

Introduction to Machine Learning

– Prof. Balaraman Ravindran | IIT Madras

Problem Solving Session (Week-11)

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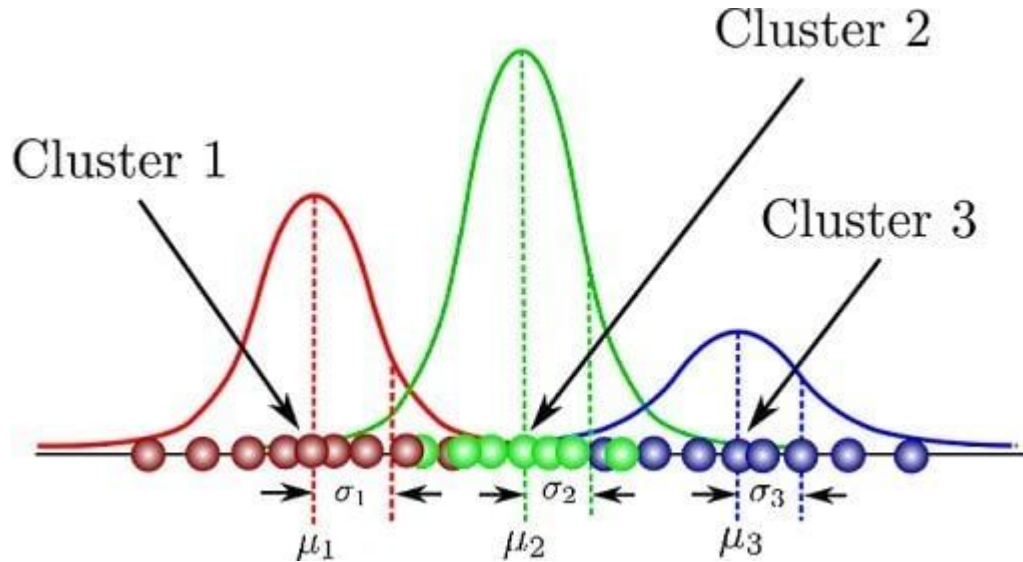
Week-11 Contents

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1. Gaussian Mixture Model
2. Expectation Maximization

Introduction to Mixture Models

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Mixture Models

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- **Definition:** A mixture model is a linear combination of probability distributions.
- **Density Formula:**

$$p(x_n) = \sum_{k=1}^K \pi_k p(x_n | \theta_k)$$

- π_k : Mixing weight ($\sum \pi_k = 1$).
- $p(x | \theta_k)$: Component distribution (e.g., Gaussian).

Gaussian Mixture Models (GMMs)

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- **Most common form: Components are Gaussian distributions.**
- **Gaussian Density:**

$$\mathcal{N}(x|\mu, \Sigma) = \frac{1}{(2\pi)^{p/2} |\Sigma|^{1/2}} \exp \left(-\frac{1}{2} (x - \mu)^T \Sigma^{-1} (x - \mu) \right)$$

- **Constraints:**

- $0 \leq \pi_k \leq 1$
- $\sum_{k=1}^K \pi_k = 1$

Gaussian Distribution Refresher

- **Univariate Gaussian:**

$$\mathcal{N}(x|\mu, \sigma^2) = \frac{1}{\sqrt{2\pi\sigma^2}} \exp\left(-\frac{(x - \mu)^2}{2\sigma^2}\right)$$

- **Multivariate Gaussian (p-dimensional):**

$$\mathcal{N}(x|\mu, \Sigma) = \frac{1}{(2\pi)^{p/2} |\Sigma|^{1/2}} \exp\left(-\frac{1}{2}(x - \mu)^T \Sigma^{-1} (x - \mu)\right)$$

Introduction to Gaussian Mixture Models

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- A Gaussian Mixture Model (GMM) is a probabilistic model that assumes data is generated from a mixture of several Gaussian distributions with unknown parameters. It is widely used for:
- Clustering
- Density estimation
- Anomaly detection
- Key Features:
- Each Gaussian component has its own mean (μ_k) and covariance (Σ_k).
- Each component has a weight (π_k) representing its contribution to the mixture.

Mathematical Formulation

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A GMM with K components is defined as:

$$p(x) = \sum_{k=1}^K \pi_k \mathcal{N}(x|\mu_k, \Sigma_k)$$

where:

- π_k = weight of the k -th Gaussian ($\sum_{k=1}^K \pi_k = 1$)
- $\mathcal{N}(x|\mu_k, \Sigma_k)$ = Gaussian PDF:

$$\mathcal{N}(x|\mu_k, \Sigma_k) = \frac{1}{(2\pi)^{d/2} |\Sigma_k|^{1/2}} \exp \left(-\frac{1}{2} (x - \mu_k)^T \Sigma_k^{-1} (x - \mu_k) \right)$$

Derivation of EM for GMM

E-Step Derivation

The responsibility $\gamma(z_{nk})$ is the posterior probability that x_n comes from the k -th Gaussian:

$$\gamma(z_{nk}) = p(z_k = 1 | x_n) = \frac{p(x_n | z_k = 1)p(z_k = 1)}{\sum_j p(x_n | z_j = 1)p(z_j = 1)}$$

M-Step Derivation

Maximize the expected log-likelihood:

$$Q(\theta, \theta^{\text{old}}) = \sum_{n=1}^N \sum_{k=1}^K \gamma(z_{nk}) [\log \pi_k + \log \mathcal{N}(x_n | \mu_k, \Sigma_k)]$$

Taking derivatives w.r.t. μ_k, Σ_k, π_k and setting to zero gives the update rules.

