=1 to n-L+1 j=i+L-1 m[i,j]=inf for k=i to j-1 q=m[i,k] + m[k+1,j]+p[i-1]*p[k]*p[j] if q<m[i,j] m[i,j]=q s[i,j]=k return m and s</pre>	<p>LCS-Length</p> <pre>m=length[X]; n=length[Y]; for i=1 to m {c[i,0]=0}; for j=0 to n {c[0,j]=0} for i=1 to m for j=1 to n if x[i]=y[j] then c[i,j] = c[i-1, j-1]+1 else if c[i-1, j] >= c[i, j-1] then c[i,j]=c[i-1, j] else c[i,j]=c[i, j-1]+1 return c[m, n]</pre>	<p>Optimal-BST(p,q,n)</p> <pre>for i=1 to n+1 {e[i,i-1]=q[i-1]; w[i,i-1]=q[i-1]} for L=1 to n for i=1 to n-L+1 j=i+L-1 e[i,j]=inf w[i,j]=w[i,j-1]+p[j]+q[j] for k=1 to j t=e[i, k-1]+e[k+1, j]+w[i, j] if t<e[i, j] then e[i, j] = t root[i, j]=k return e and root</pre>
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Optimal BST: Let $K = \langle k_1, \dots, k_n \rangle$ be n distinct keys $\ni k_1 < \dots < k_n$. We also must have $n+1$ dummy keys representing all the values not in K . $\langle d_0, \dots, d_n \rangle \ni d_0$ represents all values less than k_1 , d_1 represents all values between k_1 and k_2 , and d_n represents all values greater than k_n . Each k_i has a probability p_i and each d_i has a probability q_i . Define: $w(i, j) =$

$$\sum_{i=1}^j p_i + \sum_{i=1}^j q_i \text{ and } e[i, j] = \begin{cases} q_{i-1} & \text{if } j = i - 1 \\ \min_{i \leq k \leq j} \{e[i, k-1] + e[k+1, j] + w(i, j)\} & \text{if } i \leq j \end{cases}$$

Greedy Algorithm: Prove *Optimal substructure*, and *Greedy choice property* – (that any optimal solution may or must contain the greedy choice.)

1. Cast the optimization problem as one in which we make a choice and are left with one subproblem to solve.
2. Prove that there is always an optimal solution to the original problem that makes the greedy choice. \rightarrow Greedy choice is safe.
3. Demonstrate that, ..., what is left is a subproblem with the property that if we combine an optimal solution with the greedy choice we get an optimal solution.

Fractional Knapsack, Super increasing coin problem... Huffman Encoding	Huffman(C) n = length of C //C linked list Q=C //min priority queue for i = 1 to n-1	... allocate a new node z left[z]=x=extractMin[Q] right[z]=y=extractMin[Q] f[z]=f[x]+f[y]	... Insert(Q,z) return extractMin(Q) //return the //root of the tree
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Flow Networks: A flow network $G = (V, E)$ is a directed graph in which each edge $(u, v) \in E$ has a nonnegative capacity $c(u, v) \geq 0$. Each vertex v is on a path from (s)Source to (t)Sink \rightarrow the graph is connected and $|E| \geq |V| - 1$. **Flow** is defined by $f(u, v) \leq c(u, v)$ with the following properties: **Capacity constraint** above or $\sum_{u \in V} f(u, v) = 0$ for $v \in V - \{s, t\}$,

Skew symmetry $\forall u, v \in V, f(u, v) = -f(v, u)$, **Flow conservation** $\forall u \in V = \{s, t\} \sum_{v \in V} f(u, v) = 0$. **Value of a flow** f denoted $|f| = \sum_{v \in V} v f(s, v)$. **Implicit Sum notation** $f(X, Y) = \sum_{x \in X} \sum_{y \in Y} f(x, y)$. **Lemma 26.1:** $\forall X, Y, Z \subseteq V$ with $X \cap Y = \emptyset, f(X \cup Y, Z) = f(X, Z) + f(Y, Z)$ and $f(Z, X \cup Y) = f(Z, X) + f(Z, Y)$. **Residual networks** are those nasty looking networks with back flow arrows instead of used/capacity numbers on the original edges. residual networks are defined as: $G_r = (V, E_r)$ where $E_r = \{(u, v) \in V \times V : c_r(u, v) > 0\}$. **Augmenting Paths:** p is a simple path from s to t in a residual network G_r . A flow is maximum if there does not exist any more augmenting paths. **Max-Flow, Min-Cut:** Given all cuts – see def. (S, T) where $s \in S, t \in T$, the cut with the minimum flow $f(S, T)$ is the maximum flow. Ford-Fulkerson = $O(E|f^*|)$ where f^* is the max flow.

