Graph Algorithms II

Summer 2017 • Lecture 07/20

A Few Notes

Homework 3

Due tomorrow at 11:59 p.m. on Gradescope.

Homework 4

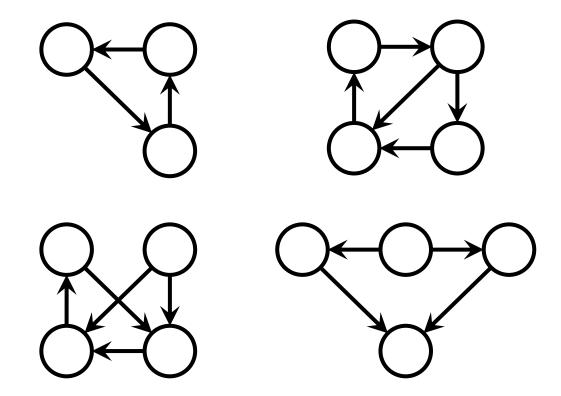
Released tomorrow.

Outline for Today

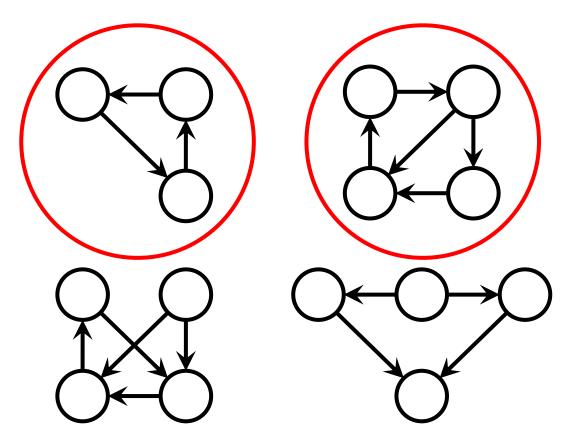
Graph algorithms

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

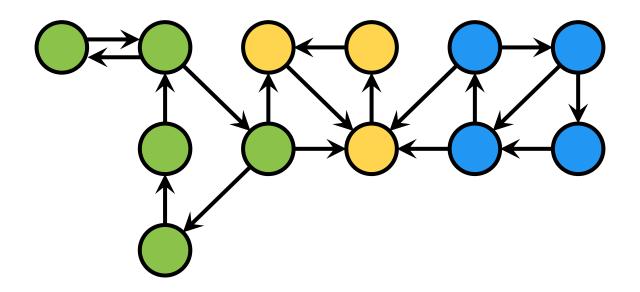
A directed graph G = (V, E) is strongly connected if, for all pairs of vertices u and v, there's a path from u to v and a path from v to u.



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?

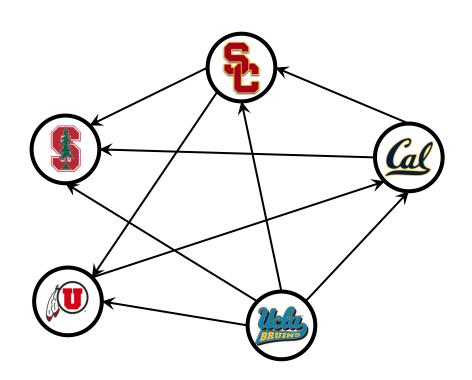
SCCs provide information about communities.

A computer scientist might want to decompose the Internet into SCCs to find related topics.

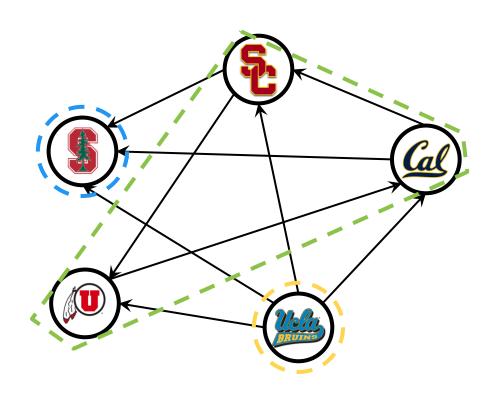
An economist might want to decompose labor market data into SCCs before making sense of it.

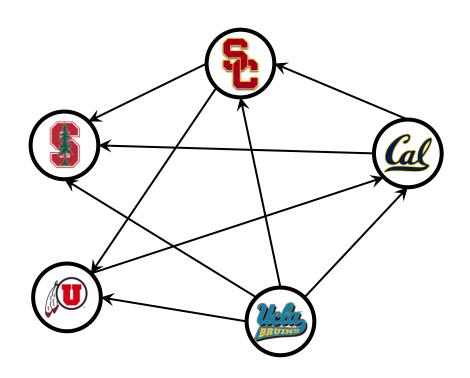
A football executive might want to determine which Pac-12 school should play in the Rose Bowl.

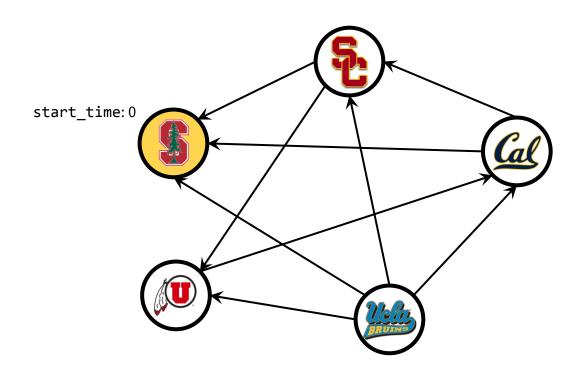
How many SCCs are in this graph? 🤔

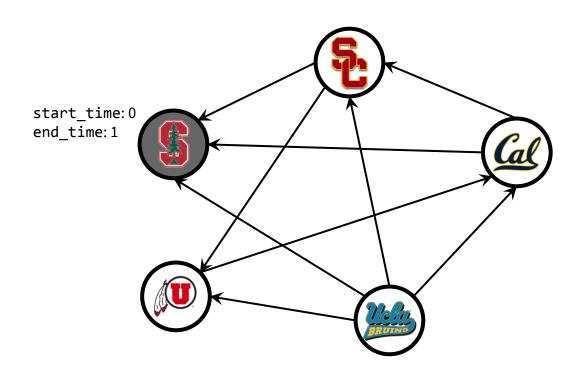


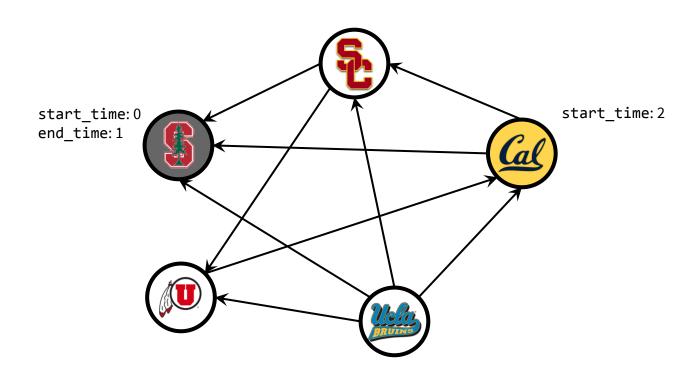
How many SCCs are in this graph? 9 3; let's find them!

