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NRun ■ C >>
                                                        :::::
                                         Code
                                                     ~
 In [3]: import numpy as np
         import matplotlib.pyplot as plt
         1a)
 In [8]: # Circle topology
         # Unweighted adjacency matrix
         # Option 1: Manually enter the entries
         Atilde = np.array(
                  [[0,1,0,0,0,0,0,1],
                   [1,0,1,0,0,0,0,0],
                   [1,1,0,1,1,0,0,0],
                   [0,0,1,0,1,0,0,0],
                   [0,0,0,1,0,1,0,0],
                   [0,0,0,0,1,0,1,0],
                   [0,0,0,0,0,1,0,1],
                   [1,0,0,0,0,0,1,0]])
         # Option 2: or you can exploit the patterns
         # Atilde = np.zeros((8,8))
         # for i in range(8): #
              Atilde[i,(i+1)\%8] = 1
               Atilde[i,(i-1)\%8] = 1
         # Atilde[2,0] = 1
         # Atilde[2,4] = 1
         print('Unweighted adjacency matrix')
         print(Atilde)
         print(' ')
         Unweighted adjacency matrix
         [[01000001]
          [10100000]
          [1 1 0 1 1 0 0 0]
          [0 0 1 0 1 0 0 0]
          [0 0 0 1 0 1 0 0]
          [0 0 0 0 1 0 1 0]
          [0 0 0 0 0 1 0 1]
          [10000010]]
         1b)
In [12]: # Find weighted adjacency matrix
         # option 1: normalize columns with a for loop
         A = np.zeros((8,8), dtype=float)
         for k in range(8):
             norm = np.sum(Atilde[:,k])
             A[:,k] = Atilde[:,k]/norm
         # option 2: normalize using numpy.sum() and broadcasting, in a single line
         #A = ???
         print('Weighted adjacency matrix')
         print(A)
         Weighted adjacency matrix
         [[0.
                      0.5
                                            0.
                                                       0.
                                                                  0.
          [0.3333333 0.
                                 0.5
                                                       0.
                                                                 0.
           0.
          [0.3333333 0.5
                                                       0.33333333 0.
                                 0.
                                            0.5
           0.
                      0.
          [0.
                      0.
                                 0.5
                                            0.
                                                       0.33333333 0.
           0.
                                                                  0.5
          [0.
                                            0.5
                                                       0.
                                 0.
           0.
                                                       0.33333333 0.
          [0.
                                 0.
                                            0.
           0.5
          [0.
                                                       0.
                                                                  0.5
                                 0.
                                            0.
                      0.5
          [0.33333333 0.
                                                       0.
                                                                  0.
                                 0.
                                            0.
           0.5
         1c) and 1d)
In [14]: # Power method
         b0 = 0.125*np.ones((8,1))
         print('b0 = ', b0)
         print(' ')
         b1 = 0.125*np.ones((8,1))
         print('b1 = ', b1)
         print(' ')
         b = b0.copy()
         for k in range(1000):
             b = 0.125*b
         print('1000 iterations')
         print('b = ',b)
         b0 = [[0.125]]
          [0.125]
          [0.125]
          [0.125]
          [0.125]
          [0.125]
          [0.125]
          [0.125]]
         b1 = [[0.125]]
          [0.125]
          [0.125]
          [0.125]
          [0.125]
          [0.125]
          [0.125]
          [0.125]]
         1000 iterations
         b = [[0.]]
          [0.]
          [0.]
          [0.]
          [0.]
          [0.]
          [0.]
          [0.]]
         1e) Explination goes here.
