## Lecture 12 notes

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Centrality<sup>1</sup>

Question: How important is a vertex in a graph?

**Definition:** A structural index (S.I.) of a graph G : (V, E) is a function  $C : V \to \mathbb{R}$  such that for isomorphic graphs  $G, H, C_G(v) = C_H(\phi(v))$ , where  $\phi(v)$  is the image of v in H.

**Definition (Matrix form)**: Let  $f : \mathbb{R}^{n \times n} \to \mathbb{R}^n$  be a function on the adjacency matrix A. f is a structural index if and only if  $P^T f(PAP^T) = f(A)$ , where P is any permutation matrix.

Example of permutation matrices:

If we have 
$$\mathbf{x} = \begin{bmatrix} x_1 \\ x_2 \\ x_3 \\ x_4 \end{bmatrix}$$
 and need  $\mathbf{P}\mathbf{x} = \begin{bmatrix} x_2 \\ x_1 \\ x_3 \\ x_4 \end{bmatrix}$   
By setting  $\mathbf{P} = \begin{bmatrix} 0 & 1 & 0 & 0 \\ 1 & 0 & 0 & 0 \\ 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 1 \end{bmatrix}$   
we have  $\mathbf{P}\mathbf{x} = \begin{bmatrix} 0 & 1 & 0 & 0 \\ 1 & 0 & 0 & 0 \\ 0 & 0 & 1 & 0 \\ 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 1 \end{bmatrix} \begin{bmatrix} x_1 \\ x_2 \\ x_3 \\ x_4 \end{bmatrix} = \begin{bmatrix} x_2 \\ x_1 \\ x_3 \\ x_4 \end{bmatrix}$ 

We also have  $\boldsymbol{P}^T \boldsymbol{P} = \boldsymbol{I}$ 

Example of P applied to adjacency matrix. Suppose we have a graph G shown below:



<sup>1</sup>Chapter 3-5 of Network Analysis

Its adjacency matrix  $\boldsymbol{A}$  is:

$$\boldsymbol{A} = \begin{bmatrix} 0 & 1 & 1 & 0 & 0 & 1 \\ 1 & 0 & 1 & 0 & 0 & 0 \\ 1 & 1 & 0 & 1 & 1 & 0 \\ 0 & 0 & 1 & 0 & 0 & 0 \\ 0 & 0 & 1 & 0 & 0 & 0 \\ 1 & 0 & 0 & 0 & 0 & 0 \end{bmatrix}$$

Suppose G is relabeled into H:



We set the permutation matrix  $\mathbf{P}$  to be a  $6 \times 6$  matrix with 1 at the index  $(label_{new}, label_{old})$ , such that  $\mathbf{P}\mathbf{x}_{old} = \mathbf{x}_{new}$  and  $\mathbf{P}^T\mathbf{x}_{new} = \mathbf{x}_{old}$ .