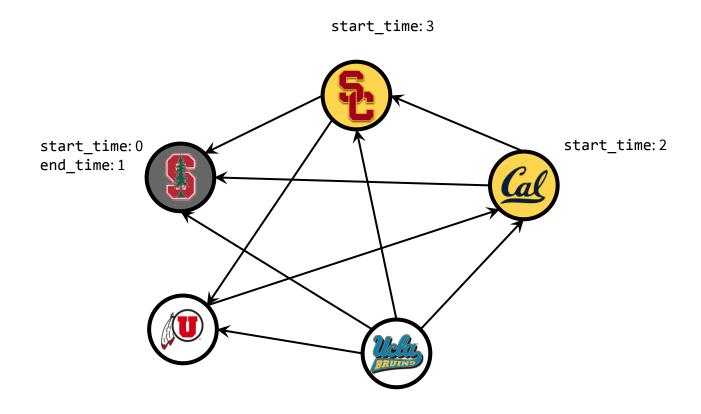


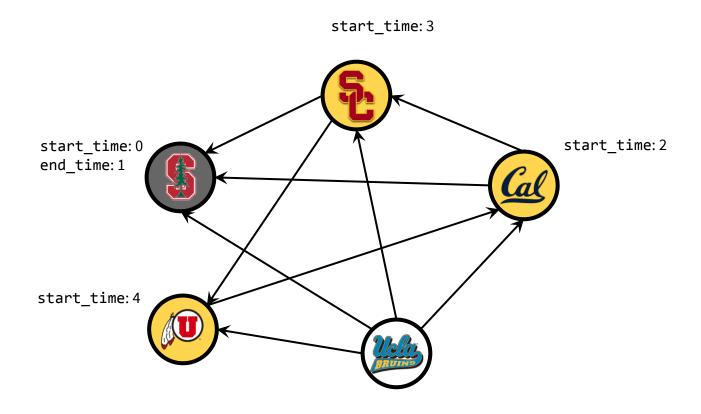


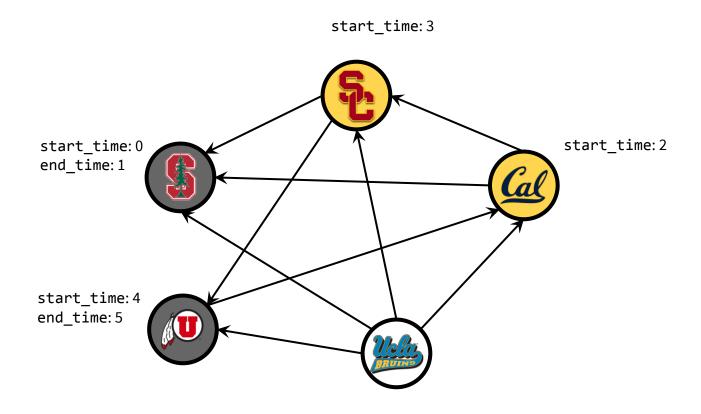


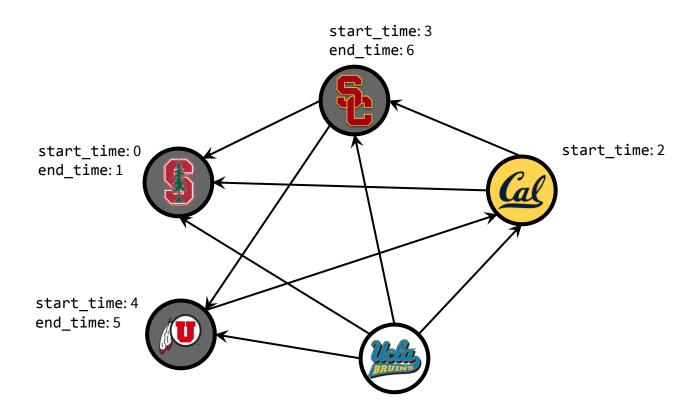


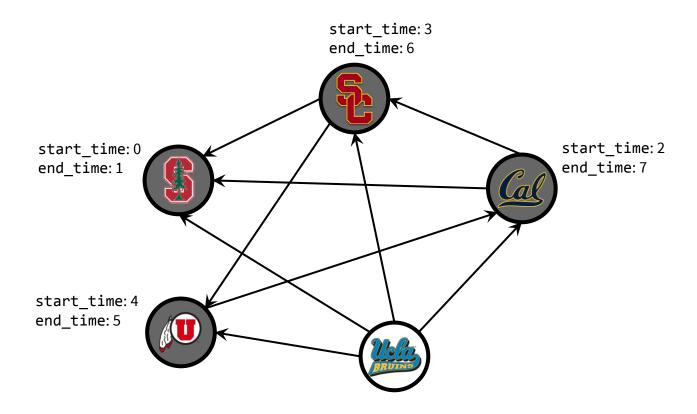


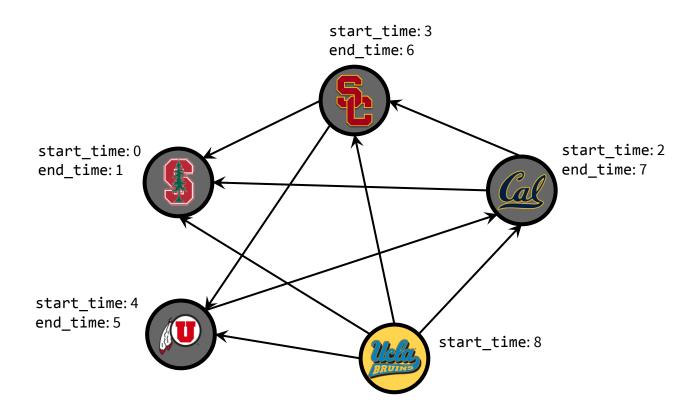


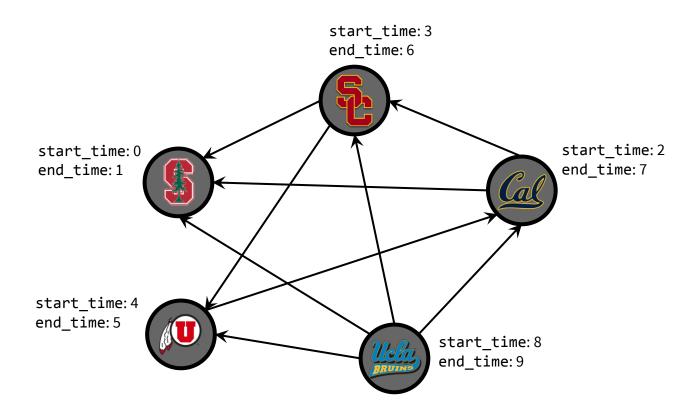




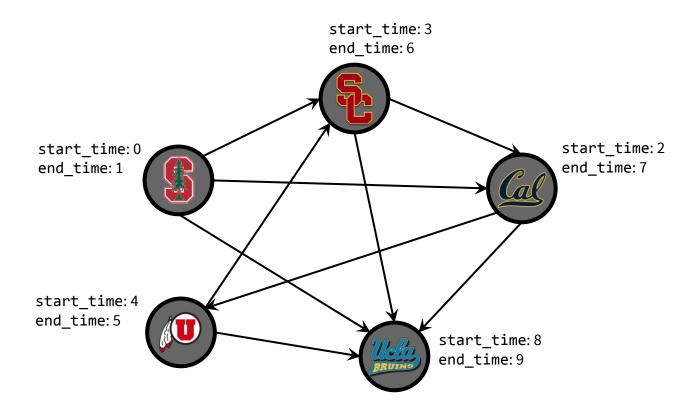


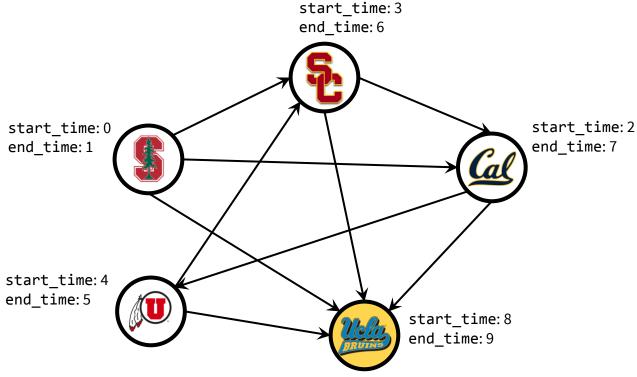


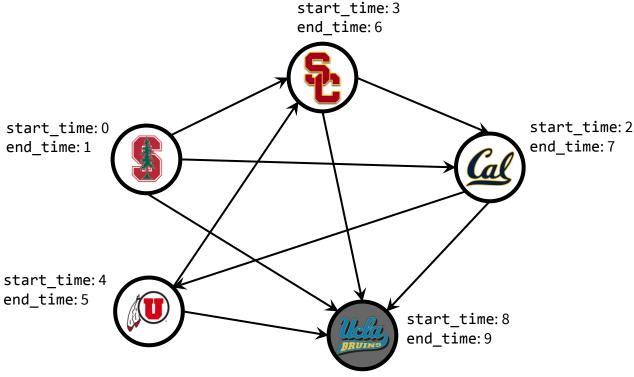


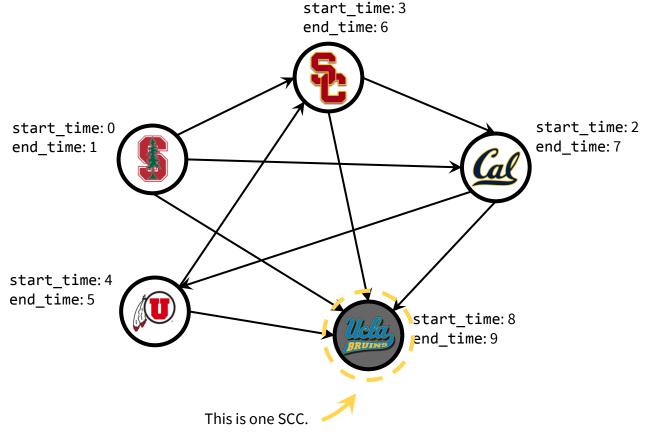


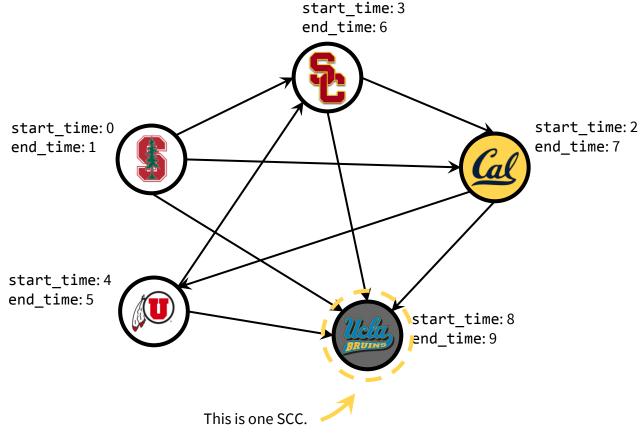
2. Reverse all of the edges.