 In [ ]: # yes, since the eigenvector corresponds to each node reflects the importance
         2a)
In [18]: # Hub topology
         Atildehub = np.array(
                  [[0,1,0,0,0,0,0,1,1],
                   [1,0,1,0,0,0,0,0,1],
                   [1,1,0,1,1,0,0,0,1],
                   [0,0,1,0,1,0,0,0,1],
                   [0,0,0,1,0,1,0,0,1],
                   [0,0,0,0,1,0,1,0,1],
                   [0,0,0,0,0,1,0,1,1],
                   [1,0,0,0,0,0,1,0,1],
                   [1,1,1,1,1,1,1,0]])
         print('Unweighted adjacency matrix')
         print(Atildehub)
         print(' ')
         Unweighted adjacency matrix
         [[010000011]
          [101000001]
          [1 1 0 1 1 0 0 0 1]
          [0 0 1 0 1 0 0 0 1]
          [0 0 0 1 0 1 0 0 1]
          [0 0 0 0 1 0 1 0 1]
          [0 0 0 0 0 1 0 1 1]
          [100000101]
          [1 1 1 1 1 1 1 0]]
         2b)
In [19]:
         # find weighted adjacency matrix
         A = np.zeros((9,9), dtype=float)
         for k in range(9):
             norm = np.sum(Atildehub[:,k])
             A[:,k] = Atildehub[:,k]/norm
         Ahub = A
         print('Weighted adjacency matrix')
         print(Ahub)
         Weighted adjacency matrix
                      0.33333333 0.
         [[0.
                                            0.
                                                       0.
                                                                  0.
           0.
                      0.33333333 0.125
          [0.25
                                 0.33333333 0.
                                                       0.
                                                                  0.
           0.
                                 0.125
          [0.25
                      0.33333333 0.
                                            0.3333333 0.25
                                                                  0.
           0.
                                 0.125
          [0.
                                 0.33333333 0.
                                                       0.25
                                                                  0.
           0.
                                 0.125
          [0.
                                 0.
                                            0.33333333 0.
                                                                  0.33333333
           0.
                                 0.125
          [0.
                                                                  0.
                                                       0.25
           0.33333333 0.
                                 0.125
          [0.
                                                       0.
                                                                  0.33333333
                      0.33333333 0.125
                                                                  0.
          [0.25
                                 0.
                                                       0.
           0.33333333 0.
                                 0.125
          [0.25
                      0.33333333 0.33333333 0.33333333 0.25
                                                                  0.33333333
           0.33333333 0.33333333 0.
         2c) and 2d)
In [20]: b0 = (1/9)*np.ones((9,1))
         print('b0 = ', b0)
         print(' ')
         bhub1 = (1/9)*np.ones((9,1))
         print('bhub1 = ', bhub1)
         print(' ')
         bhub = b0.copy()
         for k in range(1000):
             bhub = (1/9)*bhub
         print('1000 iterations')
         print('bhub = ', bhub)
         print(' ')
         bhubr = b0.copy()
         for k in range(100):
             bhubr = (1/9)*bhubr
         print('100 iterations')
         print('bhubr = ',bhubr)
         b0 = [[0.11111111]]
          [0.1111111]
          [0.1111111]
          [0.1111111]
          [0.1111111]
          [0.1111111]
          [0.1111111]
          [0.1111111]
          [0.1111111]]
         bhub1 = [[0.11111111]]
          [0.1111111]
          [0.1111111]
          [0.1111111]
          [0.1111111]
          [0.1111111]
          [0.1111111]
          [0.1111111]
          [0.1111111]]
         1000 iterations
         bhub = [[0.]]
          [0.]
          [0.]
          [0.]
          [0.]
          [0.]
          [0.]
          [0.]
          [0.]]
         100 iterations
         bhubr = [[4.18317994e-97]
          [4.18317994e-97]
          [4.18317994e-97]
          [4.18317994e-97]
          [4.18317994e-97]
          [4.18317994e-97]
          [4.18317994e-97]
          [4.18317994e-97]
          [4.18317994e-97]]
```

Complete 2e and 2f below.

In [ ]: # e) yes, since the eigenvector corresponds to each node reflects the importance

Jupyter Period14\_Activity\_start (autosaved)

Insert

Cell

Kernel

File

Edit

View

Help

Widgets

Logout

Python 3 O

Trusted