

CMSC 451: Lecture 3
Cycles and Strong Components
Tuesday, Sep 7, 2017

Reading: Our algorithm for strong components follows the presentation of Sect. 3.4 of DPV.

Applications of DFS: Last time we introduced depth-first search (DFS). Today, we discuss some applications of this powerful and efficient method of graph traversal.

Directed Acyclic Graphs (Optional): A *directed acyclic graph*, or *DAG*, is a directed graph that has no cycles. DAGs arise in many applications where there are precedence or ordering constraints. For example, if there are a series of tasks to be performed, and certain tasks must precede other tasks (e.g., in construction you have to build the walls before you install the windows). In general a *precedence constraint graph* is a DAG in which vertices are tasks and the edge (u, v) means that task u must be completed before task v begins.

It is easy to see that every DAG must have at least one vertex with no incoming edges, and at least one vertex with no outgoing edges. A vertex with no incoming edges (only outgoing) is called a *source* and a vertex with no outgoing edges (only incoming) is called a *sink*.

Acyclicity Testing (Optional): Let us consider the problem of determining whether a digraph is acyclic. We are given a directed graph $G = (V, E)$, and we wish to determine whether G contains a cycle. If so, G is not a DAG.

We will present a simple algorithm based on DFS. Recall that in addition to tree edges, a DFS forest contains three other types of edges, back edges (which go to a vertex's ancestor), forward edges (which go to a vertex's descendant), and cross edges (everything else). Observe that if the DFS forest of G contains at least one back edge, then G has a cycle. This is easy to see. If (u, v) is a back edge, then there is a path in the tree from the ancestor v to the descendant u , and the back edge from u to v completes the cycle. The following lemma shows that this condition is not only sufficient, but necessary.

Claim: If a digraph G has a cycle, then *any* DFS forest of G (i.e., no matter what order the vertices are visited) has a *back edge*.

Proof: The proof is based on a very simple observation about the various edge types and finish times. Recall from the Parenthesis Lemma (from the previous lecture) that if u is an ancestor of v then we have $[d[v], f[v]] \subset [d[u], f[u]]$. It follows that if (u, v) is a tree edge or forward edge then $f[u] > f[v]$. Also, observe that (u, v) is a cross edge, it must be u was discovered after v was finished (for otherwise, u would have made a DFSvisit call on v , implying that this would be a tree edge). Therefore $d[u] > f[v]$. Since a vertex cannot finish until after it was discovered, we have $f[u] > f[v]$.

In summary, for all these three edge types (tree, forward, and cross), the finish time of the origin is strictly greater than the finish time of the destination. It follows directly that it is impossible to complete a cycle from any combination of just these three edge types. Only for back edges do we have $f[u] < f[v]$, and therefore in order to form a cycle we need to have *at least one* back edge. Therefore, if a graph G has a cycle, in any DFS forest of G there must be at least one back edge.

The above theorem implies that in order to determine whether a graph G has a cycle, it suffices to test whether it has a back edge. How do we know whether an edge is a back edge. The proof of the above theorem provides an easy way. We can first apply DFS to G , and we then run through the edges, checking whether $d[u] > d[v]$. Can we do this on the fly as DFS is running? The answer is yes. Observe that a back edge goes from a vertex u to an ancestor v . Such an ancestor must have been discovered, but not yet finished. For the other non-tree edge types, the destination v will have already finished. The main DFS function is the same, only DFSvisit needs to be updated.

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Determining whether a graph has a cycle
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DFSvisit(u) {
    mark[u] = discovered           // perform a DFS search at u
    d[u] = ++time                 // u has been discovered
    for each (v in Adj(u)) {
        if (mark[v] == undiscovered) { // undiscovered neighbor?
            pred[v] = u
            DFSvisit(v)               // ...visit it
        }
        else if (mark[v] != finished) { // found a back edge?
            output "G has a cycle!"
        }
    }
    mark[u] = finished           // we're done with u
    f[u] = ++time
}

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Strong Components: (The following material applies *only* to directed graphs!)

