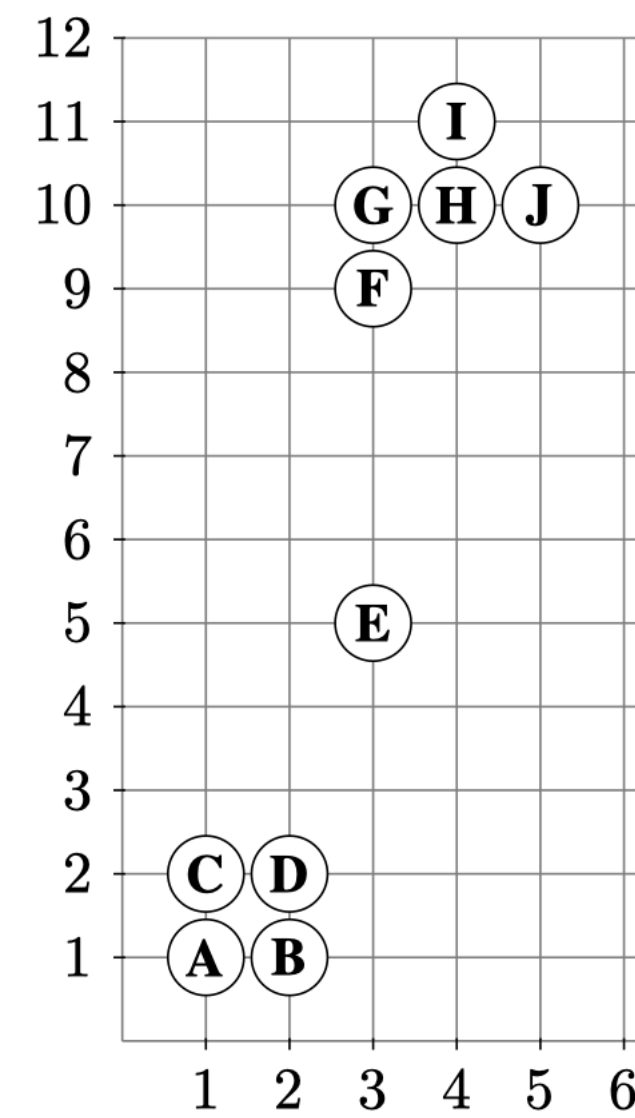
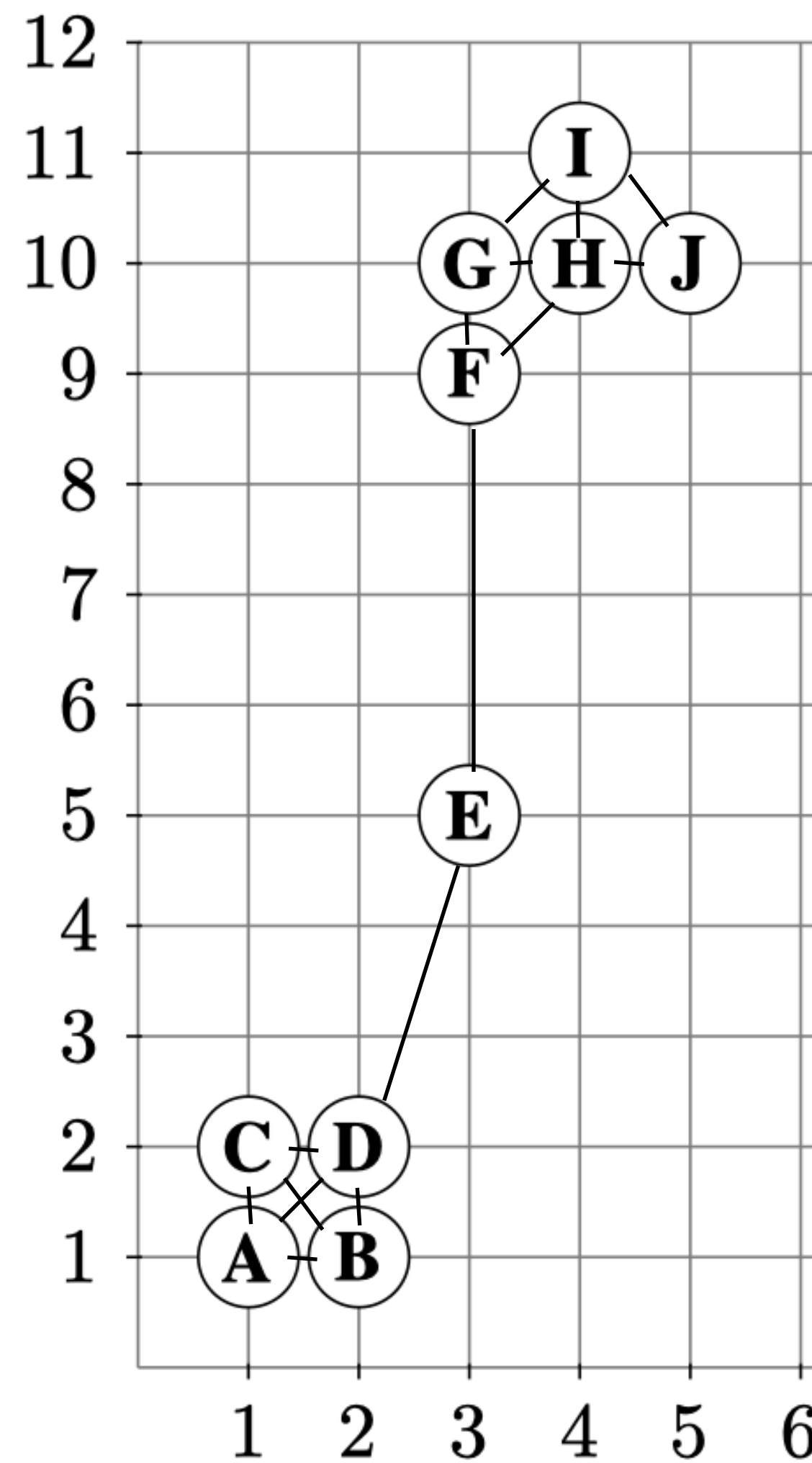


