Absorbing Random Walk Centrality

Theory and Algorithms

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Background

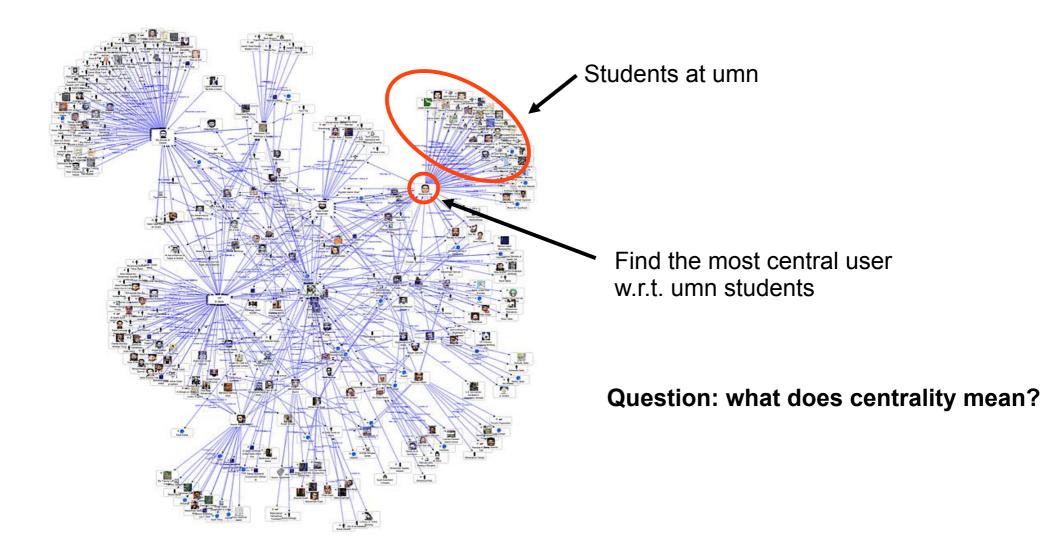


#Facebook users

#find the k most central users in this large network

#How about finding central users w.r.t particular users?

Background

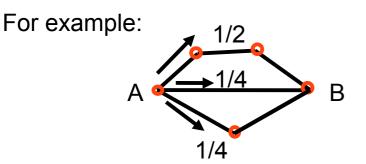


Centrality

- **Degree Centrality**: centrality of a node is simply quantified by its <u>degree</u>
- Closeness centrality: the <u>average distance</u> of a node from all other nodes in the graph
- Betweenness cerntrality: the <u>number of shortest paths</u> between pairs of nodes in the graph that pass through a given node
 Drawbacks: Above centrality can change dramatically with the insertion or deletion of a single edge.

Random-Walk Centrality (RWC)

- **Definition**: the <u>expected first passage time</u> of a random walk of a given node of the graph, when it starts from a random node of the graph.
- Strengths: robust when change edges in a graph



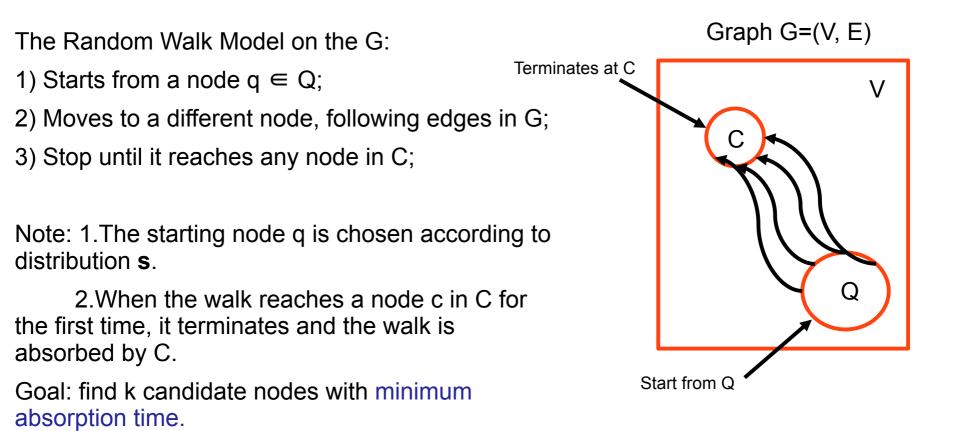
RWC of node B= 3* 1/2+1* 1/4+2* 1/4=9/4

Problem Formulation

-Given a graph G=(V, E), where V is the set of n nodes, E is the set of m undirected edges. A subset of nodes $Q \subseteq V$, referred to the query nodes.

