Efficient Algorithms for Spanning Tree Centrality

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Abstract

In a connected graph, spanning tree centralities of a vertex and an edge measure how crucial they are for the graph to be connected. In this paper, we propose efficient algorithms for estimating spanning tree centralities with theoretical guarantees on their accuracy. We experimentally demonstrate that our methods are orders of magnitude faster than previous methods. Then, we propose a novel centrality notion of a vertex, called aggregated spanning tree centrality, which also considers the number of connected components obtained by removing the vertex. We also give an efficient algorithm for estimating aggregated spanning tree centrality. Finally, we experimentally show that those spanning tree centralities are useful to identify vulnerable edges and vertices in infrastructure networks.

Introduction

Measuring the importance of edges and vertices is a fundamental task in network analysis. Many importance notions have been proposed in the literature such as degree centrality, betweenness centrality [Anthonisse, 1971; Freeman, 1977], closeness centrality [Bavelas, 1950], and Katz centrality [Katz, 1953], and they are used in various applications such as community detection [Girvan and Newman, 2002; Newman and Girvan, 2004], suppressing the epidemics [Barrat *et al.*, 2008], and maximizing the spread of influence [Kempe *et al.*, 2003].

Among those centralities, in this paper, we focus on *spanning tree centrality* [Teixeira *et al.*, 2013; Qi *et al.*, 2015]. In a connected graph G = (V, E), the spanning tree centrality of an edge $e \in E$ is defined as the probability that e is used in a uniformly sampled spanning tree. Although the spanning tree centrality was originally used to reconstruct a phylogenetic tree from evolutionary relations among species [Teixeira *et al.*, 2013], it turns out that it measures how crucial an edge is for the graph to be connected. Similar centrality notion for vertices was also introduced in [Qi *et al.*, 2015].

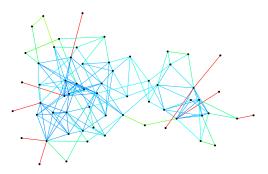


Figure 1: Dolphins network. The color of an edge represents its relative centrality value in the network. Red, green, and blue edges represent that they have high, medium, and low centrality values, respectively.

To get intuition, we depicted spanning tree centralities of edges in a social network formed by bottlenose dolphins [Lusseau *et al.*, 2003] in Figure 1. We can observe that edges on the periphery have high spanning tree centralities because the network will be disconnected by removing a few of those edges. On the other hand, edges in the middle of a community has a small spanning tree centrality because there are many paths bypassing them, and hence they can be safely removed without disconnecting the graph.

Spanning tree centrality is a useful notion for identifying vulnerable edges and vertices. For example, infrastructure networks such as power grids should be robust against physical attacks and natural disasters, and it is important to maintain its connectivity so that all the users can access the infrastructure [Bernstein *et al.*, 2014; Pagani and Aiello, 2013]. Spanning tree centrality directly measures how edges and vertices are important to keep the graph connected. Also, it is reported that edges of high spanning tree centralities are important for disseminating information through a social network [Mavroforakis *et al.*, 2015].

Although several exact and approximation methods have been proposed for computing spanning tree centralities [Teixeira *et al.*, 2013; Mavroforakis *et al.*, 2015; Qi *et al.*, 2015], those methods are not scalable enough to find vulnerable edges and vertices in large real-world networks, and resolving this issue is the main motivation of this paper.

Our Contributions

In this paper, we propose efficient algorithms for estimating spanning tree centrality with theoretical guarantees on their accuracy. Our algorithms sample spanning trees uniformly at random sufficiently many times using Wilson's algorithm [Wilson, 1996], and then estimate the spanning tree centralities of edges or vertices using those samples. Combined with several speeding-up techniques, our algorithms run orders of magnitude faster than previous methods on realworld networks. Because of the efficiency of our methods, we can now identify vulnerable edges and vertices in large networks. Then, we experimentally show that protecting edges and vertices of high spanning tree centralities makes a power grid network robust against random failures of edges and vertices.

The effect of vertex failures in infrastructure networks is often assessed by the resulting number of connected components [Bernstein et al., 2014]. We introduce a novel notion, called aggregated spanning tree centrality, which measures the importance of vertices with respect to this criteria. Then, we give an efficient algorithm for estimating aggregated spanning tree centrality with theoretical guarantee on its accuracy. Using a power grid network, we experimentally show that the failures of vertices with high aggregated spanning tree centralities indeed result in a large number of connected components compared to other centrality notions.

