## HPSC - Assignment 1

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## **Problem**

Measure the convergence rate of FDM step09.py, given the error function:

error = 
$$\sqrt{\sum_{i,j=1}^{nx,ny} \frac{(p_{exact} - p_{approx})^2}{p_{exact}^2}}$$

The exact solution is available from BEM step02.py:

$$p_{exact} = \frac{x}{4} - 4\sum_{n=odd}^{\infty} \frac{1}{(nx)^2 \sinh 2n\pi}$$

The source code for this assignment can be found at:

https://github.com/gear/HPSC/tree/master/hw

## Answer

In this assignment, I will use SNAP.PY, a network analysis developed by Stanford University, as a computational tool. All illustrations (figures, graphs ...) are drawn using Adobe Illustrator CC 2015.

Question 1: Compute eigenvector centrality and betweenness centrality of each vertex.

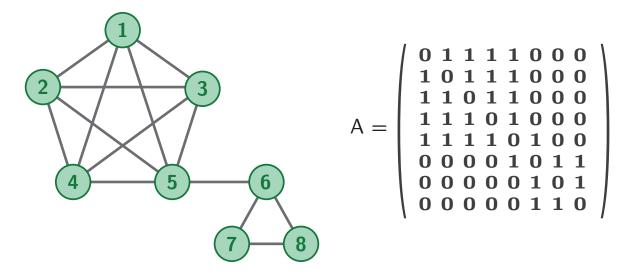


Figure 1: Simple network and its corresponding adjacency matrix.

Eigenvector centrality is given by the formula:

$$x_i^{(t)} = \sum_{j \neq i} A_{ij} \times x_i^{(t)}$$

Rewrite in final matrix form:

$$A\boldsymbol{x} = k_1 \boldsymbol{x}$$

where  $\boldsymbol{x}$  is a vector storing score of all nodes, and  $k_1$  is the largest eigenvalue of the adjacency matrix A.

Listing 1: Eigenvector centrality computation with SNAP.PY

```
# Extracted from UnweightedUndirectedGraph class - File: cn_a2_p1.py

...

import snap as sn

self._graph = sn.LoadEdgeList(sn.PUNGraph, edge_list, 0, 1, '\')

...

# Compute eigenvector centrality and store to a hash table.

def EigenvectorCentrality(self):

# Create a hash map: Int -> Float

NIdEigenH = sn.TIntFltH()

sn.GetEigenVector(self._graph, NIdEigenH)

return NIdEigenH
```

The vector result of Listing 1 is shown as follow:

Figure 2 shows the result in the graph. Eigenvector centrality of each vertex is shown by a blue decimal number next to it. As we can see, vertex number 5 has the highest eigenvector centrality means that vertex number 5 is the most *central* vertex. By looking at the graph, we can intuitively understand this fact.

Betweenness centrality is another metric to measure how important a vertex is within the network. Different from other metric, betweenness centrality measure how important a vertex is in the information flow

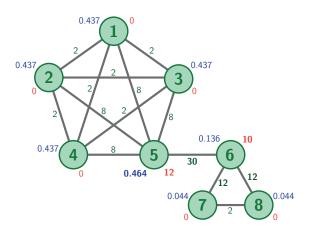


Figure 2: The given network with eigenvector centrality (blue), and betweenness centrality of vertices (red) and edges (green)

between other vertices. In [?], the author defines betweenness centrality  $x_i$  of vertex i as follow:

$$x_i = \sum_{st} n_{st}^i$$
, or  $x_i = \sum_{st} \frac{n_{st}^i}{g_{st}}$ 

where  $n_{st}^i$  is the number of geodesic paths between vertex s and vertex t that go through i, and  $g_{st}$  is the total number of geodesic paths between s and t. Besides the normal betweenness centrality, in [?] the author also mentioned 2 other types of betweenness: flow betweenness and random walk betweenness. However, in this assignment, I will only compute the standard betweenness for the given network.