A digraph $G = (V, E)$ is said to be *strongly connected* if for every vertex u and v there is a path from u to v and from v to u . It is easy that this *mutual reachability* relation between vertices is an equivalence relation. This implies that it partitions V into equivalence class, called the *strong components* (or *strongly-connected components*) of G (see Fig. 1(a) and (b)).

If the vertices within each strong component are collapsed into a single vertex, the resulting digraph is called the *component digraph* (see Fig. 1(c)). It is easy to see that the component digraph must be acyclic (since if a number of components could be joined in a cycle, they would collapse into a single larger strong component). Therefore, this graph is usually called the *component DAG*.

There exists an $O(n + m)$ -time DFS algorithm for computing strong components. It is based on the following lemma.

Claim 1: If DFSvisit is started at a vertex u , it will terminate precisely when all the vertices reachable from u have been visited.

Proof: This follows from the exhaustive nature of DFS. (Note that some of these vertices may have been reached by earlier calls to DFSvisit.)

Claim 2: If C and C' are two strong components, and there is an edge from a vertex in C to a vertex in C' , then the highest finish time in C is bigger than the highest finish time in C' .

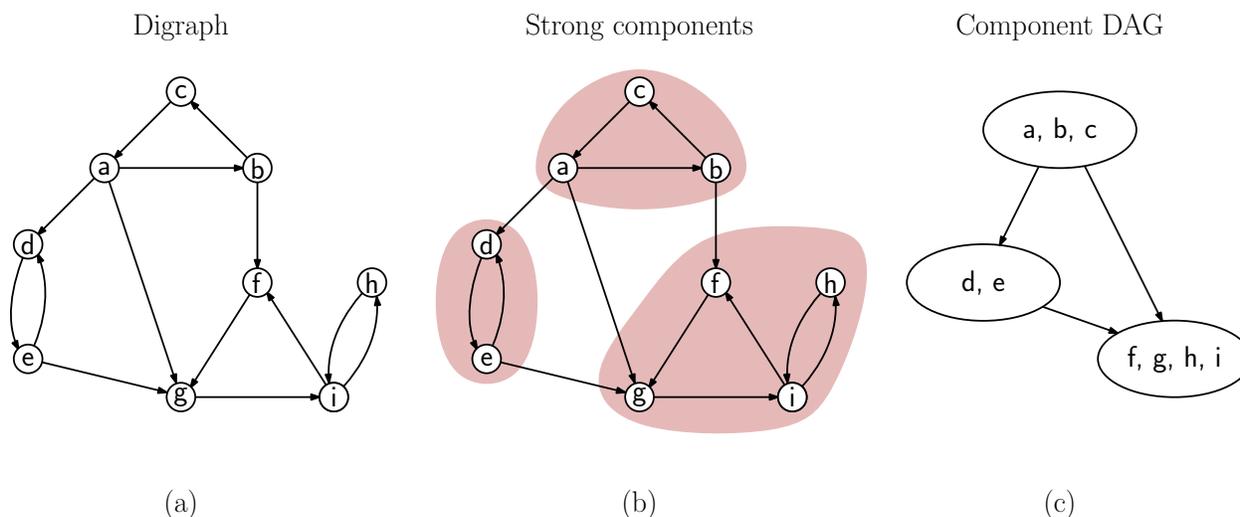


Fig. 1: Strong components and the component DAG.

Proof: There are two cases depending on whether the DFS first encounters a vertex from C or C' . If it first encounters a vertex u in C , then by Claim 1 the DFS will visit all the vertices of both C and C' before returning to u . Therefore, u will have the highest finish time of every vertex in $C \cup C'$. If it first visits a vertex in C' , then the DFS will get stuck in C' (since it is not possible to reach anything in C). It follows that all the vertices of C will have higher discovery times than those of C' , which further implies that they will have higher finish times as well.

Claim 3: The vertex that receives the highest finish time in a DFS must lie in a source vertex of the component DAG. (Recall that a vertex in a DAG is called a source if it has no incoming edges.)