Related Work

The notion of spanning tree centrality for edges was introduced in [Teixeira *et al.*, 2013] for reconstructing a phylogenetic tree from evolutionary relationships among species. Their exact method is based on Kirchhoff's matrix-tree theorem and computes determinants of matrices polynomially many times. This makes their algorithm slow, and it can only handle graphs of tens of thousands of edges within an hour.

An approximation algorithm based on dimension reduction through the Johnson-Lindenstraus lemma [Johnson and Lindenstrauss, 1984] was proposed by [Spielman and Srivastava, 2011]. A more practical version of this algorithm was given by [Mavroforakis *et al.*, 2015]. Their algorithm uses an almost linear-time algorithm for solving symmetric diagonally dominant (SDD) systems [Koutis *et al.*, 2011]. Despite these sophisticated techniques, their algorithm can only handle graphs of million edges within an hour. We will experimentally show that our algorithms can accurately estimate spanning tree centrality on graphs of millions of edges within a minute.

Spanning tree centrality for vertices was introduced in [Qi *et al.*, 2015]. Their method is also an exact method based on Kirchhoff's matrix-tree theorem and hence impractical.

A spanning centrality of an edge is often called *effective resistance* as it corresponds to effective resistance when we regard a graph as an electrical circuit. See [Lovász, 1993] for consequences of this interpretation and connections with random walk. Theoretically, the spanning tree centrality of an edge can be well approximated in almost linear time [Spielman and Srivastava, 2011] and is used as a building block of designing almost linear time algorithms for other problems

such as spectral sparsifier construction [Spielman and Srivastava, 2011] and approximate maximum flow [Kelner *et al.*, 2014].

Preliminaries

In this section, we introduce basic notions used throughout this paper and review several concentration inequalities.

For an integer $k \in \mathbb{N}$, [k] denotes the set $\{1, 2, \ldots, k\}$. For $a, b \in \mathbb{R}$ and $\epsilon > 0$, $a = b \pm \epsilon$ means $b - \epsilon \le a \le b + \epsilon$. For a probabilistic event X, [X] denotes its indicator, that is, [X] = 1 if X holds and 0 otherwise.

Let G = (V, E) be a graph. For a vertex $v \in V$, $d_G(v)$ denotes the degree of v. For a set of vertices $S \subseteq V$, G[S] denotes the subgraph of G induced by S.

We call an edge a *bridge* if the number of connected components increases by removing it. Similarly, we call a vertex an *articulation point* if the number of connected components increases by removing it. A graph is called *biconnected* if it does not have any articulation point. A *biconnected component* of a graph is its maximal biconnected subgraph.

Symbols n and m will denote the number of vertices and edges, respectively, of the input graph.

Spanning Tree Centrality

Let G=(V,E) be a connected graph. We define μ_G as the uniform distribution over spanning trees of G. By $T\sim \mu_G$, we mean that we sample a spanning tree T from μ_G . We introduce three kinds of spanning tree centralities, one for edges and the other two for vertices.

Definition 1 (Teixeira *et al.* 2013). Let G = (V, E) be a connected graph. The spanning tree centrality of an edge $e \in E$, denoted by $\operatorname{st}_G(e)$, is defined as $\operatorname{Pr}_{T \sim \mu_G}[e \in T]$.

Definition 2 (Qi *et al.* 2015). Let G=(V,E) be a connected graph. The spanning tree centrality of a vertex $v\in V$, denoted by $\operatorname{st}_G(v)$, is defined as $\Pr_{T\sim \mu_G}[d_T(v)\geq 2]$.

Definition 3. Let G=(V,E) be a connected graph. The aggregated spanning tree centrality of a vertex $v\in V$, denoted by ${\rm ast}_G(v)$, is defined as $\mathbf{E}_{T\sim\mu_G}[d_T(v)]$.

We omit subscripts when they are clear from the context. Below, we explain intuitions behind those definitions.

