220962050 Arhaan Lab11

October 18, 2024

0.1 Question

Consider the mentioned dataset and apply the hierarchical data-clustering algorithm, to identify the clusters. Write a Python function (without using the scikit-learn library) to do the following:-a. Plot a graph that displays the number of clusters on the x-axis and the Sum of Squared Errors (SSE) on the y-axis.

- b. Display the proximity matrix using Euclidean distance, Manhattan distance, and Minkowski distance.
- c. Plot the dendrogram for single, complete, average, centroid, and ward linkage methods.

```
[8]: import numpy as np
     import matplotlib.pyplot as plt
     import pandas as pd
     import scipy.cluster.hierarchy as sch
     import scipy.spatial.distance as dist
     data = np.array([
         [1, 1],
                   # p1
         [3, 2],
                  # p2
         [9, 1],
                 # p3
         [3, 7],
                 # p4
         [7, 2],
                 # p5
         [9, 7],
                  # p6
         [4, 8],
                   # p7
         [8, 3],
                   # p8
         [1, 4]
                   # p9
    ])
```

0.1.1 (B)

```
manhattan_df = pd.DataFrame(manhattan_dist, columns=[f'p{i+1}' for i in_
  →range(len(data))],
                                 index=[f'p{i+1}' for i in range(len(data))])
    minkowski_df = pd.DataFrame(minkowski_dist, columns=[f'p{i+1}' for i in_
  →range(len(data))],
                                 index=[f'p{i+1}' for i in range(len(data))])
    return euclidean_df, manhattan_df, minkowski_df
euclidean df, manhattan df, minkowski df = proximity matrices(data)
print("Euclidean Distance Matrix:\n", euclidean df)
print("\nManhattan Distance Matrix:\n", manhattan_df)
print("\nMinkowski Distance Matrix:\n", minkowski_df)
Euclidean Distance Matrix:
                     p2
                                                                       p7 \
           p1
                               рЗ
                                         p4
                                                   p5
                                                             p6
                                  6.324555 6.082763
    0.000000 2.236068 8.000000
                                                      10.000000 7.615773
p1
    2.236068 0.000000
                        6.082763
                                  5.000000
                                            4.000000
                                                       7.810250
                                                                6.082763
p2
p3
    8.000000 6.082763
                        0.000000
                                  8.485281
                                            2.236068
                                                       6.000000 8.602325
    6.324555 5.000000
                        8.485281
                                  0.000000
                                            6.403124
                                                       6.000000 1.414214
p4
    6.082763 4.000000
                        2.236068 6.403124
                                           0.000000
                                                       5.385165 6.708204
р5
p6 10.000000 7.810250 6.000000 6.000000 5.385165
                                                       0.000000 5.099020
    7.615773 6.082763
                        8.602325 1.414214 6.708204
                                                       5.099020 0.000000
р7
8q
    7.280110 5.099020
                        2.236068 6.403124 1.414214
                                                       4.123106 6.403124
p9
    3.000000 2.828427
                        8.544004 3.605551 6.324555
                                                       8.544004 5.000000
         р8
                   p9
p1 7.280110 3.000000
p2 5.099020 2.828427
p3 2.236068 8.544004
p4 6.403124 3.605551
p5 1.414214 6.324555
p6 4.123106 8.544004
p7 6.403124 5.000000
p8 0.000000 7.071068
p9 7.071068 0.000000
Manhattan Distance Matrix:
            p2
                  рЗ
      p1
                             p5
                                   p6
                                         p7
                                              p8
                                                   p9
    0.0
                          7.0
                                14.0 10.0
р1
          3.0
                8.0
                      8.0
                                            9.0
                                                  3.0
                      5.0
                                       7.0
p2
    3.0
          0.0
                7.0
                           4.0
                                11.0
                                            6.0
                                                  4.0
                     12.0
                                     12.0
рЗ
    8.0
          7.0
                0.0
                           3.0
                                 6.0
                                            3.0
                                                 11.0
p4
    8.0
          5.0
               12.0
                      0.0
                           9.0
                                 6.0
                                       2.0
                                            9.0
                                                  5.0
p5
                                            2.0
    7.0
          4.0
                3.0
                      9.0
                           0.0
                                 7.0
                                       9.0
                                                  8.0
         11.0
                6.0
                      6.0
                           7.0
                                 0.0
                                       6.0
                                           5.0
                                                 11.0
   14.0
р6
                           9.0
                                 6.0
                                           9.0
                                                  7.0
          7.0
               12.0
                      2.0
                                       0.0
р7
   10.0
```

```
9.0 0.0
               6.0
                          9.0 2.0
                                     5.0
                                                     8.0
     p8
         9.0
                     3.0
     p9
         3.0
               4.0 11.0
                          5.0 8.0 11.0
                                           7.0 8.0
                                                     0.0
     Minkowski Distance Matrix:
               p1
                        p2
                                  рЗ
                                           р4
                                                     р5
     p1 0.000000 2.080084 8.000000 6.073178 6.009245 8.995883 7.179054
     p2 2.080084 0.000000
                           6.009245 5.000000 4.000000
                                                        6.986368 6.009245
     p3 8.000000 6.009245 0.000000 7.559526 2.080084 6.000000 7.763936
     p4 6.073178 5.000000 7.559526 0.000000 5.738794 6.000000 1.259921
     p5 6.009245 4.000000 2.080084 5.738794 0.000000 5.104469 6.240251
     p6 8.995883 6.986368 6.000000 6.000000 5.104469 0.000000 5.013298
     p7 7.179054 6.009245 7.763936 1.259921 6.240251 5.013298 0.000000
     p8 7.054004 5.013298 2.080084 5.738794 1.259921 4.020726 5.738794
     p9 3.000000 2.519842 8.138223 3.271066 6.073178 8.138223 4.497941
                       p9
              p8
     p1 7.054004 3.000000
     p2 5.013298 2.519842
     p3 2.080084 8.138223
     p4 5.738794 3.271066
     p5 1.259921 6.073178
     p6 4.020726 8.138223
     p7 5.738794 4.497941
     p8 0.000000 7.006796
     p9 7.006796 0.000000
     0.1.2 (A)
[10]: def calculate_sse(data, clusters):
         sse = 0
         for cluster in clusters:
             cluster_points = data[cluster]
             cluster_center = np.mean(cluster_points, axis=0)
             sse += np.sum((cluster_points - cluster_center) ** 2)
         return sse
     def plot_sse(data):
         sse_values = []
         for k in range(1, len(data) + 1):
```