Spectral Clustering

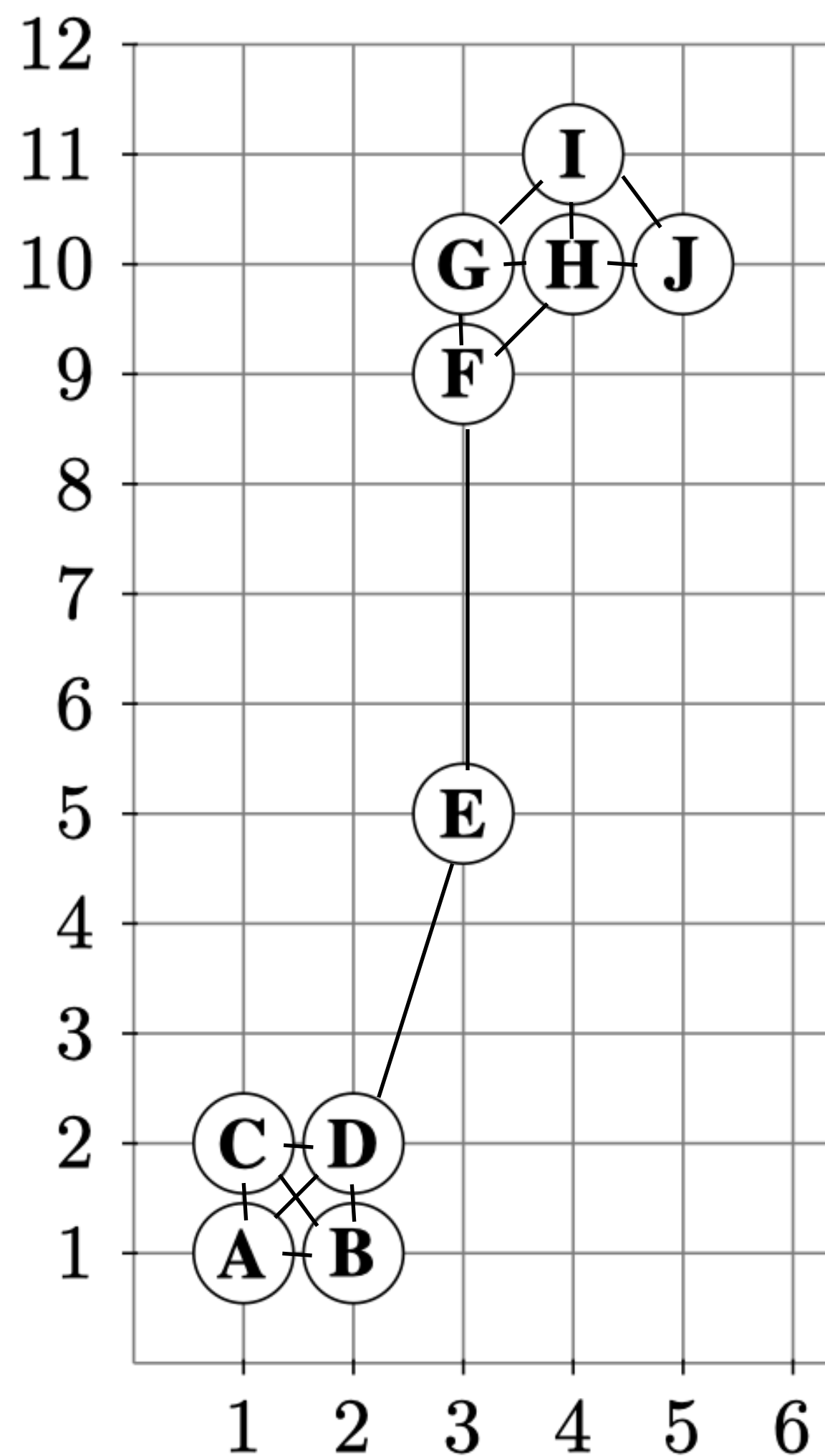
Apply spectral clustering to the first ten points of the previous dataset