The spanning tree centrality of an edge measures how it is like a bridge. Indeed, if an edge e=(u,v) is a bridge, then e is used in any spanning tree and hence $\operatorname{st}(e)=1$. On the other hand, if there are many other paths between u and v that bypass e, then e is unlikely used in a uniformly sampled spanning tree, and hence $\operatorname{st}(e)$ will be small.

The spanning tree centrality of a vertex measures how it is like an articulation point. Indeed, if a vertex v is an articulation point, then v cannot be a leaf in a spanning tree and hence $\operatorname{st}(v)=1$. On the other hand, if v is contained in a tightly connected community and is not an articulation point, then $\operatorname{st}(v)$ becomes small. For example, if v is a vertex in a complete graph of v vertices, then by Kirchhoff's matrix tree theorem, we can easily show that $\operatorname{st}(v)=1-(1-1/n)^{n-2}\to 1-1/e\ (n\to\infty)$.

The aggregated spanning tree centrality of a vertex incorporates the number of connected components obtained by removing the vertex. To see this, note that $\operatorname{ast}(v) = \sum_{i=1}^{\infty} \Pr_{T \sim \mu_G}[d_T(v) \geq i]$. Here, $\Pr_{T \sim \mu_G}[d_T(v) \geq i]$ can be seen as a generalization of $\operatorname{st}(v)$ in the sense that it measures how v is like an articulation point such that the number of connected components increases by at least i-1 by removing v.

Concentration Inequalities

We will use following Hoeffding's inequality in our analysis.

Lemma 4. Let X_1, \ldots, X_q be independent random variables in [0,1] and $X = \sum_{i \in [q]} X_i$. Then for any $0 < \epsilon < 1$, we have

$$\Pr[|X - \mathbf{E}[X]| > \epsilon q] \le 2 \exp(-2\epsilon^2 q).$$

We say that random variables X_1,\ldots,X_q are negatively correlated if $\Pr[\bigwedge_{i\in[q]}X_i]\leq\prod_{i\in[q]}\Pr[X_i]$. Even when random variables are dependent, if they are negatively correlated, then the following concentration inequality holds:

Lemma 5 (Panconesi and Srinivasan 2006). Let X_1, \ldots, X_q be random variables in $\{0,1\}$ that are negatively correlated and $X = \sum_{i \in [q]} X_i$. Then for any $0 < \epsilon < 1$, we have

$$\Pr[X > (1 + \epsilon) \mathbf{E}[X]] \le \left(\frac{e^{\epsilon}}{(1 + \epsilon)^{1 + \epsilon}}\right)^{\mathbf{E}[X]},$$
$$\Pr[X < (1 - \epsilon) \mathbf{E}[X]] \le \exp(-\epsilon^2 \mathbf{E}[X]/2).$$

For an edge $e \in E$, let X_e be the indicator corresponding to the event that e is used in a spanning tree sampled from μ_G . Then, we have the following fact.

Lemma 6 (Burton and Pemantle 1993). Let G = (V, E) be a connected graph. Random variables $\{X_e\}_{e \in F}$ for any $F \subseteq E$ are negatively correlated.

Hence, we can apply Lemma 5 to the set $\{X_e\}_{e\in F}$ for any set of edges $F\subseteq E$.

Algorithms

In this section, we explain our algorithms for estimating spanning tree centralities.

Spanning Tree Centrality of Edges

Our algorithm for estimating spanning tree centralities of edges is based on Wilson's algorithm [Wilson, 1996], which samples a spanning tree uniformly at random. We need several notions to explain Wilson's algorithm.

Let G=(V,E) be a graph and $P=(v_1,\ldots,v_\ell)$ be a path in G, that is, $\{v_i,v_{i+1}\}\in E$ for each $i\in [\ell-1]$. We define the loop erasure of P, denoted by $\mathrm{LE}(P)$, as the path obtained from P by removing cycles. More formally, we define indices i_j inductively using the following rule: $i_1=1$ and $i_{j+1}=\max\{i\in [\ell]:v_i=v_{i_j}\}+1$. The induction stops when we have $v_{i_j}=v_\ell$ for the current j. Let ℓ' be the last index j for which i_j is defined. Then, $\mathrm{LE}(P)$ is defined as the path $(v_{i_1},\ldots,v_{i_{\ell'}})$.