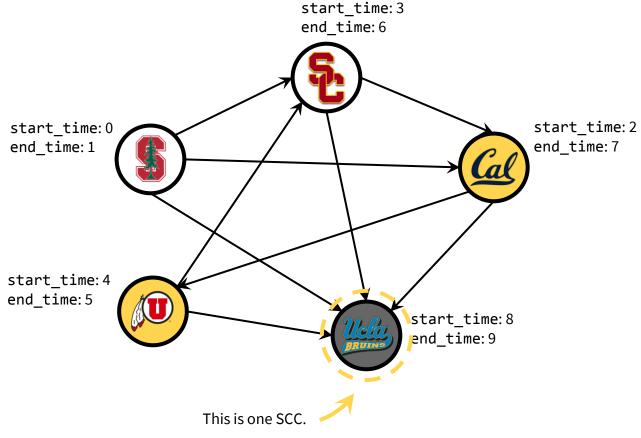


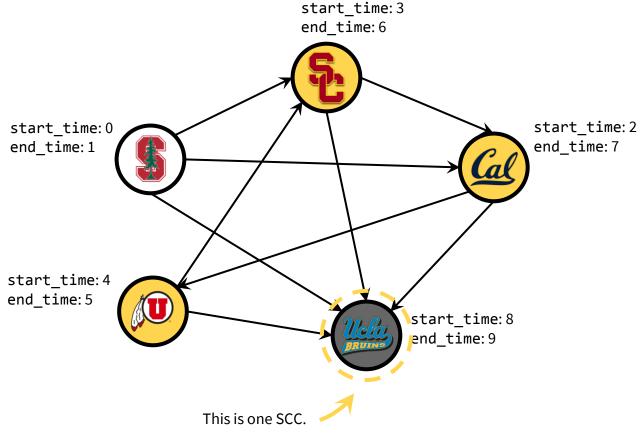


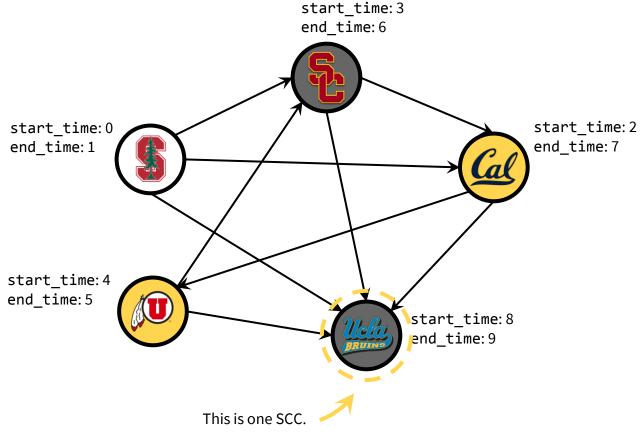


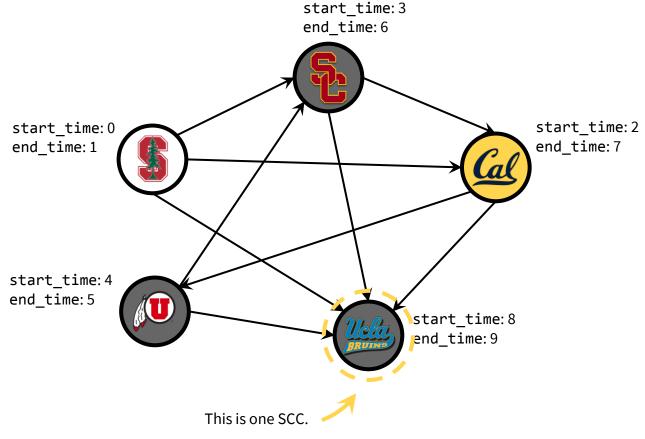


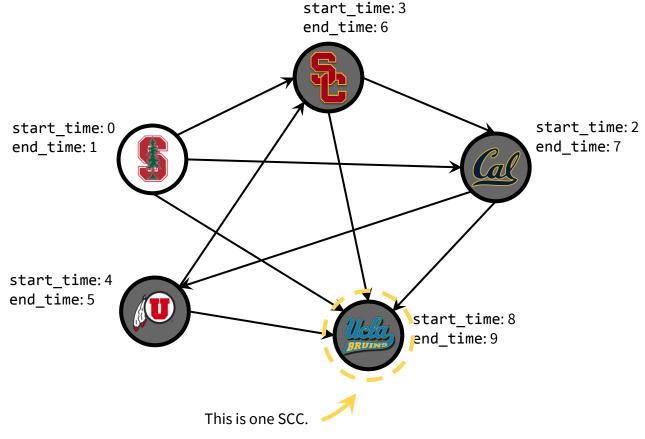


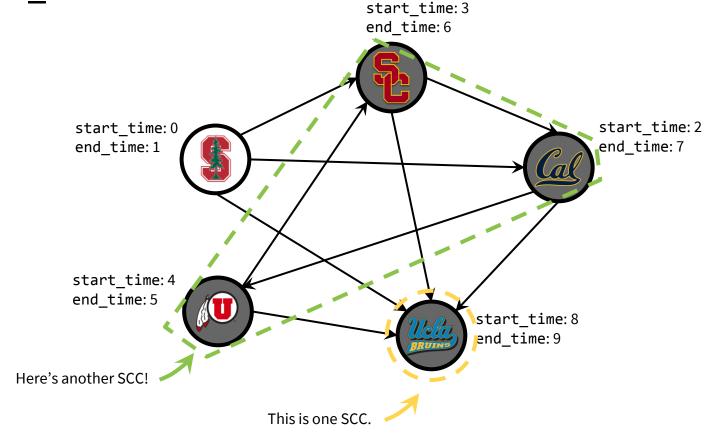


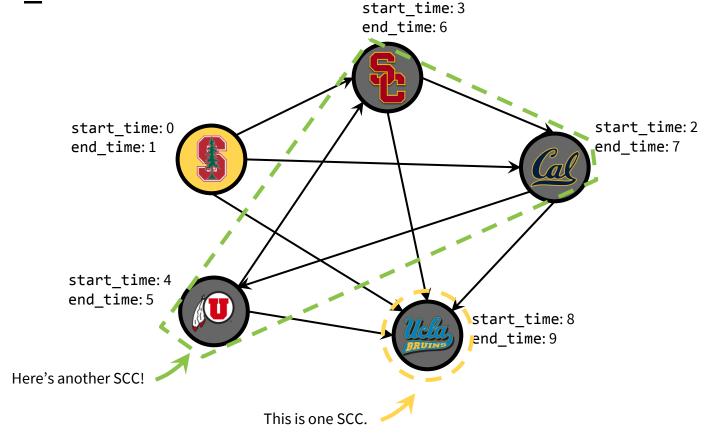


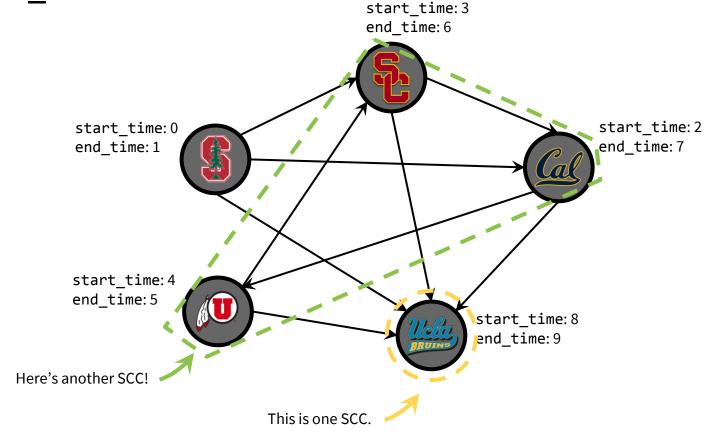


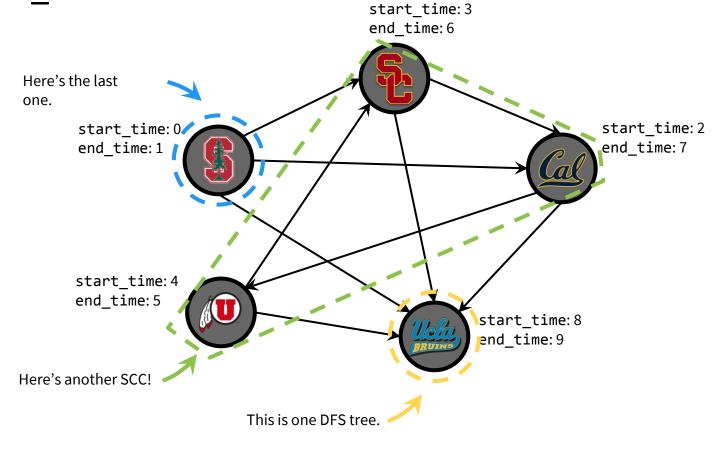








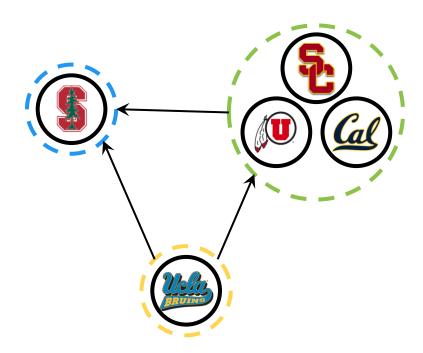




Whoa. How did that work?

