```
# Project 13: Social networks, clustering, and eigenvalue problems
import numpy as np
import matplotlib.pyplot as plt
import scipy.io
from scipy.linalg import eigh
from scipy.io import loadmat
print("\nSubtask 1\n")
# 1. Simple graph: Define adjacency matrix
AdjMatrix = np.array([[0, 1, 1, 0],
[1, 0, 0, 1],
[1, 0, 0, 1],
[0, 1, 1, 0]])
print("Adjacency Matrix:")
print(AdjMatrix)
    Subtask 1
    Adjacency Matrix:
    [[0 1 1 0]
     [1 0 0 1]
     [1 0 0 1]
     [0 1 1 0]]
print("\nSubtask 2\n")
# 2. Find the row sums of the matrix AdjMatrix
RowSums = np.sum(AdjMatrix, axis=1)
print("\nRow Sums:")
print(RowSums)
    Subtask 2
    Row Sums:
     [2 2 2 2]
```

```
print("\nSubtask 3\n")
# 3. Compute the Laplacian of the graph
LaplaceGraph = np.diag(RowSums) - AdjMatrix
print("\nLaplacian Matrix:")
print(LaplaceGraph)
# Check if LaplaceGraph is singular
test vector = np.ones(len(LaplaceGraph))
singularity_check = LaplaceGraph @ test_vector
print("\nSingularity Check (Laplacian * ones):")
print(singularity_check)
    Subtask 3
    Laplacian Matrix:
     [[2-1-1 0]
     [-1 \ 2 \ 0 \ -1]
     [-1 \ 0 \ 2 \ -1]
     [0 -1 -1 2]
    Singularity Check (Laplacian * ones):
     [0. 0. 0. 0.]
print("\nSubtask 4\n")
# 4. Find eigenvalues and eigenvectors using the eig function
D, V = np.linalg.eig(LaplaceGraph)
\rightarrow
    Subtask 4
```

```
print("\nSubtask 5\n")
d, ind = np.argsort(D), np.argsort(D)
D = np.diag(D[ind])
V = V[:, ind]
print("\nEigenvalues (sorted):")
print(np.diag(D))
print("\nEigenvectors (sorted):")
print(V)
    Subtask 5
    Eigenvalues (sorted):
     [-2.22044605e-16 \quad 2.00000000e+00 \quad 2.00000000e+00 \quad 4.00000000e+00]
    Eigenvectors (sorted):
     [[ 5.00000000e-01 4.08248290e-01 7.07106781e-01 -5.00000000e-01]
     [5.00000000e-01 -5.77350269e-01 4.80181756e-16 5.00000000e-01]
     [ 5.00000000e-01 5.77350269e-01 -1.77321568e-16 5.00000000e-01]
      [ 5.00000000e-01 -4.08248290e-01 -7.07106781e-01 -5.00000000e-01]]
print("\nSubtask 6\n")
# 6. Identify the second smallest eigenvalue and its corresponding eigenvector
second smallest eigenvalue = D[1, 1]
V2 = V[:, 1]
# Ensure V2 has a positive first entry
if V2[0] < 0:
   V2 = -V2
print("\nSecond Smallest Eigenvalue:")
print(second smallest eigenvalue)
print("\nEigenvector corresponding to the second smallest eigenvalue (V2):")
print(V2)
    Subtask 6
    Second Smallest Eigenvalue:
    1.999999999999999
    Eigenvector corresponding to the second smallest eigenvalue (V2):
     [ 0.40824829 -0.57735027  0.57735027 -0.40824829]
print("\nSubtask 7\n")
# 7. Separate the elements of the eigenvector V2
pos = []
neg = []
```

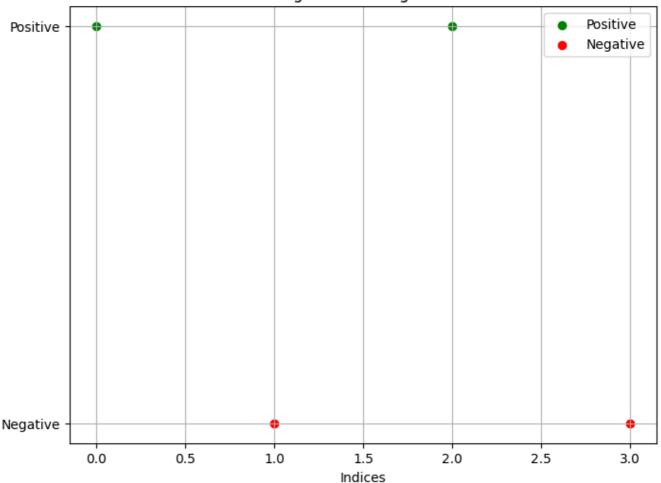
```
for j in range(len(V2)):
    if V2[j] > 0:
       pos.append(j)
    else:
        neg.append(j)
print("\nPositive Indices (V2 > 0):")
print(pos)
print("\nNegative Indices (V2 <= 0):")</pre>
print(neg)
# Optional: Visualize the clusters
plt.figure(figsize=(8, 6))
plt.scatter(pos, [1]*len(pos), color='green', label='Positive')
plt.scatter(neg, [0]*len(neg), color='red', label='Negative')
plt.yticks([0, 1], ['Negative', 'Positive'])
plt.title('Clustering based on Eigenvector V2')
plt.xlabel('Indices')
plt.legend()
plt.grid()
plt.show()
```



Subtask 7

Positive Indices (V2 > 0):
[0, 2]
Negative Indices (V2 <= 0):
[1, 3]</pre>

## Clustering based on Eigenvector V2

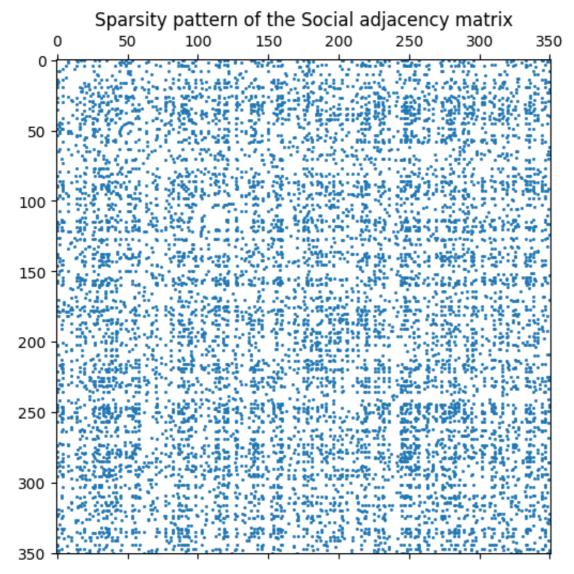


```
print("\nSubtask 8\n")
# 8. Load the data
data = loadmat("/content/social.mat")
Social = data['Social']
print("Loaded Social adjacency matrix with shape:", Social.shape)
# Spy plot of the Social matrix
plt.figure(figsize=(8, 6))
plt.spy(Social, markersize=1)
plt.title('Sparsity pattern of the Social adjacency matrix')
plt.show()
```

**→**▼

Subtask 8

Loaded Social adjacency matrix with shape: (351, 351)



```
print("\nSubtask 9\n")