-Goal: Find a set C of k nodes that are central w.r.t the query nodes Q.

-The centrality of a set of nodes C w.r.t Q is based on the notion of randomwalk centrality.



Problem Formulation

The Random Walk Model:

1) Starts from a node $q \in Q$;

2) Moves to a different node, following edges in G;

3)May restart from Q with probability α

4) Stop until it reaches any node in C;

Problem Statement:

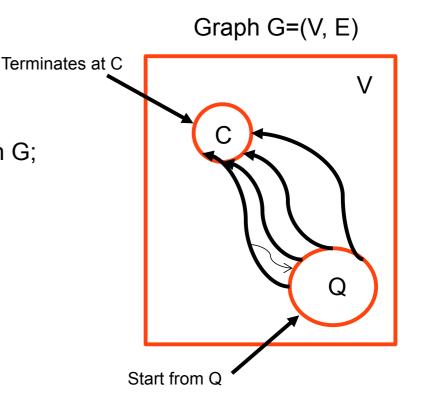
Let $\operatorname{ac}_{Q}^{q}(C)$ denote this expected length from q to C.

Define the absorbing random-walk centrality of a set of nodes C w.r.t. query nodes Q as:

 $\operatorname{ac}_Q(C) = \sum_{q \in Q} \mathbf{s}(q) \operatorname{ac}_Q^q(C).$

How to calculate the random-walk centralit of set C regarding Q?(Compute expected length from Q to C.)

Ans: Using absorbing random walks model



Goal: Calculate the expected length of a random walk from Q to C.

Define a random walk on the graph G:

-Let P be the transition matrix for a random walk, with P(i,j) to be the transition probability from node i to node j in one step.

-Define P(c,c)=1 and P(c,j)=0 if $j \neq c$, for all absorbing nodes c in C. -For the rest T=V\C of non-absorbing (transient) nodes. Define the transition probability as:

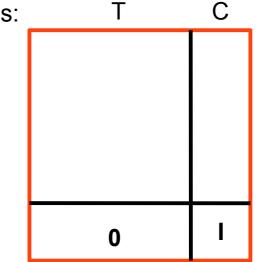
$$\mathbf{P}(i,j) = \begin{cases} \alpha \,\mathbf{s}(j) & \text{if } j \in Q \setminus N(i), \\ (1-\alpha)/d_i + \alpha \,\mathbf{s}(j) & \text{if } j \in N(i). \end{cases}$$

N(i) denotes the neighbors of node i.

So the transition matrix of the random walk is written as:

$$\mathbf{P} = \left(\begin{array}{cc} \mathbf{P}_{TT} & \mathbf{P}_{TC} \\ \mathbf{0} & \mathbf{I} \end{array} \right),$$

Nodes in V represent the states, P defines the P= transition matrix. We get the **random walk model**! In oerder to compute expected length from Q to C. We first compute the expected length from node i to node j.in the **defined random walk model**



Т

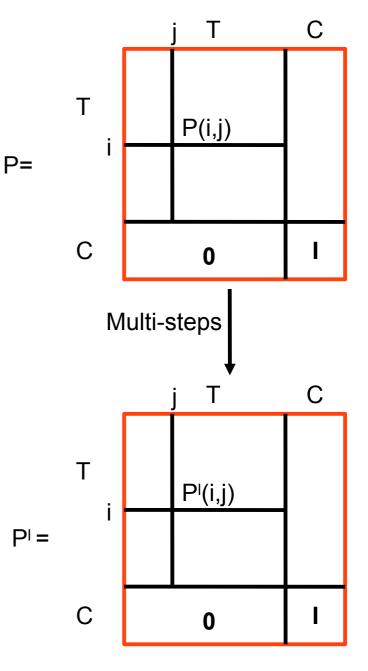
С

The probability **from i to j via I step** is given by the (i,j)-entry if matrix $P^{I}(i,j)$. So the expected length that the random walk vist node j starting from node i is given by the (i,j)-entry of the |T|x|T| matrix:

$$\mathbf{F} = \sum_{\ell=0}^{\infty} \mathbf{P}_{TT}^{\ell} = \left(\mathbf{I} - \mathbf{P}_{TT}\right)^{-1},$$

F =

F(i,j)



