K-Means and Gaussian Mixture Models

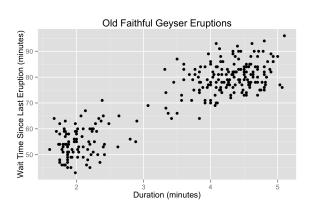
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K-Means Clustering

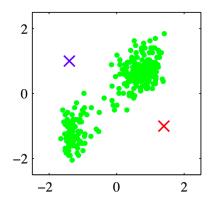
Example: Old Faithful Geyser



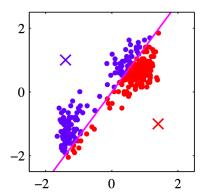
- Looks like two clusters.
- How to find these clusters algorithmically?

k-Means: By Example

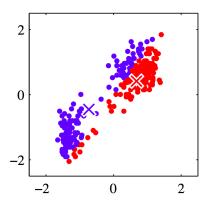
- Standardize the data.
- Choose two cluster centers.



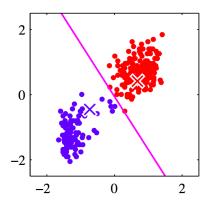
Assign each point to closest center.



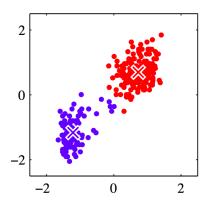
• Compute new class centers.



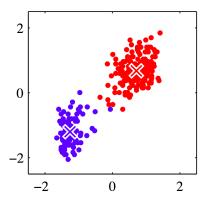
Assign points to closest center.



Compute cluster centers.

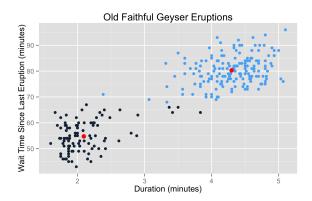


• Iterate until convergence.



k-Means Algorithm: Standardizing the data

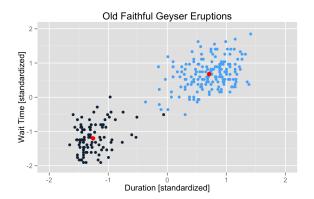
Without standardizing:



- Blue and black show results of k-means clustering
- Wait time dominates the distance metric

k-Means Algorithm: Standardizing the data

With standardizing:



• Note several points have been reassigned from black to blue cluster.

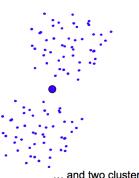
k-Means: Failure Cases

k-Means: Suboptimal Local Minimum

• The clustering for k = 3 below is a local minimum, but suboptimal:



Would be better to have one cluster here



and two clusters here

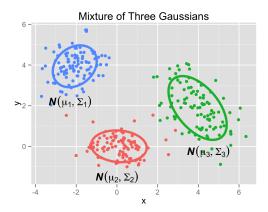
Gaussian Mixture Models

Probabilistic Model for Clustering

- Let's consider a **generative model** for the data.
- Suppose
 - There are k clusters.
 - We have a probability density for each cluster.
- Generate a point as follows
 - **1** Choose a random cluster $z \in \{1, 2, ..., k\}$.
 - 2 Choose a point from the distribution for cluster Z.

Gaussian Mixture Model (k = 3)

- **1** Choose $z \in \{1, 2, 3\}$ with $p(1) = p(2) = p(3) = \frac{1}{3}$.
- 2 Choose $x \mid z \sim \mathcal{N}(X \mid \mu_z, \Sigma_z)$.

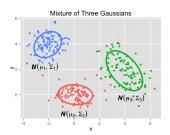


Gaussian Mixture Model Parameters (k Components)

Cluster probabilities: $\pi = (\pi_1, ..., \pi_k)$

Cluster means : $\mu = (\mu_1, ..., \mu_k)$

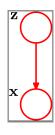
Cluster covariance matrices: $\Sigma = (\Sigma_1, \dots \Sigma_k)$



For now, suppose all these parameters are known.

