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CS4423-Networks: Week 6 (19+20 Feb 2025)

Part 3: Eigenvector Centrality - Computation

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This Jupyter notebook, and PDF and HTML versions, can be found at https://www.niallmadden.ie/2425-CS4423/#Week06

This notebook was written by Niall Madden, adapted from notebooks by Angela Carnevale.

Modules for this notebook

```
In [1]: import networkx as nx
import numpy as np
opts = { "with_labels": True, "node_color": "xkcd:sky blue"} # show labels; nodes are
np.set_printoptions(precision=3) # just display arrays to 3 decimal places
np.set_printoptions(suppress=True) # avoid scientific notation (better for matrices)
```

Eigenvector Centrality

So now we know (see CS4423-W06-Part-2.pdf)

- The adjacency matrix, A of a connected graph, G, is an irreducible non-negative matrix.
- So the F-B Therom applies to it
- So ${\cal A}$ has an eigenvalue that is real and positive, and greater than the modulus of any other.
- It has a corresponding positive eigenvector, $ec{v}.$
- v_i is the Eigenvector Centrality node i.

Normalisation

One minor issue is that any multiple of \vec{v} is also an eigenvector for the same eigenvalue. This is not a major problem: we are mainly interested in if, for example $v_i > v_j$ which would mean that Node i has greater centrality than Node j.

Nonetheless, by convection we choose $ec{v}$ so that

- + $ec{v}>0$ (already discussed)
- + $ec{v}^Tec{v}=1$ (equivalently, $\|ec{v}\|_2=\sqrt{v_1^2+v_2^2+\ldots v_n^2}=1$

We say such an eigenvector is **normalised**.

Computing Eigenvalues

Presently, we'll learn about a method called the Power Method

For now, though, we'll use the np.linalg.eig() which computes the eigenvalues and eigenvectors of a matrix:

l, V = np.linalg.eig(A) computes

- l : an array of length *n* containing the eigenvalues of *A*. (Note: we can't call this array lambda, since that is a keyword in Python.
- V : a n imes n matrix; column i of V is the eigenvector corresponding to the eigenvalues λ_i .

Example: Find the eigenvalues and corresponding eigenvectors of

$$A = \begin{pmatrix} 2 & 2 \\ 3 & 1 \end{pmatrix}$$

```
In [2]: A = np.array([[2,2],[3,1]])
l, V = np.linalg.eig(A)
print(f"The eigenvalues of A are {l[0]} and {l[1]}.")
print(f"The corresponding eigenvectors are {V[:,0]} and {V[:,1]}")
```

The eigenvalues of A are 4.0 and -1.0. The corresponding eigenvectors are [0.707 0.707] and [-0.555 0.832]

Let's check if this worked:

```
In [3]: print(A@V[:,0])
print(l[0]*V[:,0])
[2.828 2.828]
[2.828 2.828]
```

Let's check if the columns of V are normalised:

```
In [4]: print(f" ||v|| = {np.linalg.norm(V[:,1])}")
```

||v|| = 1.0

Eigenvalues of adjacency matrices

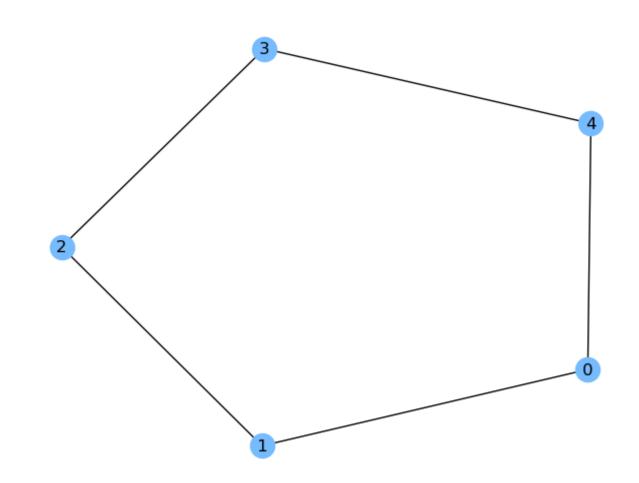
Let's look at some examples, for which we may have an intuition for the centrality. we want to check that

- there is a dominant positive eigenvalue
- there is a corresponding positive eigenvector
- the centrality values seem "sensible"

(Note: Extra details given on the white board!)