Lemma 1: The SCC metagraph is a DAG.

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

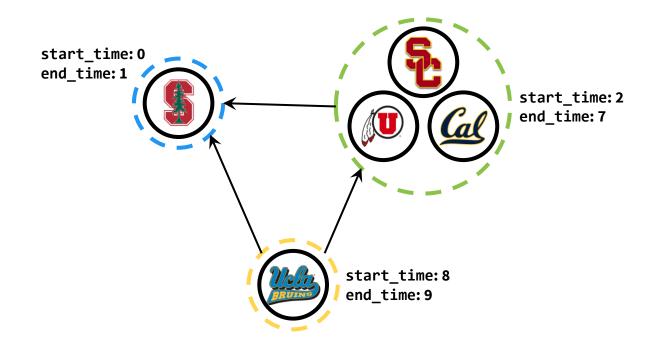


Let the **end time** of a SCC be the largest end time of any element of that SCC.

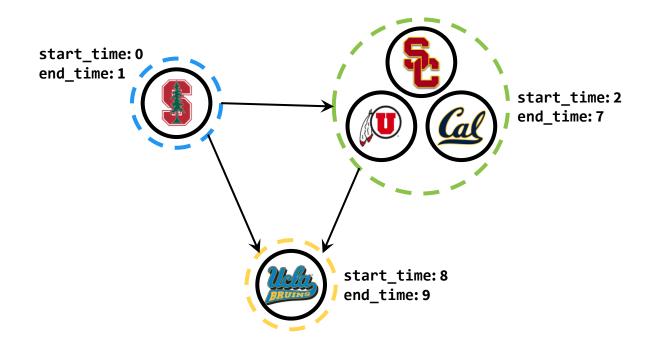
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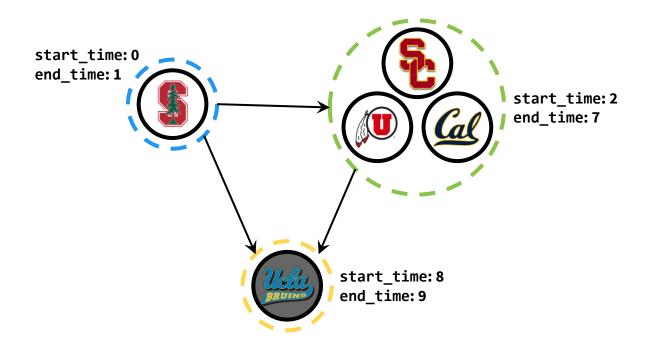
The main idea leverages the fact that vertex in the SCC metagraph with the largest end_time has no incoming edges.



