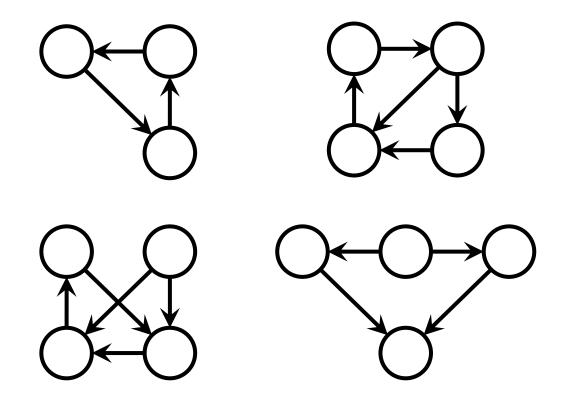
Graph Algorithms II

Outline for Today

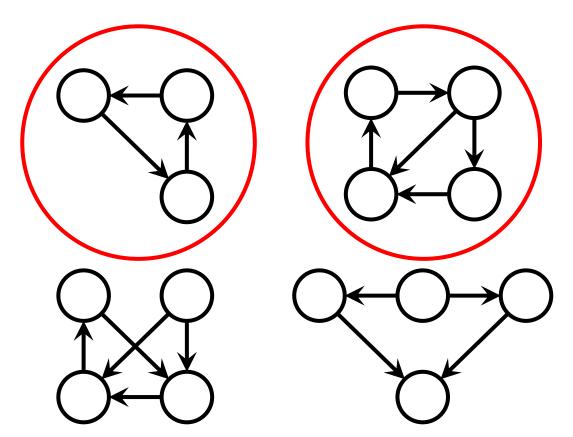
Graph algorithms

Kosaraju's Algorithm for finding strongly connected components Karger's Algorithm for finding global minimum cuts

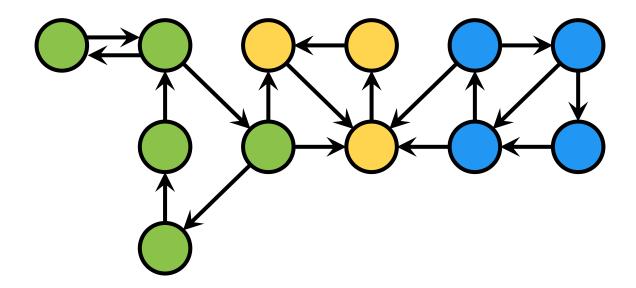
A directed graph G = (V, E) is strongly connected if, for all pairs of vertices u and v, there's a path from u to v and a path from v to u.



A directed graph G = (V, E) is strongly connected if, for all pairs of vertices u and v, there's a path from u to v and a path from v to u.



We can decompose a graph into its strongly connected components (SCCs).



Why do we care about SCCs?

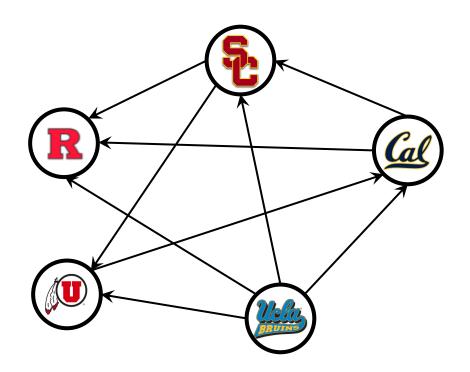
SCCs provide information about communities.

A computer scientist might want to decompose the Internet into SCCs to find connected areas.

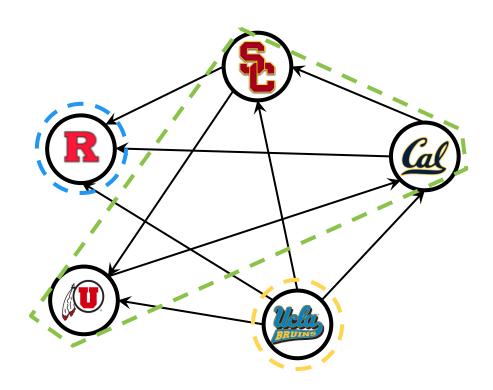
An economist might want to decompose labor market data (e.g. employeremployee graph) into SCCs to analyze labor status.

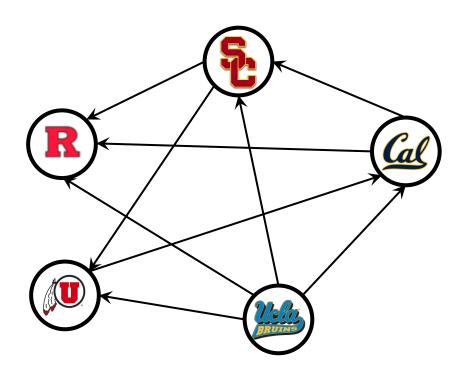
A football executive might want to determine which schools should play in NCAA.

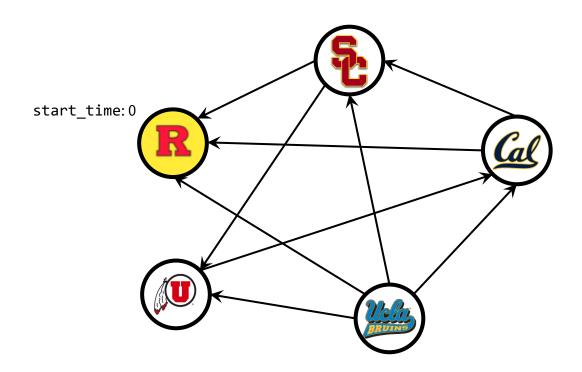
How many SCCs are in this graph? 🤔

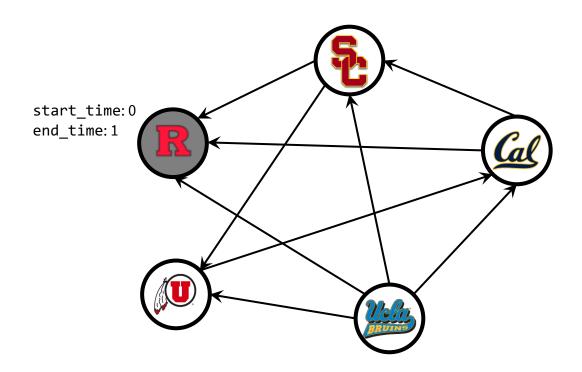


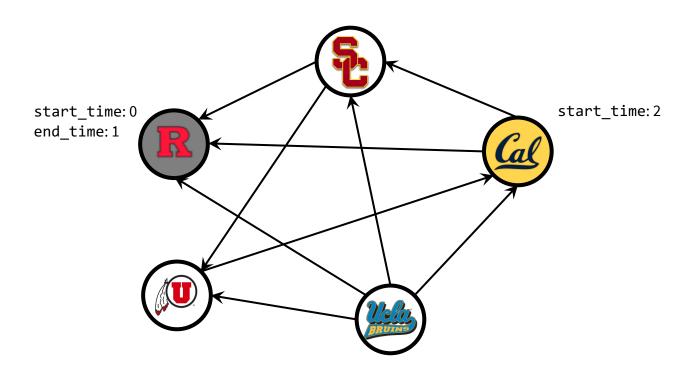
How many SCCs are in this graph? 9 3; let's find them!