# 9. Define DiagSocial and LaplaceSocial
DiagSocial = np.sum(Social, axis=1)
LaplaceSocial = np.diag(DiagSocial) - Social
print("\nDiagonal matrix DiagSocial:")
print(DiagSocial)
print("\nLaplacian matrix LaplaceSocial:")
print(LaplaceSocial)
     Subtask 9
    Diagonal matrix DiagSocial:
     [42 19 25 36 50
                             2 29
                                    5 28
                                                8 38 17 47
                                                            4 29 49 28 17
                      5 7
                                          5
                                            4
                                                      32
       8 43 25 48 30 20 43 47
                               6 39 34 46 40 50 27
                                                         15 46 33 31 24 42
       6 44 26
                5 24 47
                          6 24 50
                                   7 56 11
                                             4
                                                4 15
                                                     24
                                                          9 21
                                                                7 22 23
                                                                          5
                                                7 56 34 49
          5
             9
                   4 45
                          4 23
                                5 35 46
                                          9
                                             8
                                                             9 48 42
                                                                       6 44 33
                4
                   8
                      4 21 43 28
                                   6 21
                                            5
                                                8 23
                                                       3
                                                                  48 18
      29 24 50
                4
                                         26
                                                         27 51 42
                                                       3 34
      46 45 25
                4
                   4
                      7
                         22 52 41
                                    4
                                       5
                                         24 16
                                                8 20
                                                             4 46 46 43
                                                                         27 43
                                                                               32
       4 42 42
                4
                   5
                       4
                         26
                             7
                               20
                                    7 51 26 32 48 42
                                                      55 48
                                                             6
                                                                6
                                                                    7
                                                                      28
      29 33 19 39
                       6
                          6 47
                               32 45 45 22 30 39 19
                                                       5
                                                         27 44 30
                                                                  34
                                                                      43
                                                                          5