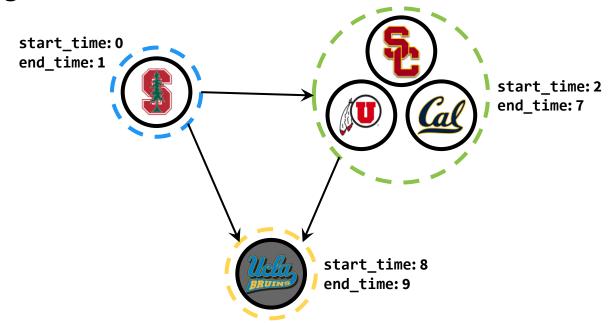
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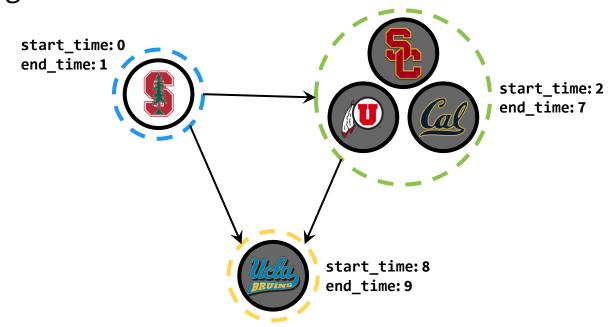
The main idea leverages the fact that vertex in the SCC metagraph with the largest end_time has no incoming edges. After reversing the edges, it has no outgoing edges. Running dfs on that vertex finds exactly that component.



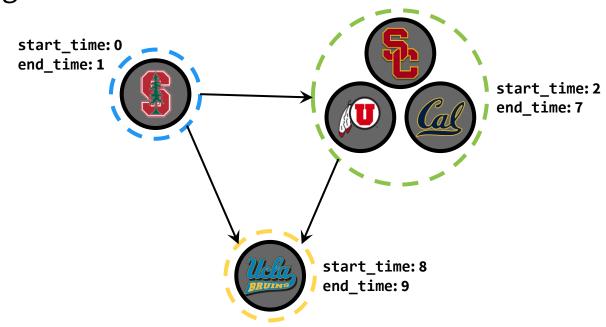
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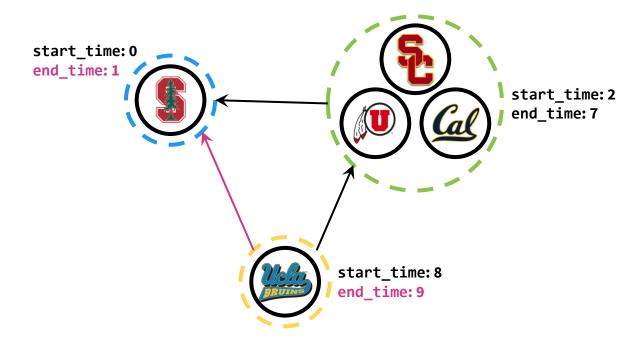


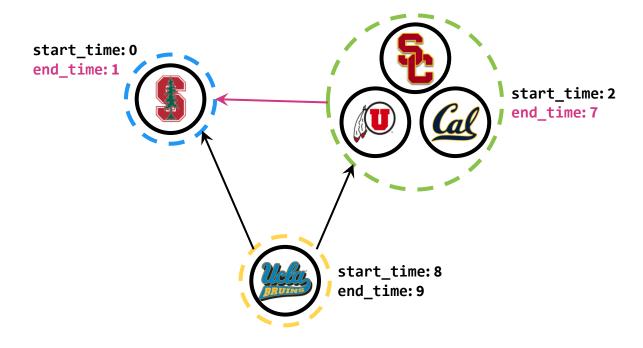
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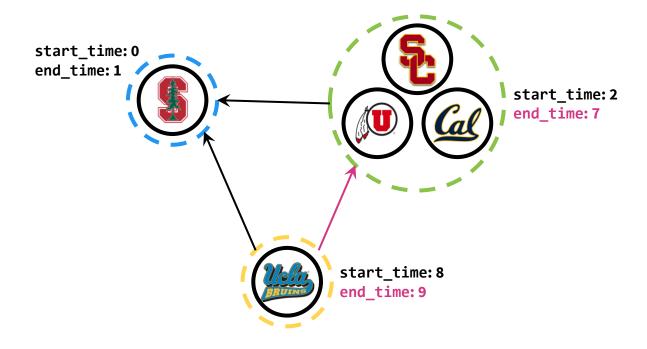


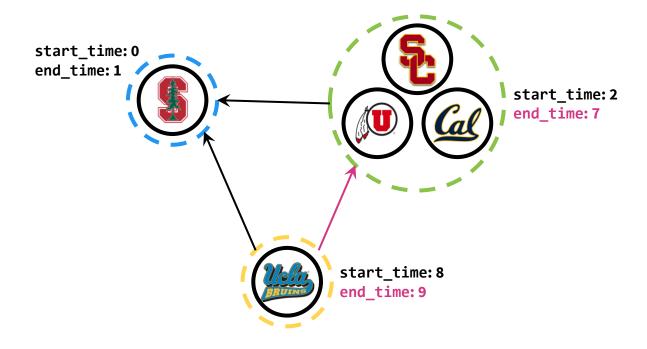
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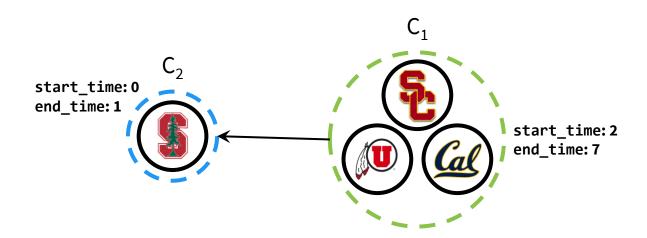






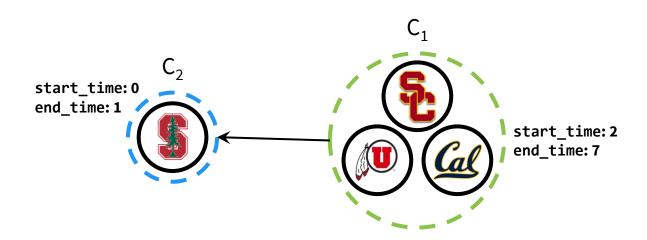






Claim: For each edge (u, v) in the SCC metagraph where $u \in C_1$ and $v \in C_2$, end_time of C_1 must be larger than end_time of C_2 .

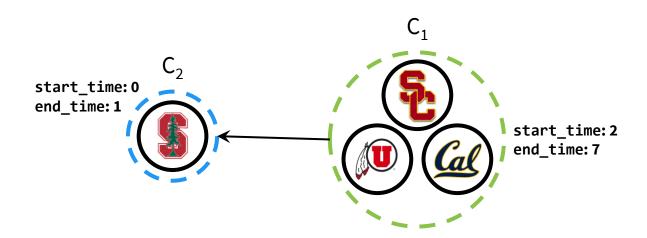
Intuition: In order for the end_time of C_1 to be smaller than the end_time of C_2 , all vertices in C_1 must have end_times smaller than at least one end_time of C_2 .



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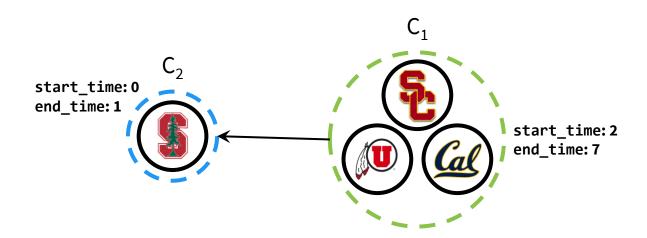
For this to occur, the first dfs must have marked all vertices in C_1 as "done" before at least one vertex in C_2 .