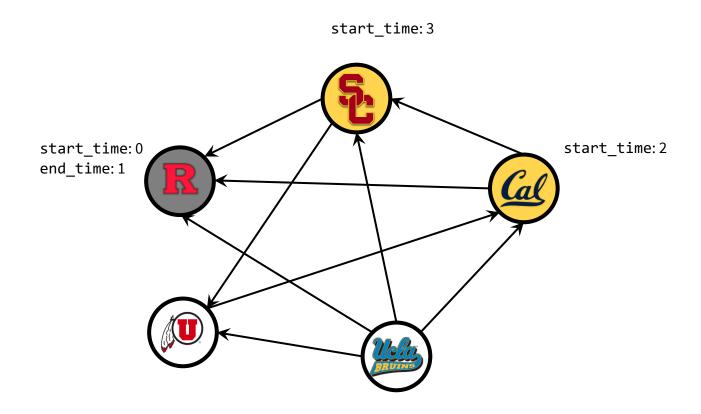


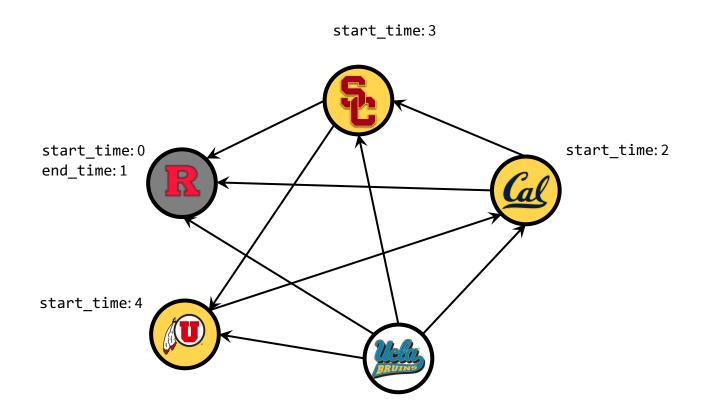


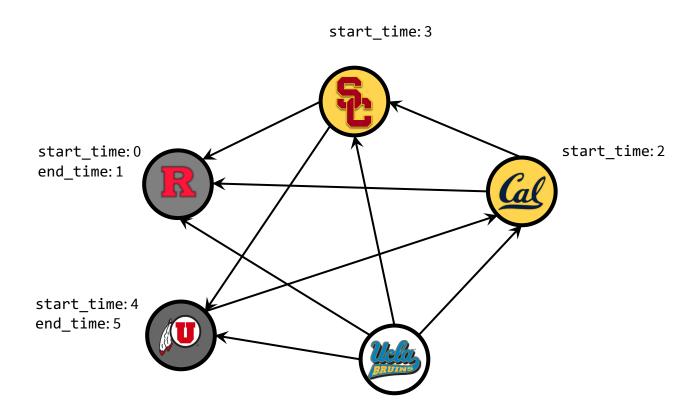


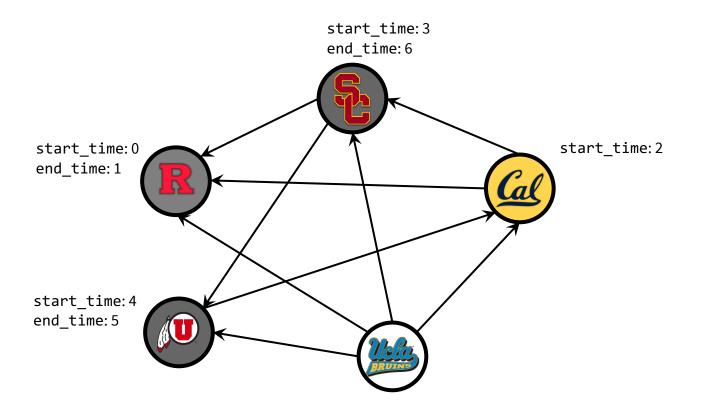


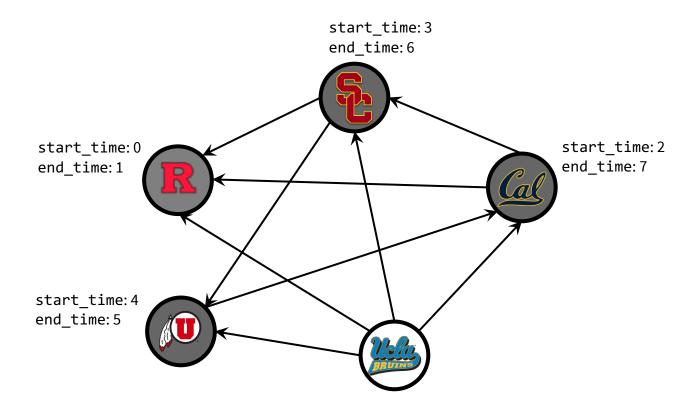


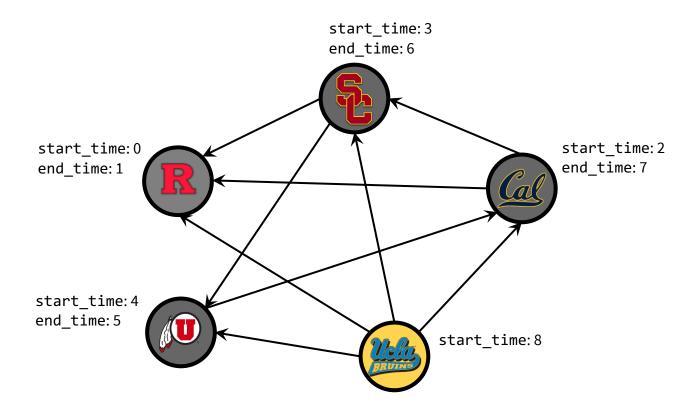


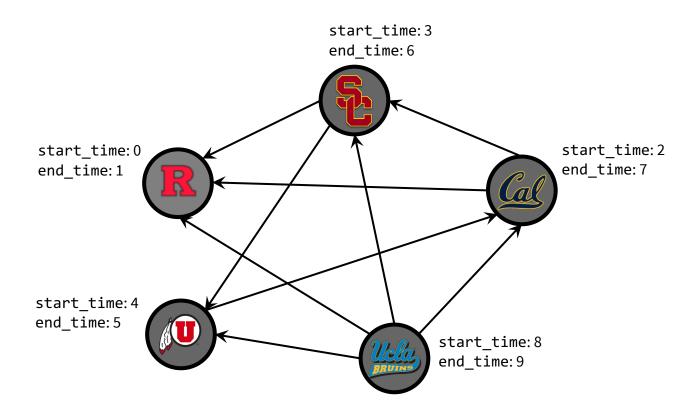




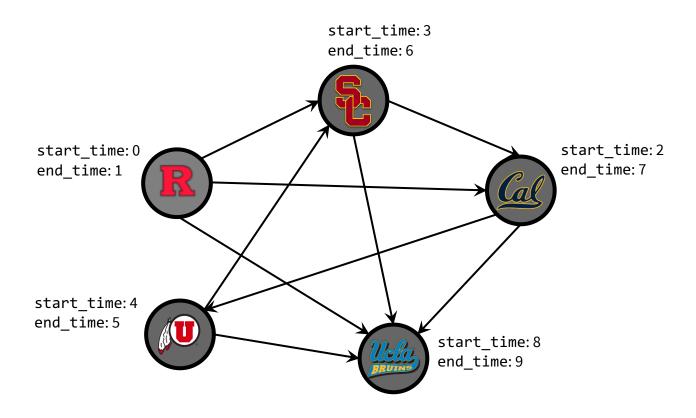


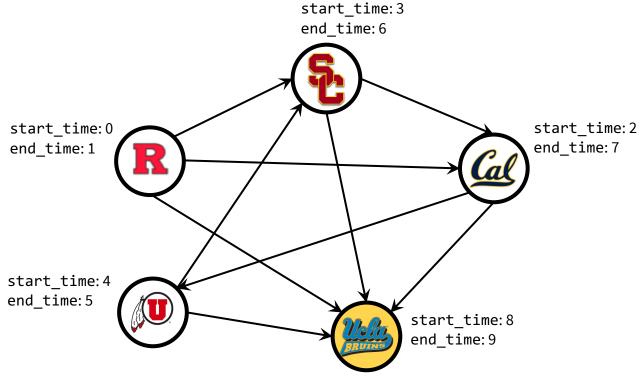


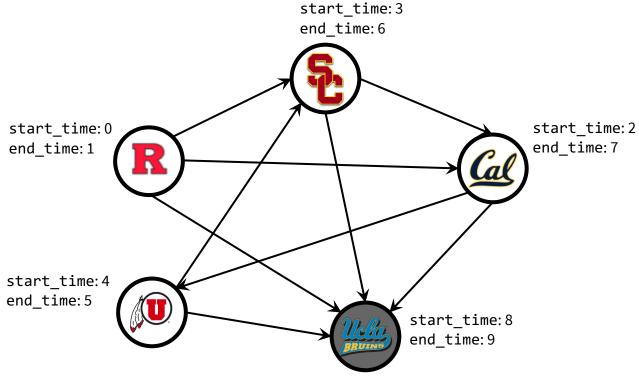


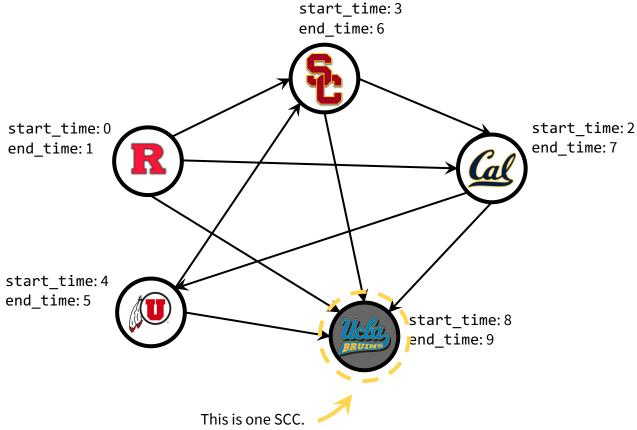


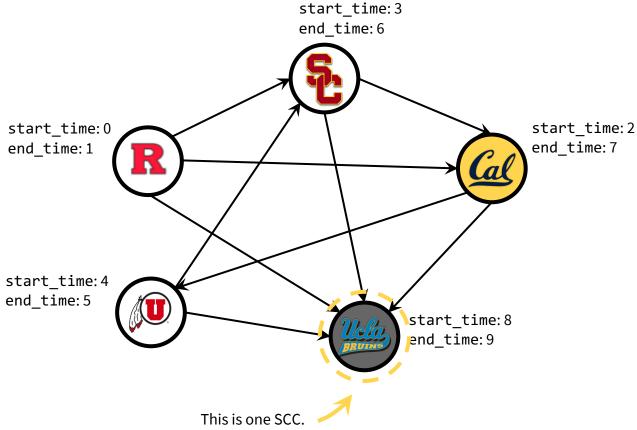
2. Reverse all of the edges.

