```
Ts'ui Pe must have said once: I am withdrawing to write a book.

And another time: I am withdrawing to construct a labyrinth.

Every one imagined two works;

to no one did it occur that the book and the maze were one and the same thing.

— Jorge Luis Borges, "El jardín de senderos que se bifurcan" (1942)

English translation ("The Garden of Forking Paths") by Donald A. Yates (1958)

How beautiful the world would be if there were a procedure for moving through labyrinths.

— Umberto Eco, Il nome della rosa (1980)

English translation (The Name of the Rose) by William Weaver (1983)
```

# 17½ Depth-First Search in Detail

### 17½.1 Depth-First Search

Recall from the previous lecture the recursive formulation of depth-first search in undirected graphs.

```
DFS(v):

if v is unmarked

mark v

for each edge vw

DFS(w)
```

We can make this algorithm slightly faster (in practice) by checking whether a node is marked *before* we recursively explore it. This modification ensures that we call DFS( $\nu$ ) only once for each vertex  $\nu$ . We can further modify the algorithm to set the parent pointers and to compute other information about the vertices. This additional computation will be performed using two black-box subroutines PREVISIT and POSTVISIT, which we leave unspecified for now.

```
\frac{\text{DFS}(v):}{\text{mark } v}
\frac{\text{PreVisit}(v)}{\text{for each edge } vw}
\text{if } w \text{ is unmarked}
\frac{parent(w) \leftarrow v}{\text{DFS}(w)}
\frac{\text{PostVisit}(v)}{\text{PostVisit}(v)}
```

We can search any *connected* graph by unmarking all vertices and then calling DFS(s) for an arbitrary start vertex s. As we argued in the previous lecture, the subgraph of all parent edges  $v \rightarrow parent(v)$  defines a spanning tree of the graph, which we consider to be rooted at the start vertex s.

**Lemma 1.** Let T be a depth-first spanning tree of a connected undirected graph G, computed by calling DFS(s). For any node v, the vertices that are marked during the execution of DFS(v) are the proper descendants of v in T.

**Proof:** T is also the recursion tree for DFS(s).

**Lemma 2.** Let T be a depth-first spanning tree of a connected undirected graph G. For every edge vw in G, either v is an ancestor of w in T, or v is a descendant of w in T.

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**Proof:** Assume without loss of generality that v is marked before w. Then w is unmarked when DFS(v) is invoked, but marked when DFS(v) returns, so the previous lemma implies that w is a proper descendant of v in T.

Lemma 2 implies that any depth-first spanning tree *T* divides the edges of *G* into two classes: *tree* edges, which appear in *T*, and *back* edges, which connect some node in *T* to one of its ancestors.

### 17½.2 Counting and Labeling Components

For graphs that might be connected, we can compute a depth-first spanning *forest* by calling the following wrapper function; again, we introduce a generic black-box subroutine PREPROCESS to perform any necessary preprocessing for the POSTVISIT and POSTVISIT functions.

```
\frac{\text{DFSALL}(G):}{\text{PREPROCESS}(G)}
for all vertices v
unmark v
for all vertices v
if v is unmarked
\text{DFS}(v)
```

With very little additional effort, we can count the components of a graph; we simply increment a counter inside the wrapper function. Moreover, we can also record which component contains each vertex in the graph by passing this counter to DFS. The single line  $comp(v) \leftarrow c$  is a trivial example of PREVISIT.<sup>1</sup>

```
COUNTANDLABEL(G):

c \leftarrow 0

for all vertices v

unmark v

for all vertices v

if v is unmarked

c \leftarrow c + 1

LABELCOMPONENT(v, c)

return c
```

```
LABELCOMPONENT(v, c):

mark v

comp(v) ← c

for each edge vw

if w is unmarked

LABELCOMPONENT(w, c)
```

## 171/2.3 Preorder and Postorder Labeling

You should already be familiar with preorder and postorder traversals of rooted trees, both of which are derived from depth-first search. Similar traversal orders can be defined for arbitrary graphs as follows:

```
LABELCOMPONENT(v, clock):

mark v

pre(v) \leftarrow clock

clock \leftarrow clock + 1

for each edge vw

if w is unmarked

clock \leftarrow LABELCOMPONENT(w, clock)

post(v) \leftarrow clock

clock \leftarrow clock + 1

return clock
```

<sup>&</sup>lt;sup>1</sup>The absence of instructions after the for loop is a vacuous example of PostVisit.