P =	0	0	0	0	0	1]
	1	0	0	0	0	0
	0	0	0	0	1	0
	0	0	1	0	0	0
	0	0	0	1	0	0
	0	1	0	0	0	0

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\boldsymbol{P}\boldsymbol{A}\boldsymbol{P}^T = (\boldsymbol{P}\boldsymbol{A})\boldsymbol{P}^T
                       0 0
                                  0 0 0
              \begin{bmatrix} 1 & 0 & 0 & 0 & 0 & 0 \\ 0 & 1 & 1 & 0 & 0 & 1 \\ 0 & 0 & 1 & 0 & 0 & 0 \\ 1 & 1 & 0 & 1 & 1 & 0 \\ 0 & 0 & 1 & 0 & 0 & 0 \\ 1 & 0 & 1 & 0 & 0 & 0 \end{bmatrix}
                                                    \boldsymbol{P}^{T} (flipping rows of \boldsymbol{A} according to \boldsymbol{P})
         =
               Γ1
                      0 0
                                  0
                                         0
                                              0]
                                                      Γ0
                                                           1
                                                                   0
                                                                         0 0
                                                                                     0
                                      0
                                                       0 \ 0 \ 0 \ 0 \ 0 \ 1
                0
                     1 \ 1
                                  0
                                              1
                     \begin{bmatrix} 1 & 0 & 0 & 1 \\ 0 & 1 & 0 & 0 & 0 \\ 1 & 0 & 1 & 1 & 0 \\ 0 & 1 & 0 & 0 & 0 \\ 0 & 1 & 0 & 0 & 0 \end{bmatrix}
                                                     0 0 0 1 0 0
                0
                                                                                           (flipping columns according to \boldsymbol{P}^T)
         =
                                                    \begin{bmatrix} 0 & 0 & 0 & 0 & 1 & 0 \\ 0 & 0 & 1 & 0 & 0 & 0 \end{bmatrix}
                1
                0
                                                     1 0 0 0 0
                                                                                     0
                1
               Γ0
                                  0
                                       0
                      1
                           0
                                              0
                1
                      0 \ 0 \ 1 \ 0
                                              1
                      0 \ 0 \ 1
                                              0
               0
                                         0
                                                      (this is the adjacency matrix of H)
          =
                0
                      1
                            1
                                   0
                                         1 1
                0 0 0
                                  1
                                        0 0
               0
                    1 \ 0 \ 1
                                        0 0
From now on, we will use a specific graph \mathcal{G} to illustrate different kinds of
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**Naming convention**: we use C(x) is centrality of vertex x or  $C_x$ . When we have  $C(x) \ge C(y)$  (i.e.  $C_x \ge C_y$ ), we say x is more important than y.

**Example 1**: out-degree and in-degree d(x).

For  $\mathcal{G}$ , we have the degree for each vertex marked as below:



To prove d(x) is an S.I.:

• By isomorphism: obviously d(x) is an S.I. because it is label independent.

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• By matrix:

$$f(\mathbf{A}) = \mathbf{A}\mathbf{e}$$
$$\mathbf{P}^{T}f(\mathbf{P}\mathbf{A}\mathbf{P}^{T}) = \mathbf{P}^{T}(\mathbf{P}\mathbf{A}\mathbf{P}^{T})\mathbf{e}$$
$$= (\mathbf{P}^{T}\mathbf{P})\mathbf{A}(\mathbf{P}^{T}\mathbf{e})$$
$$= \mathbf{I}\mathbf{A}\mathbf{e}$$
$$= \mathbf{A}\mathbf{e}$$
$$= f(\mathbf{A})$$

**Example 2**: Eccentricity.

Let  $e(u) = \max_{v} d(u, v)$ , where d(u, v) denote the distance between u and v. We have  $e_c(u) = \frac{1}{e(u)}$  is an S.I.



## Example 3: Closeness / Transmission number.

Let  $t(u) = \sum_{v} d(u, v)$ , where d(u, v) denote the distance between u and v. We have  $t_c(u) = \frac{1}{t(u)}$  is an S.I.



## Example 4: Betweeness.

 $b_c(u) = \sum_{s,t \neq u} \frac{\sigma_{st}(u)}{\sigma_{st}}$ .  $\sigma_{st}$  is the number of shortest paths between vertices s

and t.  $\sigma_{st}(u)$  is the number of shortest paths between s and t that pass u. Suppose we have a graph shown below, then vertex 4 is the most important vertex according to betweeness.



Example 5: Katz Index.