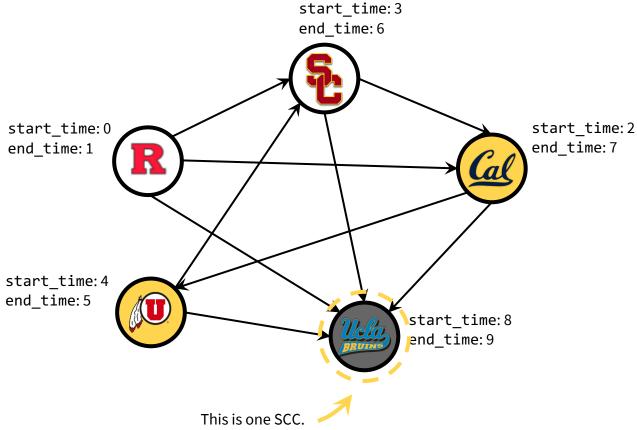


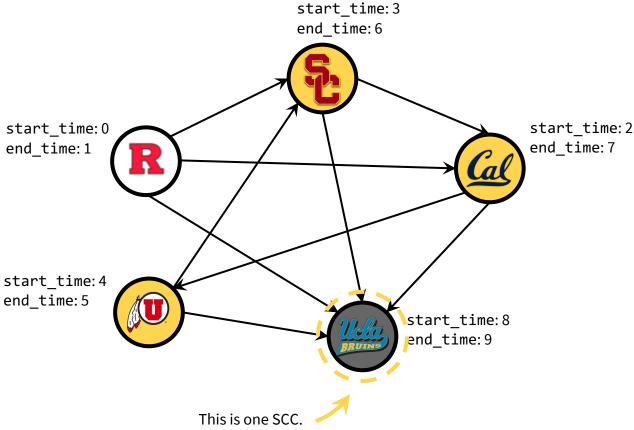


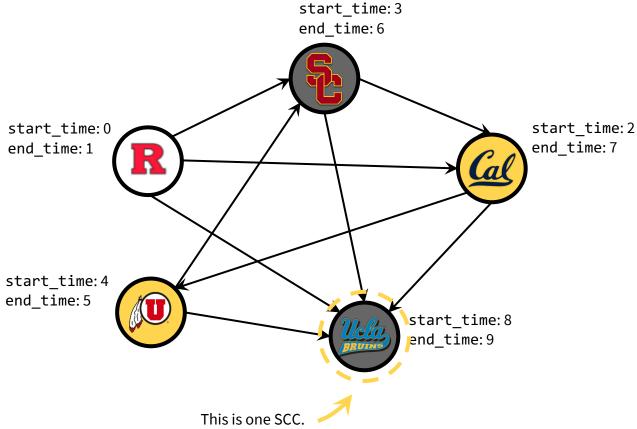


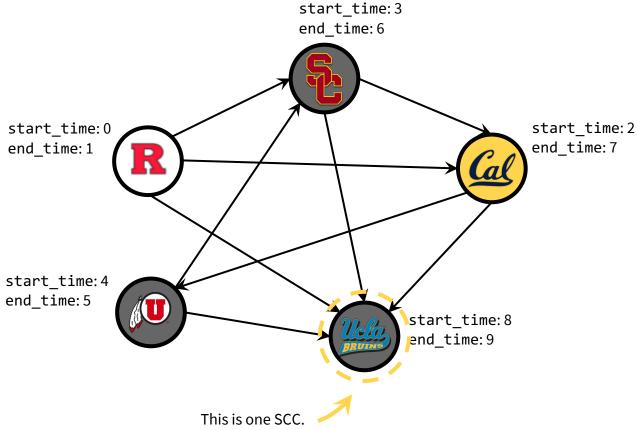


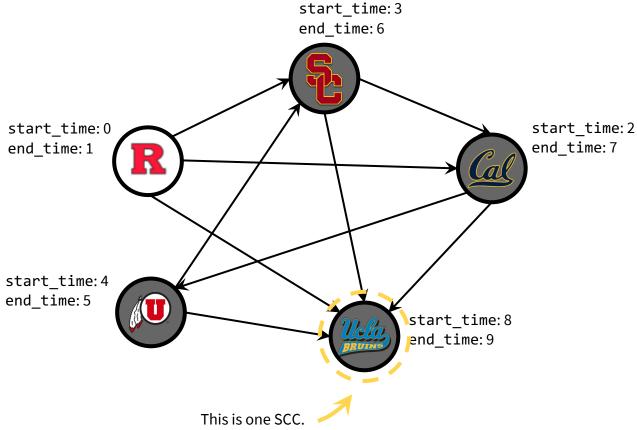


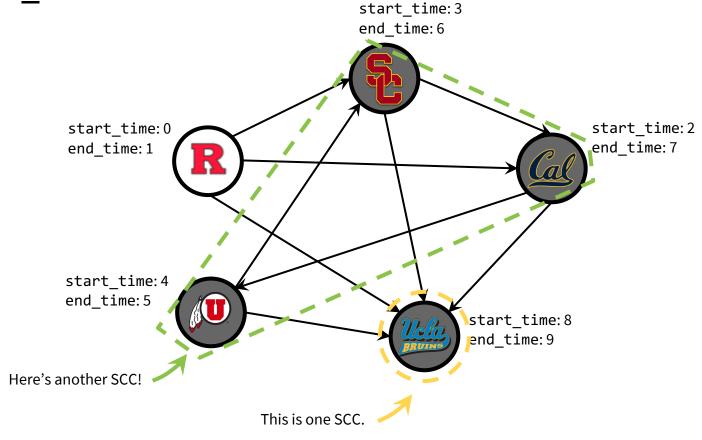


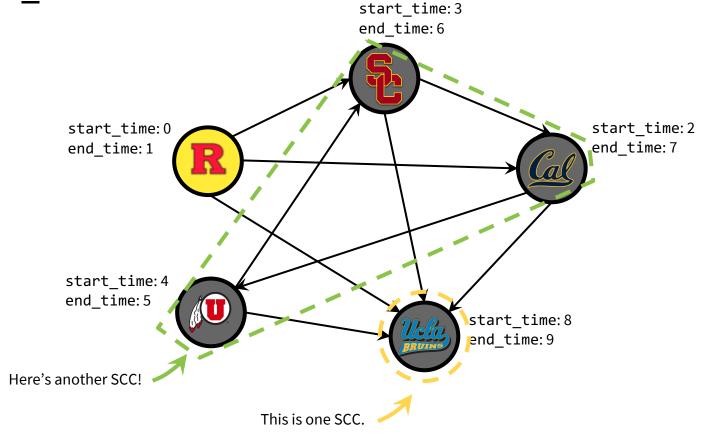


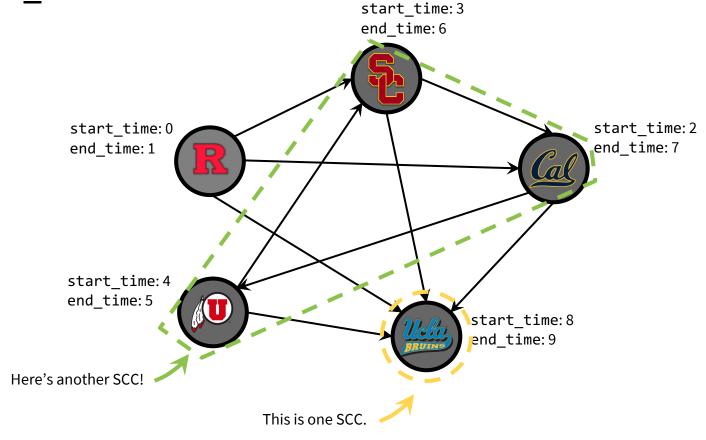


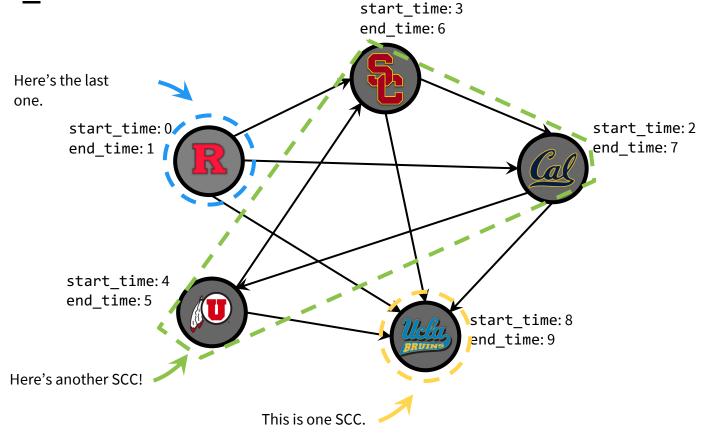












Why Kosaraju's Algorithm Works

Whoa. How did that work?

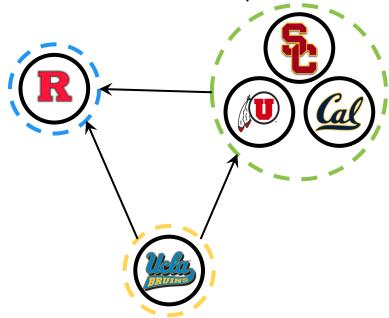
We explain by answering two questions:

- 1) Why do we use Depth First Search
- 2) In the second DFS, why do we start from vertex with largest end_time in each round.

Why do we use Depth First Search

Lemma 1: The SCC metagraph is a directed acyclic graph (DAG).

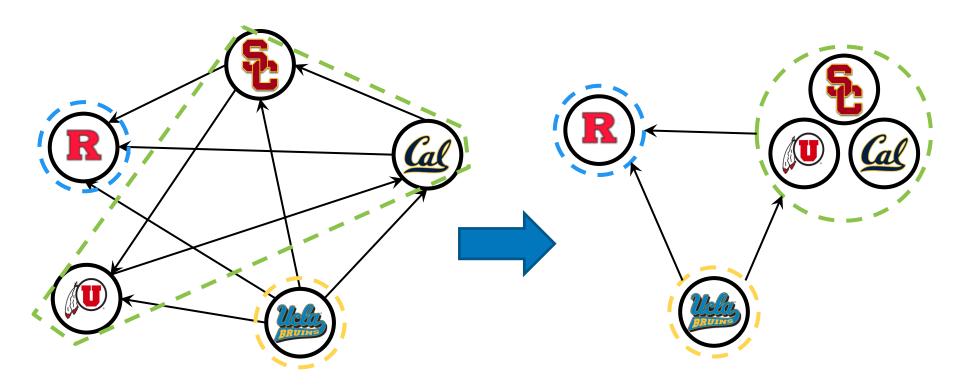