Expectation-Maximization (EM) Algorithm

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Since GMM parameters are unknown, we use the **EM algorithm** to estimate them iteratively.

EM Steps

1. E-step (Expectation):

Compute the posterior probability (responsibility) of each data point belonging to each Gaussian:

$$\gamma(z_{nk}) = \frac{\pi_k \mathcal{N}(x_n | \mu_k, \Sigma_k)}{\sum_{j=1}^K \pi_j \mathcal{N}(x_n | \mu_j, \Sigma_j)}$$

2. M-step (Maximization):

Update parameters (π_k, μ_k, Σ_k) using the responsibilities:

$$\begin{aligned}\mu_k &= \frac{\sum_{n=1}^N \gamma(z_{nk}) x_n}{\sum_{n=1}^N \gamma(z_{nk})} \\ \Sigma_k &= \frac{\sum_{n=1}^N \gamma(z_{nk}) (x_n - \mu_k)(x_n - \mu_k)^T}{\sum_{n=1}^N \gamma(z_{nk})} \\ \pi_k &= \frac{\sum_{n=1}^N \gamma(z_{nk})}{N}\end{aligned}$$

Introduction to EM

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- EM is an iterative optimization algorithm used for maximum likelihood estimation when data involves latent (hidden) variables.
- Commonly used in:
- Gaussian Mixture Models (GMM)
- Hidden Markov Models (HMM)
- Missing data problems

Why EM is Needed

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- Direct maximization of the log-likelihood is difficult due to hidden variables.
- For example, in GMM, the identity of the cluster (latent) for each data point is unknown.
- EM provides a framework to iteratively estimate the latent variables and optimize model parameters.

Basic Idea of EM

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- Start with initial guesses for parameters.
- E-Step (Expectation): Compute expected value of the log-likelihood, given current parameters.
- M-Step (Maximization): Maximize the expected log-likelihood w.r.t. parameters.
- Repeat until convergence.

EM Notation

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- Let: X : Observed data, Z : Latent variables, θ : Model parameters

- Log-likelihood:

$$\log p(X | \theta) = \log \sum_z p(X, Z | \theta)$$

- This is intractable \rightarrow EM maximizes a lower bound

E-Step (Expectation)

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Compute the **posterior distribution** over latent variables using current parameters:

$$Q(\theta, \theta^{(t)}) = \mathbb{E}_{Z|X, \theta^{(t)}} [\log p(X, Z | \theta)]$$

This is the **expected complete-data log-likelihood**.

M-Step (Maximization)

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Maximize the expected log-likelihood w.r.t. parameters:

$$\theta^{(t+1)} = \arg \max_{\theta} Q(\theta, \theta^{(t)})$$

This gives updated parameter estimates that improve the log-likelihood.

EM Algorithm Summary

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Repeat until convergence:

1. **E-step:** Compute posterior $p(Z \mid X, \theta^{(t)})$
2. **M-step:** Maximize $\mathbb{E}_Z[\log p(X, Z \mid \theta)]$

EM guarantees that:

$$\log p(X \mid \theta^{(t+1)}) \geq \log p(X \mid \theta^{(t)})$$

EM - Intuition

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- EM alternates between guessing the missing data (E-step) and optimizing parameters assuming the guess is correct (M-step).
- The E-step fills in the missing data.
- The M-step finds best parameters given the completed data.

Applications of EM

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- Clustering (GMM)
- Missing data imputation
- Topic modeling (LDA)
- Image segmentation
- Speech recognition

Advantages and Limitations

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- **Advantages:**

- **Handles incomplete data**
- **Has theoretical guarantees**
- **Simple to implement**

- **Limitations:**

- **Can get stuck in local optima**
- **Requires careful initialization**
- **Convergence can be slow**

Example- 1D

Problem Statement

Given the following 1D data points:

$$X = \{1, 2, 6, 7\}$$

Assume a **GMM with 2 components** ($K = 2$):

- Initial parameters:
 - Means: $\mu_1 = 1, \mu_2 = 6$
 - Variances: $\sigma_1^2 = 1, \sigma_2^2 = 1$
 - Mixing coefficients (priors): $\pi_1 = 0.5, \pi_2 = 0.5$

Task: Perform **one full iteration of the EM algorithm** to update the parameters.

Example- 1D

— — —

Step 1: E-Step (Compute Responsibilities)

The responsibility $\gamma(z_{ik})$ measures how much component k is responsible for data point x_i .