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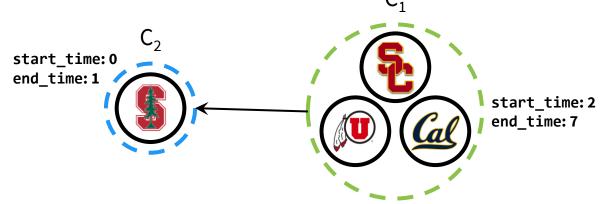
For this to occur, the first dfs must have marked all vertices in C₁ as "done" before at least one vertex in C₂. But this is impossible since the first dfs must have explored edge (u, v) before marking all vertices in C₁ as "done."



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 .

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For this to occur, the first dfs must have marked all vertices in C_1 as "done" before at least one vertex in C_2 . But this is impossible since the first dfs must have explored edge (u, v) before marking all vertices in C_1 as "done." Since C_2 is an SCC, all vertices in it are reachable from v; therefore, all must have an end_time smaller than u.



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|).

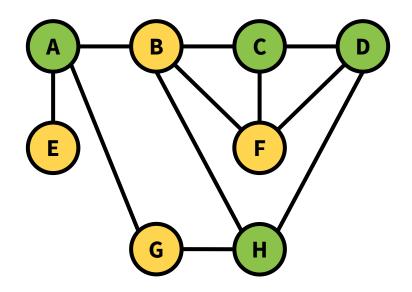
Runtime for the second DFS is O(|V| + |E|).

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

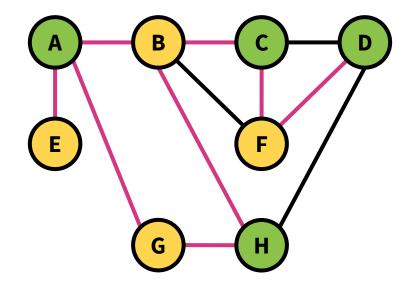
3 min break

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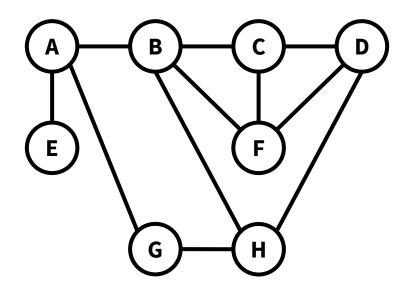


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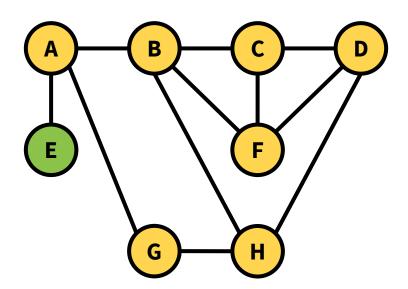
Edges that **cross the cut** go from one part to the other. e.g. These edges cross the cut.

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



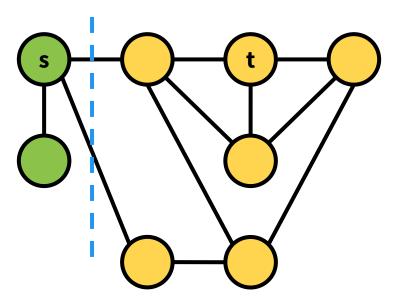
A **global minimum cut** is a cut that has the fewest edges possible crossing it.

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



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

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



Global Minimum Cuts

Why might we care about global minimum cuts?

Application: Image segmentation

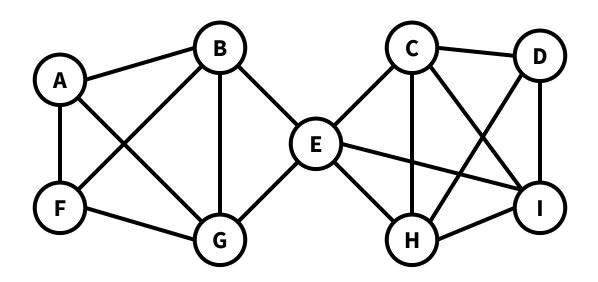
Karger's Algorithm finds global minimum cuts.

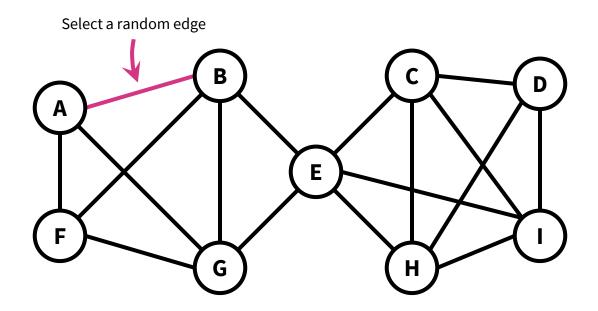
It's a Monte Carlo randomized algorithm! Unlike quicksort, which is always correct but sometimes slow, Karger's algorithm is always fast but sometimes Incorrect.

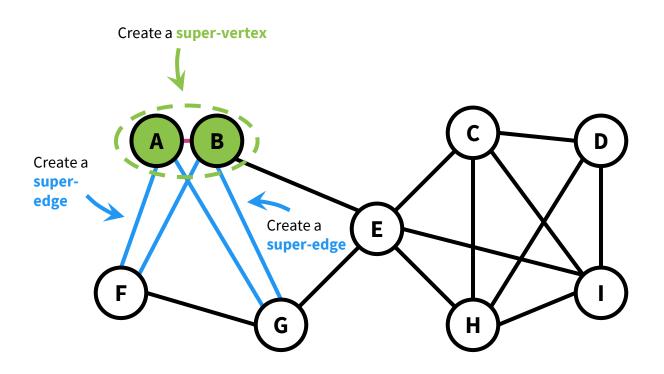
For all inputs A, quicksort returns a sorted list. For all inputs A, with high probability over the choice of pivots, quicksort runs fast.

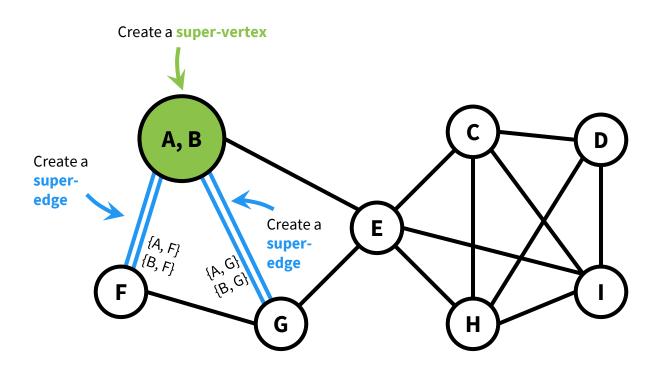
For all inputs G, karger runs fast. For all inputs G, with high probability over the randomness in the algorithm, karger returns a minimum cut.

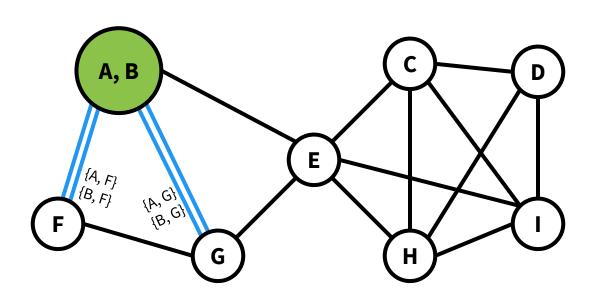
The general idea is to pick random edges to "contract" until there are a minimal number of vertices and edges left.