We'll discuss how to learn or estimate them later.

Gaussian Mixture Model: Joint Distribution



Factorize the joint distribution:

$$p(x,z) = p(z)p(x \mid z)$$

= $\pi_z \mathcal{N}(x \mid \mu_z, \Sigma_z)$

- π_z is probability of choosing cluster z.
- $x \mid z$ has distribution $\mathcal{N}(\mu_z, \Sigma_z)$.
- z corresponding to x is the true cluster assignment.
- Suppose we know the model parameters π_z, μ_z, Σ_z .
- Then we can easily compute the joint p(x,z).

Latent Variable Model

- We observe x.
- We don't observe z. (Cluster assignment).
- ullet Cluster assignment z is called a **hidden variable** or **latent variable**.

Definition

A **latent variable model** is a probability model for which certain variables are never observed.

e.g. The Gaussian mixture model is a latent variable model.

The GMM "Inference" Problem

- We observe x. We want to know z.
- The conditional distribution of the cluster z given x is

$$p(z \mid x) = p(x, z)/p(x)$$

- The conditional distribution is a soft assignment to clusters.
- A hard assignment is

$$z^* = \underset{z \in \{1, \dots, k\}}{\operatorname{arg\,min}} p(z \mid x).$$

• So if we have the model, clustering is trival.

Mixture Models

Gaussian Mixture Model: Marginal Distribution

• The marginal distribution for a single observation x is

$$p(x) = \sum_{z=1}^{k} p(x, z)$$
$$= \sum_{z=1}^{k} \pi_{z} \mathcal{N}(x \mid \mu_{z}, \Sigma_{z})$$

- Note that p(x) is a convex combination of probability densities.
- This is a common form for a probability model...

Mixture Distributions (or Mixture Models)

Definition

A probability density p(x) represents a mixture distribution or mixture model, if we can write it as a convex combination of probability densities. That is,

$$p(x) = \sum_{i=1}^{k} w_i p_i(x),$$

where $w_i \ge 0$, $\sum_{i=1}^k w_i = 1$, and each p_i is a probability density.

- In our Gaussian mixture model, x has a mixture distribution.
- \bullet More constructively, let S be a set of probability distributions:
 - $oldsymbol{0}$ Choose a distribution randomly from S.
 - Sample x from the chosen distribution.
- Then x has a mixture distribution.

Learning in Gaussian Mixture Models

The GMM "Learning" Problem

- Given data x_1, \ldots, x_n drawn from a GMM,
- Estimate the parameters:

Cluster probabilities:
$$\pi = (\pi_1, ..., \pi_k)$$

Cluster means:
$$\mu = (\mu_1, ..., \mu_k)$$

Cluster covariance matrices:
$$\Sigma = (\Sigma_1, \dots \Sigma_k)$$

- Once we have the parameters, we're done.
- Just do "inference" to get cluster assignments.

Estimating/Learning the Gaussian Mixture Model

- One approach to learning is maximum likelihood
 - find parameter values that give **observed data** the **highest likelihood**.
- The model likelihood for $\mathcal{D} = \{x_1, \dots, x_n\}$ is

$$L(\pi, \mu, \Sigma) = \prod_{i=1}^{n} p(x_i)$$
$$= \prod_{i=1}^{n} \sum_{z=1}^{k} \pi_z \mathcal{N}(x_i \mid \mu_z, \Sigma_z).$$

As usual, we'll take our objective function to be the log of this:

$$J(\pi, \mu, \Sigma) = \sum_{i=1}^{n} \log \left\{ \sum_{z=1}^{k} \pi_{z} \mathcal{N}(x_{i} \mid \mu_{z}, \Sigma_{z}) \right\}$$