For each x_i , compute:

$$\gamma(z_{ik}) = \frac{\pi_k \cdot \mathcal{N}(x_i | \mu_k, \sigma_k^2)}{\sum_{j=1}^2 \pi_j \cdot \mathcal{N}(x_i | \mu_j, \sigma_j^2)}$$

where $\mathcal{N}(x | \mu, \sigma^2) = \frac{1}{\sqrt{2\pi\sigma^2}} \exp\left(-\frac{(x-\mu)^2}{2\sigma^2}\right)$.

Example- 1D

Compute for each data point:

1. $x_1 = 1$:

- $\mathcal{N}(1|\mu_1 = 1, \sigma_1^2 = 1) = \frac{1}{\sqrt{2\pi}}e^0 = 0.3989$
- $\mathcal{N}(1|\mu_2 = 6, \sigma_2^2 = 1) = \frac{1}{\sqrt{2\pi}}e^{-\frac{(1-6)^2}{2}} = 0.0015$
- $\gamma(z_{11}) = \frac{0.5 \times 0.3989}{0.5 \times 0.3989 + 0.5 \times 0.0015} = \frac{0.1995}{0.2009} \approx 0.993$
- $\gamma(z_{12}) = 1 - 0.993 = 0.007$

2. $x_2 = 2$:

- $\mathcal{N}(2|\mu_1 = 1, \sigma_1^2 = 1) = \frac{1}{\sqrt{2\pi}}e^{-\frac{(2-1)^2}{2}} = 0.2420$
- $\mathcal{N}(2|\mu_2 = 6, \sigma_2^2 = 1) = \frac{1}{\sqrt{2\pi}}e^{-\frac{(2-6)^2}{2}} = 0.0003$
- $\gamma(z_{21}) = \frac{0.5 \times 0.2420}{0.5 \times 0.2420 + 0.5 \times 0.0003} \approx 0.998$
- $\gamma(z_{22}) \approx 0.002$

3. $x_3 = 6$:

- $\mathcal{N}(6|\mu_1 = 1, \sigma_1^2 = 1) = 0.0003$
- $\mathcal{N}(6|\mu_2 = 6, \sigma_2^2 = 1) = 0.3989$
- $\gamma(z_{31}) \approx 0.002$
- $\gamma(z_{32}) \approx 0.998$

4. $x_4 = 7$:

- $\mathcal{N}(7|\mu_1 = 1, \sigma_1^2 = 1) = 0.0001$
- $\mathcal{N}(7|\mu_2 = 6, \sigma_2^2 = 1) = 0.2420$
- $\gamma(z_{41}) \approx 0.0004$
- $\gamma(z_{42}) \approx 0.9996$

Example- 1D

— — —

Responsibilities Summary:

x_i	$\gamma(z_{i1})$	$\gamma(z_{i2})$
1	0.993	0.007
2	0.998	0.002
6	0.002	0.998
7	0.0004	0.9996

Example- 1D

Step 2: M-Step (Update Parameters)

Now, update π_k, μ_k, σ_k^2 using the responsibilities.

1. Update Mixing Coefficients (Priors) π_k :

$$\pi_k = \frac{\sum_{i=1}^N \gamma(z_{ik})}{N}$$

- $\pi_1 = \frac{0.993+0.998+0.002+0.0004}{4} = \frac{1.9934}{4} \approx 0.498$
- $\pi_2 = \frac{0.007+0.002+0.998+0.9996}{4} = \frac{2.0066}{4} \approx 0.502$

2. Update Means μ_k :

$$\mu_k = \frac{\sum_{i=1}^N \gamma(z_{ik}) x_i}{\sum_{i=1}^N \gamma(z_{ik})}$$

- $\mu_1 = \frac{(0.993 \times 1) + (0.998 \times 2) + (0.002 \times 6) + (0.0004 \times 7)}{1.9934} \approx \frac{2.987}{1.9934} \approx 1.498$
- $\mu_2 = \frac{(0.007 \times 1) + (0.002 \times 2) + (0.998 \times 6) + (0.9996 \times 7)}{2.0066} \approx \frac{12.98}{2.0066} \approx 6.47$

3. Update Variances σ_k^2 :

$$\sigma_k^2 = \frac{\sum_{i=1}^N \gamma(z_{ik}) (x_i - \mu_k)^2}{\sum_{i=1}^N \gamma(z_{ik})}$$

- $\sigma_1^2 = \frac{0.993(1-1.498)^2 + 0.998(2-1.498)^2 + 0.002(6-1.498)^2 + 0.0004(7-1.498)^2}{1.9934} \approx 0.25$
- $\sigma_2^2 = \frac{0.007(1-6.47)^2 + 0.002(2-6.47)^2 + 0.998(6-6.47)^2 + 0.9996(7-6.47)^2}{2.0066} \approx 0.25$

Example- 1D

— — —

Final Updated Parameters After 1 EM Iteration:

Parameter	Component 1 ($k = 1$)	Component 2 ($k = 2$)
π_k	0.498	0.502
μ_k	1.498	6.47
σ_k^2	0.25	0.25

Example- 1D

— — —

Interpretation

- The means shifted slightly towards the data points they are most responsible for.
- The variances decreased because the model is becoming more confident in cluster assignments.
- The priors π_k adjusted slightly based on cluster responsibilities.