Listing 2: Betweenness centrality computation with SNAP.PY

```
# Extracted from UnweightedUndirectedGraph class - File: cn_a2_p1.py
2
   import snap as sn
   self._graph = sn.LoadEdgeList(sn.PUNGraph, edge_list, 0, 1, 'u')
   # Compute betweenness centrality of vertices and edges and store to a hash table.
   def BetweennessCentrality(self, isNode = true):
       # Create 2 hash maps: Int -> Float
       NodeScore = sn.TIntFltH()
       EdgeScore = sn.TIntFltH()
10
       \verb|sn.GetBetweennessCentr(self.\_graph, NodeScore, EdgeScore, 1.0)| \\
11
        if (isNode):
            return NodeScore
13
14
       else:
           return EdgeScore
```

The vector result of Listing 2 is shown as follow:

Figure 2 shows the result in the graph. Betweenness centrality of each vertex is shown by a red integer next to it. As we can see, vertex number 5 has the highest betweenness centrality since it is the connection between two cliques.

**Question 2:** Construct Laplacian matrix and Modularity matrix for the given network. Laplacian matrix and Modularity matrix of the graph is computed with *numpy* as follow:

Listing 3: Laplacian matrix and Modularity matrix computation with Numpy

```
1 # Extracted from UnweightedUndirectedGraph class - File: cn_a2_p1.py
2
  import numpy as np
4 ...
   self._adj_matrix = edge_list_to_np(edge_list)
  # Compute and return Laplacian matrix
   def LaplacianMatrix(self):
      D = np.diag(np.sum(self._adj_matrix, 0))
       return D - self._adj_matrix
10
11
12 # Compute and return Modularity matrix
13 def ModularityMatrix(self):
       B = np.zeros(self._adj_matrix.shape)
14
       d = np.sum(self._adj_matrix, 0)
15
       m = sum(d)
16
       for i in range(B.shape[0]):
17
18
           for j in range(B.shape[1]):
19
               B[i,j] = A[i,j] - (d[i] * d[j]) / float(m)
       return B
20
```

The result of Listing 3 is:

$$\mathcal{L} = \begin{pmatrix} 4 & -1 & -1 & -1 & -1 & 0 & 0 & 0 \\ -1 & 4 & -1 & -1 & -1 & 0 & 0 & 0 \\ -1 & -1 & -4 & -1 & -1 & 0 & 0 & 0 \\ -1 & -1 & -1 & -4 & -1 & 0 & 0 & 0 \\ -1 & -1 & -1 & -1 & 5 & -1 & 0 & 0 \\ 0 & 0 & 0 & 0 & -1 & 3 & -1 & -1 \\ 0 & 0 & 0 & 0 & 0 & -1 & 2 & -1 \\ 0 & 0 & 0 & 0 & 0 & -1 & -1 & 2 \end{pmatrix}$$

$$\mathcal{B} = \begin{pmatrix} -0.571 & 0.428 & 0.428 & 0.428 & 0.281 & -0.443 & -0.429 & -0.289 \\ 0.428 & -0.571 & 0.428 & 0.428 & 0.285 & -0.423 & -0.229 & -0.229 \\ 0.428 & 0.428 & -0.571 & 0.428 & 0.285 & -0.423 & -0.229 & -0.229 \\ 0.428 & 0.428 & 0.428 & -0.571 & 0.285 & -0.423 & -0.229 & -0.229 \\ 0.285 & 0.285 & 0.285 & 0.285 & -0.892 & 0.464 & -0.356 & -0.356 \\ 0.428 & -0.428 & -0.428 & -0.428 & 0.464 & -0.327 & 0.729 & 0.729 \\ 0.285 & -0.285 & -0.285 & -0.285 & -0.357 & 0.785 & -0.144 & 0.856 \\ 0.285 & -0.285 & -0.285 & -0.285 & -0.357 & 0.785 & 0.856 & -0.144 \end{pmatrix}$$

(a) Compute the eigenvalue of the Laplacian Matrix, deduce the second smallest eigenvalue, from there perform a spectral bisection on the graph into 2 equal parts.

