PageRank Algorithm

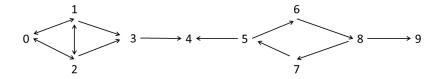
The Page rank algorithm was the initial Google webpage ranking algorithm. It assumes 100% page rank in a directed network. The algorithm computes the probability that a person clicking at random on links will arrive at a particular page. The probability, at any step, that the person will continue is given by the damping factor $\alpha \in (0,1)$.

Suppose there are N nodes in the network.

Algorithm:

- Let $\alpha \in (0,1)$ be the damping factor.
- Use discrete uniform distribution to initialize the rank of each data point. That is, at time 0, each node is assigned rank $\frac{1}{N}$.
- At each step, α of a node's rank is evenly distributed among its outgoing links and the remainder of its rank gets distributed to all the nodes (including itself).
- Iterate until the resulting rank vector stabilizes.

For our example with 10 nodes, let's add directions to edges:



Let $\alpha = 0.84$ and initialize the nodes with rank $r_0(i) = \frac{1}{10}$ for $0 \le i \le 9$. Let's look at step t + 1. Each node will update to rank of at least

$$\sum_{i=1}^{N} \frac{(1-\alpha)r_t(i)}{N} = \frac{1-\alpha}{N} = .016$$

$$r_{t+1}(0) = \frac{\alpha}{3}r_t(1) + \frac{\alpha}{3}r_t(2) + (.016)$$

$$r_{t+1}(1) = \frac{\alpha}{2}r_t(0) + \frac{\alpha}{3}r_t(2) + (.016)$$

$$r_{t+1}(2) = \frac{\alpha}{2}r_t(0) + \frac{\alpha}{3}r_t(1) + (.016)$$

$$r_{t+1}(3) = \frac{\alpha}{3}r_t(1) + \frac{\alpha}{3}r_t(2) + (.016)$$

$$r_{t+1}(4) = \frac{\alpha}{1}r_t(3) + \frac{\alpha}{2}r_t(5) + \frac{\alpha}{1}r_t(4) + (.016)$$

$$r_{t+1}(5) = \frac{\alpha}{1}r_t(7) + (.016)$$

$$r_{t+1}(6) = \frac{\alpha}{2}r_t(5) + (.016)$$

$$r_{t+1}(7) = \frac{\alpha}{2}r_t(8) + (.016)$$

$$r_{t+1}(8) = \frac{\alpha}{1}r_t(6) + (.016)$$

$$r_{t+1}(9) = \frac{\alpha}{2}r_t(8) + \frac{\alpha}{1}r_t(9) + (.016)$$

Note that the ranks add up to 1:

$$\sum_{i=0}^{9} r_{t+1}(i) = \left[\alpha \sum_{i=0}^{9} r_t(i)\right] + (1 - \alpha) = 1$$

For example, at step 1,

$$r_1(0) = .028 + .028 + .016 = .072$$
 $r_1(5) = .084 + .016 = .1$ $r_1(1) = .042 + .028 + .016 = .086$ $r_1(6) = .042 + .016 = .058$ $r_1(2) = .042 + .028 + .016 = .086$ $r_1(7) = .042 + .016 = .058$ $r_1(3) = .028 + .028 + .016 = .072$ $r_1(8) = .084 + .016 = .1$ $r_1(4) = .084 + .042 + .016 + .084 = .226$ $r_1(9) = .042 + .016 + .084 = .142$

The algorithm runs until there is no change in the ranks.

So what are we really finding? If you have encountered Markov chains, then we can interpret the network as a Markov chain with transitions given by weights depending on α . Looking for stability in the network translates to finding a stationary distribution for the Markov chain. For our example, the transition matrix associated with the Markov chain is