Edmonds-Karp algorithm uses a depth first search to find the shortest path when adding augmenting paths to the residual network and it runs in $O(V^2E)$

Classification problems: **P** (polynomial time solvable) **NP** (non-deterministic polynomial), **NPC**, **NP-Hard**
Proving NP: A language $L \in NP$ means that given a certificate we can verify it in polynomial time. So: Precisely define the certificate and the verification algorithm, show that the algorithm verifies in polynomial time and cannot be fooled.
Proving NP-Hard: Given a language L and every language $L' \in NP$ $L' \leq_p L$ and possibly $L \notin NP$. Thus we must only prove that a known NP-hard problem or NP-complete problem reduces to this one in polynomial time.
Proving NP-Complete: Given a language L , prove that $L \in NP$ and prove that some known language $L' \in NPC$ reduces to this language. NOTE: the reduction may seem totally arbitrary! what you have to do is show that $(L'(x) \rightarrow \text{yes}) \Leftrightarrow (L(x) \rightarrow \text{yes})$, that is find a polynomial time algorithm to transform L' into L . Don't worry about anything except that \Leftrightarrow condition! Not all instances of your problem will cover the NPC problem! Don't worry about it!

LIST OF NPC Problems and sketches of the reductions:
 Circuit-Sat : original problem we don't do this one!
 Sat (Boolean formula sat) : label the wires and create formulas for each gate like

$$x_n \wedge (x_3 \Leftrightarrow x_1 \wedge x_2) \wedge \dots \wedge (x_7 \Leftrightarrow x_8 \vee x_9)$$

 3-CNF-SAT :
 Clique : A complete subgraph of G – thus a K -Clique is a complete subgraph of G with $k=|V|$. We reduce by creating a graph that has 3 vertice sets (same number as clauses) and put in an edge from each vertex to each vertex in other clauses that don't contradict it. If there is an n -clique where n is the number of clauses, it is satisfiable. Draw and convince yourself you should be able to reproduce it.
 Vertex-Cover : If there is a k -clique in G , then there is a vertex cover of size $n-k$ in the complement of G .
 Subset-Sum : Evil and we don't have to do it!
 Ham-Cycle: Evil and we don't have to do it!
 TSP(Traveling Salesperson): $TSP = \{ \langle G, c, k \rangle : G = (V, E) \text{ is a complete graph, } c \text{ is a function from } V \times V \rightarrow Z, k \in Z, \text{ and } G \text{ has a traveling-salesman tour with cost at most } k \}$. The reduction is simple: Take an instance of ham-cycle $G(V, E)$ and map it to a complete graph $G'(V', E')$ where if $(u, v) \in E, c(u, v) = 0$, otherwise $c(u, v) = 1$. Is there a TSP($G', c, 0$)?

For any graph $G=(V,E)$ and subset $V' \subseteq V$, the following statements are equivalent: (1) V' is a vertex cover for G . (2) $V-V'$ is an independent set for G . (3) $V-V'$ is a clique in the complement of $G \setminus \{c\}$ of G where $G^c=(V, E^c)$ and $E^c = \{(u, v) : u, v \in V \text{ and } \{u, v\} \notin E\}$

Connected graph: An undirected graph that has a path between every pair of vertices. **Strongly connected graph**: A directed graph that has a path from each vertex to every other vertex. **Degree** of a vertex: the number of edges connected to it. **Degree** of a graph: the maximum degree of any vertex. **Residual network**: Instead of using flow/capacity flow is denoted by an arrow in the opposite direction and capacity is reduced.

Augmenting paths: A path from s to t such that for each edge in the path $c(u,v) > 0$. **Cuts** (graphs/networks): A nonempty, proper subset of vertices of a graph. **Bipartite graph**: An undirected graph where vertices can be divided into two sets such that no edge connects vertices in the same set. **Trees**: Are zero based! **Bijection**: Given two non-empty sets X, Y , a function $f: X \rightarrow Y$ is a bijection if it is one-to-one (injective) and onto (surjective). **Vertex Cover**: of an undirected graph $G=(V,E)$ is a subset $V' \subseteq V$ such that if $(u,v) \in E$, then $u \in V'$, or $v \in V'$ or both. That is a vertex cover is a set of vertices that cover the edges. Running time for BFS is $O(V+E)$ BFS finds only those vertices that are reachable from S . The graph created from a BFS has no cycles. DFS may result in different trees based on the order in which adjacent vertices are visited, but this will not change the effectiveness of the results. **Running Time** $O(V+E)$. The Predecessor subgraph forms a **depth-first forest** composed of several depth-first trees. A **topological sort** of a DAG $G=(V,E)$ is a linear ordering of all its vertices such that if G contains an edge (u,v) , then u appears before v in the ordering. (If the graph is not acyclic, then no linear ordering is possible.) A **strongly connected component** of a directed graph $G=(V,E)$ is a **maximal set of vertices** $C \subseteq V$ such that " $u, v \in C$, we have both a path from u to v and v to u ."