Wilson's algorithm is given in Algorithm 1. It takes a graph G=(V,E) and a vertex $r\in V$, called the root, as its input. Let v_1,\ldots,v_{n-1} be an arbitrary ordering of $V\setminus\{r\}$. We start with a tree T consisting of a single vertex r. Then, for each

Algorithm 1 Wilson's algorithm

```
1: procedure WILSON(G, r)

2: T \leftarrow the tree consisting of a single vertex r.

3: Let v_1, \ldots, v_{n-1} be an arbitrary ordering of V \setminus \{r\}.

4: for i = 1 to n - 1 do

5: Let P be a random walk from v_i to T.

6: Add the loop erasure \mathrm{LE}(P) to T.

7: return T.
```

Algorithm 2 Spanning Tree Centrality for Edges

```
1: procedure ST-EDGE(G, r, \epsilon, \delta)

2: q \leftarrow \lceil \log(2m/\delta)/(2\epsilon^2) \rceil.

3: for i = 1 to q do

4: T_i \leftarrow \text{WILSON}(G, r).

5: for each e \in E do

6: \widetilde{\text{st}}(e) \leftarrow \frac{1}{q} \sum_{i \in [q]} [e \in T_i].

7: return \{\widetilde{\text{st}}(e)\}_{e \in E}.
```

 $i \in [n-1]$, we perform a random walk from v_i until it reaches some vertex in T, and then add to T the loop erasure $\mathrm{LE}(P)$ of the obtained random walk P. Note that, if v_i is already a vertex in T, then P is a walk consisting of a single vertex v_i , and hence T does not change. It is clear that Wilson's algorithm always returns a spanning tree when G is connected. Indeed the following holds.

Theorem 7 (Wilson 1996). For a connected graph G = (V, E) and a vertex $r \in V$, WILSON(G, r) uniformly samples a spanning tree of G. The expected running time is $\sum_{u \in V \setminus \{r\}} \pi_G(u) \kappa_G(u, r)$.

Here, $\pi_G(u)$ denotes the probability of staying at a vertex $u \in V$ in the stationary distribution of a random walk on G, and $\kappa_G(u,v)$ denotes the commute time between two vertices $u,v \in V$, that is, the expected number of steps in a random walk starting at u, before v is visited and then u is reached again. Note that $\pi_G(u) = d_G(u)/2m$, where m is the number of edges, as we consider undirected connected graph.

The choice of the root r and the vertex ordering v_1,\ldots,v_{n-1} much affect the practical performance. We will discuss several strategies later.

With the aid of Wilson's algorithm, we can efficiently estimate spanning edge centralities of edges by sampling a sufficient number of spanning trees and then counting how many of them use a particular edge (Algorithm 2). We have the following guarantee on its accuracy and time complexity:

Theorem 8. Let G=(V,E) be a connected graph, $r\in V$ be a vertex, and $0<\epsilon,\delta<1$. With probability at least $1-\delta$, ST-EDGE (G,r,ϵ,δ) outputs $\{\widetilde{\operatorname{st}}(e)\}_{e\in E}$ with $\widetilde{\operatorname{st}}(e)=\operatorname{st}(e)\pm\epsilon$ for every $e\in E$. The expected running time is $\lceil\log(2m/\delta)/(2\epsilon^2)\rceil\cdot\sum_{u\in V}\pi_G(u)\kappa_G(u,r)$.

Proof. By Hoeffding's inequality, for each edge $e \in E$, the probability that $\widetilde{\operatorname{st}}(e) = \operatorname{st}(e) \pm \epsilon$ holds is at least $1 - \delta/m$. By union bound, the probability that $\widetilde{\operatorname{st}}(e) = \operatorname{st}(e) \pm \epsilon$ holds for every $e \in E$ is at least $1 - \delta$.