Step 1: Create a similarity Graph



We can now get the Adjacency/Degree Matrix



	A	B	C	D	E	F	G	H	I	J
A	0	1	1	1	0	0	0	0	0	0
B	1	0	1	1	0	0	0	0	0	0
C	1	1	0	1	0	0	0	0	0	0
D	1	1	1	0	1	0	0	0	0	0
E	0	0	0	1	0	1	0	0	0	0
F	0	0	0	0	1	0	1	1	0	0
G	0	0	0	0	0	1	0	1	1	1
H	0	0	0	0	0	1	1	0	1	1
I	0	0	0	0	0	0	1	1	0	1
J	0	0	0	0	0	0	1	1	1	0

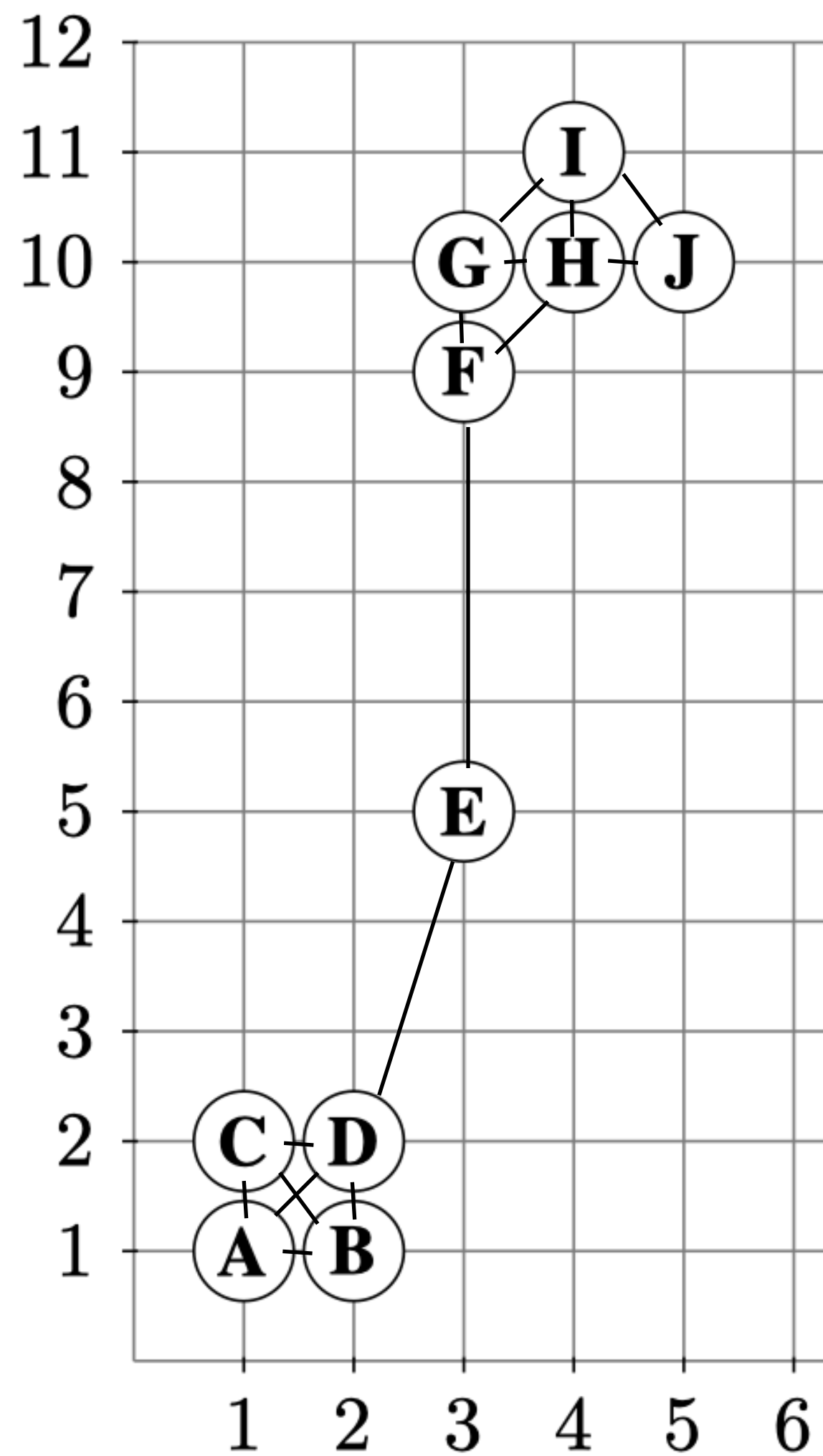
A

	A	B	C	D	E	F	G	H	I	J
A	3	0	0	0	0	0	0	0	0	0
B	0	3	0	0	0	0	0	0	0	0
C	0	0	3	0	0	0	0	0	0	0
D	0	0	0	4	0	0	0	0	0	0
E	0	0	0	0	2	0	0	0	0	0
F	0	0	0	0	0	3	0	0	0	0
G	0	0	0	0	0	0	4	0	0	0
H	0	0	0	0	0	0	0	4	0	0
I	0	0	0	0	0	0	0	0	3	0
J	0	0	0	0	0	0	0	0	0	3

D

The non-zero values of the Degree Matrix are the rows of the Adjacency Matrix added up

Now we can calculate the Laplace Matrix $L = D - A$



$$\begin{pmatrix} 3 & -1 & -1 & -1 & 0 & 0 & 0 & 0 & 0 & 0 \\ -1 & 3 & -1 & -1 & 0 & 0 & 0 & 0 & 0 & 0 \\ -1 & -1 & 3 & -1 & 0 & 0 & 0 & 0 & 0 & 0 \\ -1 & -1 & -1 & 4 & -1 & 0 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & -1 & 2 & -1 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 & -1 & 3 & -1 & -1 & 0 & 0 \\ 0 & 0 & 0 & 0 & 0 & -1 & 4 & -1 & -1 & -1 \\ 0 & 0 & 0 & 0 & 0 & -1 & -1 & 4 & -1 & -1 \\ 0 & 0 & 0 & 0 & 0 & 0 & -1 & -1 & 3 & -1 \\ 0 & 0 & 0 & 0 & 0 & 0 & -1 & -1 & -1 & 3 \end{pmatrix}$$