                                6 34 51
                                               29 32
                                                       5
                                                                       5
      19 26 35
                7 50 31
                          6 31
                                          7
                                             6
                                                          4
                                                            18 29
                                                                         38 32
                                5 18 49
      43 24 46 54 34 48
                            5
                                         14 47
                                               22 28
                                                            22
                                                                6
                                                                    7
                                                                      45
                                                                          5
                          6
                                                     46 46
       6 16 23
               6 41 43 49 50 40 48 41 44
                                            23 31 27 40 46 52
                                                                5
                                                                   5
                                                                       5 27
                                                                            25
             4 36 16 43
                            5
                                7 44
                                     55 41
                                             3 29 50 34 39
                                                            25 44 50
                                                                       8 47 19
      47
                          8
                                          3
      21 46 40 34 22
                       6 48 28 41 29 18
                                             5
                                                6 52 44
                                                         8
                                                             6
                                                                5 18
                                                                       6 42 44 45
                                          5 49 45 15 17 23
                                                                5
          6 25 33 51 20
                          5 32 41
                                    5
                                      20
                                                             4
                                                                   9
                                                                       7 50 36 51
                          5
                                    9
                   6 30
                             5 48
                                       8 53 21 19 221
     Laplacian matrix LaplaceSocial:
                                                  0
     [ [
                          42
                                                                        0
                           0
                                                  0
                                                                        01
                           0
                                                19
                                                                        0 ...
       18446744073709551615 18446744073709551615
                                                                        01
                                                                       25 ...