The time complexity is immediate from Theorem 7. \Box

Algorithm 3 Spanning Tree Centrality for Vertices

```
1: procedure ST-VERTEX(G, r, \epsilon, \delta)

2: q \leftarrow \lceil \log(2n/\delta)/(2\epsilon^2) \rceil.

3: for i = 1 to q do

4: T_i \leftarrow \text{WILSON}(G, r).

5: for each u \in V do

6: \widetilde{\text{st}}(u) \leftarrow \frac{1}{q} \sum_{i \in [q]} [d_{T_i}(u) \geq 2].

7: return \{\widetilde{\text{st}}(u)\}_{u \in V}.
```

Algorithm 4 Aggregated Spanning Tree Centrality

```
1: procedure AST-VERTEX(G, r, \epsilon, \delta)

2: \epsilon' \leftarrow \widetilde{\Theta}(\epsilon/\log(n/\delta)). (\widetilde{\Theta}(\cdot)) hides the log factor.)

3: q \leftarrow \widetilde{\Theta}(\log(n/\delta)/\epsilon'^2).

4: for i=1 to q do

5: T_i \leftarrow \text{WILSON}(G, r).

6: for each u \in V do

7: \widetilde{\text{ast}}(u) \leftarrow \frac{1}{q} \sum_{i \in [q]} d_{T_i}(u).

8: return \{\widetilde{\text{ast}}(u)\}_{u \in V}.
```

We will see that, in real-world networks, most edges have large spanning tree centralities, say, more than 0.1. Hence, we can detect important edges by choosing ϵ a small constant, that is, a value independent of the graph size.

Spanning Tree Centrality of Vertices

Our algorithm for estimating spanning tree centralities of vertices is very similar to the one for edges (Algorithm 3). We have the following guarantee:

Theorem 9. Let G=(V,E) be a connected graph, $r\in V$ be a vertex, and $\epsilon,\delta>0$. Then with probability at least $1-\delta$, ST-VERTEX (G,r,ϵ,δ) outputs $\{\widetilde{\operatorname{st}}(u)\}_{u\in V}$ with $\widetilde{\operatorname{st}}(u)=\operatorname{st}(u)\pm\epsilon$ for every $u\in V$. The expected running time is $\lceil\log(2n/\delta)/(2\epsilon^2)\rceil\cdot\sum_{u\in V}\pi_G(u)\kappa_G(u,r)$.

As the proof of Theorem 9 is quite similar to that of Theorem 8, we omit the details.

Aggregated Spanning Tree Centrality of Vertices

In this section, we explain our algorithm for estimating aggregated spanning tree centralities of vertices. Our algorithm is given in Algorithm 4. It simply samples a sufficient number of spanning trees, and then computes the average degree of each vertex in those spanning trees.

If we simply apply Hoeffding's inequality, then we can only guarantee that, with high probability, $\widetilde{\operatorname{ast}}(u)$ approximates $\operatorname{ast}(u)$ to within $\epsilon d_G(u)$ for every $u \in V$. We want a stronger guarantee because $\operatorname{ast}(u)$ is typically much smaller than $d_G(u)$. To this end, we use the fact that edges in a uniformly sampled spanning tree are negatively correlated, and we get the following guarantee:

Theorem 10. Let G = (V, E) be a connected graph, $r \in V$ be a vertex, and $\epsilon, \delta > 0$. Then with probability at least $1 - \delta$, AST-VERTEX (G, r, ϵ, δ) outputs $\{\widetilde{\operatorname{ast}}(u)\}_{u \in V}$ with $\widetilde{\operatorname{ast}}(u) = \operatorname{ast}(u) \pm (\epsilon(1 + \operatorname{ast}(u)) + \delta)$ for every

 $u \in V$. The expected running time is $\widetilde{O}(\log^3(n/\delta)/\epsilon^2 \cdot \sum_{u \in V} \pi_G(u) \kappa_G(u,r))$.

Proof. We fix a vertex $u \in V$. We show that the algorithm outputs $\widetilde{\operatorname{ast}}(u)$ with $\widetilde{\operatorname{ast}}(u) = \operatorname{ast}(u) \pm \left(\epsilon(1+\operatorname{ast}(u)) + \delta\right)$ with probability $1-\delta/n$. Then, the first claim follows by union bound.