This process is repeated until convergence (when parameters change negligibly).

Example- 3D

— Problem Statement

Given the following **3D data points**:

$$X = \left\{ \begin{bmatrix} 1 \\ 0 \\ 2 \end{bmatrix}, \begin{bmatrix} 2 \\ 1 \\ 1 \end{bmatrix}, \begin{bmatrix} 5 \\ 4 \\ 6 \end{bmatrix}, \begin{bmatrix} 6 \\ 5 \\ 5 \end{bmatrix} \right\}$$

Assume:

- **Initial means:**

$$\mu_1 = \begin{bmatrix} 1 \\ 0 \\ 2 \end{bmatrix}, \quad \mu_2 = \begin{bmatrix} 5 \\ 4 \\ 6 \end{bmatrix}$$

- **Initial covariances** (isotropic):

$$\Sigma_1 = \Sigma_2 = \begin{bmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & 1 \end{bmatrix}$$

- **Initial mixing coefficients (priors):**

$$\pi_1 = 0.5, \quad \pi_2 = 0.5$$

Task: Perform **one full EM iteration** to update the parameters.

Example- 3D

— — —

Step 1: E-Step (Compute Responsibilities)

The **responsibility** $\gamma(z_{ik})$ measures how likely point x_i belongs to cluster k :

$$\gamma(z_{ik}) = \frac{\pi_k \mathcal{N}(x_i | \mu_k, \Sigma_k)}{\sum_{j=1}^K \pi_j \mathcal{N}(x_i | \mu_j, \Sigma_j)}$$

where the **multivariate Gaussian PDF** is:

$$\mathcal{N}(x | \mu, \Sigma) = \frac{1}{(2\pi)^{D/2} |\Sigma|^{1/2}} \exp \left(-\frac{1}{2} (x - \mu)^T \Sigma^{-1} (x - \mu) \right)$$

Example- 3D

Compute for each point:

1. For $x_1 = \begin{bmatrix} 1 \\ 0 \\ 2 \end{bmatrix}$:

- $\mathcal{N}(x_1|\mu_1, \Sigma_1) = 0.0635$ (high, since $x_1 = \mu_1$)
- $\mathcal{N}(x_1|\mu_2, \Sigma_2) \approx 0.0000$ (very low)
- $\gamma(z_{11}) = \frac{0.5 \times 0.0635}{0.5 \times 0.0635 + 0.5 \times 0.0000} \approx 1.0$
- $\gamma(z_{12}) \approx 0.0$

2. For $x_2 = \begin{bmatrix} 2 \\ 1 \\ 1 \end{bmatrix}$:

- $\mathcal{N}(x_2|\mu_1, \Sigma_1) \approx 0.0299$
- $\mathcal{N}(x_2|\mu_2, \Sigma_2) \approx 0.0000$
- $\gamma(z_{21}) \approx 1.0$
- $\gamma(z_{22}) \approx 0.0$

3. For $x_3 = \begin{bmatrix} 5 \\ 4 \\ 6 \end{bmatrix}$:

- $\mathcal{N}(x_3|\mu_1, \Sigma_1) \approx 0.0000$
- $\mathcal{N}(x_3|\mu_2, \Sigma_2) = 0.0635$ (since $x_3 = \mu_2$)
- $\gamma(z_{31}) \approx 0.0$
- $\gamma(z_{32}) \approx 1.0$

4. For $x_4 = \begin{bmatrix} 6 \\ 5 \\ 5 \end{bmatrix}$:

- $\mathcal{N}(x_4|\mu_1, \Sigma_1) \approx 0.0000$
- $\mathcal{N}(x_4|\mu_2, \Sigma_2) \approx 0.0299$
- $\gamma(z_{41}) \approx 0.0$
- $\gamma(z_{42}) \approx 1.0$

Example- 3D

— — —

Responsibilities Summary:

x_i	$\gamma(z_{i1})$	$\gamma(z_{i2})$
$[1, 0, 2]^T$	1.0	0.0
$[2, 1, 1]^T$	1.0	0.0
$[5, 4, 6]^T$	0.0	1.0
$[6, 5, 5]^T$	0.0	1.0

Example- 3D

Step 2: M-Step (Update Parameters)

Now, update π_k, μ_k, Σ_k using the responsibilities.

