CIRCLE Dataset k=20

```
clear
clc
```

Load the file

```
circle_mat = load('Circle.mat');

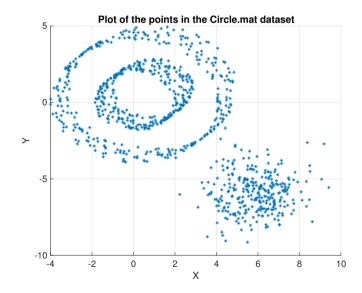
% Display the structure of the file
disp(circle_mat);
```

X: [900x2 double]

```
% Extract the matrix of points
X = circle_mat.X;
```

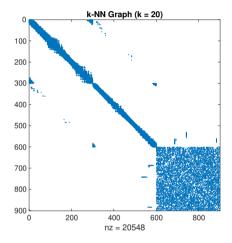
Plot the points

```
figure;
scatter(X(:,1), X(:,2), 10, Marker='*');
xlabel('X');
ylabel('Y');
title('Plot of the points in the Circle.mat dataset');
grid on;
```



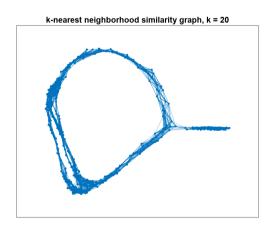
1. Similarity matrix and adjacency matrix

```
k_values = [10, 20, 40];
k = 20;
% Construct the k-nearest neighborhood similarity graph and its adjacency matrix W
W = knn_graph(X, k);
% Visualize the graph using its similarity matrix
figure;
spy(W);
title(['k-NN Graph (k = ', num2str(k), ')']);
```



```
% Store W as a sparse matrix
W = sparse(W);

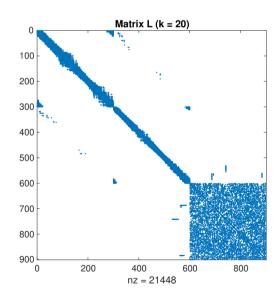
% Visualize the graph G corresponding to the adjacency matrix W
G = graph(W);
figure;
plot(G);
title(['k-nearest neighborhood similarity graph, k = ', num2str(k)]);
```



2. Construct the degree matrix D and the Laplacian matrix L

The matrix L is stored in a sparse format

```
% Plot the Laplacian matrix L
figure;
spy(L);
title(['Matrix L (k = ', num2str(k), ')']);
```



3. Compute the number of connected components of the similarity graph

```
% The points with the same number belong to the same connected component
bins = conncomp(G);
% Number of connected components
```

```
num_components = max(bins);

% Display the result
disp(['Number of connected components: ', num2str(num_components)]);
```

Number of connected components: 1

4 - 5. Compute eigenvalues and eigenvectors

```
% Set a number M of values to be computed (later it will be changed)
M = 10;
% Inizialize the eigenvalues vector and the eigenvectors matrix
eigvalues = zeros(M, 1);
eigvects = zeros(N, M);
% Choose the vector v that will be used for the inverse power method or deflation_method
v = 0.5 * ones(N, 1);
v(1:2:N) = -0.5;
% Max iterations in the power method
maxIter = 1000:
% Relative tolerance
relTol = 1e-10;
% A known fact from theory is that L is semi pos def and has at least one
% eigenvalue = 0 and that the vector of all ones is a corresponding
\% eigvalues(1) = 0;
eigvects(:, 1) = ones(N,1)/ norm(ones(N,1));
% [eigvalues(1), eigvects(:, 1)] = invpower_method(L, v(1:end), maxIter, relTol);
% Compute the reamining eigenvetors and eigenvalues
[eigvalues, eigvects, residualnorms] =
   deflation_method(L, v, eigvects, eigvalues, M, maxIter, relTol);
% Check how good the approximation is by comparing with eigs function of
% Matlab
[mat_eigvects, mat_eigs] = eigs(L, M, 'smallestabs');
norm(eigvalues - diag(mat_eigs))
```

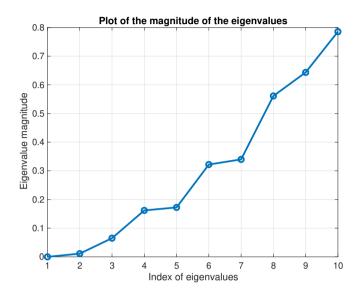
ans = 4.7619e-10

Now, find the actual number M of eigenvalues that will be used for the clustering algorithm

```
% Plot the computed eigenvalues
x = 1:M;

figure
plot(x, eigvalues, '-o', 'LineWidth', 2);
xlabel('Index of eigenvalues');
```

```
ylabel('Eigenvalue magnitude');
title('Plot of the magnitude of the eigenvalues');
grid on;
```