$$P = \begin{bmatrix} .016 & .436 & .436 & .016 & .016 & .016 & .016 & .016 & .016 \\ .296 & .016 & .296 & .296 & .016 & .016 & .016 & .016 & .016 \\ .296 & .296 & .016 & .296 & .016 & .016 & .016 & .016 & .016 \\ .016 & .016 & .016 & .016 & .016 & .016 & .016 & .016 & .016 \\ .016 & .016 & .016 & .016 & .0856 & .016 & .016 & .016 & .016 \\ .016 & .016 & .016 & .016 & .436 & .016 & .016 & .016 & .016 \\ .016 & .016 & .016 & .016 & .016 & .016 & .016 & .016 & .016 \\ .016 & .016 & .016 & .016 & .016 & .016 & .016 & .016 & .016 \\ .016 & .016 & .016 & .016 & .016 & .016 & .016 & .016 & .016 \\ .016 & .016 & .016 & .016 & .016 & .016 & .016 & .016 & .016 & .016 \\ .016 & .016 & .016 & .016 & .016 & .016 & .016 & .016 & .016 & .016 \\ .016 & .016 & .016 & .016 & .016 & .016 & .016 & .016 & .016 & .016 \\ .016 & .016 & .016 & .016 & .016 & .016 & .016 & .016 & .016 & .016 \\ .016 & .016 & .016 & .016 & .016 & .016 & .016 & .016 & .016 & .016 \\ .016 & .016 & .016 & .016 & .016 & .016 & .016 & .016 & .016 & .016 \\ .016 & .016 & .016 & .016 & .016 & .016 & .016 & .016 & .016 \\ .016 & .016 & .016 & .016 & .016 & .016 & .016 & .016 & .016 \\ .016 & .016 & .016 & .016 & .016 & .016 & .016 & .016 & .016 \\ .016 & .016 & .016 & .016 & .016 & .016 & .016 & .016 \\ .016 & .016 & .016 & .016 & .016 & .016 & .016 & .016 \\ .016 & .016 & .016 & .016 & .016 & .016 & .016 \\ .016 & .016 & .016 & .016 & .016 & .016 & .016 & .016 \\ .016 & .016 & .016 & .016 & .016 & .016 & .016 \\ .016 & .016 & .016 & .016 & .016 & .016 & .016 \\ .016 & .016 & .016 & .016 & .016 & .016 \\ .016 & .016 & .016 & .016 & .016 & .016 \\ .016 & .016 & .016 & .016 & .016 \\ .016 & .016 & .016 & .016 & .016 \\ .016 & .016 & .016 & .016 \\ .016 & .016 & .016 & .016 \\ .016 & .016 & .016 & .016 \\ .016 & .016 & .016 & .016 \\ .016 & .016 & .016 & .016 \\ .016 & .016 & .016 \\ .016 & .016 & .016 \\ .016 & .016 & .016 \\ .016 & .016 & .016 \\ .016 & .016 & .016 \\ .016 & .016 & .016 \\ .016 & .016 & .016 \\ .016 & .016 & .016 \\ .016 & .016 \\ .016 & .016 \\ .016 & .016 \\ .016 & .016 \\ .016 & .016 \\ .016 & .016 \\$$

P is a stochastic matrix, meaning that each row adds up to 1. In the language of Markov chains, we are looking for a unit length row vector π such that

$$\pi P = \pi$$
.

If you would rather work with column vectors, then transpose both sides of the equation and find a unit length column vector $\mathbf{v} = \pi^T$ such that

$$P^T \mathbf{v} = \mathbf{v}$$

This means that \mathbf{v} is the eigenvector of P^T associated to the largest eigenvalue 1 (this is the largest eigenvalue due to the fact that P is stochastic.) We need to ensure that \mathbf{v} is of length one, so we might have to rescale if we use a calculator! Using computational software, or a calculator, we can find that

$$\mathbf{v} = \begin{bmatrix} 0.042244 \\ 0.046865 \\ 0.046865 \\ 0.042244 \\ 0.441189 \\ 0.045488 \\ 0.035105 \\ 0.035105 \\ 0.045488 \\ 0.219407 \end{bmatrix}$$

Another way to find π is to look at the rows of $\lim_{n\to\infty} P^n$. Under certain circumstances, each row of this limiting matrix will be π , such as in our example.

On the other hand, the PageRank Algorithm should also converge to this probability distribution vector, which we computed here exactly.

Remark: Note the similarity with the eigenvector centrality discussed last lecture. The difference is that here we are working in a directed graph and only α of a node's weight is distributed to its (directed) neighbors.