1. Update Mixing Coefficients (Priors) π_k :

$$\pi_k = \frac{\sum_{i=1}^N \gamma(z_{ik})}{N}$$

- $\pi_1 = \frac{1.0+1.0+0.0+0.0}{4} = 0.5$
- $\pi_2 = \frac{0.0+0.0+1.0+1.0}{4} = 0.5$

2. Update Means μ_k :

$$\mu_k = \frac{\sum_{i=1}^N \gamma(z_{ik}) x_i}{\sum_{i=1}^N \gamma(z_{ik})}$$

- $\mu_1 = \frac{1.0 \times [1, 0, 2]^T + 1.0 \times [2, 1, 1]^T}{2} = \frac{[3, 1, 3]^T}{2} = [1.5, 0.5, 1.5]^T$
- $\mu_2 = \frac{1.0 \times [5, 4, 6]^T + 1.0 \times [6, 5, 5]^T}{2} = \frac{[11, 9, 11]^T}{2} = [5.5, 4.5, 5.5]^T$

Example- 3D

- 3. Update Covariance Matrices Σ_k :

$$\Sigma_k = \frac{\sum_{i=1}^N \gamma(z_{ik})(x_i - \mu_k)(x_i - \mu_k)^T}{\sum_{i=1}^N \gamma(z_{ik})}$$

• For Σ_1 :

$$\Sigma_1 = \frac{1.0([1, 0, 2]^T - [1.5, 0.5, 1.5]^T)([1, 0, 2]^T - [1.5, 0.5, 1.5]^T)^T + 1.0([2, 1, 1]^T - [1.5, 0.5, 1.5]^T)([2, 1, 1]^T - [1.5, 0.5, 1.5]^T)^T}{2}$$

$$= \frac{1.0 \begin{bmatrix} 0.25 & 0.25 & -0.25 \\ 0.25 & 0.25 & -0.25 \\ -0.25 & -0.25 & 0.25 \end{bmatrix} + 1.0 \begin{bmatrix} 0.25 & -0.25 & -0.25 \\ -0.25 & 0.25 & 0.25 \\ -0.25 & 0.25 & 0.25 \end{bmatrix}}{2} = \begin{bmatrix} 0.25 & 0.0 & -0.25 \\ 0.0 & 0.25 & 0.0 \\ -0.25 & 0.0 & 0.25 \end{bmatrix}$$

• For Σ_2 :

$$\Sigma_2 = \frac{1.0([5, 4, 6]^T - [5.5, 4.5, 5.5]^T)([5, 4, 6]^T - [5.5, 4.5, 5.5]^T)^T + 1.0([6, 5, 5]^T - [5.5, 4.5, 5.5]^T)([6, 5, 5]^T - [5.5, 4.5, 5.5]^T)^T}{2}$$

$$= \begin{bmatrix} 0.25 & 0.25 & -0.25 \\ 0.25 & 0.25 & -0.25 \\ -0.25 & -0.25 & 0.25 \end{bmatrix}$$

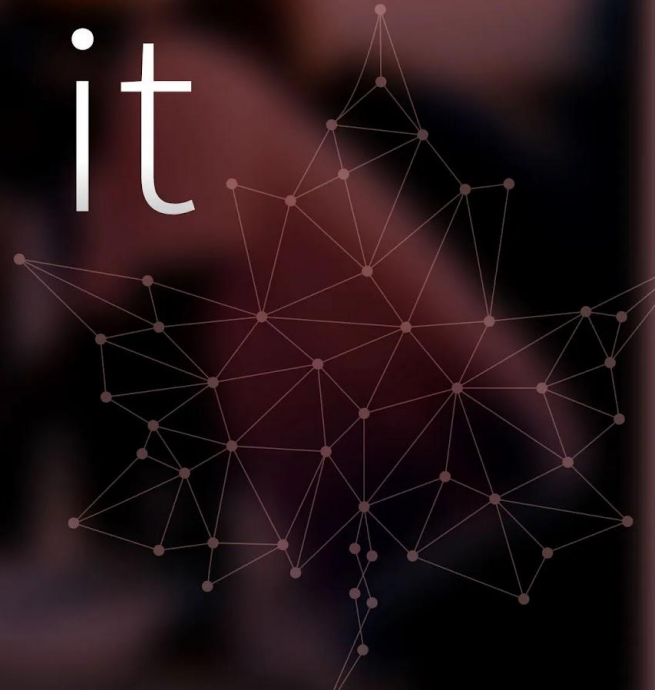
Example- 3D

Final Updated Parameters After 1 EM Iteration:

Parameter	Cluster 1 ($k = 1$)	Cluster 2 ($k = 2$)
π_k	0.5	0.5
μ_k	$[1.5, 0.5, 1.5]^T$	$[5.5, 4.5, 5.5]^T$
Σ_k	$\begin{bmatrix} 0.25 & 0.0 & -0.25 \\ 0.0 & 0.25 & 0.0 \\ -0.25 & 0.0 & 0.25 \end{bmatrix}$	$\begin{bmatrix} 0.25 & 0.25 & -0.25 \\ 0.25 & 0.25 & -0.25 \\ -0.25 & -0.25 & 0.25 \end{bmatrix}$

Assignment-11 (Cs-101- 2024) (Week-11)

Let's ^{SOLVE} = it



Question-1

— — —

01:00

What constraint must be satisfied by the mixing coefficients (π_k) in a GMM?

- a) $\pi_k > 0 \forall k$
- b) $\sum_k \pi_k = 1$
- c) $\pi_k < 1 \forall k$
- d) $\sum_k \pi_k = 0$

Question-1- Correct answer

What constraint must be satisfied by the mixing coefficients (π_k) in a GMM?