The suitable number of eigenvalues is 2 since eig3 is much larger than eig2 for k=20

```
M = 2;
% Define the matrix U that will be used for the spectral clustering
U = eigvects(:, 1:M);
```

6 - 7 - 8. Spectral clustering, k means

```
% Clusterize using k means and obtain the indices (and the centroids)
% inside the clusters of each point
[idx, C] = kmeans(U, M);

% Assing the original data to the corresponding clusters
A = cell(M, 1);

for i = 1:N
    % Find the cluster of y_i
    cluster_idx = idx(i);

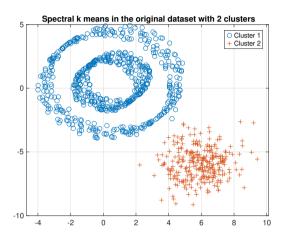
    % Assing it to x_i
        A{cluster_idx} = [A{cluster_idx}; X(i, :)];
end

% Plot of the clusterized data in the original space
X_spect_clust = zeros(N, 3);
X_spect_clust(:, 1: 2) = X;
X_spect_clust(:, 3) = idx;
```

```
figure;
markers = ['o', '+', 's'];
gscatter(X_spect_clust(:,1), X_spect_clust(:,2), X_spect_clust(:, 3), [], markers, [], 5);
title(['Spectral k means in the original dataset with ', num2str(M), ' clusters']);
legend('Cluster 1', 'Cluster 2', 'Cluster 3');
```

Warning: Ignoring extra legend entries.

grid on;



9.a K MEANS TO THE ORIGINAL DATA

```
k_value = 2; % 2 , 3

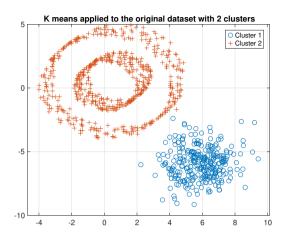
% Clusterize the original data
[idx_k, C_k] = kmeans(X, k_value);

% Add the index to X_kmeans
X_kmeans = zeros(N, 3);
X_kmeans(:, 1: 2) = X;
X_kmeans(:, 3) = idx_k;

figure;
markers = ['o', '+', 's'];
gscatter(X_kmeans(:,1), X_kmeans(:,2), X_kmeans(:, 3), [], markers, [], 5);
title(['K means applied to the original dataset with ', num2str(k_value), ' clusters']);
legend('Cluster 1', 'Cluster 2', 'Cluster 3');
```

Warning: Ignoring extra legend entries.

```
grid on;
```

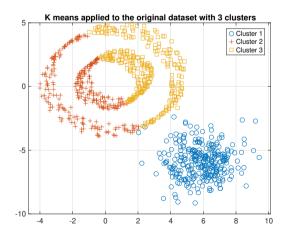


```
k_value = 3;

% Clusterize the original data
[idx_k, C_k] = kmeans(X, k_value);

% Add the index to X_kmeans
X_kmeans = zeros(N, 3);
X_kmeans(:, 1: 2) = X;
X_kmeans(:, 3) = idx_k;

figure;
markers = ['o', '+', 's'];
gscatter(X_kmeans(:,1), X_kmeans(:,2), X_kmeans(:, 3), [], markers, [], 5);
title(['K means applied to the original dataset with ', num2str(k_value), ' clusters']);
legend('Cluster 1', 'Cluster 2', 'Cluster 3');
grid on;
```



9.b DBSCAN TO THE ORIGINAL DATA

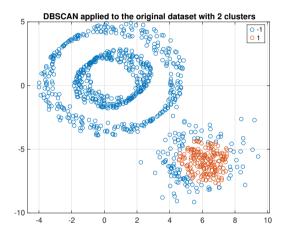
Apply the DBSCAN algorithm

```
M = 2;
% Neighborhood radius
epsilon = 0.8; % 1, if M = 3
% Minimum points for a cluster
minPts = 45;

idx_d = dbscan(X, epsilon, minPts);

X_dbscan = zeros(N, 3);
X_dbscan(:, 1: 2) = X;
X_dbscan(:, 3) = idx_d;

figure;
gscatter(X_dbscan(:,1), X_dbscan(:,2), X_dbscan(:, 3), [], 'o', 5);
title(['DBSCAN applied to the original dataset with ', num2str(M), ' clusters']);
grid on;
```



```
M = 3;
% Neighborhood radius
epsilon = 1;
% Minimum points for a cluster
minPts = 45;

idx_d = dbscan(X, epsilon, minPts);

X_dbscan = zeros(N, 3);
X_dbscan(:, 1: 2) = X;
X_dbscan(:, 3) = idx_d;

figure;
gscatter(X_dbscan(:,1), X_dbscan(:,2), X_dbscan(:, 3), [], 'o', 5);
```

title(['DBSCAN applied to the original dataset with ', num2str(M), ' clusters']);
grid on;