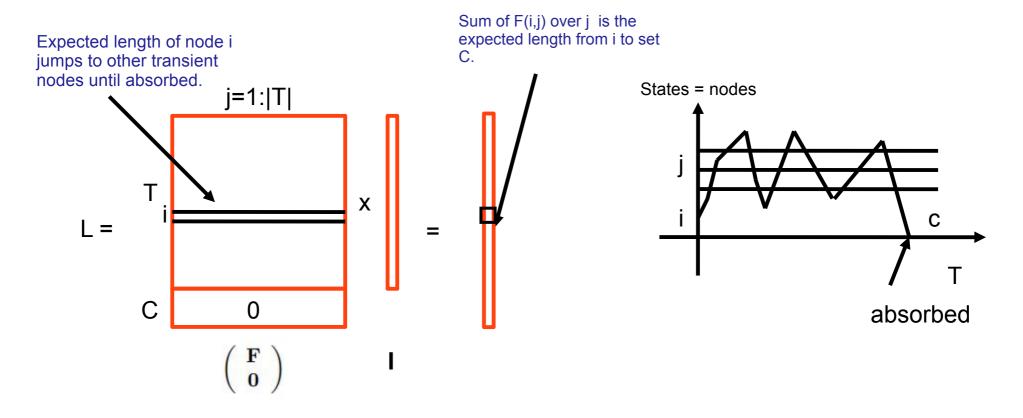
The (i,j)-entry of F is the expected length of the random walk from state i to state j until it is absorbed by C.

Next State i to absorption set C

The expected length of a random walk that starts from node i and reaches set C is given by the i-th element of the following nx1 vector:

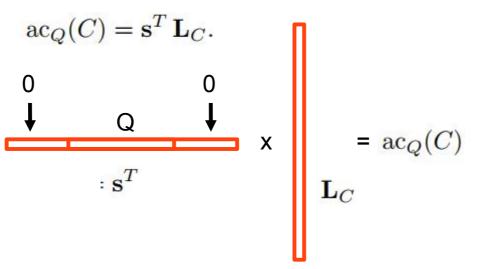
$$\mathbf{L} = \mathbf{L}_C = \left(\begin{array}{c} \mathbf{F} \\ \mathbf{0} \end{array}\right) \mathbf{1},$$

Where I is an Tx1 vector of all 1s.



 $F = i \frac{F(i,j)}{F(i,j)}$

The expected number of steps when starting from Q until being absorbed by C is obtained by summing over all query nodes:



Difficulties:

- Computing objective functions for candidate C requries an expensive matrix inversion.
- Searching for the optimal set C involves considering an exponential number of candidate sets.

How to effciently compute the random walk centrality?

ApproximateAC Algorithm

Compute $ac_Q(C)$ via **AproximateAC** algorithm, which follows from the infinite-sum expansion as:

$$\operatorname{ac}_{Q}(C) = \mathbf{s}^{T} \mathbf{L}_{C} = \mathbf{s}^{T} \begin{pmatrix} \mathbf{F} \\ \mathbf{0} \end{pmatrix} \mathbf{1} = \sum_{\ell=0}^{\infty} \mathbf{x}_{\ell} \mathbf{1},$$

for $\mathbf{x}_{0} = \mathbf{s}^{T}$ and $\mathbf{x}_{\ell+1} = \mathbf{x}_{\ell} \begin{pmatrix} \mathbf{P}_{TT} \\ \mathbf{0} \end{pmatrix}.$

Algorithm 1 ApproximateAC

Input: Transition matrix \mathbf{P}_{TT} , threshold ϵ , starting probabilities \mathbf{s} Output: Absorbing centrality ac_Q $\mathbf{x}_0 \leftarrow \mathbf{s}^T$; $\delta \leftarrow \mathbf{x}_0 \cdot \mathbf{1}$; $\operatorname{ac} \leftarrow \delta$; $\ell \leftarrow 0$ while $\delta < \epsilon$ do $\mathbf{x}_{\ell+1} \leftarrow \mathbf{x}_\ell \begin{pmatrix} \mathbf{P}_{TT} \\ \mathbf{0} \end{pmatrix}$ $\delta \leftarrow \mathbf{x}_{\ell+1} \cdot \mathbf{1}$ $\operatorname{ac} \leftarrow \operatorname{ac} + \delta$ $\ell \leftarrow \ell + 1$ return ac

How to select the k nodes for set C that has lowest ac?