- a) $\pi_k > 0 \forall k$
- b) $\sum_k \pi_k = 1$
- c) $\pi_k < 1 \forall k$
- d) $\sum_k \pi_k = 0$

Correct options: (b) - Summation of priors probability=1

Question-2

— — —

01:00

The EM algorithm is guaranteed to decrease the value of its objective function on any iteration

- a) True
- b) False

Question-2- Correct answer

— — —

The EM algorithm is guaranteed to decrease the value of its objective function on any iteration

a) True

b) False

Correct options: (b).

Question-3

— — —

03:00

Why might the EM algorithm for GMMs converge to a local maximum rather than the global maximum of the likelihood function?

- a) The algorithm is not guaranteed to increase the likelihood at each iteration
- b) The likelihood function is non-convex
- c) The responsibilities are incorrectly calculated
- d) The number of components K is too small

Question-3 - Correct answer

— — —

Why might the EM algorithm for GMMs converge to a local maximum rather than the global maximum of the likelihood function?

- a) The algorithm is not guaranteed to increase the likelihood at each iteration
- b) The likelihood function is non-convex
- c) The responsibilities are incorrectly calculated
- d) The number of components K is too small

Correct options: (b)

Question-4

— — —

03:00

What does soft clustering mean in GMMs?

- a) There may be samples that are outside of any cluster boundary.
- b) The updates during maximum likelihood are taken in small steps, to guarantee convergence.
- c) It restricts the underlying distribution to be gaussian.
- d) Samples are assigned probabilities of belonging to a cluster

Question-4 - Correct answer

What does soft clustering mean in GMMs?

- a) There may be samples that are outside of any cluster boundary.
- b) The updates during maximum likelihood are taken in small steps, to guarantee convergence.
- c) It restricts the underlying distribution to be gaussian.
- d) Samples are assigned probabilities of belonging to a cluster

Correct options: (d)

Question-5

— — —

01:00

KNN is a special case of GMM with the following properties: (Multiple Correct)

(a) $\gamma_i = \frac{i}{(2\pi\epsilon)^{1/2}} e^{-\frac{1}{2\epsilon}}$

(b) Covariance = $\epsilon \mathbb{I}$

(c) $\mu_i = \mu_j \forall i, j$

(d) $\pi_k = \frac{1}{k}$

Question-5 - Correct answer

KNN is a special case of GMM with the following properties: (Multiple Correct)

(a) $\gamma_i = \frac{i}{(2\pi\epsilon)^{1/2}} e^{-\frac{1}{2\epsilon}}$

(b) Covariance = $\epsilon\mathbb{I}$

(c) $\mu_i = \mu_j \forall i, j$

(d) $\pi_k = \frac{1}{k}$

Correct options: (b) (d)

Question-6

— — —

We apply the Expectation Maximization algorithm to $f(D, Z, \theta)$ where D denotes the data, Z denotes the hidden variables and θ the variables we seek to optimize. Which of the following are correct?

- a) EM will always return the same solution which may not be optimal
- b) EM will always return the same solution which must be optimal
- c) The solution depends on the initialization

Question-6 - Correct answer

— — —

We apply the Expectation Maximization algorithm to $f(D, Z, \theta)$ where D denotes the data, Z denotes the hidden variables and θ the variables we seek to optimize. Which of the following are correct?

- a) EM will always return the same solution which may not be optimal
- b) EM will always return the same solution which must be optimal
- c) The solution depends on the initialization

Correct options: (c)

Question-7

— — —

03:00

True or False: Iterating between the E-step and M-step of EM algorithms always converges to a local optimum of the likelihood.

- a) True
- b) False

Question-7 - Correct answer

— — —

True or False: Iterating between the E-step and M-step of EM algorithms always converges to a local optimum of the likelihood.

a) True

b) False

Correct options: (a)

Question-8

— — —

03:00

The number of parameters needed to specify a Gaussian Mixture Model with 4 clusters, data of dimension 5, and diagonal covariances is:

- a) Lesser than 21
- b) Between 21 and 30
- c) Between 31 and 40
- d) Between 41 and 50

Question-8

— — —

03:00

The number of parameters needed to specify a Gaussian Mixture Model with 4 clusters, data of dimension 5, and diagonal covariances is:

1. Mixing Coefficients (Priors) π_k :

- There are $K = 4$ mixing coefficients, but they must sum to 1, so only $K - 1 = 3$ are free parameters.

2. Means μ_k :

- Each cluster has a mean vector of dimension $D = 5$.
- Total parameters: $K \times D = 4 \times 5 = 20$.

3. Diagonal Covariance Matrices Σ_k :

- Since the covariance is diagonal, each cluster has $D = 5$ variance parameters (one per dimension).
- Total parameters: $K \times D = 4 \times 5 = 20$.