Let $Y_i=d_{T_i}(u)$ for each $i\in[q]$. Let $N\geq \operatorname{ast}(u)$ be a parameter determined later, and we define $Y_{i,\leq N}=Y_i\cdot [Y_i\leq N]$ and $Y_{i,>N}=Y_i\cdot [Y_i>N]$ for each $i\in[q]$. Note that $Y_i=Y_{i,\leq N}+Y_{i,>N}$ holds. Hence, $\operatorname{ast}(u)=\frac{1}{q}\sum_i Y_i=\frac{1}{q}\sum_i Y_{i,\leq N}+\frac{1}{q}\sum_i Y_{i,>N}$ holds. By Hoeffding's bound,

$$\frac{1}{q} \sum_{i} Y_{i, \leq N} = \frac{1}{q} \sum_{i} \mathbf{E}[Y_{i, \leq N}] \pm \epsilon' N$$

holds with probability at least $1 - \delta/(2n)$.

Let $\epsilon'' = N/\operatorname{ast}(u) - 1$. Note that $\epsilon'' \ge 0$ because we chose $N \ge \operatorname{ast}(u)$. For any $i \in [q]$, by Lemma 5, we have

$$\Pr[Y_i > N] \le \frac{e^{\epsilon'' \operatorname{ast}(u)}}{(1 + \epsilon'')^{(1 + \epsilon'') \operatorname{ast}(u)}} \le \frac{e^N}{(N/\operatorname{ast}(u))^N}.$$

By choosing $N=\max\{2e\cdot \operatorname{ast}(u), \log(2qn/\delta)\}$, the probability of having $Y_i>N$ becomes at most $\delta/(2qn)$. Then by union bound, with probability at least $1-\delta/(2n)$, we have $Y_i\leq N$ for every $i\in [q]$. In other words, $Y_{i,>N}=0$ for every $i\in [q]$.

Since $Y_i \leq n$, we have $\mathbf{E}[Y_{i,>N}] \leq n \cdot \delta/(2qn) = \delta/(2q)$, which means that $\mathbf{E}[Y_{i,\leq N}] = \mathrm{ast}(u) - \mathbf{E}[Y_{i,>N}] \geq \mathrm{ast}(u) - \delta/(2q)$. Also, it is clear that $\mathbf{E}[Y_{i,\leq N}] = \mathrm{ast}(u) - \mathbf{E}[Y_{i,>N}] \leq \mathrm{ast}(u)$.

Combining arguments above, with probability at least $1 - \delta/n$, we have

$$\widetilde{\operatorname{ast}}(u) = \frac{1}{q} \sum_{i} Y_i = \frac{1}{q} \sum_{i} Y_{i, \leq N} = \frac{1}{q} \sum_{i} \mathbf{E}[Y_{i, \leq N}] \pm \epsilon' N$$
$$= \operatorname{ast}(u) \pm \left(\delta + \epsilon' \max\{2e \cdot \operatorname{ast}(u), \log(2qn/\delta)\}\right).$$

By choosing $\epsilon' = \epsilon/\big(2e\log(2qn/\delta)\big)$, we have $\widetilde{\mathrm{ast}}(u) = \mathrm{ast}(u) \pm \big(\delta + \epsilon(1 + \mathrm{ast}(u))\big)$. Note that the dependency between ϵ' and q is cyclic because we need $q = \Theta(\log(n/\delta)/\epsilon'^2)$. However, this recurrence has a solution $\epsilon' = \widetilde{\Theta}(\epsilon/\log(n/\delta))$ and $q = \widetilde{\Theta}(\log(n/\delta)/\epsilon^2)$.

The analysis of the time complexity is obvious.

Speeding-up techniques

In this section, we discuss several techniques for speeding up our algorithms.

Biconnected-component decomposition

Let G=(V,E) be a connected graph. Suppose that biconnected components C_1,\ldots,C_k are attached at an articulation point $u\in V$. Then, we can uniformly sample a spanning tree T of G by uniformly sampling spanning trees T_1,\ldots,T_k of $G[C_1],\ldots,G[C_k]$, respectively, and attaching them at u. This means that $\operatorname{st}_G(e)=\operatorname{st}_{G[C_i]}(e)$ for any edge $e\in G[C_i]$. Hence, it suffices to process smaller graphs $G[C_1],\ldots,G[C_k]$ separately.

[Mavroforakis *et al.*, 2015] proposed to reduce the input graph to its 2-*core*, a maximal subgraph of the minimum degree at least two. Our decomposition technique surpasses the reduction to the 2-core because after splitting the graph at articulation points, we get bridges and biconnected components, which are subgraphs of the 2-core.