L

Calculate the Eigenvectors and Eigenvalues of L

The first eigenvalue will always be 0 for a connected graph

The first non-zero (smallest) eigenvalue is our spectral gap, which gives us a notion of the density of our graph

$$\begin{pmatrix} 3 & -1 & -1 & -1 & 0 & 0 & 0 & 0 & 0 & 0 \\ -1 & 3 & -1 & -1 & 0 & 0 & 0 & 0 & 0 & 0 \\ -1 & -1 & 3 & -1 & 0 & 0 & 0 & 0 & 0 & 0 \\ -1 & -1 & -1 & 4 & -1 & 0 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & -1 & 2 & -1 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 & -1 & 3 & -1 & -1 & 0 & 0 \\ 0 & 0 & 0 & 0 & 0 & -1 & 4 & -1 & -1 & -1 \\ 0 & 0 & 0 & 0 & 0 & -1 & -1 & 4 & -1 & -1 \\ 0 & 0 & 0 & 0 & 0 & 0 & -1 & -1 & 3 & -1 \\ 0 & 0 & 0 & 0 & 0 & 0 & -1 & -1 & -1 & 3 \end{pmatrix}$$

L

The Eigenvalues

- 0
- 0.17316160468804823
- 1.3998043064038725
- 3.094787587743074
- 3.9999999999999956
- 3.999999999999999
- 4.0
- 4.999999999999997
- 5.00000000000000036
- 5.332246501165005

Finding Clusters

**We can now perform K-Means clustering on our results. Reference the python file.
If you want to run it at home, make sure to install numpy and sklearn in your corresponding venv**

**For further reading: <https://towardsdatascience.com/spectral-clustering-aba2640c0d5b>
Get back to me if you run into a paywall**