Question-8 - Correct answer

— — —

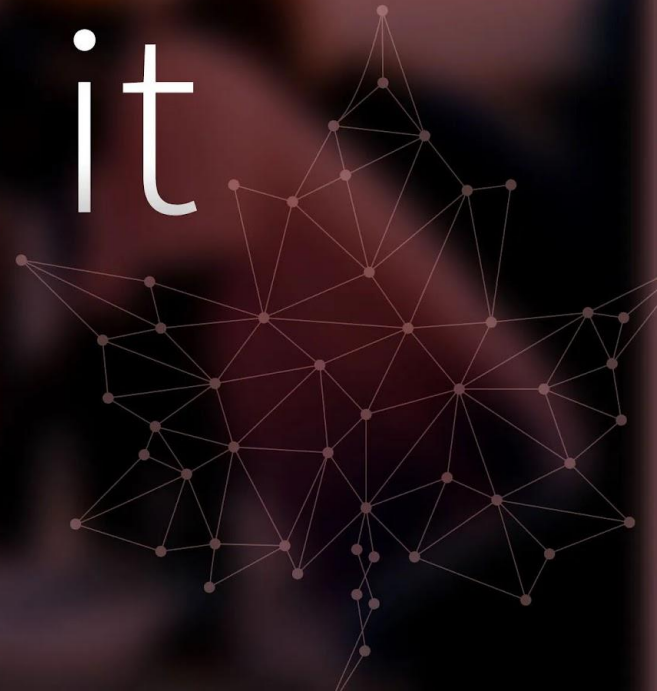
The number of parameters needed to specify a Gaussian Mixture Model with 4 clusters, data of dimension 5, and diagonal covariances is:

- a) Lesser than 21
- b) Between 21 and 30
- c) Between 31 and 40
- d) Between 41 and 50

Correct options: (d)

Assignment-11 (Cs-46- 2025) (Week-11)

Let's ^{SOLVE} = it



Question-1

— — —

01:00

Which of the following is/are estimated by the Expectation Maximization (EM) algorithm for a Gaussian Mixture Model (GMM)?

- a) K (number of components)
- b) π_k (mixing coefficient of each component)
- c) μ_k (mean vector of each component)
- d) Σ_k (covariance matrix of each component)
- e) None of the above

Question-1- Correct answer

— — —

Which of the following is/are estimated by the Expectation Maximization (EM) algorithm for a Gaussian Mixture Model (GMM)?

- a) K (number of components)
- b) π_k (mixing coefficient of each component)
- c) μ_k (mean vector of each component)
- d) Σ_k (covariance matrix of each component)
- e) None of the above

Correct options: (b) (c) (d)

Question-2

— — —

01:00

Which of the following is/are true about the responsibility terms in GMMs?
Assume the standard notation used in the lectures.

- a) $\sum_k \gamma(z_{nk}) = 1 \forall n$
- b) $\sum_n \gamma(z_{nk}) = 1 \forall k$
- c) $\gamma(z_{nk}) \in \{0,1\} \forall n,k$
- d) $\gamma(z_{nk}) \in [0,1] \forall n,k$
- e) $\pi_j > \pi_k \Rightarrow \gamma(z_{nj}) > \gamma(z_{nk}) \forall n$

Question-2- Correct answer

Which of the following is/are true about the responsibility terms in GMMs?
Assume the standard notation used in the lectures.

- a) $\sum_k \gamma(z_{nk}) = 1 \forall n$
- b) $\sum_n \gamma(z_{nk}) = 1 \forall k$
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- d) $\gamma(z_{nk}) \in [0,1] \forall n,k$
- e) $\pi_j > \pi_k \Rightarrow \gamma(z_{nj}) > \gamma(z_{nk}) \forall n$

Correct options: (a) (d)

Question-3

03:00

What is the update equation for μ_k in the EM algorithm for GMM?

- a) $\mu^{(m)}_k = \sum_{n=1}^N \gamma(z_{nk}) |v^{(m)} x_n \sum_{n=1}^N \gamma(z_{nk}) |v^{(m-1)}$
- b) $\mu^{(m)}_k = \sum_{n=1}^N \gamma(z_{nk}) |v^{(m-1)} x_n \sum_{n=1}^N \gamma(z_{nk}) |v^{(m-1)}$
- c) $\mu^{(m)}_k = \sum_{n=1}^N \gamma(z_{nk}) |v^{(m-1)} x_n N$
- d) $\mu^{(m)}_k = \sum_{n=1}^N \gamma(z_{nk}) |v^{(m)} x_n N$

Question-3 - Correct answer

What is the update equation for μ_k in the EM algorithm for GMM?

- a) $\mu^{(m)}_k = \sum_{n=1}^N \gamma(z_{nk}) |v^{(m)}_k x_n| \sum_{n=1}^N \gamma(z_{nk}) |v^{(m-1)}_k|$
- b) $\mu^{(m)}_k = \sum_{n=1}^N \gamma(z_{nk}) |v^{(m-1)}_k x_n| \sum_{n=1}^N \gamma(z_{nk}) |v^{(m-1)}_k|$
- c) $\mu^{(m)}_k = \sum_{n=1}^N \gamma(z_{nk}) |v^{(m-1)}_k x_n| N$
- d) $\mu^{(m)}_k = \sum_{n=1}^N \gamma(z_{nk}) |v^{(m)}_k x_n| N$

Correct options: (b)

Question-4

03:00

Select the correct statement(s) about the EM algorithm for GMMs.