Equivalently, if we're willing to use a (shudder) global variable, we can use our generic depth-first-search algorithm with the following subroutines PREPROCESS, PREVISIT, and POSTVISIT.

```
\frac{\text{Preprocess}(G):}{\text{clock} \leftarrow 0}
```

```
\frac{\text{PReVisit}(v):}{\text{pre}(v) \leftarrow \text{clock}}\text{clock} \leftarrow \text{clock} + 1
```

```
\frac{\text{PostVisit}(v):}{post(v) \leftarrow clock}clock \leftarrow clock + 1
```

Consider two vertices u and v, where u is marked after v. Then we must have pre(u) < pre(v). Moreover, Lemma 1 implies that if v is a descendant of u, then post(u) > post(v), and otherwise, pre(v) > post(u). Thus, for any two vertices u and v, the intervals [pre(u), post(u)] and [pre(v), post(v)] are either disjoint or nested; in particular, if uv is an edge, Lemma 2 implies that the intervals must be nested.

## 171/2.4 Directed Graphs and Reachability

The recursive algorithm requires only one minor change to handle directed graphs:

 $\frac{\text{DFSALL}(G):}{\text{for all vertices } v}$  unmark v for all vertices v if v is unmarked DFS(v)

```
\frac{\text{DFS}(v):}{\text{mark } v}
\text{PREVISIT}(v)
\textit{for each edge } v \rightarrow w
\text{if } w \text{ is unmarked}
\text{DFS}(w)
\text{PostVISIT}(v)
```

However, it's less clear now that we can use this modified algorithm to count components. Suppose G is a single directed path. Depending on the order that we choose the visit the nodes in DFSALL, we may discover any number of "components" between 1 and n; all that we can be guaranteed is that the "component" numbers are non-decreasing as we traverse the path. In fact, the real problem is that the definition of "component" is only suitable for *undirected* graphs.

Instead, for directed graphs we rely on a more subtle notion of *reachability*. We say that a node v is *reachable* from another node u in a directed graph G—or more simply, that u can reach v—if and only if there is a directed path in G from u to v. Let Reach(u) denote the set of vertices that are reachable from u (including u itself). A simple inductive argument proves that Reach(u) is precisely the subset of nodes that are marked by calling DFS(u).

## 17½.5 Directed Acyclic Graphs

A *directed acyclic graph* or *dag* is a directed graph with no directed cycles. Any vertex in a dag that has no incoming vertices is called a *source*; any vertex with no outgoing edges is called a *sink*. Every dag has at least one source and one sink (Do you see why?), but may have more than one of each. For example, in the graph with *n* vertices but no edges, every vertex is a source and every vertex is a sink.

We can check whether a given directed graph G is a dag in O(V+E) time as follows. First, to simplify the algorithm, we add a single artificial source s, with edges from s to every other vertex. Because s has no outgoing edges, no directed cycle in G+s goes through s, which implies that G+s is a dag if and only if G is a dag. Then we preform a depth-first search of G+s starting at the new source vertex s; by construction every other vertex is reachable from s, so this search visits every node in the graph.

Instead of vertices being merely marked or unmarked, each vertex has one of three statuses—New, Active, or Done—which depend on whether we have started or finished the recursive depth-first search at that vertex. (Since this algorithm never uses parent pointers, I've removed the line " $parent(w) \leftarrow v$ ".)

```
ISACYCLIC(G):

add vertex s

for all vertices v \neq s

add edge s \rightarrow v

status(v) \leftarrow New

return IsACYCLICDFS(s)
```

```
IsAcyclicDFS(\nu):

status(\nu) \leftarrow Active

for each edge \nu \rightarrow w

if status(w) = Active

return False

else if status(w) = New

if IsAcyclicDFS(w) = False

return False

status(\nu) \leftarrow Done

return True
```

Suppose the algorithm returns False. Then the algorithm must discover an edge  $v \rightarrow w$  such that status(w) = Active. The active vertices are precisely the vertices currently on the recursion stack, which are all ancestors of the current vertex v. Thus, there is a directed path from w to v, and so the graph has a directed cycle.

On the other hand, suppose G has a directed cycle. Let w be the first vertex in this cycle that we visit, and let  $v \rightarrow w$  be the edge leading into v in the same cycle. Because there is a directed path from w to v, we must call IsAcyclicDFS(v) during the execution of IsAcyclicDFS(w), unless we discover some other cycle first. During the execution of IsAcyclicDFS(v), we consider the edge  $v \rightarrow w$ , discover that status(w) = Active. The return value False bubbles up through all the recursive calls to the top level.