                           0
                                                  0
                                                                        01
                           0 18446744073709551615
                          21
                                                  0 18446744073709551615
                           0 18446744073709551615
                           0
                                                19
                                                    18446744073709551615]
                           0
       18446744073709551615 18446744073709551615
                                                                       2211
print("\nSubtask 10\n")
# 10. Compute eigenvalues and eigenvectors
D, V = np.linalg.eig(LaplaceSocial)
print("\nEigenvalues (D):")
print(D)
```

```
print("\nEigenvectors (V):")
print(V)
# Check the shapes
print("Shape of V (eigenvectors):", V.shape)
print("Shape of D (eigenvalues):", D.shape)
d, ind = np.argsort(D), np.argsort(D)
D = np.diag(D[ind])
V = V[:, ind]
print("\nEigenvalues (sorted):")
print(np.diag(D))
print("\nEigenvectors (sorted):")
print(V)
     -7.93309180e+18 -7.30039076e+18 -6.60029818e+18 -6.19182146e+18
     -5.42242603e+18 -5.09404014e+18 -4.64429621e+18 -4.04194737e+18
     -3.42093068e+18 -2.67495781e+18 -2.22908528e+18 -1.67326547e+18
     -1.35945445e+18 -1.10252185e+18 -7.33223777e+17 -1.50655293e+16
      3.66850571e+17 8.38595118e+17 1.13289697e+18 1.34061746e+18
      2.00955967e+18 2.39555123e+18 2.74286959e+18 3.22248239e+18
      3.69741316e+18 4.26543390e+18 4.78141228e+18 5.66915645e+18
      5.96810177e+18 6.61028109e+18 6.86736139e+18 7.26178903e+18
      7.41826324e+18 8.17536984e+18 8.43627167e+18 8.60545879e+18
      9.28826591e+18 1.02872225e+19 1.14035216e+19 1.16610911e+19
      1.21049666e+19
                      1.27651677e+19
                                      1.32045922e+19 1.34443840e+19
      1.42145110e+19
                      1.44559151e+19 1.51891359e+19 1.54114602e+19
      1.62568233e+19
                      1.66780272e+19
                                      1.70721347e+19
                                                     1.76953777e+19
                      1.85117454e+19
                                     1.88675545e+19
                                                     1.91554788e+19
      1.78689735e+19
      2.01182608e+19
                      2.10114103e+19 2.14614933e+19
                                                      2.23804992e+19
      2.31046764e+19
                      2.35318090e+19
                                      2.39152487e+19
                                                      2.45434280e+19
      2.59155943e+19
                      2.62691177e+19
                                      2.67046286e+19 2.69038645e+19
                      2.89546542e+19
      2.75392242e+19
                                      2.90832364e+19
                                                      2.97080121e+19
      3.01780849e+19
                      3.04988154e+19 3.07283472e+19
                                                     3.12584236e+19
      3.19791971e+19
                      3.24261364e+19
                                      3.29021359e+19