- BFS**: Given a Graph G and a starting node s
1. Color all the nodes white
 2. Distance for all nodes u to be $d[u]=\infty$
 3. Parent of each node u to be $p[u]=nil$
 4. Color s grey
 5. Enqueue $s \rightarrow Q$
 6. while Q is not empty
 7. $u = dequeue(Q)$
 8. find each white neighbor v of u do
 9. $d[v]=d[u]+1$
 10. $p[v]=u$
 11. enqueue(v)
 12. color[u]=black

- Topological Sort(G)**
1. call DFS(G) to compute the finish times $f[v]$ for each vertex v
 2. as each vertex is finished, insert it on the front of a linked list
 3. Return the linked list of vertices
- (1) $O(V+E)$ (2-3) $O(1) = O(V+E)$

Strongly-Connected-Components(G)

- 1) call DFS(G) to compute finishing times $f[u]$ for each vertex u
- 2) compute G^T
- 3) call DFS(G^T), but in the main loop of DFS, consider the vertices in order of decreasing $f[u]$ (as computed in line 1)
- 4) output the vertices of each tree in the depth-first forest formed in line 3 as separate strongly connected components.

DFS(G)

- 1) for each vertex $u \in V[G]$
 - 2) do color[u]=white
 - 3) $p[u]=nil$
 - 4) time=0
 - 5) for each vertex $u \in V[G]$
 - 6) do if color[u]=white
 - 7) then DFS_Visit(u)
- DFS_Visit(u)**
- 1) color[u]=gray
 - 2) time++
 - 3) discover[u]=time
 - 4) for each $v \in Adj[u]$ //explore edges
 - 5) do if color[v]=white
 - 6) then $p[v]=u$
 - 7) DFS_Visit(v)
 - 8) color[u]=black
 - 9) time++
 - 10) finish[u]=time

MST - Kruskal(G, w)

1. $A = ?$ // A is the set of edges that define the MST
2. for each vertex $v \in V$
3. do Make-Set(v)
4. sort the edges of E into nondecreasing order by weight
5. for each edge $(u, v) \in E$, taken in nondecreasing order by weight
6. do if FindSet(u) \neq FindSet(v)
7. Then $A = A \cup \{(u, v)\}$ //include (u, v) safe edge in MST
8. Union(u, v) //Union the sets u and v
9. return A

Running Time: where $a(V) = O(\log|V|) = O(\log|E|)$ (because $E < V^2$) is a slowly growing function defined on Cp511. So we get $O(V \log E + E \log E) = O(E \log E)$ or using the above logic $O(E \log V)$

MST - Prim's($G(V, E), w(e \in E), r$) // graphweight, start

- 1 for each $u \in V$
 2. $key[u] = \infty$
 3. $p[u] = Nil$
 - 4 $key[r] = 0$ //decrease key
 5. Insert V into Q //amin priority queue
 6. While $Q \neq ?$
 7. $u = Extract-Min(Q)$
 8. for each $v \in Adj[u]$
 9. if $v \in Q$ and $w(u, v) < key[v]$
 10. $p[v] = u$
 11. $key[v] = w(u, v)$ //decrease key operation
- Running Time** = $O(E \log V)$

However MST running time can be decreased by using a Fibonacci heap where Extract-Min in $O(\log v)$ and Decrease-Key in $O(1)$, thus we can reduce line 11, and get a running time of $O(E + V \log V)$. **Shortest Paths**: Adjacency list representation is used for all these algorithms.

B-F Allows: negative weight edges, but not negative cycles on the shortest path.

The running

Relax(u, v, w)

- if $d[v] > d[u] + w(u, v)$
 $d[v] = d[u] + w(u, v)$
 $p[v] = u$
 $O(1)$

time for Dijkstra's algorithm is quite complex

because it depends on how the Min-Queue Q is implemented. if we use an array and take advantage of the numbered list of vertices we have \rightarrow $Insert(u) = O(1)$
 Thus the **running time** is $O(V^2 + E) = O(V^2)$. This can be reduced to $Decrease Key(u) = O(1)$

Dijkstra(G, w, s)

