Exercise 6b Date: 29/01/2025

Gausian Mixture Model - Multidimensional dataset

<u>Aim:</u> To apply the Gaussian Mixture Model (GMM) clustering algorithm to a multidimensional dataset, to identify clusters within the data.

Algorithm:

The Gaussian Mixture Model (GMM) is a probabilistic clustering algorithm that assumes the data is generated from a mixture of several Gaussian distributions with unknown parameters. Unlike K-Means, which assigns points to the nearest cluster centroid, GMM assigns probabilities to each data point belonging to a particular cluster.

The probability of a data point x_i belonging to cluster k is given by:

$$P(z_i = k | x_i) = rac{\pi_k \mathcal{N}(x_i | \mu_k, \Sigma_k)}{\sum_{j=1}^K \pi_j \mathcal{N}(x_i | \mu_j, \Sigma_j)}$$

Where:

- $\mathcal{N}(x_i|\mu_k,\Sigma_k)$ is the Gaussian distribution with mean μ_k and covariance Σ_k .
- π_k is the mixing coefficient for cluster k.
- *K* is the total number of clusters.

Step 1: Import Libraries

• Import necessary Python libraries such as NumPy, Matplotlib, and Scikit-learn.

Step 2: Load the Dataset

• Load the Iris dataset and extract all four features (multidimensional dataset).

Step 3: Scale the Data

• Normalize the dataset using StandardScaler to ensure that features contribute equally to clustering.

Step 4: Dimensionality Reduction using PCA

Apply Principal Component Analysis (PCA) to reduce the dimensionality to 2D for visualization purposes.

Step 5: Apply GMM

- Initialize the Gaussian Mixture Model with n_components=3 to find three clusters.
- Fit the model to the scaled data and predict cluster labels.

Step 6: Visualize the Clusters

- Plot the clustered data points in 2D using PCA-reduced features.
- Display a 3D scatter plot using three original scaled features.
- Label the axes and provide a title for the graphs.

Import the libraries

```
import numpy as np
import matplotlib.pyplot as plt
from sklearn import datasets
import warnings
from sklearn.mixture import GaussianMixture
from sklearn.preprocessing import StandardScaler
from sklearn.decomposition import PCA
warnings.filterwarnings("ignore")
```

Load the Dataset

```
In [47]: iris = datasets.load_iris()
X = iris.data
```

Scale the data

```
In [48]: scaler = StandardScaler()
X_scaled = scaler.fit_transform(X)
```

Dimensionality Reduction using PCA

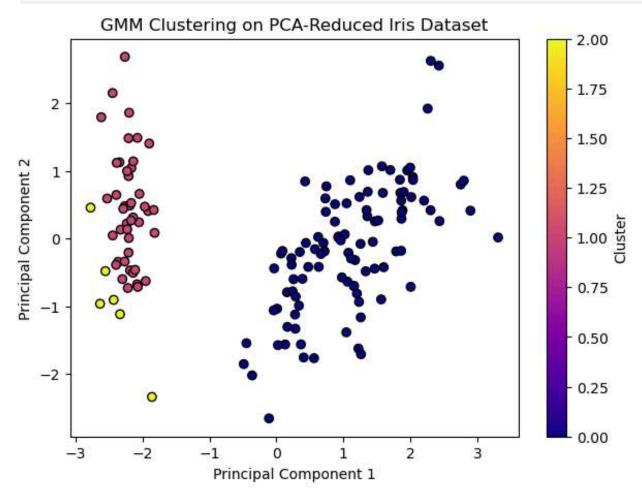
```
In [42]: pca = PCA(n_components=2)
X_pca = pca.fit_transform(X_scaled)
```

Apply Gaussian Mixture Model

```
In [49]: gmm = GaussianMixture(n_components=3, random_state=42)
gmm_labels = gmm.fit_predict(X_pca)
```

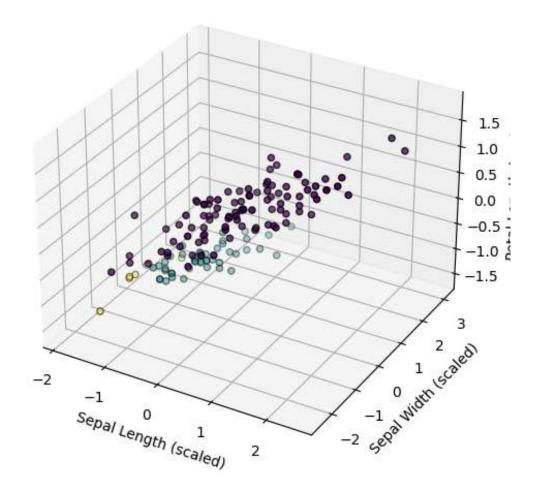
Plot the graphs

```
In [50]: plt.figure(figsize=(7,5))
    plt.scatter(X_pca[:, 0], X_pca[:, 1], c=gmm_labels, cmap='plasma', edgecolors='k')
    plt.colorbar(label='Cluster')
    plt.xlabel('Principal Component 1')
    plt.ylabel('Principal Component 2')
    plt.title('GMM Clustering on PCA-Reduced Iris Dataset')
    plt.show()
```



```
In [51]: fig = plt.figure(figsize=(8,6))
    ax = fig.add_subplot(111, projection='3d')
    ax.scatter(X_scaled[:, 0], X_scaled[:, 1], X_scaled[:, 2], c=gmm_labels, cmap='viridis', edgecolors='k')
    ax.set_xlabel('Sepal Length (scaled)')
    ax.set_ylabel('Sepal Width (scaled)')
    ax.set_zlabel('Petal Length (scaled)')
    ax.set_title('GMM Clustering on Iris Dataset (Multivariate: All Features)')
    plt.show()
```

GMM Clustering on Iris Dataset (Multivariate: All Features)



Result

The Gaussian Mixture Model successfully clustered the multidimensional dataset into three clusters based on probability distributions.

Comparison with K-Means

- Soft Clustering: Unlike K-Means, which assigns each point to a single cluster, GMM assigns probabilities, allowing for overlapping clusters.
- Elliptical Clusters: K-Means assumes clusters are spherical, whereas GMM models them as ellipses using covariance matrices.
- More Flexible Boundaries: GMM results in more flexible decision boundaries, making it more suitable for complex datasets where clusters are not well separated.
- Computation Time: GMM is generally more computationally expensive due to iterative probability updates, unlike K-Means which uses direct centroid assignments.