Claim 3 is equivalent to saying that the strong components can be linearly arranged in decreasing order of their highest finish times. By doing so, every edge in the component DAG will go from an earlier component in the linear order to a later one. How can we exploit this to obtain an efficient algorithm to find the strong components.

Claim 3 allows us to identify a vertex in some *source* of the component DAG. Unfortunately, this is not all that useful. What *would* be useful is to identify a *sink* of the component DAG. If we could do this, we could start a DFS at this vertex, with the knowledge (by Claim 2) that no other strong components would be visited. We could then delete all these vertices (or equivalently, mark them as visited), and repeat the process. Eventually, all the strong components will be identified, each one arising as a separate subtree of the DFS forest.

So how do we convert an algorithm that identifies sources to one that identifies sinks in the component DAG? The trick is to reverse all the edges of G . Let G^R denote the directed graph that has the same vertex set as G , but every edge (u, v) is replaced by its reverse (v, u) . Note that the strong components of G^R are the same as G , but the direction of edges in the component DAG have all be *reversed*. Thus, the sources in the component DAG of G^R are

sinks in the component DAG of G . This leads to the following (insanely clever) algorithm for computing strong components.

- (1) Given G , compute G^R (see Fig. 2(a)). (Note: This is a small programming exercise, which involves a simple traversal of G 's adjacency list. It can be done in $O(n + m)$ time.)
- (2) Run $\text{DFS}(G^R)$ and label each vertex of G with the finish time of the corresponding vertex of G^R (see Fig. 2(b)).
- (3) Sort the vertices of G in decreasing order of finish times. (Note: Since finish times are integers in the range $[1, 2n]$, this can be done in $O(n)$ time through Bucket Sort.)
- (4) Run $\text{DFS}(G)$, but in the outermost loop, whenever we need to find a new vertex to start DFSvisit , take the vertex with the highest finish time (using the above sorted order).
- (5) Each subtree of the DFS forest will be a strong component (see Fig. 2(c)).

The correctness of the above algorithm follows from the remarks made earlier. The running time is $O(n + m)$, dominated by the time to compute G^R and the times for the two DFS's.

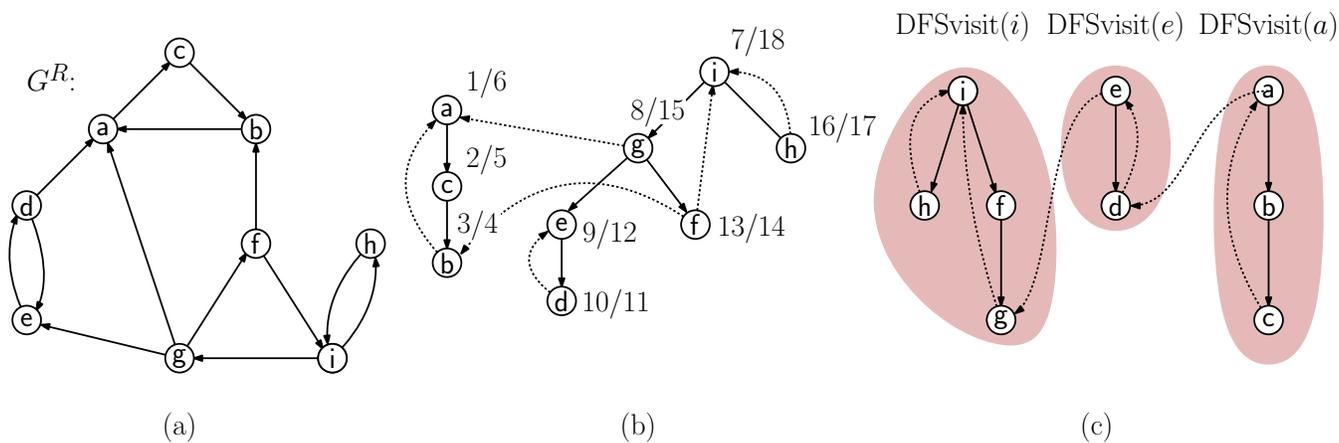


Fig. 2: Strong components and DFS.