Intuition: If not, then two SCCs would collapse into one.



Why do we use Depth First Search

Review: DAG, a directed graph with no directed cycles.

Metagraph: connect a pair of meta vertex as long as there exists an edge between any vertex pair from each meta vertex.

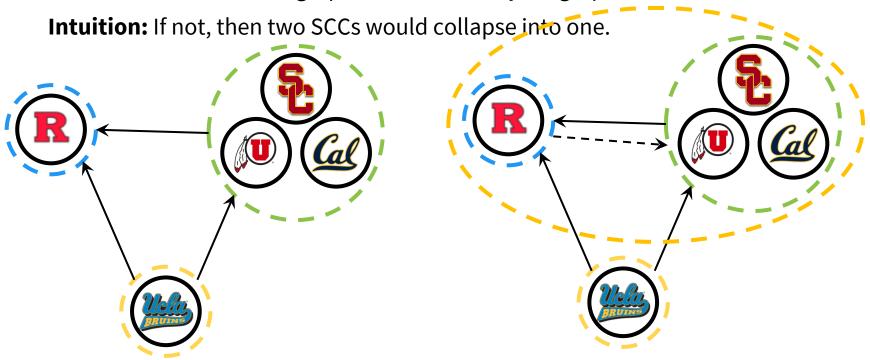


Original Graph SCC MetaGraph

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Why do we use Depth First Search

Lemma 1: The SCC metagraph is a directed acyclic graph (DAG).



Proof: assume there exist a cycle path between two meta vertex, then they can be merged into a bigger metavertex.

Why do we use Depth First Search

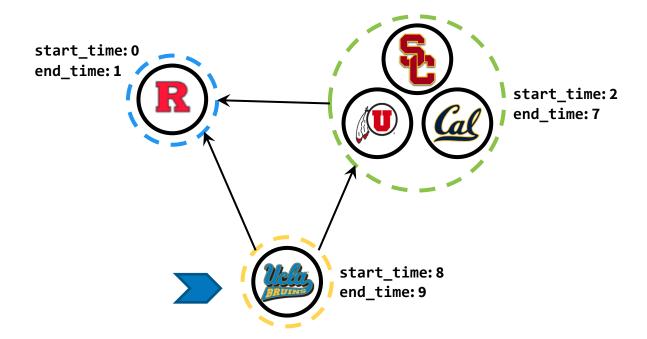
Let the **end time** of a SCC be the largest end time of any element of that SCC.