- a) In the m th iteration, the $\gamma(z_{nk})$ values are computed using the parameter estimates computed in the same iteration.
- b) In the m th iteration, the $\gamma(z_{nk})$ values are computed using the parameter estimates computed in the $(m-1)$ th iteration.
- c) The Σ_k parameter estimates are computed during the E step.
- d) The π_k parameter estimates are computed during the M step.

Question-4 - Correct answer

— — —

Select the correct statement(s) about the EM algorithm for GMMs.

- a) In the m th iteration, the $\gamma(z_{nk})$ values are computed using the parameter estimates computed in the same iteration.
- b) In the m th iteration, the $\gamma(z_{nk})$ values are computed using the parameter estimates computed in the $(m-1)$ th iteration.
- c) The Σ_k parameter estimates are computed during the E step.
- d) The π_k parameter estimates are computed during the M step.

Correct options: (b)(d)

Question-5-7

01:00

For questions 5 to 7, use the following data consisting of 8 points (x_i, y_i)

x_i	1.0	1.8	0.9	8.2	9.1	7.8	7.6	8.8
y_i	1.5	1.2	1.6	7.3	7.2	9.5	8.0	7.5

Question-5

— — —

03:00

Fit a GMM with 2 components for this data. What are the mixing coefficients of the learned components? (Note: Use the sklearn implementation of GMM with random state = 0. Do not change the other default parameters).

- a) (0.791, 0.209)
- b) (0.538, 0.462)
- c) (0.714, 0.286)
- d) (0.625, 0.375)

Question-5 - Correct answer

— — —

Fit a GMM with 2 components for this data. What are the mixing coefficients of the learned components? (Note: Use the sklearn implementation of GMM with random state = 0. Do not change the other default parameters).

- a) (0.791, 0.209)
- b) (0.538, 0.462)
- c) (0.714, 0.286)
- d) (0.625, 0.375)

Correct options: (d)

Question-6

— — —

03:00

Using the model trained in question 5, compute the log-likelihood of the following points. Which of these points has the highest likelihood of being sampled from the model?

- a) (2.0, 0.5)
- b) (-1.0, -0.5)
- c) (7.5, 8.0)
- d) (5.0, 5.5)

Question-6 - Correct answer

— — —

Using the model trained in question 5, compute the log-likelihood of the following points. Which of these points has the highest likelihood of being sampled from the model?

- a) (2.0, 0.5)
- b) (-1.0, -0.5)
- c) (7.5, 8.0)
- d) (5.0, 5.5)

Correct options: (c)

Question-7

— — —

03:00

Let Model A be the GMM with 2 components that was trained in question 5. Using the same data from question 5, estimate a GMM with 3 components (Model B). (Note: Use the sklearn implementation of GMM with random state = 0 and all the other default parameters.)

Select the pair(s) of points that have the same label in Model A but different labels in Model B.

- a) (1.0, 1.5) and (0.9, 1.6)
- b) (1.8, 1.2) and (0.9, 1.6)
- c) (7.8, 9.5) and (8.8, 7.5)
- d) (7.8, 9.5) and (7.6, 8.0)
- e) (8.2, 7.3) and (7.6, 8.0)

Question-7 – Correct answer

— — —

Let Model A be the GMM with 2 components that was trained in question 5. Using the same data from question 5, estimate a GMM with 3 components (Model B). (Note: Use the sklearn implementation of GMM with random state = 0 and all the other default parameters.)
Select the pair(s) of points that have the same label in Model A but different labels in Model B.

- a) (1.0, 1.5) and (0.9, 1.6)
- b) (1.8, 1.2) and (0.9, 1.6)
- c) (7.8, 9.5) and (8.8, 7.5)
- d) (7.8, 9.5) and (7.6, 8.0)
- e) (8.2, 7.3) and (7.6, 8.0)

Correct options: (c) (e)

Question-8

— — —

03:00

Consider the following two statements.

Statement A: In a GMM with two or more components, the likelihood can attain arbitrarily high values.

Statement B: The likelihood increases monotonically with each iteration of EM.

- a) Both the statements are correct and Statement B is the correct explanation for Statement A.
- b) Both the statements are correct, but Statement B is not the correct explanation for Statement A.
- c) Statement A is correct and Statement B is incorrect.
- d) Statement A is incorrect and Statement B is correct.
- e) Both the statements are incorrect.

Question-8 - Correct answer

— — —

Consider the following two statements.

Statement A: In a GMM with two or more components, the likelihood can attain arbitrarily high values.

Statement B: The likelihood increases monotonically with each iteration of EM.

- a) Both the statements are correct and Statement B is the correct explanation for Statement A.
- b) Both the statements are correct, but Statement B is not the correct explanation for Statement A.
- c) Statement A is correct and Statement B is incorrect.
- d) Statement A is incorrect and Statement B is correct.
- e) Both the statements are incorrect.

Correct options: (b)

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THANK YOU

Suggestions and Feedback



Next Session:

**Tuesday:
15-Apr-2025
6:00 - 8:00 PM**