

Mathematics for Machine Learning

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Lecture 10: Density Estimation with GMMs

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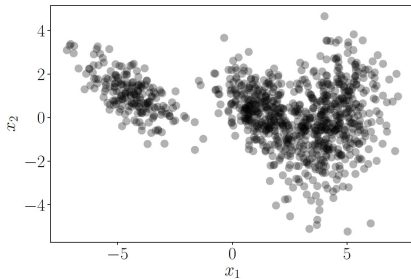
Introduction

The aim here is to represent a given dataset compactly using a probability density function from some parametric family (e.g. a Gaussian distribution).

Useful especially for large datasets.

From a density we can sample, that is, generate new data.

We can also compute the likelihood that a new point comes from the same distribution.



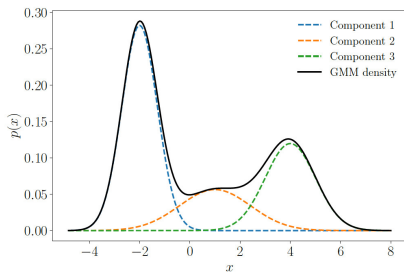
11.1 Gaussian mixture model

A **Gaussian mixture model** is a linear (convex) combination of K Gaussian distributions:

$$p(\mathbf{x}|\boldsymbol{\theta}) = \sum_{k=1}^K \pi_k \mathcal{N}(\mathbf{x}|\boldsymbol{\mu}_k, \boldsymbol{\Sigma}_k) \quad \text{with } \pi_k \in [0, 1] \quad \text{and} \quad \sum_{k=1}^K \pi_k = 1$$

where $\boldsymbol{\theta} = \{\mathbf{x}_k, \boldsymbol{\Sigma}_k, \pi_k : k = 1, \dots, K\}$ contains all the parameters of the model.

This gives us significantly more flexibility than a single unimodal Gaussian distribution.



Example

$$\begin{aligned} p(x|\boldsymbol{\theta}) &= 0.5\mathcal{N}(x|-2, \frac{1}{2}) \\ &\quad + 0.2\mathcal{N}(x|1, 2) \\ &\quad + 0.3\mathcal{N}(x|4, 1) \end{aligned}$$

11.2 Parameter estimation via maximum likelihood

Given a dataset $\mathcal{X} = \{\mathbf{x}_1, \dots, \mathbf{x}_N\}$ of points i.i.d. sampled from some distribution $p(\mathbf{x})$, our task is to **represent the unknown $p(\mathbf{x})$ by a GMM** with K mixture components.

The idea is to find the maximum likelihood estimate θ_{ML} of the GMM parameters.

The data is i.i.d., so
$$p(\mathcal{X}|\theta) = \prod_{n=1}^N p(\mathbf{x}_n|\theta) = \prod_{n=1}^N \left(\sum_{k=1}^K \pi_k \mathcal{N}(\mathbf{x}_n|\boldsymbol{\mu}_k, \boldsymbol{\Sigma}_k) \right)$$

The log-likelihood is
$$\mathcal{L}(\theta) = \sum_{n=1}^N \log p(\mathbf{x}_n|\theta) = \sum_{n=1}^N \log \sum_{k=1}^K \pi_k \mathcal{N}(\mathbf{x}_n|\boldsymbol{\mu}_k, \boldsymbol{\Sigma}_k)$$

We'd like to find the gradient of \mathcal{L} w.r.t. the model parameters θ , set it to $\mathbf{0}$, and solve for θ . Unfortunately, here we cannot obtain a closed-form solution. Instead we will use an **iterative scheme**, where the idea is to update one parameter at a time while keeping the others fixed.

The **responsibility** of the k -th mixture component for the n -th data point is defined as

$$r_{n,k} = \frac{\pi_k \mathcal{N}(\mathbf{x}_n | \boldsymbol{\mu}_k, \boldsymbol{\Sigma}_k)}{\sum_{j=1}^K \pi_j \mathcal{N}(\mathbf{x}_n | \boldsymbol{\mu}_j, \boldsymbol{\Sigma}_j)}$$

$(r_{n,1}, \dots, r_{n,K})$ is a probability vector; a “soft assignment” of \mathbf{x}_n to the K components.

Updating the GMM **means**: $\boldsymbol{\mu}_k^{\text{new}} = \frac{1}{N_k} \sum_{n=1}^N r_{n,k} \mathbf{x}_n$ with $N_k = \sum_{n=1}^N r_{n,k}$

Updating the GMM **covariances**: $\boldsymbol{\Sigma}_k^{\text{new}} = \frac{1}{N_k} \sum_{n=1}^N r_{n,k} (\mathbf{x}_n - \boldsymbol{\mu}_k)(\mathbf{x}_n - \boldsymbol{\mu}_k)^\top$

Updating the GMM **mixture weights**: $\pi_k = \frac{N_k}{N}$

11.3 EM algorithm

Unfortunately, the updates on the previous slide are not a closed-form solution for the parameters of the GMM, because the responsibilities $r_{n,k}$ depend on those parameters.

But they do suggest a simple iterative scheme, called **expectation maximisation**.

Choose initial values for μ_k , Σ_k , π_k , and alternate until convergence between:

E-step: evaluate the responsibilities $r_{n,k}$

M-step: use the updated responsibilities to re-estimate μ_k , Σ_k , π_k