Vertex ordering

We consider the following strategies for the vertex ordering strategies used in Wilson's algorithm: Random, Degree, Distance, and Reverse. In the first three strategies, we choose a vertex with the highest degree as the root. Other vertices are randomly shuffled, sorted in decreasing order of degrees, and sorted in increasing order of distances from the root in Random, Degree, and Distance, respectively. In Reverse, a vertex with the lowest degree is chosen as the root and other vertices are sorted in increasing order of degrees.

Experiments

In this section, we demonstrate the efficiency of our methods and the effectiveness of spanning tree centralities for finding vulnerable edges and vertices in real-world networks.

Environment All the proposed methods were implemented in C++11 and compiled with gcc 4.8.3 with -O3 option. Unless stated otherwise, we used biconnected-component decomposition and Distance strategy for vertex ordering. The implementation of the algorithm proposed by [Mavroforakis *et al.*, 2015], which we refer to the MGKT algorithm, was obtained from authors' website¹. All experiments were conducted with a single thread on a Linux server with Intel Xeon X5675 3.07GHz CPU and 288GB Memory.

Datasets We conducted all experiments on real-world social, road, and infrastructure networks. All datasets used in our experiments can be downloaded from Stanford Network Analysis Project [Leskovec and Krevl, 2014]. Table 1 shows the information of datasets.

Spanning Tree Sampling Performance

We first evaluate the runtime of sampling spanning trees with our speed-up techniques. Note that the runtime of our centrality estimation algorithms is dominated by the time for sampling spanning trees. Therefore, the overall running time is easily obtained by multiplying the time for sampling a spanning tree and the number of required spanning trees, which is determined by the error parameter ϵ .

Speed-up by the biconnected-component decomposition

Table 2 shows the average runtime of sampling a spanning tree, where we sampled 1,000 spanning trees in total. Note that we can sample a spanning trees in less than 2 seconds on Orkut, which is a large social network with over one million edges. Biconnected components decomposition consistently reduces the runtime of sampling a spanning tree. In particular, we can sample spanning trees more than 40% faster on sparse networks such as DBLP and Youtube.

Table 1: Statistics of datasets. We denote the number of vertices and edges in the largest biconnected component of the original graph by n' and m', respectively. ST_e denotes the spanning edge centrality of the vertex with 100,000-th highest centrality value except US-PowerGrid. For US-PowerGrid, ST_e denotes that of the vertex with 3,000-th highest centrality value since US-PowerGrid has only 6.6K edges.

| Dataset | n | m | n' | m' | ST_e |
|--------------|------|------|------|------|-----------------|
| US-PowerGrid | 4.9K | 6.6K | 3.0K | 4.6K | 0.77 |
| Gnutella | 63K | 148K | 34K | 119K | 0.20 |
| Epinions | 76K | 509K | 36K | 365K | 0.17 |
| Slashdot | 77K | 905K | 47K | 439K | 0.13 |
| DBLP | 426K | 1.0M | 211K | 884K | 0.60 |
| Youtube | 1.2M | 3.0M | 452K | 2.3M | 1.00 |
| RoadNet-TX | 1.4M | 1.9M | 1.1M | 1.6M | 1.00 |
| Skitter | 1.7M | 11M | 1.4M | 11M | 1.00 |
| Pokec | 1.6M | 22M | 1.5M | 22M | 1.00 |
| Orkut | 3.1M | 117M | 3.0M | 117M | 0.51 |
| | | | | | |

Table 2: The average runtime of sampling a spanning tree. T_G and T_{BC} denote the runtime in milliseconds when we sample spanning trees from the whole graph and from each biconnected component, respectively.

| Dataset | T_G | T_{BC} | T_{BC}/T_G |
|------------|--------|----------|--------------|
| DBLP | 132.1 | 78.7 | 0.60 |
| Youtube | 297.1 | 133.1 | 0.45 |
| RoadNet-TX | 1288.2 | 997.9 | 0.77 |
| Skitter | 494.0 | 371.2 | 0.75 |
| Pokec | 652.0 | 574.5 | 0.88 |
| Orkut | 1462.9 | 1380.6 | 0.94 |

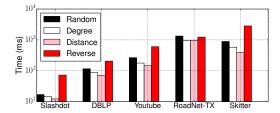


Figure 2: Effect of vertex ordering strategies.