We conclude that IsAcyclic(G) returns True if and only if G is a dag.

## 17½.6 Topological Sort

A **topological ordering** of a directed graph G is a total order  $\prec$  on the vertices such that  $u \prec v$  for every edge  $u \rightarrow v$ . Less formally, a topological ordering arranges the vertices along a horizontal line so that all edges point from left to right. A topological ordering is clearly impossible if the graph G has a directed cycle—the rightmost vertex of the cycle would have an edge pointing to the left! On the other hand, every dag has a topological order, which can be computed by either of the following algorithms.

```
TOPOLOGICALSORT(G):
n \leftarrow |V|
for i \leftarrow n down to 1
v \leftarrow \text{ any sink in } G
S[1] \leftarrow v
remove v and all edges entering v
return S[1..n]
```

The correctness of these algorithms follow inductively from the observation that *deleting* a vertex cannot *create* a cycle.

This simple algorithm has two major disadvantages. First, the algorithm actually destroys the input graph. This destruction can be avoided by simply marking the "deleted" vertices, instead of actually deleting them, and defining a vertex to be a source (sink) if none of its incoming (outgoing) edges come from (lead to) an unmarked vertex. The more serious problem is that finding a source vertex seems to require  $\Theta(V)$  time in the worst case, which makes the running time of this algorithm  $\Theta(V^2)$ . In fact, a careful implementation of this algorithm computes a topological ordering in O(V+E) time without removing any edges.

But there is a simpler linear-time algorithm based on our earlier algorithm for deciding if a directed graph is acyclic. The new algorithm is based on the following observation:

**Lemma 3.** For any directed acyclic graph G, the first vertex marked Done by IsAcyclic(G) must be a sink.

**Proof:** Let v be the first vertex marked Done during an execution of IsAcyclic. For the sake of argument, suppose v has an outgoing edge  $v \rightarrow w$ . When IsAcyclicDFS considers the edge  $v \rightarrow w$ , there are three cases to consider.

- If status(w) = Done, then w is marked Done before v, which contradicts the definition of v.
- If status(w) = New, the algorithm calls TopoSortDFS(w), which (among other computation) marks w. Thus, w is marked Done before v, which contradicts the definition of v.
- If status(w) = Active, then G has a directed cycle, contradicting our assumption that G is acyclic.

In all three cases, we have a contradiction, so  $\nu$  must be a sink.

Thus, to topologically sort a dag *G*, it suffice to list the vertices in the *reverse* order of being marked Done. For example, we could push each vertex onto a stack when we mark it Done, and then pop every vertex off the stack.

```
TOPOLOGICALSORT(G):
add vertex s
for all vertices v \neq s
add edge s \rightarrow v
status(v) \leftarrow \text{New}
if TOPOSORTDFS(s) = FALSE
return \text{ FAILURE}
for \ i \leftarrow 1 \ to \ V
S[i] \leftarrow \text{POP}
return \ S[1..V]
```

```
TopoSortDFS(v):

status(v) \leftarrow Active

for each edge v \rightarrow w

if status(w) = Active

return False

else if status(w) = New

if TopoSortDFS(w) = False

return False

status(v) \leftarrow Done

Push(v)

return True
```

But maintaining a separate data structure is really overkill. In most applications of topological sort, out goal is not to compute an explicit sorted list of the vertices; instead, we need to process the graph either in topological order or in reverse topological order, performing some fixed computation at each vertex. In this case, it is not necessary to *record* the topological order. To process the graph in *reverse* topological order, we can just process each vertex at the end of its recursive depth-first search. (The following algorithm omits the cycle detection; the input graph is assumed to be a dag.)

```
PROCESS DAGBACKWARD (G):
add vertex s
for all vertices v \neq s
add edge s \rightarrow v
status(v) \leftarrow \text{New}
PROCESS BACKWARD DFS (s)
```

```
\frac{\text{PROCESSBACKWARDDFS}(v):}{status(v) \leftarrow \text{ACTIVE}}
for each edge v \rightarrow w
if status(w) = \text{NEW}
\text{PROCESSBACKWARDDFS}(w)
else if status(w) = \text{ACTIVE}
fail gracefully
status(v) \leftarrow \text{Done}
\text{PROCESS}(v)
```

If we already *know* that the input graph is acyclic, we can simplify this algorithm a bit. In particular, we no longer need a separate 'active' status for vertices.

```
PROCESSBACKWARDDFS(v):
mark v
for each edge v \rightarrow w
if w is unmarked
PROCESSBACKWARDDFS(w)
PROCESS(v)
```

Except for the addition of the artificial source vertex *s*, which we need to ensure that every vertex is visited, this is just the standard depth-first search algorithm, with PostVisit renamed to Process!