Listing 4: Perform Spectral Bisection on the graph

```
# Extracted from UnweightedUndirectedGraph class - File: cn_a2_p1.py
2
   import numpy as np
3
4
   self._adj_matrix = edge_list_to_np(edge_list)
5
   # Compute spectral bisection
   def SpectralBisection(self):
       L = self.LaplacianMatrix()
       # Get eigenvectors of \L
10
       u , v = np.linalg.eig(L)
11
       # Get the second smallest eigenvalue and its eigenvector
12
       i = np.argsort(u)[1]
13
       eigv = v[:,i]
14
       partition = np.ones(eigv.shape[0])
15
16
       index_sorted_eigv = np.argsort(eigv)
       for i in range(partition.shape[0] / 2, partition.shape[0]):
17
18
            partition[index_sorted_eigv[i]] = -1
       return partition
```

The vector result for this bisection is demonstrated as follow:

(b) Compute the eigenvalue of the Modularity Matrix, deduce the largest eigenvector, from there divide the network into 2 communities. In spectral modularity, we decide groups by possitiveness of each element in the eigenvector. Maximum modularity is archieved when there is a densense connection within a community and sparse between communities. The computation is performed as follow:

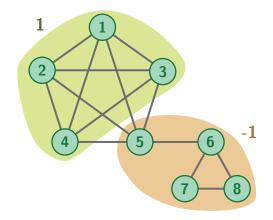


Figure 3: Illustration of spectral bisection.

Listing 5: Perform Spectral Modularity on the graph

```
# Extracted from UnweightedUndirectedGraph class - File: cn_a2_p1.py
2
   import numpy as np
   self._adj_matrix = edge_list_to_np(edge_list)
6
   # Compute spectral modularity
   def SpectralBisection(self):
       B = self.LaplacianMatrix()
       # Get eigenvectors of \L
10
       u , v = np.linalg.eig(L)
11
       # Get the second smallest eigenvalue and its eigenvector
12
       i = np.argsort(u)[1]
13
       eigv = v[:,i]
       partition = np.ones(eigv.shape[0])
15
       index_sorted_eigv = np.argsort(eigv)
16
17
       for i in range(partition.shape[0] / 2, partition.shape[0]):
            partition[index_sorted_eigv[i]] = -1
18
       return partition
```

The vector result for this bisection is demonstrated as follow:

$$partition = \begin{pmatrix} -1 & -1 & -1 & -1 & 1 & 1 \end{pmatrix}$$

Comparing result between Figure 3 and Figure 4, we can see that the spectral modularity gives us a more appropriate result since it divided the given network into two clique. Both technique I used here minimizes the number of connection between communities. However, by forcing spectral bisection to divide network into 2 equal parts, we obtained an unoptimized result. In conclusion, it is clear that modularity

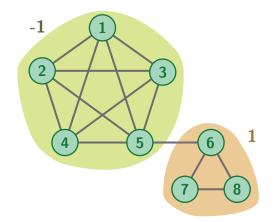


Figure 4: Illustration of spectral modularity.

method gives us a better clustering in both computational value and common sense.

Question 3: Explain quantitatively why "your friends have more friends than you do" in the configuration model. Proof: Suppose we are given a vector  $k = \{k_0, k_1, ..., k_{n-1}\}$  contains the degree of all vertices in the configuration model with n vertices and m edges. Without changing the generality of the statement, we assume that  $k_i > 0 \,\forall k_i \in k$  and n > 0. In this context, the word "friend" of a vertex i implies the adjacency vertices of vertex i. Also,  $k_i$  is the number of friend that vertex i has. Denote  $F_i$  as the number of friends that friends of vertex i has. Quantitatively, the average number of friend that an arbitary vertex i has is:

$$\mathbb{E}(F_i) = \frac{1}{n} \sum_{i=0}^{n-1} k_i = \frac{2m}{n}$$

Therefore, in this configuration model, the average friend that "you" have is 2m/n. On the other hand, the probability that there is a connection between vertex i and vertex j (i and j are friends) is:

$$p(i \leftrightarrow j) = \frac{k_i k_j}{2m - 1} \approx \frac{k_i k_j}{2m}$$

The average number of friends that a friend of vertex i has is:

$$\mathbb{E}(FF_i) = \frac{1}{k_i} \sum_{j=0}^{n-1} p(i \leftrightarrow j) \times k_j = \frac{1}{k_i} \times k_i \sum_{j=0}^{n-1} \frac{k_j}{2m} \times k_j$$
$$= \frac{1}{2m} \sum_{j=0}^{n-1} k_j^2$$