                                                      3.32991321e+19
      3.37137747e+19
                      3.53635056e+19
                                      3.57735661e+19
                                                      3.60231511e+19
      3.66045702e+19
                      3.74247446e+19
                                      3.76807250e+19 3.85223460e+19
                                                     4.10046768e+19
      3.86551930e+19
                      3.96194703e+19
                                     4.01591963e+19
      4.13572777e+19
                      4.18206269e+19 4.29225349e+19
                                                     4.34763987e+19
      4.37777443e+19
                      4.45337445e+19 4.49655765e+19
                                                     4.61948643e+19
                      4.74903802e+19 4.81613153e+19
      4.65683594e+19
                                                     4.85664149e+19
      4.95902084e+19
                      5.08369335e+19
                                      5.14957088e+19
                                                      5.24100864e+19
      5.36300416e+19
                      5.43992704e+19
                                     5.56434325e+19
                                                     5.61031079e+19
      5.66253756e+19
                      5.75744520e+19 5.78253400e+19
                                                     5.81711581e+19
      5.87009950e+19
                      5.88330137e+19
                                      5.99211317e+19
                                                      6.11436509e+19
      6.21840746e+19
                      6.27850796e+19 6.36190789e+19 6.43909357e+19
      6.48545829e+19
                      6.66507944e+19
                                      6.75559282e+19
                                                     6.90258886e+19
      6.97450187e+19
                      6.99694147e+19
                                     7.10174515e+19
                                                     7.23210111e+19
                                     7.51349827e+19
                                                      7.71211873e+19
      7.34244853e+19
                      7.41688396e+19
      7.85607743e+19
                      7.93338019e+19
                                     7.97817119e+19
                                                      8.01801141e+19
      8.23006362e+19
                      8.28036393e+19 8.52640489e+19 8.69930358e+19
      8.82524068e+19
                      8.88683300e+19
                                     9.01311417e+19 9.18096401e+19
      9.27589905e+19
                      9.33279231e+19 9.49128050e+19 9.57261108e+19
      9.78420558e+19
                      9.88690551e+19
                                      1.01403950e+20
                                                      1.03063596e+20
```

```
1.04124596e+20
                       1.05052168e+20 1.06299208e+20 1.07829813e+20
      1.10547598e+20
                      1.11356675e+20 1.12100301e+20 1.14938035e+20
      1.16908138e+20 1.18005302e+20 1.22794790e+20 1.23581014e+20