Consider an adjacency matrix  $\mathbf{A}$  for representing a voting result, where if  $\mathbf{A}_{ij} = 1$ , we say *i* voted for *j*.  $[\mathbf{A}^T \mathbf{e}]$  is the number of votes for *j*. Suppose people there were a set of people who voted for *i* and then *i* voted for *j*. We wanted to count the votes from all of these people who voted for *i* as well. Then the count of votes becomes:

$$\begin{bmatrix} \mathbf{A}^T \mathbf{e} \end{bmatrix} + \begin{bmatrix} (\mathbf{A}^T)^2 \mathbf{e} \end{bmatrix}$$

Then following this logic, why cannot we count the votes in an infinite order:

$$\begin{bmatrix} \mathbf{A}^T \mathbf{e} \end{bmatrix} + \begin{bmatrix} (\mathbf{A}^T)^2 \mathbf{e} \end{bmatrix} + \dots + \begin{bmatrix} (\mathbf{A}^T)^k \mathbf{e} \end{bmatrix} + \dots$$

This scheme has a problem that it'll generate infinite counts. We can modify the counting scheme a little bit by dampening the weight of vote as the order becomes higher. This is accomplished by multiplying  $\alpha$  to  $\boldsymbol{A}$  where  $0 < \alpha < 1$ . Then we have the count of votes as:

$$\left[\alpha \mathbf{A}^{T}\mathbf{e}\right] + \left[(\alpha \mathbf{A}^{T})^{2}\mathbf{e}\right] + \dots + \left[(\alpha \mathbf{A}^{T})^{k}\mathbf{e}\right] + \dots$$

Katz index is then defined as:

$$\mathbf{k} = \sum_{l=1}^{\infty} (\alpha \mathbf{A}^T)^l \mathbf{e}$$

When **A** is 1-by-1 matrix (i.e. scalar 1),  $k_1 = 1 + \alpha + \alpha^2 + \cdots = \frac{1}{1-\alpha}$ , if  $|\alpha| < 1$ . That is,  $k_1$  is a geometric series.

If we generalize geometric series to matrices, we have the Neumann series, which is named after Carl Gottfried Neumann.

**Neumann series:**  $\sum_{l=0}^{\infty} \mathbf{A}^l \to (\mathbf{I} - \mathbf{A})^{-1}$  if  $\rho(\mathbf{A}) < 1$ .  $\rho(\mathbf{A})$  is the spectral radius of  $\mathbf{A}$ . (recall  $\rho(\mathbf{A}) = \max_i(|\lambda_i|)$ ).

Then we can write Katz index as:

$$\mathbf{k} = ((\mathbf{I} - \alpha \mathbf{A}^T)^{-1} - \mathbf{I})\mathbf{e}, \text{ if } \rho(\alpha \mathbf{A}^T) < 1$$
$$(\mathbf{I} - \alpha \mathbf{A}^T)\mathbf{k} = (\mathbf{I} - (\mathbf{I} - \alpha \mathbf{A}^T))\mathbf{e}$$
$$(\mathbf{I} - \alpha \mathbf{A}^T)\mathbf{k} = \alpha \mathbf{A}^T \mathbf{e}$$

We can solve this linear system to get Katz index  ${\bf k}.$  The Richardson method gives:

$$(\mathbf{I} - \alpha \mathbf{A}^{T})\mathbf{x} = \alpha \mathbf{A}^{T}\mathbf{e}$$
$$\mathbf{r}^{(t)} = \mathbf{f} - (\mathbf{I} - \alpha \mathbf{A}^{T})\mathbf{x}^{(t)}$$
$$\mathbf{r}^{(t)} = \mathbf{f} - \mathbf{x}^{(t)} + \alpha \mathbf{A}^{T}\mathbf{x}^{(t)}$$
$$\mathbf{x}^{(t+1)} = \mathbf{x}^{(t)} + \mathbf{r}^{(t)}$$
$$= \mathbf{f} + \alpha \mathbf{A}^{T}\mathbf{x}^{(t)}$$

Let  $\mathbf{k}^{(\ell)} = \sum_{i=1}^{\ell} (\alpha \mathbf{A}^T)^i \mathbf{e}$ . In the homework, we'll see that  $\mathbf{x}^{(t)} = \mathbf{k}^{(t)}$ , i.e. the Richardson method produces a truncated sum of the Neumann series.