Let the **starting time** of a SCC be the smallest starting time of any element of that SCC.



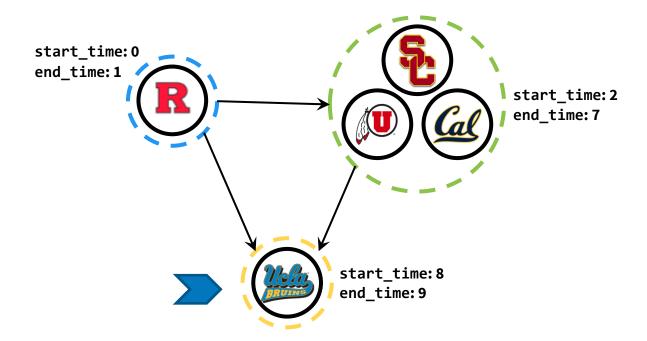
Why do we use Depth First Search

The main idea leverages the fact that vertex in the **SCC** metagraph with the largest end_time has no incoming edges.



Why do we use Depth First Search

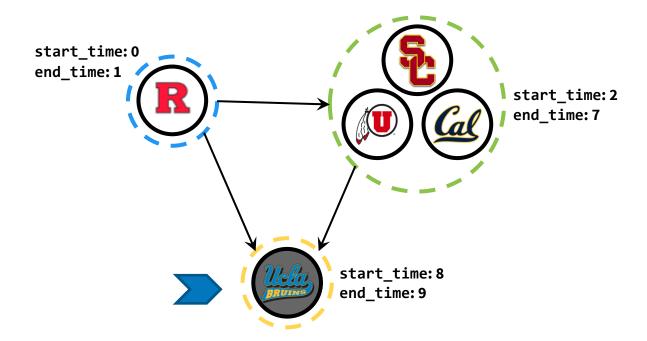
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Running dfs on that vertex finds exactly that component.

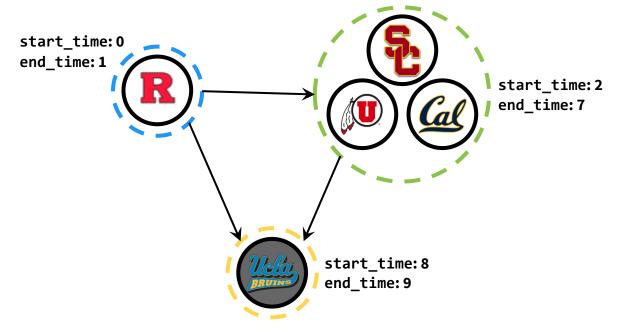


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Same argument for the rest.

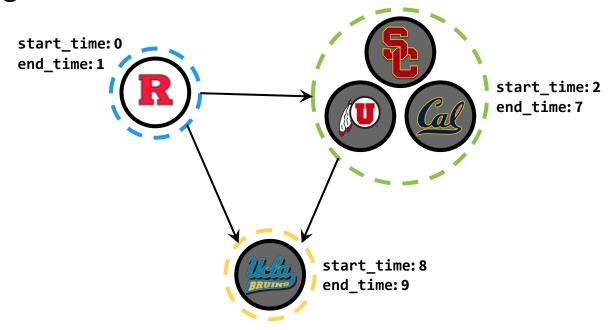


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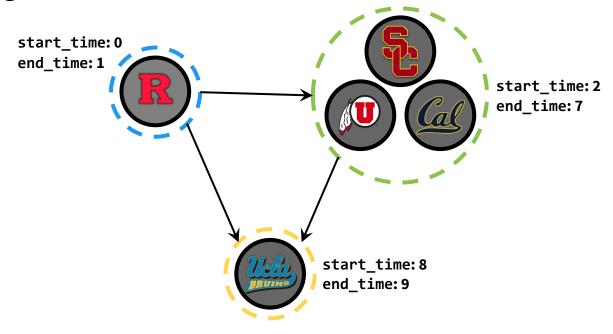


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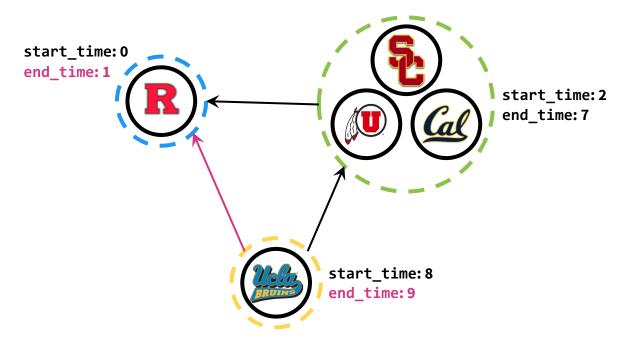
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Why do we use start with the largest end_time vertex

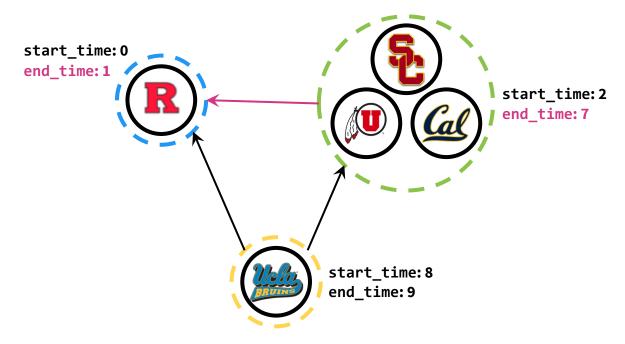
Claim: For each edge (u, v) in the SCC metagraph where $u \in C_1$ and $v \in C_2$, end_time of C_1 must be larger than end_time of C_2 .



In this way, the second DFS will converge with an SCC before entering another SCC (because all edges are inversed in the second DFS).

Why do we use start with the largest end_time vertex

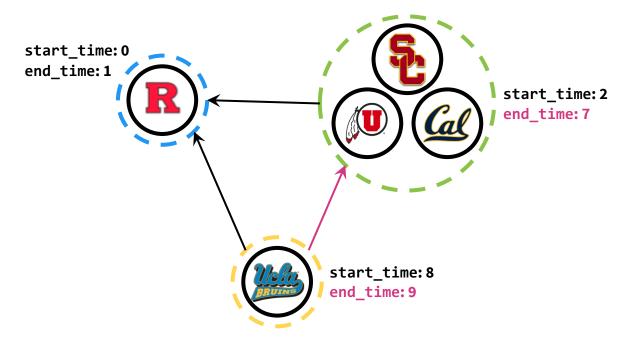
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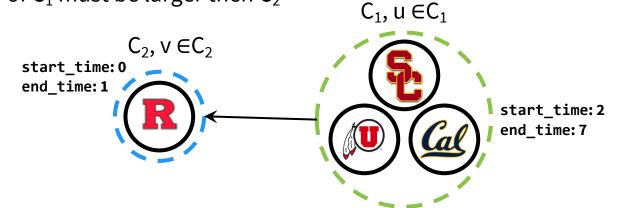
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Why do we use start with the largest end_time vertex

Claim: For each edge (u, v) in the SCC metagraph where $u \in C_1$ and $v \in C_2$, end_time of C_1 must be larger than end_time of C_2 .

Intuition: Suppose the first DFS started with C_2 , it will only finish C_2 before entering C_1 (because there can be no edge from C_2 to C_1 , or else C_2 and C_1 can be merged as a single SCC), so the end_time of C_1 must be larger then C_2 .

