

## UNIT - V ADVANCED LEARNING

### 1) K-means clustering:

K-means is a partitional clustering algorithm that divides a dataset into  $k$ -cluster based on similarity.

Steps:

- 1) Choose number of clusters  $k$ .
- 2) Randomly initialize  $k$  centroids.
- 3) Assign each data point to the nearest centroid (using distance, usually Euclidean).
- 4) Recompute centroids as the mean of points in each cluster.
- 5) Repeat steps 3-4 until centroids do not change significantly.

Objective: Minimize 
$$J = \sum_{i=1}^k \sum_{n \in C_i} \|x_n - \mu_i\|^2$$

where:  $C_i$  = cluster  $i$

$\mu_i$  = centroid of cluster  $i$

Applications:

- Customer segmentation
- Image compression.

## 2) Principal Component Analysis (PCA)

Principal Component Analysis (PCA) is a dimensionality reduction technique used to transform high-dimensional data into fewer dimensions while preserving maximum variance.

Steps:

- 1) Standardizing the data.
- 2) Compute covariance matrix.
- 3) Find eigenvalues and eigenvectors.
- 4) Select the top  $k$  eigenvectors (principal components)
- 5) Transform data into new reduced feature space.

Example:

Dataset with 2 features

$x_1$	$x_2$
2	0
0	2
3	1
1	3

Step 1: Standardizing (subtract mean = 1.5)



Step 2: Covariance Matrix

$$\Sigma = \begin{bmatrix} 1.67 & -1.0 \\ -1.0 & 1.67 \end{bmatrix}$$

First principal component aligns along direction  $[1, -1]$ .

The data can be represented along one axis  $(x_1, -x_2)$  capturing most of the variance.

Advantages:

- Removes noise and redundancy.
- Reduces computation.
- Improves visualization.

Applications:

- Face and handwriting recognition
- Image compression
- Exploratory data analysis.

### 3. GAUSSIAN MIXTURE MODELS (GMM)

Introduction:

A Gaussian Mixture Model (GMM) is a probabilistic model for representing normally distributed subpopulations within an overall population. Unlike  $k$ -means, it allows clusters of different shapes.

Model definition,

$$P(x) = \sum_{i=1}^k \pi_i N(x | \mu_i, \Sigma_i)$$

\*  $\pi_i$  - mixing coefficient (weights)

\*  $\mu_i$  - mean vector

\*  $\Sigma_i$  - covariance matrix

Learning Parameters:

Expectation-Maximization (EM) Algorithm:

1. Initialize means ( $\mu$ ), covariance ( $\Sigma$ ) and weights ( $\pi$ )
2. E-step: compute probability each point belongs to each Gaussian.
3. M-step: Update  $\mu, \Sigma, \pi$  based on these probabilities.
4. Repeat: until log-likelihood converges.

Example:

Consider 1D data: [1.0, 1.2, 1.4, 5.0, 5.2, 5.4]

we assume 2 Gaussians.

\* Initial means  $\mu_1 = 1.0$   $\mu_2 = 5.0$

\* EM will adjust  $\mu_1 = 1.2$ ,  $\mu_2 = 5.2$  and compute  $\sigma, \pi$

\* Result: two overlapping normal curves fit the two clusters better than K-Means cluster.



Advantages:

- Handles overlapping clusters.
- Provides probability of membership.

Applications:

- Speaker recognition
- Object detection
- Anomaly detection.

#### 4. Q-learning Algorithm [Reinforcement Learning]

Introduction:

Reinforcement learning [RL] is learning through interaction. An agent learns to make a sequence of decisions by receiving rewards from the environment. Q-learning is a model-free off policy RL algorithm that learns the optimal action-value function.

Components:

- State : environment's situation
- Action : what the agent can do
- Reward : numerical feedback
- Policy : mapping from state to actions.
- Q-value : expected future reward for  $(s, a)$ .

## Q-learning Update Equation:

$$Q(s, a) \leftarrow Q(s, a) + \alpha [r, \gamma \max_{a'} Q(s', a') - Q(s, a)]$$

## Example:

Grid world:

Agent starts at bottom-left and must reach top-right goal.

Actions: [Up, Down, Left, Right]

Reward: +10 for goal, -1 per step.

Initially  $Q(s, a) = 0$

When the agent moves and gets reward  $r$ , the  $Q$ -table updates using the formula.

After many episodes, the agent learns optimal actions (shortest path to goal).

## Advantages:

- learns without environment model.
- works for stochastic tasks.

## Applications:

- Game AI (chess, Go)
- Robotic navigation.
- Traffic signal optimization.

## PROBLEMS BASED ON K-MEANS CLUSTERING:

cluster these points into  $K=2$  clusters:

$(2,3), (3,3), (6,6), (8,7)$ .

step 1: Initialize centroids

$$C_1 = (2,3), C_2 = (6,6)$$

step 2: Assign points

point	$d(C_1)$	$d(C_2)$	Assigned
$(2,3)$	0	5	$C_1$
$(3,3)$	1	4.24	$C_1$
$(6,6)$	4.24	0	$C_2$
$(8,7)$	6.4	2.24	$C_2$

clusters :  $C_1 = [(2,3), (3,3)]$   
 $C_2 = [(6,6), (8,7)]$

step 3: Recompute centroids:

$$C_1 = (2.5, 3.0)$$

$$C_2 = (7.0, 6.5)$$

step 4: Reassign points  $\rightarrow$  same clusters

$$C_1 = [(2,3), (3,3)]$$

$$C_2 = [(6,6), (8,7)]$$