Example 1: $G = C_5$

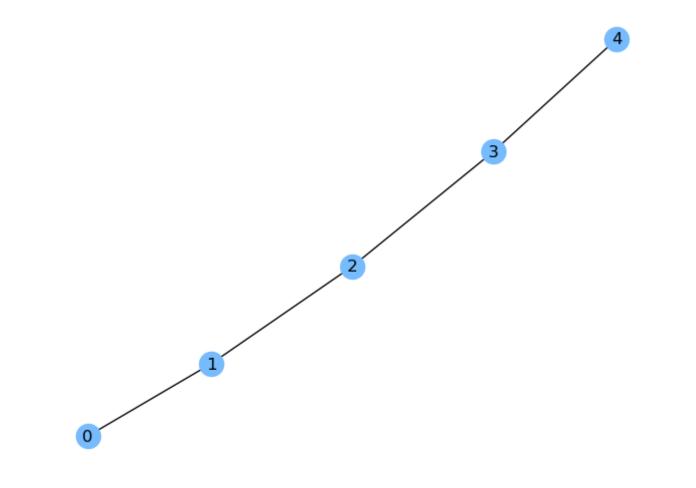
In [5]: G = nx.cycle_graph(5)
nx.draw(G, **opts)



```
In [6]: A = nx.adjacency_matrix(G).toarray()
l, V = np.linalg.eig(A)
for i in range(G.order()):
    print(f"eigenvalue {l[i]:6.3f} has eigenvectors are {V[:,i]}")
eigenvalue -1.618 has eigenvectors are [ 0.632 -0.512 0.195 0.195 -0.512]
eigenvalue 0.618 has eigenvectors are [-0.632 -0.195 0.512 0.512 -0.195]
eigenvalue 2.000 has eigenvectors are [-0.447 -0.447 -0.447 -0.447]
eigenvalue -1.618 has eigenvectors are [-0.032 0.397 -0.611 0.591 -0.345]
eigenvalue 0.618 has eigenvectors are [ 0.074 0.62 0.309 -0.429 -0.575]
```

Example 1: $G = P_5$

In [7]: G = nx.path_graph(5)
nx.draw(G, **opts)



```
In [8]: A = nx.adjacency_matrix(G).toarray()
        l, V = np.linalg.eig(A)
        for i in range(G.order()):
            print(f"eigenvalue {l[i]:6.3f} has eigenvectors are {V[:,i]}")
       eigenvalue 1.732 has eigenvectors are [0.289 0.5
                                                          0.577 0.5
                                                                      0.2891
       eigenvalue -1.732 has eigenvectors are [-0.289 0.5
                                                            -0.577 0.5
                                                                          -0.289]
       eigenvalue -1.000 has eigenvectors are [-0.5 0.5 -0. -0.5 0.5]
       eigenvalue -0.000 has eigenvectors are [ 0.577 -0.
                                                            -0.577 0.
                                                                           0.577]
       eigenvalue 1.000 has eigenvectors are [-0.5 -0.5 -0.
                                                              0.5 0.51
```

Example 3

Let's look at the eigenvalues of an adjacency matrix of a graph. It is constructed so that node 0 is more "central" than any of the others, node 5 is the least "central".

```
In [9]: G = nx.Graph(["01", "02", "03", "04", "05", "12", "13", "14"])
nx.draw(G, **opts)
```

<pre>In [10]: A = nx.adjacency_matrix(G).toarray() print(A)</pre>
$\begin{bmatrix} [0 & 1 & 1 & 1 & 1 \\ 1 & 0 & 1 & 1 & 1 \\ 1 & 1 & 0 & 0 & 0 \\ 1 & 1 & 0 & 0 & 0 \end{bmatrix}$ $\begin{bmatrix} 1 & 1 & 0 & 0 & 0 \\ 1 & 1 & 0 & 0 & 0 \\ 1 & 1 & 0 & 0 & 0 \end{bmatrix}$ $\begin{bmatrix} 1 & 0 & 0 & 0 \\ 1 & 0 & 0 & 0 \end{bmatrix}$
<pre>In [11]: l, V = np.linalg.eig(A) print(f"The eigenpairs of A are:") for i in range(5): print(f"Eigenvalue {l[i]:8.3f} with eigenvector {V[:,i]}")</pre>
The eigenpairs of A are:Eigenvalue3.102 with eigenvector [0.568 0.523 0.352 0.352 0.352 0.183]Eigenvalue0.344 with eigenvector [0.296 -0.343 -0.138 -0.138 -0.138 0.859]Eigenvalue-2.123 with eigenvector [-0.557 -0.324 0.415 0.415 0.415 0.262]Eigenvalue-1.323 with eigenvector [0.529 -0.71 0.137 0.137 0.137 -0.4]Eigenvalue-0.000 with eigenvector [-000.509 0.807 -0.298 -0.]

The Power Method

There are subfields in the *Numerical Linerar Algebra* dedicated to computing estimates for eigenvalues and eigenvectors. When we only need one eigenvalue, and it is the largest, use the **Power method**:

- 1. start with any $u=(1,1,\ldots,1)$, say;
- 2. keep replacing $u \leftarrow Au$ until $u/\|u\|$ becomes stable ...

Questions Does this work? Meaning:

- Does the sequence actually converge?
- Does it return the correct values?

```
In [12]: n = G.order()
u = np.ones((n,1)); u=u/np.linalg.norm(u)
for i in range(10):
    u = A @ u
    u = u/np.linalg.norm(u)
```

In [13]: print(u)

[[0.564] [0.521]

[0.354]

[0.354]

[0.354]

[0.185]]

In [14]: print(V)

[[0.568	0.296	-0.557	0.529	-0.	0.]
[0.523	-0.343	-0.324	-0.71	-0.	0.]
[0.352	-0.138	0.415	0.137	-0.509	0.8	316]
[0.352	-0.138	0.415	0.137	0.807	-0.4	108]
[0.352	-0.138	0.415	0.137	-0.298	-0.4	108]
[0.183	0.859	0.262	-0.4	-0.	0.]]

Finished here Thursday