Vertex ordering strategies

We examined how vertex ordering strategies affect the efficiency of sampling spanning trees. Figure 2 shows the average runtime for each strategy to sample a spanning tree, where we sampled 1,000 spanning trees in total. We can observe that Distance is the fastest among the four strategies on almost all datasets. Apparently, this is because we can avoid long unnecessary random walks in the early stage of Wilson's algorithm by starting random walks from vertices close to the root.

Edge Centrality Estimation Performance

Among the three spanning tree centralities, we take the spanning tree centrality for edges as an example, and compare our method with the previous MGKT algorithm. We obtained

http://cs-people.bu.edu/cmav/centralities

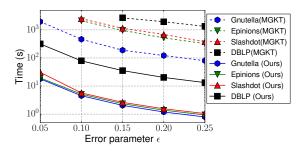


Figure 3: The runtime of our method and the MGKT algorithm under a given error parameter ϵ .

similar results for other centralities.

We compare the runtime of our method and the MGKT algorithm on several small networks. We chose the failure probability $\delta = 1/n$. For an error parameter ϵ , our method samples $\lceil \log(2m/\delta)/(2\epsilon^2) \rceil$ spanning trees and guarantees $\operatorname{st}(e) = \operatorname{st}(e) \pm \epsilon$ for every edge with probability at least $1 - \delta$ and the MGKT algorithm guarantees $(1-\epsilon)^2$ st $(e) < \widetilde{st}(e) <$ $(1+\epsilon)^2$ st(e) for every edge with probability at least $1-\delta$. We cannot directly compare those algorithms because our method guarantees absolute error whereas the MGKT algorithm guarantees relative error. However, as shown in Table 1, the important edges have large centrality values, say, more than 0.1, and hence we can claim that our method with $\epsilon = 0.05$ gives a sufficient accuracy in comparison with the MGKT algorithm with $\epsilon = 0.25$. Under this setting, our method runs orders of magnitude faster than the MGKT algorithm as shown in Figure 3.

Effectiveness of Spanning Tree Centralities

We now demonstrate the effectiveness of (aggregated) spanning tree centralities of edges and vertices.

Connectivity

We first show that, using US-PowerGrid, which is a real-world infrastructure network, spanning tree centralities of edges/vertices are useful to identify important edges/vertices for maintaining connectivity. In our experiments, we protected a certain fraction of edges/vertices with the highest centrality values in terms of spanning tree centrality or betweenness centrality, and removed each unprotected edge/vertex with a certain probability. Then, we examined the probability that the network remained connected. The results are shown in Figure 4 and 5. We can see that show spanning tree centrality is a better measure for identifying important edges and vertices to maintain connectivity.

Number of connected components

Finally, we examine the effectiveness of aggregated spanning tree centrality by considering the number of connected components. For clarity, this time, instead of protecting vertices, we removed vertices from the ones with higher centrality values. Figure 6 illustrates the result. We used the largest connected component of US-PowerGrid as the initial graph. For each different centrality measure, we removed the vertices with highest centrality values from the graph, and then computed the numbers of connected components. Larger result-

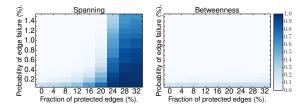


Figure 4: The probability that US-PowerGrid remained connected after edge removal among 500 trials.

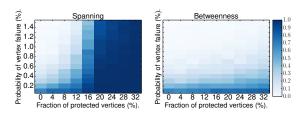


Figure 5: The probability that US-PowerGrid remained connected after vertex removal among 500 trials.

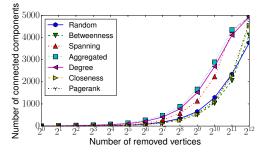


Figure 6: Number of connected components after vertex removals. Each removed vertex is considered as one connected component.

ing numbers of connected components mean that removed vertices are key vertices to connect many different parts of the graph. From Figure 6, we clearly see that our aggregated spanning tree centrality performs best in this experiment.

Conclusions

In this paper, we gave efficient algorithms for estimating spanning tree centrality. Our algorithms are orders of magnitude faster than previous methods. By experiment, we confirmed that spanning tree centralities are to identify vulnerable edges and vertices in infrastructure networks.

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