The simplest way to process a dag in *forward* topological order is to construct the *reversal* of the input graph, which is obtained by replacing each each  $v \rightarrow w$  with its reversal  $w \rightarrow v$ . Reversing a directed cycle gives us another directed cycle with the opposite orientation; thus, the reversal of a dag is another dag. Every source in G becomes a sink in the reversal of G and vice versa; it follows inductively that every topological ordering for the reversal of G is the reversal of a topological ordering of G. The reversal of any directed graph can be computed in O(V+E) time; the details of this construction are left as an easy exercise.

### 17½.7 Every Dynamic Programming Algorithm!

Our topological sort algorithm is actually the model for *every* dynamic programming algorithm! Recall that the *dependency graph* of a recurrence has a vertex for every recursive subproblem and an edge from one subproblem to another if evaluating the first subproblem requires a recursive evaluation of the second. Evaluating any recurrence with memoization is exactly the same as performing a depth-first search of the dependency graph. In particular, a vertex of the dependency graph is 'marked' if the value of the corresponding subproblem has already been computed, and the black-box subroutine Process is a placeholder for the actual value computation.

Conversely, we can use depth-first search to build dynamic programming algorithms for problems defined over dags. For example, consider the *longest path* problem, which asks for the path of *maximum* total weight from one node s to another node t in a directed graph G with weighted edges. The longest path problem is NP-hard in general directed graphs, by an easy reduction from the traveling salesman problem, but it is easy to solve in linear time if the input graph G is acyclic. For any node s, let LLP(s,t) denote the length of longest path in G from s to t. If G is a dag, this function satisfies the recurrence

$$LLP(s,t) = \begin{cases} 0 & \text{if } s = t, \\ \max_{s \to v} (\ell(s \to v) + LLP(v,t)) & \text{otherwise,} \end{cases}$$

where  $\ell(v \to w)$  is the given weight ('length') of edge  $v \to w$ . In particular, if s is a sink but not equal to t, then  $LLP(s,t) = \infty$ . We can now easily evaluate this recursive function in O(V+E) time by performing a depth-first search of G starting at s.

```
\frac{\text{LongestPath}(s,t):}{\text{if } s = t}
\text{return 0}
\text{if } LLP(v) \text{ is undefined}
LLP(v) \leftarrow \infty
\text{for each edge } s \rightarrow v
LLP(v) \leftarrow \max \{LLP(v), \ \ell(s \rightarrow v) + \text{LongestPath}(v,t)\}
\text{return } LLP(v)
```

A surprisingly large number of dynamic programming problems can be recast as either longest or shortest path problems in the associated dependency graph.

### 17½.8 Strong Connectivity

Let's go back to the proper definition of connectivity in directed graphs. Recall that one vertex u can reach another vertex v in a graph G if there is a directed path in G from u to v, and that Reach(u) denotes the set of all vertices that u can reach. Two vertices u and v are strongly strongly strongly strongly to nuclear than the equivalence relation over the set of vertices of any directed graph, just as connectivity is for undirected graphs. The equivalence classes of this relation are called the <math>strongly strongly to nuclear than the equivalence components (or more simply, the <math>to nuclear than the equivalence classes of this relation are called the <math>to nuclear than the equivalence classes of this relation are called the <math>to nuclear than the equivalence classes of this relation are called the <math>to nuclear than the equivalence classes of this relation are called the <math>to nuclear than the equivalence components (or more simply, the <math>to nuclear than the equivalence components) of to nuclear than the equivalence components (or more simply, the <math>to nuclear than the equivalence classes of this relation are called the <math>to nuclear than the equivalence classes of this relation are called the <math>to nuclear than the equivalence classes of this relation are called the <math>to nuclear than the equivalence classes of this relation are called the <math>to nuclear than the equivalence classes of this relation are called the <math>to nuclear than the equivalence classes of this relation of <math>to nuclear than the equivalence classes of this relation of <math>to nuclear than the equivalence classes of this relation of <math>to nuclear than the equivalence classes of this relation of <math>to nuclear than the equivalence classes of the equivalence classes of this relation of <math>to nuclear than the equivalence classes of the equivalence classes of this relation of <math>to nuclear than the equivalence classes of the equivalenc