      1.25564038e+20 1.30358148e+20 1.31062981e+20 1.34602091e+20
      1.38083299e+20 1.38683216e+20 1.40489284e+20 1.43213235e+20
      1.47636037e+20 1.53891311e+20 1.64450621e+20 1.65739742e+20
      1.67965098e+20 1.70182927e+20 1.80362825e+20 2.63133350e+20
      3.60858037e+20 4.69452742e+20 7.83726185e+20]
    Eigenvectors (sorted):
     [ 0.04876828  0.00286912  -0.05629825  ...  0.0046746  0.01327095
      -0.083481421
     [ 0.00317129  0.00181178  0.00862197  ...  -0.02741646  -0.01983797
      -0.003379621
      [0.00605985 \quad 0.00550859 \quad -0.00155231 \quad ... \quad -0.19655827 \quad -0.02129895
      -0.00717926
      0.02613452 0.01312016 -0.00454248 ... -0.00699816 -0.02059704
      -0.010860721
      [ \ 0.00597751 \ \ 0.02107703 \ -0.01764149 \ \dots \ -0.00769263 \ -0.01119686
print("\nSubtask 11\n")
second_smallest_eigenvalue = D[1, 1]
V2 = V[:, 1]
# Ensure V2 has a positive first entry
if V2[0] < 0:
   V2 = -V2
print("\nSecond Smallest Eigenvalue:")
print(second smallest eigenvalue)
print("\nEigenvector corresponding to the second smallest eigenvalue (V2):")
print(V2)
pos = []
neg = []
for j in range(len(V2)):
    if V2[j] > 0:
        pos append(j)
    else:
        neg.append(j)
print("\nPositive Indices (V2 > 0):")
print(pos)
print("\nNegative Indices (V2 <= 0):")</pre>
print(neg)
     -8.00193565e-02 1.67132254e-02 6.58769852e-03 1.90532286e-02
     -8.42781828e-03 2.98076253e-02 1.38280567e-01 -2.55031076e-03
      1.43493838e-02 7.13194444e-03 -1.56960824e-02 -8.25623379e-03
     -5.28984414e-03 -6.02086042e-03 1.15289001e-01 -3.82802699e-02
      4.21358191e-03 -5.04844412e-02 -6.50681138e-02 1.27276714e-01
      1.58196899e-04 1.30689236e-02 -2.76892967e-03 1.68498889e-02
     -6.74437137e-03 1.43908755e-03 -2.17921627e-02 -5.33130610e-04
     -1.84834730e-02 4.76846435e-02 -1.76915842e-03 -2.00026447e-01
```

```
7.57723128e-03 -1.04989228e-02 -2.42339790e-03 2.57064319e-01
  5.03004120e-02 -7.36586479e-02 1.11355910e-01 -5.68024422e-03
  5.33395769e-03 -1.35349559e-01 1.28771500e-02 -4.06662676e-03
 -4.04452836e-02 -7.73172302e-02 9.08812695e-03 -4.54141260e-02
 -2.91207970e-02 -1.46461800e-02 1.25450476e-02 2.81965722e-03
  8.00182029e-03 1.25771127e-02 -1.68009721e-03 -3.14141204e-03
  4.73560320e-02 1.47719070e-02 -1.03177489e-02 -2.29552328e-02
  1.55357218e-02 4.84842405e-03 -1.03687765e-02 7.47511522e-03
 -9.98732324e-03 2.69046225e-02 -8.10422890e-03 1.86000940e-02
 -5.36268616e-04 -1.56820682e-02 1.81993043e-03 -6.20376626e-03
  2.50129912e-02 -7.65756138e-02 1.88069494e-02 -1.28771303e-03
 1.41782500e-01 -8.27481775e-03 -1.64385553e-01 1.37329982e-01
  1.50010173e-02 1.08937930e-01 2.96139976e-03 1.17622786e-02
 -9.82234623e-03 1.00439358e-02 -1.09461152e-01 -1.47911596e-02
 -1.23072035e-01 -1.62957990e-02 -2.63807608e-04 -7.76195096e-02
  4.94955375e-02 1.65924989e-02 -1.42940293e-03 2.02409003e-03
 1.30568233e-01 -7.64818856e-03 -5.83725684e-05 -7.45797771e-03
  3.91360824e-03 4.29502672e-03 -5.97656877e-04 1.61091455e-03
 -1.74934784e-01 -1.53610164e-02 3.55765208e-02 -9.99878545e-02
  1.65110621e-01 -1.78460951e-01 1.20287851e-01 -3.98525368e-02
 -1.86027286e-03 -9.36484794e-03 -1.71984971e-02 6.24179773e-02