Greedy Algorithm

Greedy Algorithm

- The problem of finding k nodes with minimum random walk centrality is NP-hard.
- Approximation method:

Centrality gain function:

```
mc: mincentrality for k=1
```

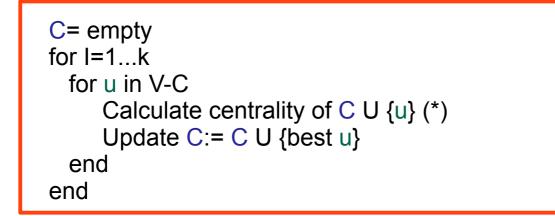
```
gain = mc – centrality, k>1
```

Note: maximize gain equals to minimize centrality.

• Greedy algorithm can guarantee (1-1/e)-appriximation for maximizing gain.

Greedy Algorithm

greedy



The complexity of greedy is O(kn³)

Related Methods

- Personalized Pagerank(PPR). This is the Pagerank alorithm with a damping factor equal to the restart probability and personalization probabilities s(q). It returns the k nodes with the highest PageRank values.
- **Degree centrality(Degree)**. Degree returns the k highest-degree nodes, being oblivious to the query nodes.
- **Distance centrality(Distance)**. Distance returns the k nodes with the highest distance centrality w.r.t. Q.
- **SpectralQ, SpectralC, SpectalD**: Project the original graph into a low-dimenstional space so that distances between nodes in the graph correspond to distance between corresponding projected points.

SpectralC performs k-means clustering on the embedding of thecandidates nodes.

SpectralD & SpectralQperforms k-means clustering on the embedding of the query nodes

data

Sillali	small		large		
Dataset	V	E	Dataset	V	E
		$\frac{ L }{78}$	kddCoauthors	2891	2891
karate	34		livejournal	3645	4141
dolphins	62	159	ca-GrQc	5242	14496
lesmis	77	254	ca-HepTh	9877	25998
	112	425	roadnet	10199	13932
football 1	115	613	oregon-1	11174	23409

cannot run greedy on these

input

Graphs from previous datasets

Query nodes:

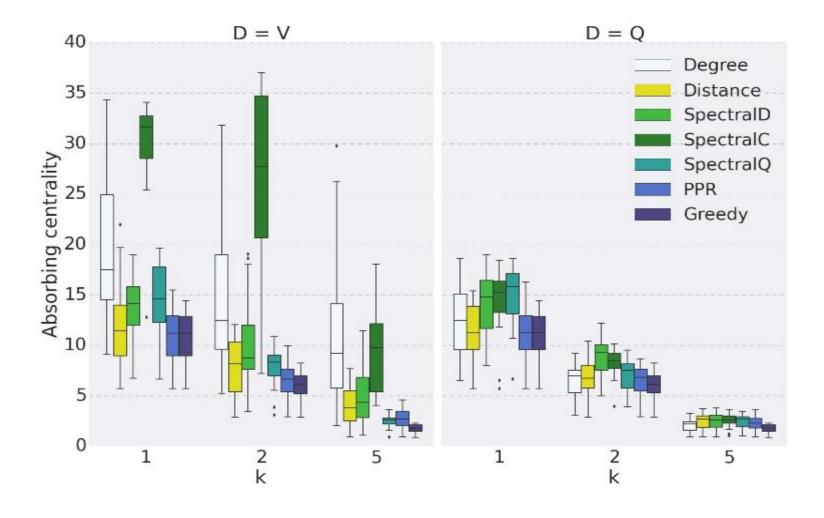
-step1. Select s seed nodes uniformly at random.

-step2. Select a ball B(s,r) around each seed nodes s, with radius r=2.

Step3. From all balls, select a set of query nodes with size q=10 and q=20, respectively for small and large datasets.