It is straightforward to compute the strong component containing a single vertex v in O(V+E) time. First we compute Reach(v) by calling DFS(v). Then we compute  $Reach^{-1}(v) = \{u \mid v \in Reach(u)\}$  by searching the reversal of G. Finally, the strong component of v is the intersection  $Reach(v) \cap Reach^{-1}(v)$ . In particular, we can determine whether the entire graph is strongly connected in O(V+E) time.

We can compute *all* the strong components in a directed graph by wrapping the single-strong-component algorithm in a wrapper function, just as we did for depth-first search in undirected graphs. However, the resulting algorithm runs in O(VE) time; there are at most V strong components, and each requires O(E) time to discover. Surely we can do better! After all, we only need O(V+E) time to decide whether every strong component is a single vertex.

### 17½.9 Strong Components in Linear Time

For any directed graph G, the **strong component graph scc(G)** is another directed graph obtained by contracting each strong component of G to a single (meta-)vertex and collapsing parallel edges. The strong component graph is sometimes also called the *meta-graph* or *condensation* of G. It's not hard to prove (hint, hint) that scc(G) is always a dag. Thus, in principle, it is possible to topologically order the strong components of G; that is, the vertices can be ordered so that every *backward* edge joins two edges in the same strong component.

Let C be any strong component of G that is a sink in scc(G); we call C a sink component. Every vertex in C can reach every other vertex in C, so a depth-first search from any vertex in C visits every vertex in C. On the other hand, because C is a sink component, there is no edge from C to any other strong component, so a depth-first search starting in C visits only vertices in C. So if we can compute all the strong components as follows:

At first glance, finding a vertex in a sink component *quickly* seems quite hard. However, we can quickly find a vertex in a *source* component using the standard depth-first search. A source component is a strong component of G that corresponds to a source in scc(G). Specifically, we compute *finishing times* for the vertices of G as follows:

```
\frac{\mathrm{DFS}(v, clock):}{\mathrm{mark}\ v}
for each edge v \rightarrow w
if w is unmarked
clock \leftarrow \mathrm{DFS}(w, clock)
clock \leftarrow clock + 1
finish(v) \leftarrow clock
return\ clock
```

**Lemma 4.** The vertex with largest finishing time lies in a source component of *G*.

**Proof:** Let v be the vertex with largest finishing time. Then DFS(v, clock) must be the last direct call to DFS made by the wrapper algorithm DFSALL.

Let *C* be the strong component of *G* that contains v. For the sake of argument, suppose there is an edge  $x \rightarrow y$  such that  $x \notin C$  and  $y \in C$ . Because v and y are strongly connected, y can reach v, and therefore x can reach v. There are two cases to consider.

- If x is already marked when DFS( $\nu$ ) begins, then  $\nu$  must have been marked during the execution of DFS(x), because x can reach  $\nu$ . But then  $\nu$  was already marked when DFS( $\nu$ ) was called, which is impossible.
- If x is not marked when DFS( $\nu$ ) begins, then x must be marked during the execution of DFS( $\nu$ ), which implies that  $\nu$  can reach x. Since x can also reach  $\nu$ , we must have  $x \in C$ , contradicting the definition of x.

We conclude that *C* is a source component of *G*.

Essentially the same argument implies the following more general result.

**Lemma 5.** For any edge  $v \rightarrow w$  in G, if finish(v) > finish(w), then v and w are strongly connected in G.

This observation is consistent with our earlier topological sorting algorithm; for *every* edge  $v \rightarrow w$  in a directed acyclic graph, we have finish(v) < finish(w).

It is easy to check (hint, hint) that any directed G has exactly the same strong components as its reversal rev(G); in fact, we have rev(scc(G)) = scc(rev(G)). Thus, if we order the vertices of G by their finishing times in DFSALL(rev(G)), the *last* vertex in this order lies in a sink component of G. Thus, if we run DFSALL(G), visiting vertices in reverse order of their finishing times in DFSALL(rev(G)), then each call to DFS visits exactly one strong component of G.