Suppose the first DFS started with C_1 , because of the existence of the edge (u,v), the algorithm will enter C_2 from u and reach v at some time during DFS; because C_2 is an SCC, all nodes in C_2 are reachable from v, so the DFS will visit all node in C_2 ; because there is no edge from C_2 to C_1 (same as above), so the DFS will only finish all nodes in C_2 and then trace back to C_1 . As a result, the end_time of C_1 must be larger then C_2



The runtime of Kosaraju's algorithm is O(V + E).

```
Runtime for the first DFS is O(|V| + |E|).
```

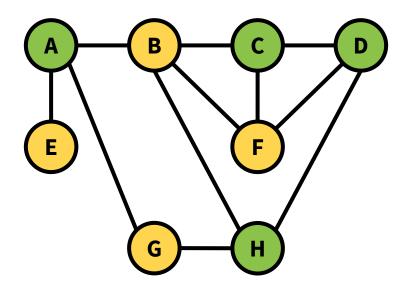
Runtime for reversing the graph is O(|V| + |E|).

Runtime for the second DFS is O(|V| + |E|).

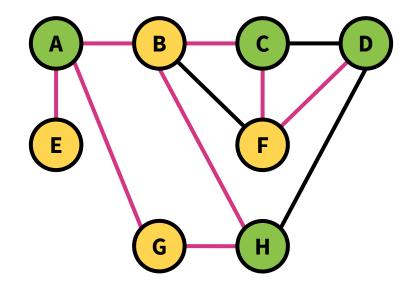
We can find connected components and SCCs in the same (asymptotic) runtime!

A **cut** is a partition of the vertices into two nonempty parts.

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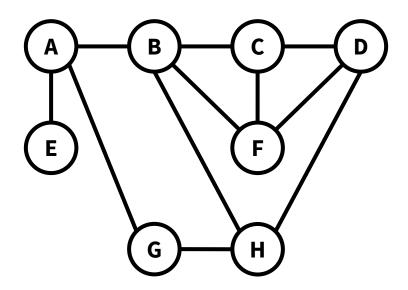
A **cut** is a partition of the vertices into two nonempty parts. e.g. This is the cut "{A, C, D, H} and {B, E, F, G}".



Edges that **cross the cut** go from one part to the other.

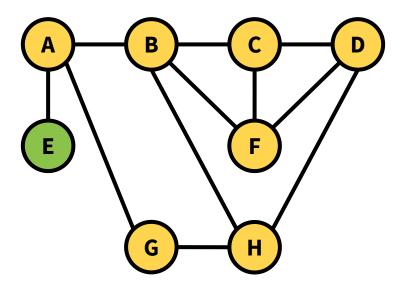
e.g. These edges cross the cut.

A **global minimum cut** is a cut that has the fewest edges possible crossing it.



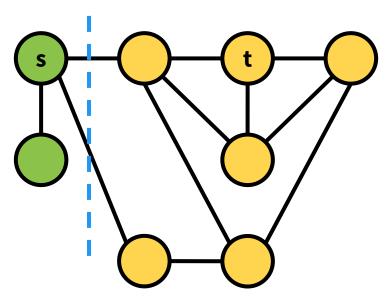
A **global minimum cut** is a cut that has the fewest edges possible crossing it.

e.g. The global minimum cut is "{A, B, C, D, F, G, H} and {E}".



Later this semester, we'll talk about **minimum s-t cuts**, which separate specific vertices **s** and **t**.

e.g. The minimum s-t cut is this cut.



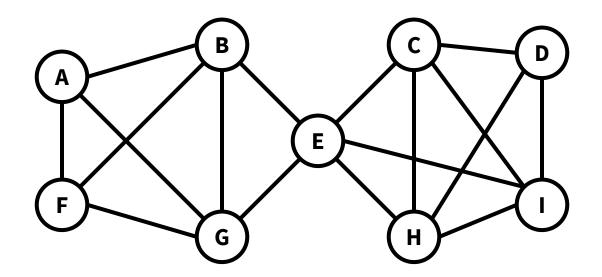
Karger's Algorithm finds global minimum cuts.

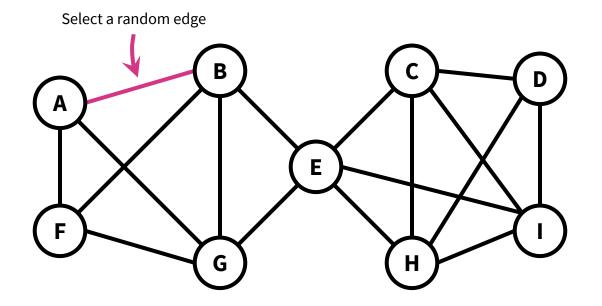
It's a Monte Carlo randomized algorithm! Unlike quicksort, which is always correct but sometimes slow, Karger's algorithm is always fast but sometimes Incorrect.

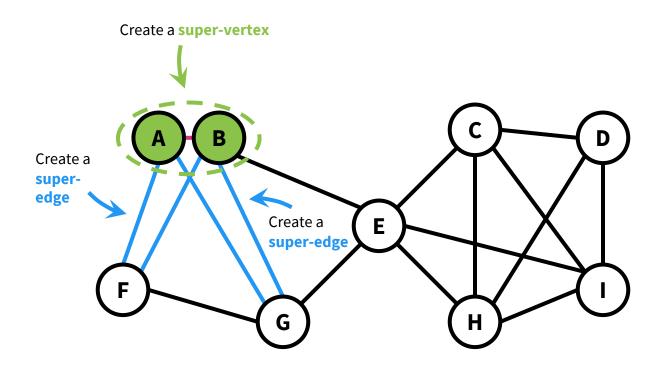
For all inputs A, quicksort returns a sorted list. For all inputs A, with high probability over the choice of pivots, quicksort runs fast.

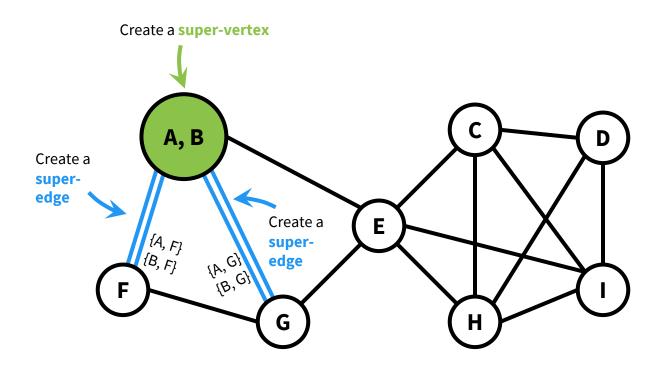
For all inputs G, karger runs fast. For all inputs G, with high probability over the randomness in the algorithm, karger returns a minimum cut.

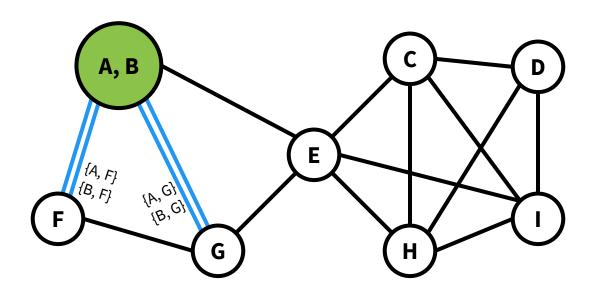
The general idea is to pick random edges to "contract" until there are a minimal number of vertices and edges left.