clusters = sch.fcluster(linkage_matrix, k, criterion='maxclust')

sse = calculate_sse(data, [np.where(clusters == i)[0] for i in range(1,__

linkage_matrix = sch.linkage(data, method='ward')

plt.plot(range(1, len(data) + 1), sse_values, marker='o')

 $\rightarrow k + 1)$

sse_values.append(sse)

plt.figure(figsize=(10, 6))

```
plt.title('Number of Clusters vs. Sum of Squared Errors (SSE)')
  plt.xlabel('Number of Clusters')
  plt.ylabel('SSE')
  plt.grid(True)
  plt.show()

plot_sse(data)
```

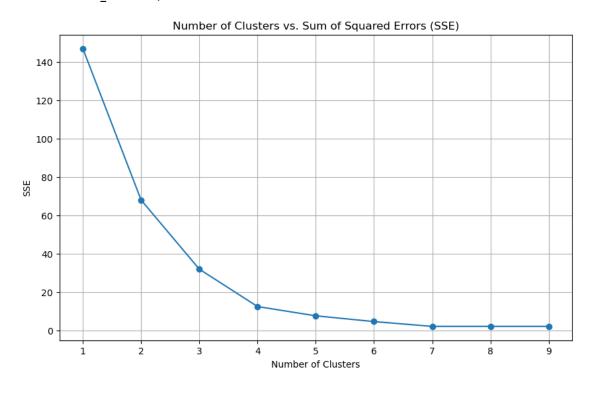
/usr/lib/python3/dist-packages/numpy/core/fromnumeric.py:3504: RuntimeWarning: Mean of empty slice.

return _methods._mean(a, axis=axis, dtype=dtype,

 $/usr/lib/python 3/dist-packages/numpy/core/_methods.py: 121: Runtime Warning: \\$

invalid value encountered in divide

ret = um.true_divide(



0.1.3 (C)

```
[11]: def plot_dendrogram(data):
    methods = ['single', 'complete', 'average', 'centroid', 'ward']

    plt.figure(figsize=(15, 10))
    for i, method in enumerate(methods):
        plt.subplot(3, 2, i + 1)
        linkage_matrix = sch.linkage(data, method=method)
```

```
sch.dendrogram(linkage_matrix)
    plt.title(f'Dendrogram ({method.capitalize()} Linkage)')
    plt.xlabel('Sample Index')
    plt.ylabel('Distance')

plt.tight_layout()
    plt.show()

plot_dendrogram(data)
```