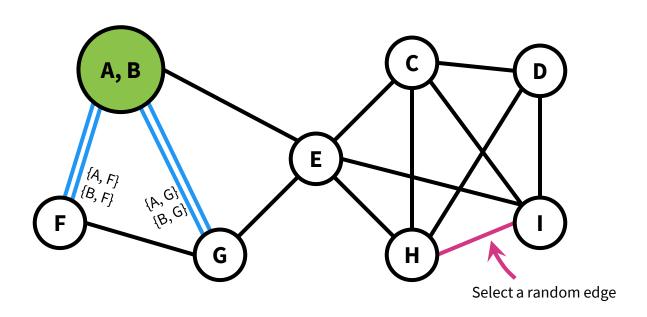


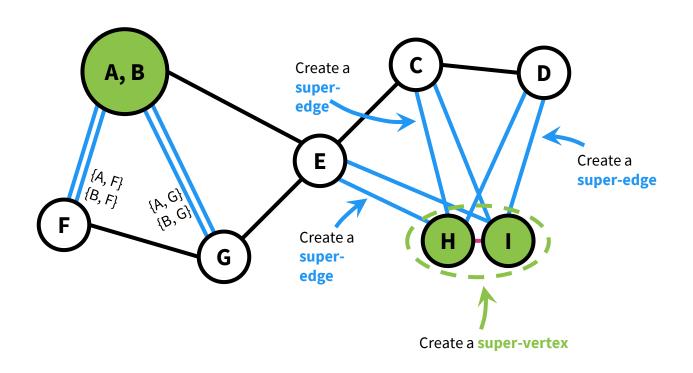


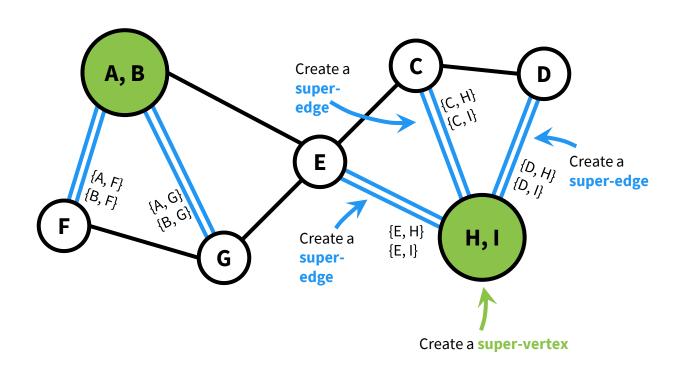


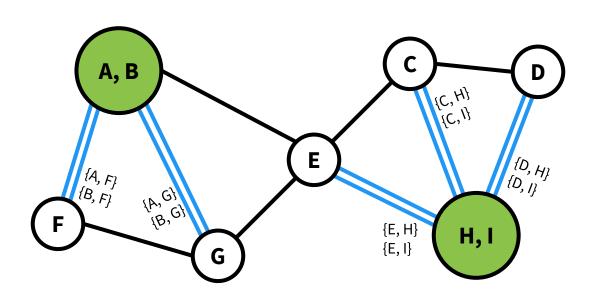


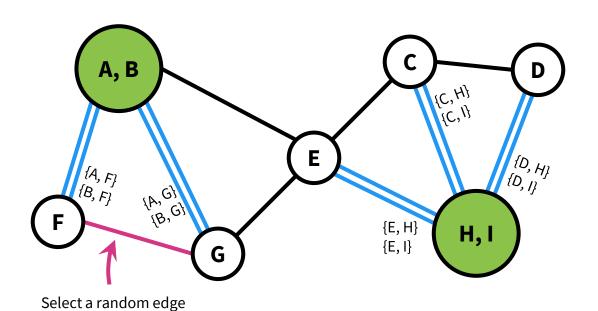


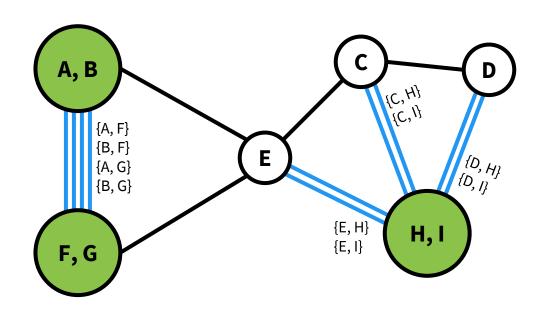


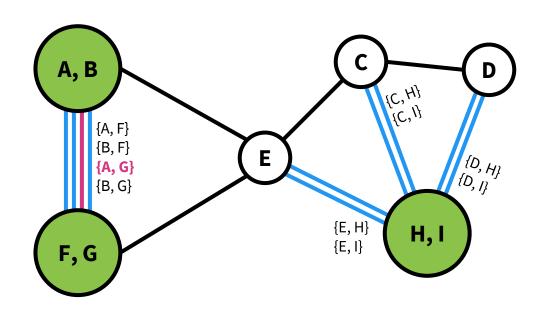


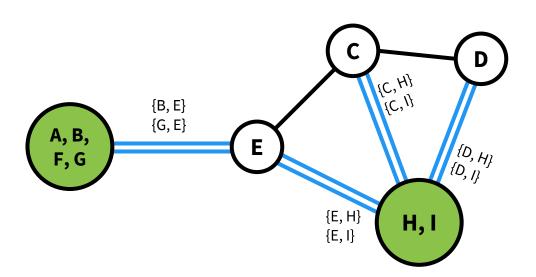


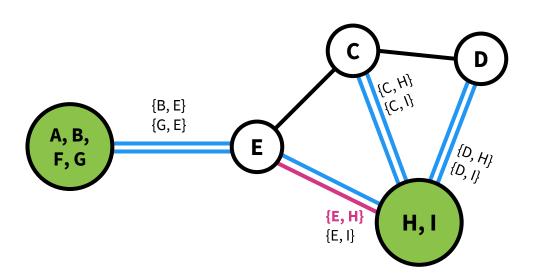


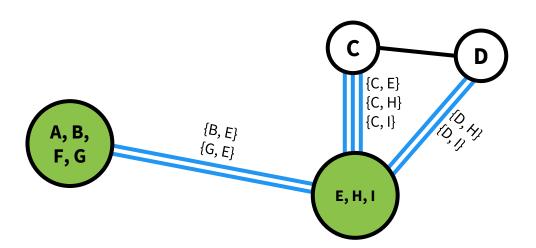


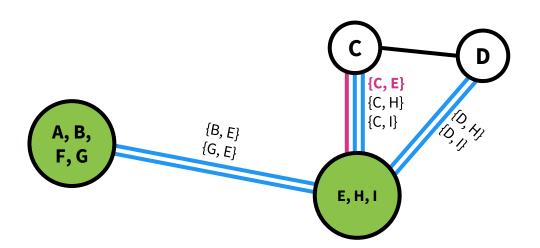


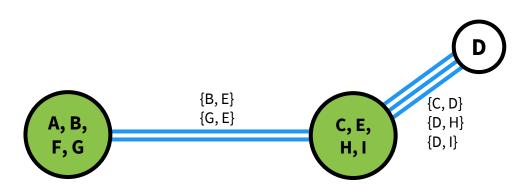


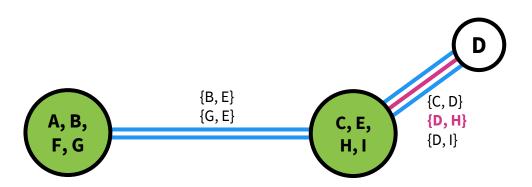






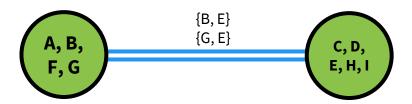






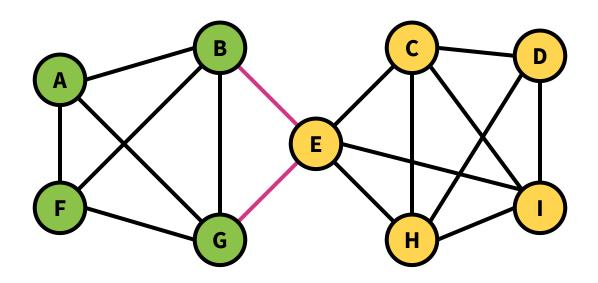
The minimum cut is given by the remaining super-vertices.