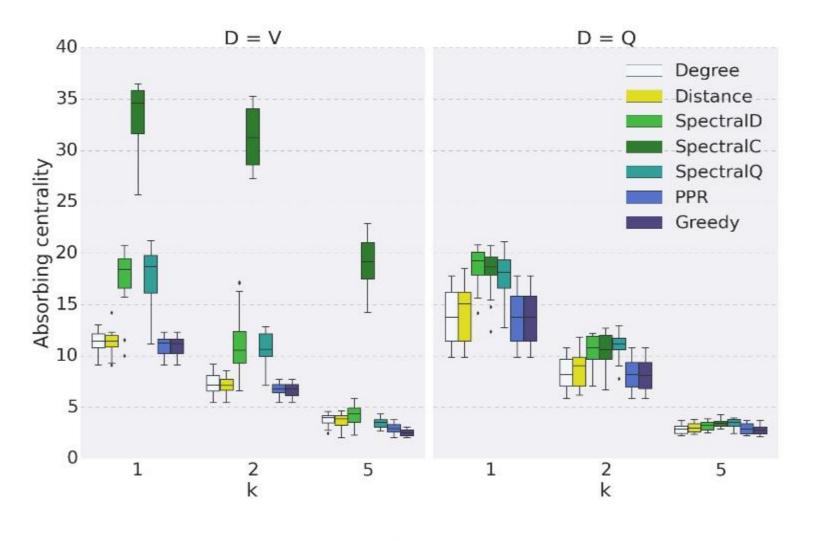
Restart probability is 0.15, and starting probability **s** are uniform over Q.

small graphs



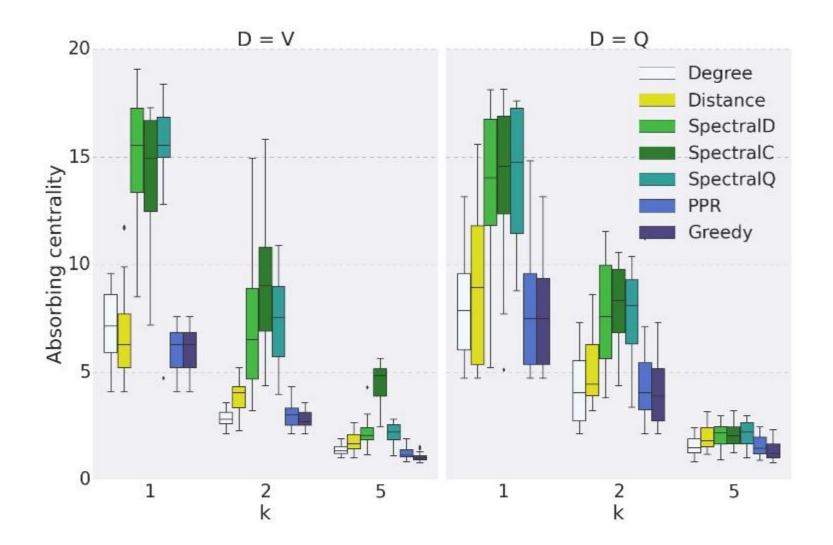
dolphins

small graphs



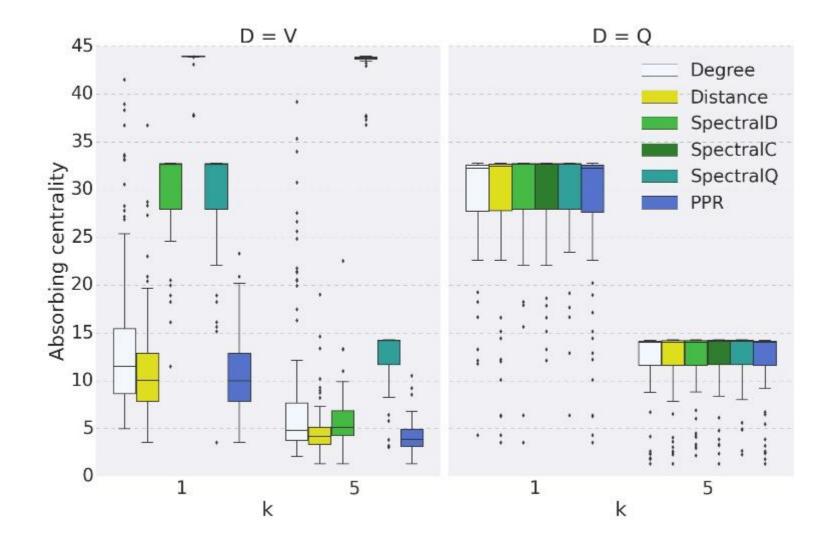
adjnoun

small graphs



karate

large graphs



oregon

Conclusion

- Adressed the problem of finding central nodes in a graph w.r.t. a set of query nodes.
- The centrality measure is based on absorbing random walks: seek to compute k nodes that minimize the expected number of steps that a random walk will need to reach at when starts from the query nodes.
- Show the problem is NP-hard and proposed a Greedy algorithm with complexity O(kn³)
- Compare related algorithms in experiments show that Greedy works well in small datasets.

The end

Thank You!