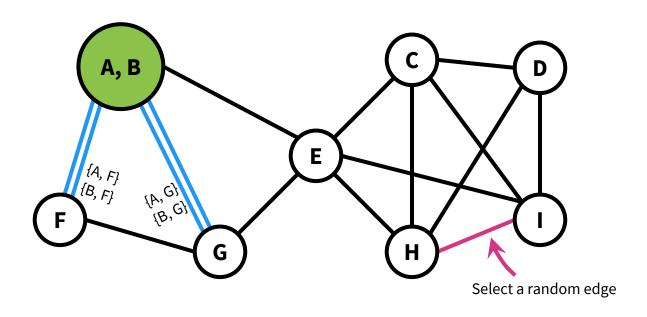


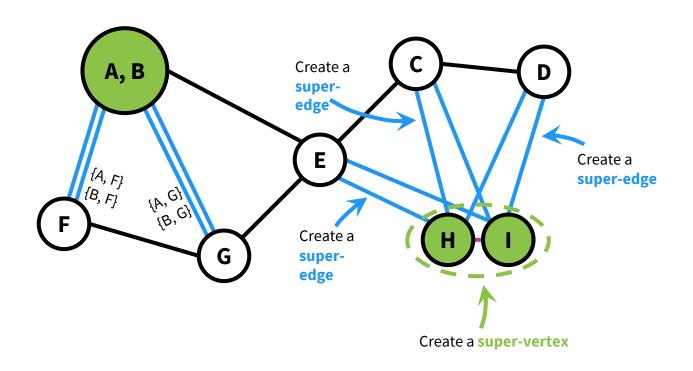


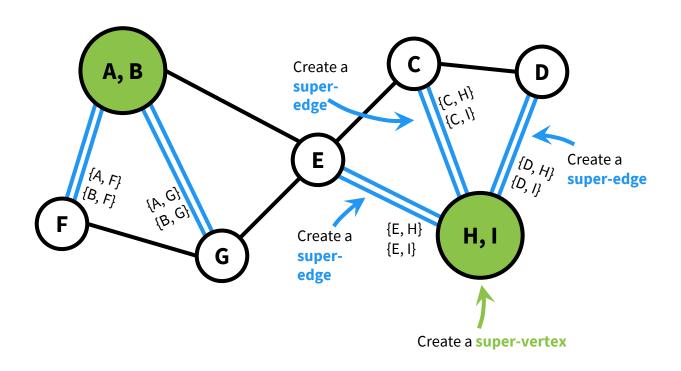


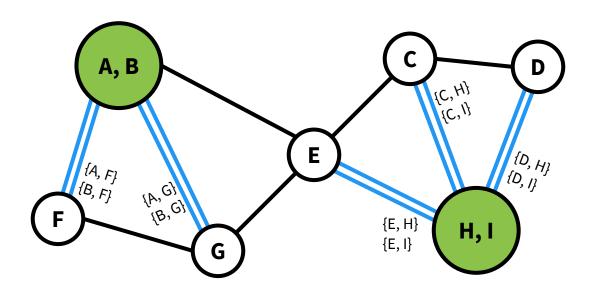


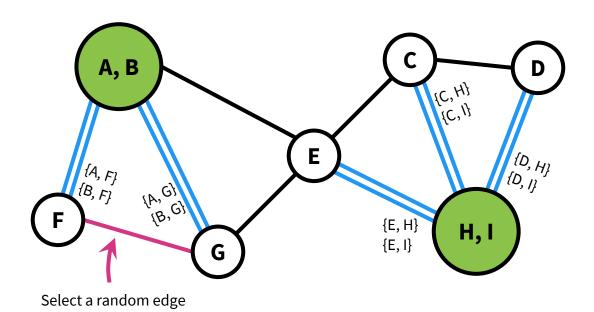


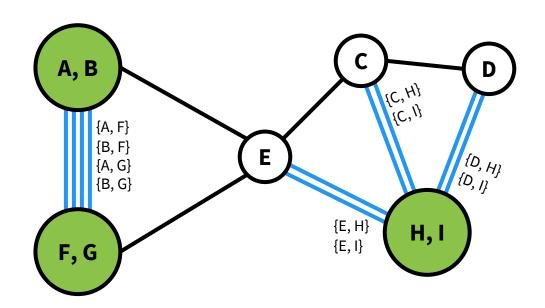


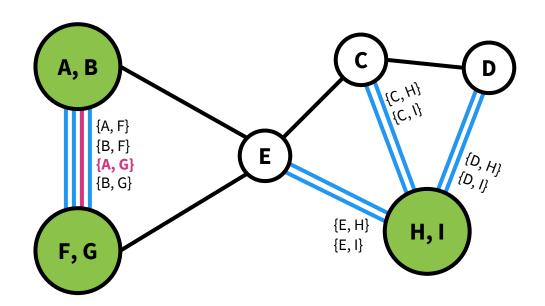


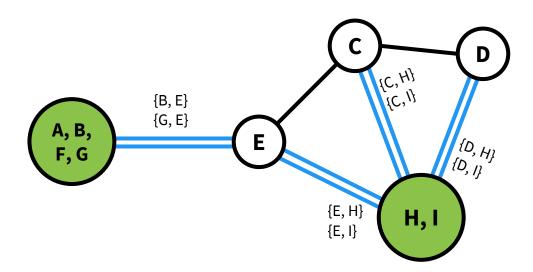


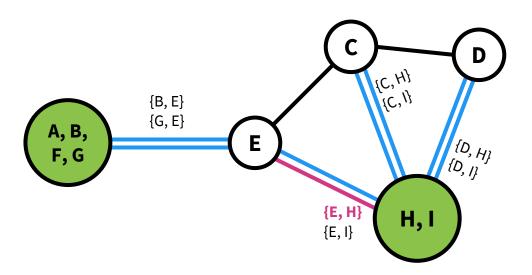


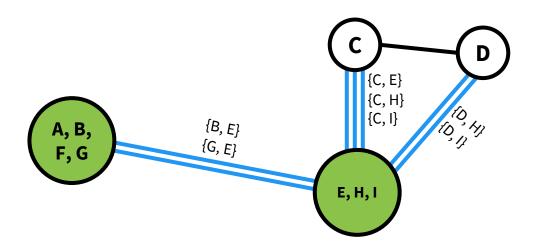


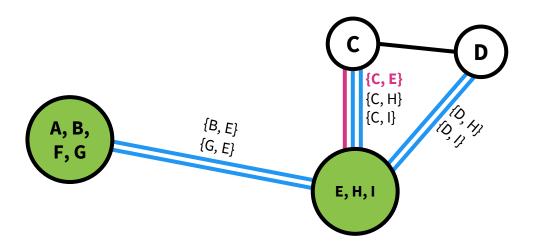


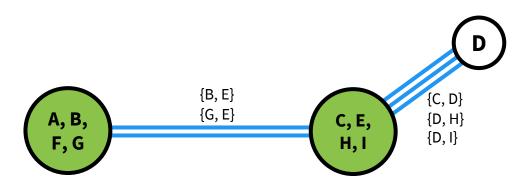


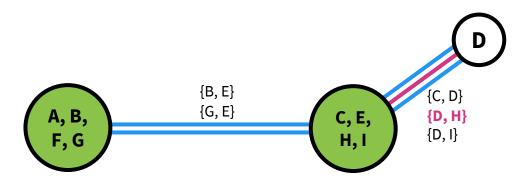






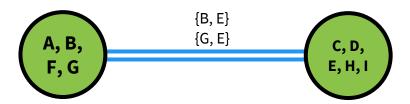






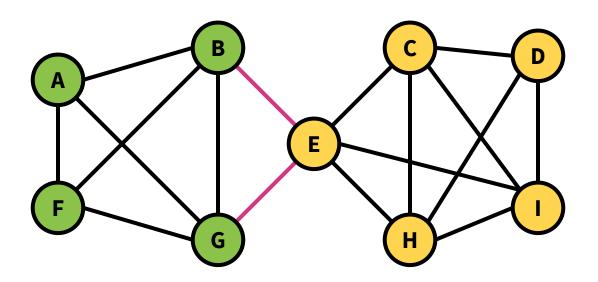
The minimum cut is given by the remaining super-vertices.

e.g. The cut is "{A, B, F, G} and {C, D, E, H, I}"; the edges that cross this cut are {B, E} and {G, E}.