1. **Initialize-Single-Source(G, s)** $O(V)$
2. $S = ?$ $O(1)$
3. Build Q from V $O(V)$
4. **While $Q \neq ?$** $O(V)$
5. $u = ExtractMin(Q)$ $O(V)$
6. $S = S \cup \{u\}$ shortest
7. **for each vertex $v \in Adj[u]$** $O(E)$
8. **Relax(u, v, w)** $O(1)$

observation: **This algorithm is an operation (call it \circ) on matrices that is associative. We have** $L^{(1)} = w$,
 $L^{(2)} = L^{(1)} \circ w = w^2$
 $L^{(3)} = L^{(2)} \circ w = w^3$
 ... running

time to be $O(n^3)$. Note: w is the adjacency matrix! Note: Dropping the superscripts allows us to diminish the space requirement from $Q(n^3) \rightarrow Q(n^2)$ without disrupting the algorithm. To construct the actual shortest paths, we can use the p 's below.

$$P_{ij}^{(k)} = \begin{cases} P_{ij}^{(k-1)} & \text{if } d_{ij}^{(k-1)} \leq d_{ik}^{(k-1)} + d_{kj}^{(k-1)} \\ P_{kj}^{(k-1)} & \text{if } d_{ij}^{(k-1)} > d_{ik}^{(k-1)} + d_{kj}^{(k-1)} \end{cases} \quad P_{ij}^{(0)} = \begin{cases} NIL & \text{if } i = j \text{ or } w_{ij} = \infty \\ i & \text{if } i \neq j \text{ and } w_{ij} < \infty \end{cases}$$

DAG-Shortest Path(G, w, s)

1. Topologically sort the vertices of G $O(V+E)$
 2. Initialize-Single-Source(G, s) $O(V)$
 3. For each vertex u , taken in Topo. sorted order $O(V)$
 4. for each vertex $v \in Adj[u]$ $O(E)$ agg analysis
 5. Relax(u, v, w) $O(1)$
- Running Time**: $O(V+E)$

Initialize-Single-Source(G, s)

1. " $v \in V(G)$
2. $d[v] = \infty$
3. $p[v] = NIL$
4. $d[s] = 0$ $O(V)$

$O(V \log V + E)$ by using a Fibonacci heap from chapter 20.

All-Pairs Shortest Paths: Given a weight matrix w and a $L^{(0)}$ matrix where $L^{(0)} = [l_{ij}^{(0)}] = \begin{cases} 0 & \text{if } i = j \\ \infty & \text{if } i \neq j \end{cases}$ Notice that the

recursive formula $l_{ij}^{(m)} = \min(l_{ik}^{(m-1)} + w_{kj})$ requires the pre-calculation of $L^{(m-1)}$ and we have three indices that go from 1 to n leading us to believe that the running time will be $O(n^4)$. Obviously $L^{(0)}$ contains the shortest path from every vertex to itself if there are no negative weight cycles. We extend this using the following algorithm:

Extend-Shortest-Paths (L, w)

- $n = \#$ of rows in L $O(1)$
 int $L'[n][n]$ is an $n \times n$ matrix $O(1)$
 for $i=1$ to n $O(n)$
 for $j=1$ to n $O(n)$
 $l_{ij} = \infty$ $O(1)$
 for $k=1$ to n $O(n)$
 $l_{ij} = \min(l_{ij}, l_{ik} + w_{kj})$ $O(1)$
 return L' $O(1)$
Total Running Time $O(n^3) = O(V^3)$

After running Extend-Shortest-Paths i times, we find the shortest path between vertices of length i . To make sure that we find all paths we must run this $|V|-1$ times. Thus the overall running time is going to be $O(V^3)$. We can improve this by the following

observation: **This algorithm is an operation (call it \circ) on matrices that is associative. We have** $L^{(1)} = w$,
 $L^{(2)} = L^{(1)} \circ w = w^2$
 $L^{(3)} = L^{(2)} \circ w = w^3$
 ... running

time to be $O(n^3)$. Note: w is the adjacency matrix! Note: Dropping the superscripts allows us to diminish the space requirement from $Q(n^3) \rightarrow Q(n^2)$ without disrupting the algorithm. To construct the actual shortest paths, we can use the p 's below.

Floyd-Warshall(w)

- $n = \#$ rows in w
 $D^{(0)} = w$
 for $k = 1$ to n
 for $i = 1$ to n
 for $j = 1$ to n
 $d_{ij}^k = \min(d_{ij}^{(k-1)}, d_{ik}^{(k-1)} + d_{kj}^{(k-1)})$
 return $D^{(n)}$
Running Time $O(n^3) = O(V^3)$

Ford-Fulkerson-Method

- 1 initialize flow to 0
- 2 while $\$p$, an augmenting path
- 3 augment flow f along p
- 4 return f