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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```
algorithm karger(G=(V,E)):
  G' = {supervertex(v) for v in V}
  E_{u'v'} = \{(u,v)\} \text{ for } (u,v) \text{ E}
  E_{u'v'} = \{\} for (u,v) not in E
  F = \{\{(u,v)\} \text{ for } (u,v) \text{ in } E\}
  while |G'| >= 2:
    \{(u,v)\}\ = uniform random edge in F
    merge supervertices(u, v)
     F = F \setminus E_{u'v'}
  return cut of the remaining super-vertices
algorithm merge supervertices(u, v):
  x' = supervertex(u' U v')
  for w' in G' \ \{u',v'\}:
    E_{v,w} = E_{u,w} U E_{v,w}
    Remove u' and v' from G' and add x'
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The while loop
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times.
                    merge_supervertices(u, v) This takes O(|V|).
                    F = F \setminus E_{\mu'\nu'}
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                    E_{v,w} = E_{u,w} U E_{v,w}
                    Remove u' and v' from G' and add x'
```

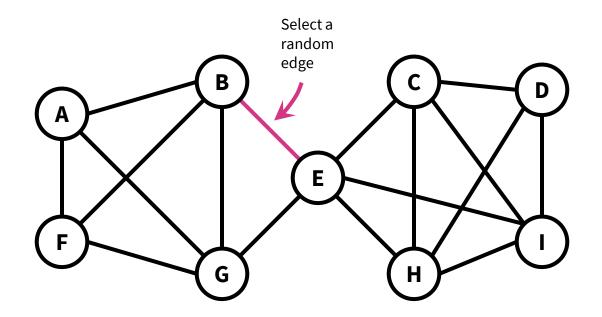
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algorithm karger(G=(V,E)):
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              E<sub>u'v'</sub> = {(u,v)} for (u,v) E
super-edges.
                 E_{u'v'} = \{\} for (u,v) not in E
                 F = \{\{(u,v)\} \text{ for } (u,v) \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
               F = F \setminus E_{u'v'}
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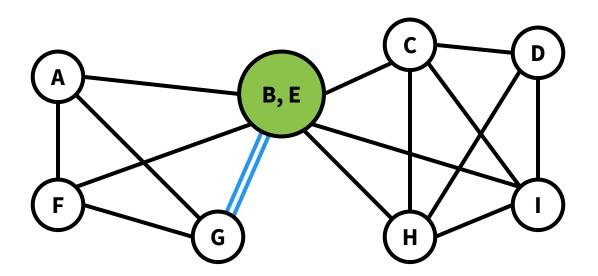
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e.g. Suppose we had chosen this edge.

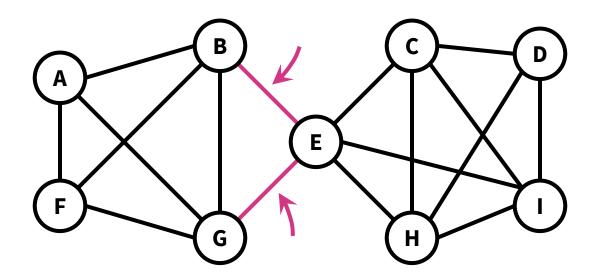


We got really lucky!

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.



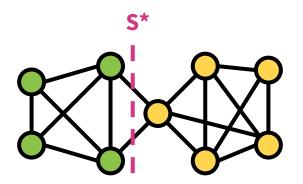
The probability that Karger's algorithm returns a minimum cut is ...

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

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*)



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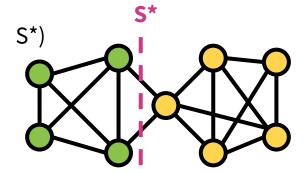
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= P(e_1 doesn't cross S*)
 \times P(e_2 doesn't cross S* | e_1 doesn't cross S*)

• • •

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

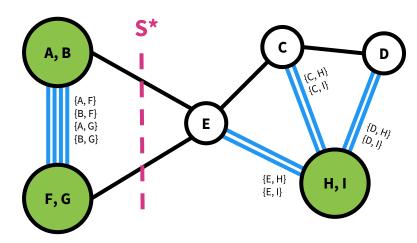


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Suppose, after j-1 iterations, karger hasn't messed up yet! What's the probability of messing up now?



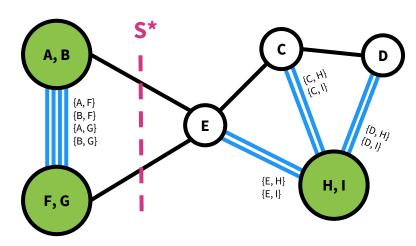
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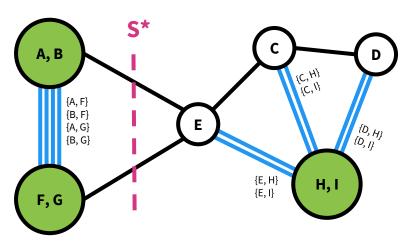
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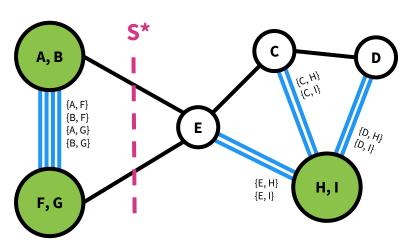
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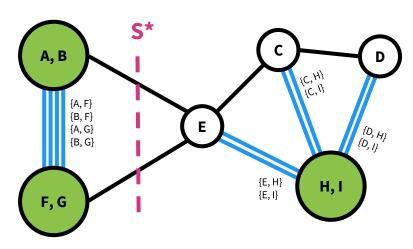
Suppose there are k edges that cross **S***.

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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

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



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Proof, cont.:

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

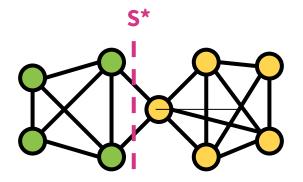
Then $P(karger returns S^*) = P(e_1 doesn't cross S^*)$

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

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

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

$$= 1/(nC2)$$

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

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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 $\leq 1 - 1/(nC2)$

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

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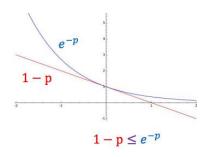
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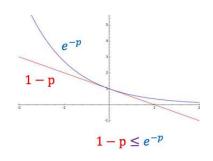
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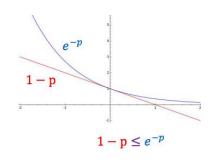
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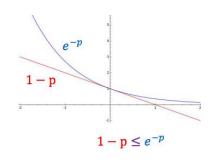
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T = (nC2) ln (1/0.1) times

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)) times = $O(|V|^2)$ 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>
```

Upshot: 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!