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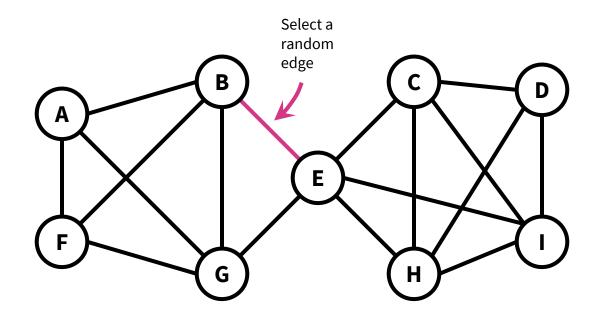
e.g. The cut is "{A, B, F, G} and {C, D, E, H, I}"; the edges that cross this cut are {B, E} and {G, E}.



```
algorithm karger(G=(V,E)):
                  G' = {supervertex(v) for v in V}
These are the
                 'E<sub>u'v'</sub> = {(u,v)} for (u,v) in E
super-edges.
                  E_{u'v'} = \{\} for (u,v) not in E
                  F = \{\{(u,v)\} \text{ for } (u,v) \text{ in } E\}
The while loop
                  while |G'| >= 2:
executes |V| - 2
                     \{(u,v)\}\ = uniform random edge in F
times.
                     merge_supervertices(u, v) This takes O(|V|).
Removes all edges
                 \rightarrow F = F \ E_{\mu'\nu'}
in the super-edge
                  return cut of the remaining super-vertices
between super-
vertices u' and v'.
               algorithm merge_supervertices(u, v):
                  x' = supervertex(u' U v')
                                                             u' is the super-vertex containing u;
                  for w' in G' \setminus {u',v'}:
                                                             v' is the super-vertex containing v.
                     E_{x'w'} = E_{u'w'} \cup E_{v'w'}
                     Remove u' and v' from G' and add x'
```

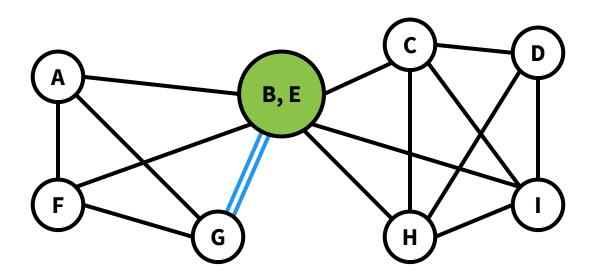
We got really lucky!

e.g. Suppose we had chosen this edge.

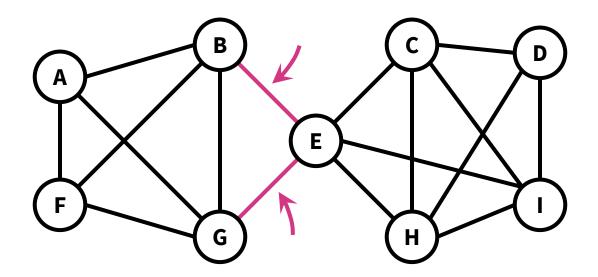


We got really lucky!

e.g. Suppose we had chosen this edge. Now there's no way to return a cut that separates B and E.



If fact, if Karger's algorithm ever randomly selects **edges in the min-cut**, then it will be incorrect.



The probability that Karger's algorithm returns a minimum cut is ...

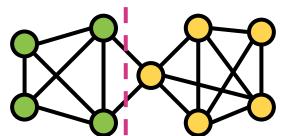
$$\geq 1/\binom{n}{2}$$

where n is the number of nodes.

Proof:

Suppose S^* is a min-cut and suppose we select edges $e_1, e_2, ..., e_{n-2}$.

Then P(karger returns S*) = P(no e_i crosses S*) = P(e_1 doesn't cross S*) S* \times P(e_2 doesn't cross S* | e_1 doesn't cross S*)



 \times P(e_{n-2} doesn't cross S* | e₁, ..., e_{n-3} doesn't cross S*)

The probability that Karger's algorithm returns a minimum cut is ...

$$\geq 1/\binom{n}{2}$$

Proof, cont.:

Suppose, after j-1 iterations, karger hasn't messed up yet! What's the probability of messing up now?

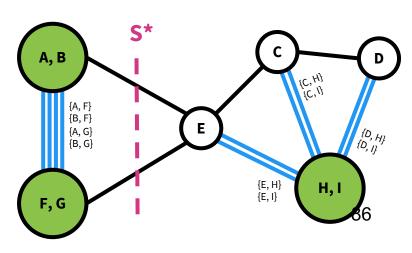
Suppose there are k edges that cross **S***.

All remaining vertices must have degree at least k (otherwise there would be a smaller cut).

So there are at least (n-j+1)k/2 total edges.

So the probability that karger chooses one of the k edges crossing **S*** at step j is at most

$$\frac{k}{\frac{(n-j+1)k}{2}} = \frac{2}{n-j+1}$$



The probability that Karger's algorithm returns a minimum cut is ...

$$\geq 1/\binom{n}{2}$$

Proof, cont.:

Suppose S^* is a min-cut and suppose we select edges $e_1, e_2, ..., e_{n-2}$.

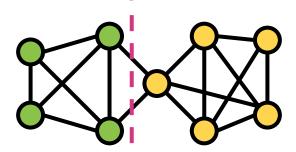
Then P(karger returns
$$S^*$$
) = P(e_1 doesn't cross S^*)

$$\times$$
 P(e₂ doesn't cross S* | e₁ doesn't cross S*)

• •

$$\times$$
 P(e_{n-2} doesn't cross S* | e₁, ..., e_{n-3} doesn't cross S*)

$$\geq \frac{(n-2) \quad (n-3) \quad (n-4) \quad (n-5) \quad (n-6)}{n \quad (n-1) \quad (n-2) \quad (n-3) \quad (n-4) \quad \dots \quad -\frac{4}{6} \quad \frac{3}{5} \quad \frac{2}{4} \quad \frac{1}{3}$$



$$= 2/(n(n-1))$$

$$= 1/(nC2)$$

1/(nC2) isn't all that great ...

For our example of n = 9, 1/(9C2) = 0.028.

Suppose we want to find the min-cut with probability 0.9.

What can we do? 🧐

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For our example of n = 9, 1/(9C2) = 0.028.

Suppose we want to find the min-cut with probability 0.9. What can we do? 🤔

How many times T do we need to repeat karger to obtain this probability?

Note that if P(find the min-cut after 1 time) $\geq 1/(nC2)$, then P(don't find the min-cut after 1 time $\leq 1 - 1/(nC2)$

P(find the min-cut after T times) ≥ 0.9

 \Leftrightarrow P(don't find the min-cut after T times) \leq 0.1.

P(don't find the min-cut after T times) = $(1 - 1/(nC2))^T$

$$\leq (e^{-1/(nC2)})^T = 0.1$$

T = (nC2) ln (1/0.1) times

Suppose we want to find the min-cut with probability p. Then we must repeat Karger $T = (nC2) \ln (1/(1-p))$ times.

 $T = (nC2) \ln (1/(1-p)) \text{ times} = O(|V|^2) \text{ times, so the overall runtime is } O(|V|^4).$

Treating 1-p as a constant.

If we use union-find data structures, then we can do better.

This might seem lousy, but then consider that enumerating over all possible cuts to find the min-cut requires $O(2^{|V|})$.

This is a huge improvement!

```
algorithm karger_loop(G=(V,E), threshold):
    cur_min_cut = None
    n = V.length, p = threshold
    for t = 1 to (nC2)ln(1/(1-p)) :
        candidate_cut = karger(G)
        if candidate_cut.size < cur_min_cut.size:
            cur_min_cut = candidate_cut
    return cur_min_cut</pre>
```

Runtime: O(|V|4)

Key point: Whenever we have a Monte-Carlo algorithm with a small probability of success, we can boost the probability of success by repeating it a bunch of times and taking the best solution!

Many statistical machine learning algorithms work in this way!