 -1.48071549e-03 -1.77973811e-04 -9.73044740e-03 -5.22832816e-03
 4.14751951e-04 7.21082587e-03 4.17500617e-02 2.69802606e-02
  5.51200467e-02 1.00573239e-02 -1.80029188e-03 5.61833973e-02
 -1.07784977e-04 3.93835087e-02 1.05435434e-02 -1.20862939e-03
 -2.14263191e-03 -2.75859101e-02 -3.51048189e-03 -1.58736574e-02
  7.78455289e-03 7.54854038e-03 -3.30893174e-02 3.82565904e-02
  3.52091027e-02 -1.48831160e-02 -1.35660455e-01 1.65122159e-01
  2.12385721e-02 4.50022337e-02 -2.16444918e-03 -1.58638892e-03
 -9.30378218e-03 -3.19200283e-02 -1.06027682e-01 -2.44161991e-02
  3.01969308e-03 1.24471287e-02 1.15639655e-02 5.49638809e-04
  3.87002943e-02 -1.23755560e-02 -1.94826448e-02 7.58025778e-05
 -1.62376724e-02 -9.12963729e-03 -5.75154258e-02 2.48648417e-02
  9.55660105e-03 2.05776586e-02 3.75088223e-03 7.34015698e-03
 -1.78854472e-02 2.95332799e-02 8.13913474e-02 9.16544786e-02
 -1.13632801e-02 -2.19693798e-03 -3.40663822e-02 -1.53527031e-02
 -2.39900672e-02 4.32027638e-03 -8.61716185e-03 2.88902614e-02
 -1.25357363e-01 2.33821546e-03 -2.96881386e-03 -7.97177269e-03
  1.29967945e-01 5.22468449e-02 -4.79550478e-03 2.09966169e-02
 3.61501335e-03 -9.19280170e-04 8.46881952e-03 9.84037223e-03
 -2.05589766e-02 -1.14310728e-01 5.25258325e-02 4.20871227e-02
 -2.08598012e-02 -2.16309642e-02 4.29919330e-02 -1.59271404e-02
 4.61564103e-03 2.48893553e-02 -1.15637571e-02 -1.47082027e-04
 -4.04863924e-02 1.24979459e-02 -2.43299556e-02 -1.54106924e-02
  1.31201612e-02 2.10770301e-02 2.15915906e-031
Positive Indices (V2 > 0):
[0, 1, 2, 3, 6, 8, 9, 11, 14, 16, 20, 21, 22, 24, 26, 28, 29, 30, 31, 33, 3
Negative Indices (V2 <= 0):
[4, 5, 7, 10, 12, 13, 15, 17, 18, 19, 23, 25, 27, 32, 35, 37, 39, 42, 43, 4
```

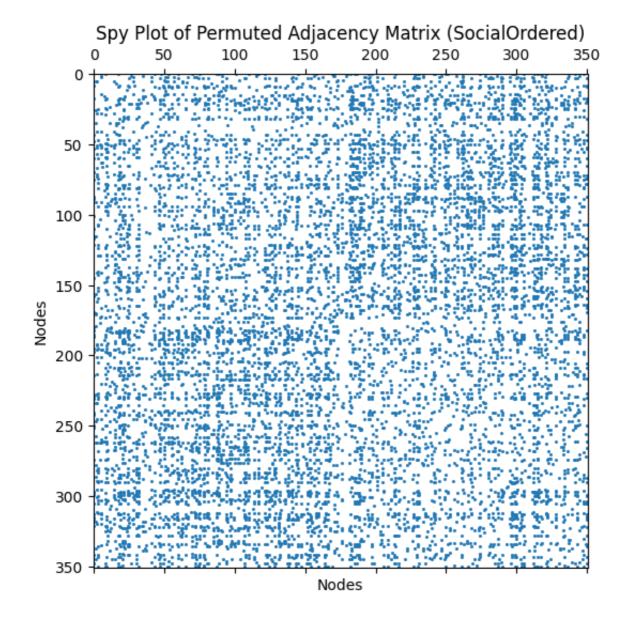
```
print("\nSubtask 12\n")
# Create the order based on positive and negative indices
order = pos + neg # Combine the positive and negative indices
m, n = Social.shape # Get the shape of the Social matrix
iden = np.eye(m) # Identity matrix of size m
# Create the permutation matrix P
P = np.zeros((m, m))
for j in range(m):
    for k in range(m):
        P[j, k] = iden[order[j], k]
# Permute the adjacency matrix
SocialOrdered = P @ Social @ P.T # Using matrix multiplication
print("Shape of SocialOrdered:", SocialOrdered.shape)

Subtask 12
Shape of SocialOrdered: (351, 351)
```

```
print("\nSubtask 13\n")
# Plot the permuted adjacency matrix
plt.figure(figsize=(8, 6))
plt.spy(SocialOrdered, markersize=1) # Using a smaller marker size for better \( \)
plt.title("Spy Plot of Permuted Adjacency Matrix (SocialOrdered)")
plt.xlabel("Nodes")
plt.ylabel("Nodes")
plt.grid(False) # Disable the grid
plt.show()
```



Subtask 13

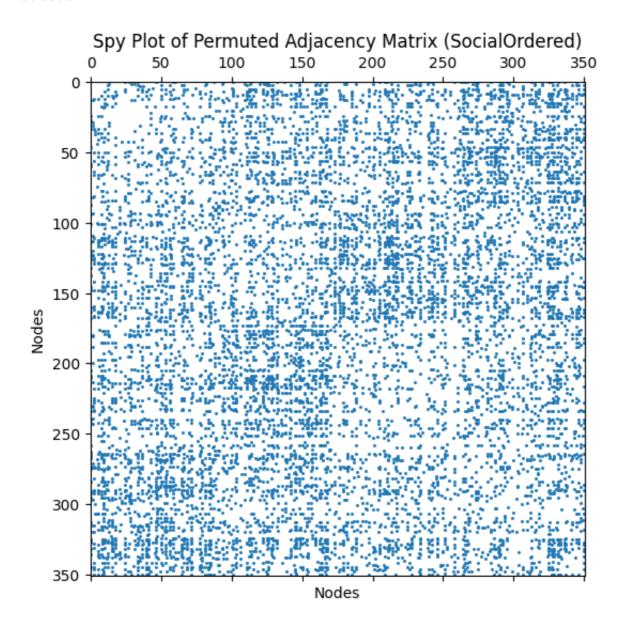


print("\nSubtask 14\n")
# Explore the third smallest eigenvalue for clustering
V3 = V[:, 2] # Get the third eigenvector
if V3[0] < 0: # Ensure V3 has a positive first entry</pre>

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V3 = -V3
# Initialize lists for the groups
pp = [] # ++ group
pn = [] # +- group
NP = [] # -+ group
nn = [] # -- group
# Grouping based on the signs of V2 and V3
for j in range(len(V2)):
    if V2[j] > 0:
        if V3[i] > 0:
            pp.append(j)
        else:
            pn.append(j)
