#### Lecture 16

### Graph Sketches: Dynamic Spanning Forest

Course: Algorithms for Big Data

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# Spanning forest



Given a graph G = (V, E), a spanning forest of G is a forest F = (V, E')

•  $E' \subseteq E$ .

Adding an edge e ∈ E/E' to F does not change the number of connected components in F.

Fact: a graph can have multiple spanning forests.

# Spanning forest: applications

- Algorithm design: intermediate step in designing graph algorithm (connectivity and path finding)
- Graph analysis: spanning forest gives the connected components of a graph.
- Network design: Minimum number of links to keep the nodes connected.

### Computing a spanning forest

Given a graph on n vertices and m edges, we can compute a spanning forest in time O(m) and O(m) space (BFS/DFS graph traversal).

What if the graph is dynamic (edges are inserted and deleted) and we have to <u>maintain</u> a spanning forest?



### Spanning forest: insert-only streams

When the stream is series of edge insertions, maintaining a spanning forest is easy.

If the new inserted edge e does not create a cycle it is added to the forest otherwise it is ignored.



Space usage:  $O(n \log n)$  bits

per-edge time: ?

### Spanning forest: dynamic graphs

When the edges, in addition to being inserted, are deleted as well it is not clear how to maintain a spanning forest without storing all existing edges.



### Dynamic spanning forest via $\ell_0$ sampling

[Ahn, Guha, McGregor, 2012] There is a randomized algorithm for maintaining a spanning forest under insertion/deletion of edges that uses  $O(n \log^3 n)$  bits of space. The algorithm uses  $\ell_0$  sampling as a subroutine.

 $\ell_0$  sampling: Given a stream of positive and negative updates on a vector  $x \in \mathbb{R}^n$ , a  $\ell_0$  sampler is a randomized algorithm that returns a random non-zero coordinate  $i \in [n]$  where the probability of returning each non-zero coordinates is

$$\frac{1}{\|\boldsymbol{x}\|_0} \pm \frac{1}{n^c}$$

With probability at most  $\delta$ , the algorithm might declare failure and return no sample. The algorithm uses  $O(\log^2 n \log(\frac{1}{\delta}))$  bits of space.

#### Computing the connected components

Lets consider a simpler problem: report the connected components after all edge-insertion/deletions are done.

The algorithm is based on a simple strategy:

Each time pick a random edge and merge its two endpoints into a super-node. Continue this process until no edge remains. In the end, the isolated super-nodes represent the connected components.











In the end, the connected components remain as isolated super-nodes.



We can pick a random edge by  $\ell_0$  sampling the adjacency matrix A.



However, in addition to this, we want to be able to sample an edge from the cut (S, V/S) when S is a super-node. How can we do this?



Suppose the vertices are labeled by numbers in  $\{1, 2, ..., n\}$ .

For each node  $i \in V$ , we define a vector  $u_i \in \{-1, 0, +1\}^{\binom{n}{2}}$  as follows.

- If the edge (i, j) exists and i < j then we set the coordinate u<sub>i</sub>(i, j) = +1
- If the edge (i, j) exists and i > j then we set the coordinate u<sub>i</sub>(i, j) = −1



The vector  $u_{i_1} + \ldots + u_{i_r}$  corresponds to the super-node  $S = \{u_{i_1}, \ldots, u_{i_r}\}.$ 



For each vector  $u_i$ , we maintain an  $\ell_0$  sampling sketch  $sk(u_i)$ .

$$sk(\boldsymbol{u}_1), sk(\boldsymbol{u}_2), \ldots, sk(\boldsymbol{u}_n)$$

If we want to sample an edge from the cut (S, V/S) where  $S = \{i_1, i_2, \ldots, i_r\}$ , we use the sketch

$$sk(\boldsymbol{u}_{i_1} + \boldsymbol{u}_{i_2} + \ldots + \boldsymbol{u}_{i_r})$$



There is one problem: if we contract the edges, one edge at a time, we may end up using the sketch  $sk(u_1)$  multiple times (n-1 times!)

$$\ell_0 \text{ sample } \boldsymbol{u}_1 : sk(\boldsymbol{u}_1)$$
  
 $\ell_0 \text{ sample } \boldsymbol{u}_1 + \boldsymbol{u}_2 : sk(\boldsymbol{u}_1 + \boldsymbol{u}_2)$   
 $\ell_0 \text{ sample } \boldsymbol{u}_1 + \boldsymbol{u}_2 + \boldsymbol{u}_3 : sk(\boldsymbol{u}_1 + \boldsymbol{u}_2 + \boldsymbol{u}_3)$   
...  
 $\ell_0 \text{ sample } \boldsymbol{u}_1 + \boldsymbol{u}_2 + \ldots + \boldsymbol{u}_{n-1} : sk(\boldsymbol{u}_1 + \boldsymbol{u}_2 + \ldots + \boldsymbol{u}_{n-1})$ 



The  $\ell_0$  sample drawn from  $u_1 + u_2$  will depend on the  $\ell_0$  sample drawn from  $u_1$ . Dependency!!

We should not use the sketch  $sk(u_i)$  multiple times because it will cause dependency issues.

If we could query a sketch multiple times we could find all neighbors of a node by using only  $O(\log^3 n)$  bits of space! This is impossible because one cannot compact  $\Omega(n)$  bits of information in  $\log^3 n$  bits of space.



How to avoid using a sketch multiple times?

Lets assume there is no isolated vertex in the input graph G = (V, E).

The algorithm works in multiple rounds. In the first round we do the following:

- For each vertex i ∈ V, we maintain an independent l<sub>0</sub> sampling sketch sk<sub>1</sub>(u<sub>i</sub>).
- We ℓ<sub>0</sub> sample the vector u<sub>i</sub> using the sketch sk<sub>1</sub>(u<sub>i</sub>). As result, for each vertex i ∈ V, we find a random neighbor of i.
- We find at least  $\frac{n}{2}$  random edges in the first round.
- We contract the random edges and create the super-nodes.



In each round, number of nodes drops by a factor of  $\frac{1}{2}$ .

number of nodes in the first round = nnumber of nodes in the second round  $\leq \frac{n}{2}$ 

As result, the algorithm finishes in at most  $\log n$  rounds.

In each round we use fresh  $\ell_0$  sampling sketches for all vertices.

Since there are at most  $\log n$  rounds, for each vertex we need to maintain  $\log n$  number of independent  $\ell_0$  sampling sketches.

In each round, we pick one of the sketches that are not used previously.

In total, we use  $n \log n$  number of  $\ell_0$  sketches. Each sketch takes  $O(\log^2 n)$  bits of space. Therefore the space complexity is  $O(n \log^3 n)$  bits.