Consider the different  $\mathcal{D}$  between  $\mathbb{E}(F_i)$  and  $\mathbb{E}(FF_i)$ . Note that  $2m = \sum_i k_i$ :

$$\mathcal{D} = \mathbb{E}(FF_i) - \mathbb{E}(F_i) = \frac{\sum_{j=0}^{n-1} k_j^2}{2m} - \frac{2m}{n} = \frac{n \sum_{j=0}^{n-1} k_j^2 - 2\left(\sum_{j=0}^{n-1} k_j\right)^2}{2mn}$$

Expand the square term  $2(...)^2$ , we have the following result:

$$\mathcal{D} = \frac{(n-1)\sum_{j=0}^{n-1} k_j^2 - 2\sum_{i < j} k_i k_j}{2mn}$$

Using the Cauchy-Schwarz inequality, we have:

$$k_i^2 + k_j^2 \ge 2k_i k_j$$

Therefore, using Cauchy-Schwarz for n(n-1)/2 pairs of  $k_i, k_j$  we have:

$$(n-1)\sum_{j=0}^{n-1} k_j^2 \ge 2\sum_{i < j} k_i k_j$$

Hence,

$$\mathcal{D} = \frac{(n-1)\sum_{j=0}^{n-1} k_j^2 - 2\sum_{i < j} k_i k_j}{2mn} \ge 0$$

In conclusion, the different between average number of a friend's friend and your friend is  $\mathcal{D} \geq 0$ , therefore the statement "(on average) your friends have more friends than you do" holds true. The equal sign happens when everyone has exactly 1 friend. QED.

**Question 4:** Example of parameter  $\beta$  and  $\gamma$  of SIR model. In the SIR epidemic model,  $\beta$  is the infection rate and  $\gamma$  is recover (or death) rate. The model is represented as a system of differential equation:

$$\frac{ds}{dt} = -\beta sx; \quad \frac{dx}{dt} = \beta sx - \gamma x; \quad \frac{dr}{dt} = \gamma x$$

where (s, x, r) are fraction of population that are *subceptible*, *recovered*, and *infectionous* respectively. Since the solution of these diffirential equation is not analytical, numerical method is applied to analyze this SIR model. In real-life application, the parameter  $\beta$  and  $\gamma$  is determined empirically by trying to fit known data with some trial value of  $\beta$  and  $\gamma$ . The result for each value (s, x, r) in this exercise is stored in an array, which is indexed my time step. According to [?], the epidemic happens when  $\beta > \gamma$ , which means the disease spreads faster than human recovery. I plotted the time graph for 4 cases:

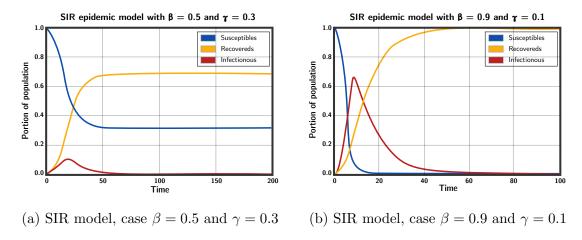


Figure 5: SIR model in case of epidemic happens.

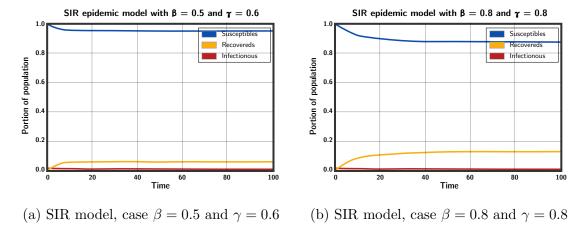


Figure 6: SIR model in case of no epidemic happens.