    else:
        if V3[j] > 0:
           NP.append(j)
        else:
            nn.append(j)
# Combine the orders of the groups
order = pp + pn + NP + nn
m = len(Social) # Get the size of Social
iden = np.eye(m) # Identity matrix of size m
P = np.zeros((m, m)) # Initialize permutation matrix
# Create the permutation matrix
for j in range(m):
    P[j, :] = iden[order[j], :]
# Permute the adjacency matrix
SocialOrdered = P @ Social @ P.T
# Plot the permuted adjacency matrix
plt.figure(figsize=(8, 6))
plt.spy(SocialOrdered, markersize=1)
plt.title("Spy Plot of Permuted Adjacency Matrix (SocialOrdered)")
plt.xlabel("Nodes")
plt.ylabel("Nodes")
plt.grid(False) # Disable the grid
plt.show()
```



Subtask 14



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print("\nSubtask 15\n")
# Step 15: Fiedler vector procedure iteratively for clusters
import numpy as np
import matplotlib.pyplot as plt
# Assuming 'Social' is your adjacency matrix, and 'pos' and 'neg' are your posit
# Define SocialPos and SocialNeg based on the positive and negative indices
SocialPos = Social[np.ix_(pos, pos)]
SocialNeg = Social[np.ix_(neg, neg)]
# Calculate the Laplacian for the positive group
rowsumpos = np.sum(SocialPos, axis=1)
DiagSocialPos = np.diag(rowsumpos)
LaplaceSocialPos = DiagSocialPos - SocialPos
# Eigen decomposition for positive group
DPos , VPos = np.linalg.eig(LaplaceSocialPos)
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DPos = np.diag(DPos[ind])
VPos = VPos[:, ind]
V2Pos = VPos[:, 1] # Second smallest eigenvector for positive group
# Group positive nodes
posp = [] # Positive group
posn = [] # Negative group
for j in range(len(V2Pos)):
    if V2Pos[i] > 0:
        posp.append(pos[j]) # Append original index
    else:
        posn.append(pos[j]) # Append original index
# Calculate the Laplacian for the negative group
rowsumneg = np.sum(SocialNeg, axis=1)
DiagSocialNeg = np.diag(rowsumneg)
LaplaceSocialNeg = DiagSocialNeg - SocialNeg
# Eigen decomposition for negative group
DNeg , VNeg = np.linalg.eig(LaplaceSocialNeg)
d, ind = np.argsort(DNeg), np.argsort(DNeg)
DNeg = np.diag(DNeg[ind])
VNeg = VNeg[:, ind]
V2Neg = VNeg[:, 1] # Second smallest eigenvector for negative group
# Group negative nodes
negp = [] # Positive group
negn = [] # Negative group
for j in range(len(V2Neg)):
    if V2Neg[j] > 0:
        negp.append(neg[j]) # Append original index
    else:
        negn.append(neg[j]) # Append original index
# Generate the final order for the permutation
ordergen = posp + posn + negp + negn
# Create the permutation matrix
m = len(Social) # Assuming the size of Social
iden = np.eye(m) # Identity matrix of size m
P = np.zeros((m, m)) # Initialize permutation matrix
# Create the permutation matrix
for j in range(m):
    P[j, :] = iden[ordergen[j], :] # Filling the permutation matrix based on ord
# Permute the adjacency matrix
SocialOrderedGen = P @ Social @ P.T # Permutation of the Social matrix
# Plot the permuted adjacency matrix
plt.figure(figsize=(10, 8))
plt.spv(SocialOrderedGen, markersize=1)
plt.title("Spy Plot of Permuted Adjacency Matrix (SocialOrderedGen)")
plt.xlabel("Nodes")
plt.ylabel("Nodes")
plt.grid(False) # Disable grid for clarity
plt.show()
```



Subtask 15