Putting everything together, we obtain the following algorithm to count and label the strong components of a directed graph in O(V+E) time, first discovered (but never published) by Rao Kosaraju in 1978. The same algorithm was independently rediscovered in 1981 by Micha Sharir. The Kosaraju-Sharir algorithm has two phases. The first phase performs a depth-first search of the reversal of G, pushing each vertex onto a stack when it is finished. In the second phase, we perform another depth-first search of the original graph G, considering vertices in the order they appear on the stack.

```
KosarajuSharir(G):
  ⟨⟨Phase 1: Push in finishing order⟩⟩
  for all vertices \nu
       unmark v
  for all vertices v
       if v is unmarked
            clock \leftarrow RevPushDFS(v)
  ⟨⟨Phase 2: DFS in stack order⟩⟩
  for all vertices \nu
       unmark v
  count \leftarrow 0
  while the stack is non-empty
       \nu \leftarrow Pop
       if v is unmarked
            count \leftarrow count + 1
            LABELONEDFS(v, count)
```

```
REVFINISHDFS(v):

mark v

for each edge v \rightarrow u in rev(G)

if u is unmarked

REVFINISHDFS(u)

PUSH(v)
```

```
LabelOneDFS(v, count):

mark v

label(v) \leftarrow count

for each edge v \rightarrow w in G

if w is unmarked

LabelOneDFS(w, count)
```

## **Exercises**

- 0. (a) Describe an algorithm to compute the reversal rev(G) of a directed graph in O(V + E) time.
  - (b) Prove that for any directed graph G, the strong component graph scc(G) is a dag.
  - (c) Prove that for any directed graph G, we have scc(rev(G)) = rev(scc(G)) is a dag.
- 1. Let *G* be a directed acyclic graph with a unique source *s* and a unique sink *t*.
  - (a) A *Hamiltonian path* in *G* is a directed path in *G* that contains every vertex in *G*. Describe an algorithm to determine whether *G* has a Hamiltonian path.
  - (b) Suppose the *vertices* of *G* have weights. Describe an efficient algorithm to find the path from *s* to *t* with maximum total weight.
  - (c) Suppose the vertices of G have integer labels, where  $label(s) = -\infty$  and  $label(t) = \infty$ . Describe an algorithm to find the longest path from s to t whose vertex labels define an increasing sequence.
  - (d) Describe an algorithm to compute the number of distinct paths from s to t in G. (Assume that you can add arbitrarily large integers in O(1) time.)
- 2. Let G be a directed acyclic graph, whose vertices have labels from some fixed finite alphabet, and let  $A[1..\ell]$  be a string over the same alphabet. Any directed path in G has a label, which is obtained by concatenating the labels of its vertices.
  - (a) Describe an algorithm to find the *number* of paths in G whose label is A. (Assume that you can add arbitrarily large integers in O(1) time.)
  - (b) Describe an algorithm to find the longest path in G whose label is a subsequence of A.
  - (c) Describe an algorithm to find the *shortest* path in *G* whose label is a *supersequence* of *A*.
- 3. Suppose two players are playing a turn-based game on a directed acyclic graph G with a unique source s. Each vertex v of G is labeled with a real number  $\ell(v)$ , which could be positive, negative,

or zero. The game starts with three tokens at s. In each turn, the current player moves one of the tokens along a directed edge from its current node to another node, and the current player's score is increased by  $\ell(u) \cdot \ell(v)$ , where u and v are the locations of the two tokens that did *not* move. At most one token is allowed on any node except s at any time. The game ends when the current player is unable to move (for example, when all three tokens lie on sinks); at that point, the player with the higher score is the winner.

Describe an efficient algorithm to determine who wins this game on a given labeled graph, assuming both players play optimally.

- \*4. Let  $x = x_1 x_2 ... x_n$  be a given *n*-character string over some finite alphabet  $\Sigma$ , and let *A* be a deterministic finite-state machine with *m* states.
  - (a) Describe and analyze an algorithm to compute the longest subsequence of x that is accepted by A. For example, if A accepts the language  $(ar)^*$  and  $x = \underline{a}\underline{b}\underline{r}\underline{a}\underline{c}\underline{a}\underline{d}\underline{a}\underline{b}\underline{r}\underline{a}$ , your algorithm should output arar.
  - (b) Describe and analyze an algorithm to compute the shortest (not necessarily contiguous) supersequence of x that is accepted by A. For example, if A accepts the language (abcdr)\* and x = abracadabra, your algorithm should output abcdrabcdrabcdrabcdrabcdr.