Properties of the GMM Log-Likelihood

GMM log-likelihood:

$$J(\pi, \mu, \Sigma) = \sum_{i=1}^{n} \log \left\{ \sum_{z=1}^{k} \pi_{z} \mathcal{N}(x_{i} \mid \mu_{z}, \Sigma_{z}) \right\}$$

Let's compare to the log-likelihood for a single Gaussian:

$$\sum_{i=1}^{n} \log \mathcal{N}(x_i \mid \mu, \Sigma)$$

$$= -\frac{nd}{2} \log (2\pi) - \frac{n}{2} \log |\Sigma| - \frac{1}{2} \sum_{i=1}^{n} (x_i - \mu)' \Sigma^{-1}(x_i - \mu)$$

- For a single Gaussian, the log cancels the exp in the Gaussian density.
 - ⇒ Things simplify a lot.
- For the GMM, the sum inside the log prevents this cancellation.
 - Expression more complicated. No closed form expression for MLE.

Issues with MLE for GMM

Identifiability Issues for GMM

Suppose we have found parameters

Cluster probabilities : $\pi = (\pi_1, ..., \pi_k)$

Cluster means: $\mu = (\mu_1, \dots, \mu_k)$

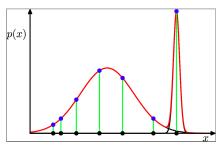
Cluster covariance matrices: $\Sigma = (\Sigma_1, \dots \Sigma_k)$

that are at a local minimum.

- What happens if we shuffle the clusters? e.g. Switch the labels for clusters 1 and 2.
- We'll get the same likelihood. How many such equivalent settings are there?
- Assuming all clusters are distinct, there are k! equivalent solutions.
- Not a problem per se, but something to be aware of.

Singularities for GMM

Consider the following GMM for 7 data points:



- Let σ^2 be the variance of the skinny component.
- What happens to the likelihood as $\sigma^2 \rightarrow 0$?
- In practice, we end up in local minima that do not have this problem.
 - Or keep restarting optimization until we do.
- Bayesian approach or regularization will also solve the problem.

From Bishop's Pattern recognition and machine learning, Figure 9.7.

Gradient Descent / SGD for GMM

What about running gradient descent or SGD on

$$J(\pi, \mu, \Sigma) = -\sum_{i=1}^{n} \log \left\{ \sum_{z=1}^{k} \pi_{z} \mathcal{N}(x_{i} \mid \mu_{z}, \Sigma_{z}) \right\}?$$

- Can be done but need to be clever about it.
- Each matrix $\Sigma_1, \ldots, \Sigma_k$ has to be positive semidefinite.
- How to maintain that constraint?
 - Rewrite $\Sigma_i = M_i M_i^T$, where M_i is an unconstrained matrix.
 - Then Σ_i is positive semidefinite.
- But we actually prefer positive definite, to avoid singularities.

Cholesky Decomposition for SPD Matrices

Theorem

Every symmetric positive definite matrix $A \in \mathbb{R}^{d \times d}$ has a unique **Cholesky** decomposition:

$$A = LL^T$$
,

where L a lower triangular matrix with positive diagonal elements.

- A lower triangular matrix has half the number of parameters.
- Symmetric positive definite is better because avoids singularities.
- Requires a non-negativity constraint on diagonal elements.
 - e.g. Use projected SGD method like we did for the Lasso.

The EM Algorithm for GMM

MLE for Gaussian Model

- Let's start by considering the MLE for the Gaussian model.
- For data $\mathcal{D} = \{x_1, \dots, x_n\}$, the log likelihood is given by

$$\sum_{i=1}^{n} \log \mathcal{N}(x_i \mid \mu, \Sigma) = -\frac{nd}{2} \log (2\pi) - \frac{n}{2} \log |\Sigma| - \frac{1}{2} \sum_{i=1}^{n} (x_i - \mu)' \Sigma^{-1}(x_i - \mu).$$

With some calculus, we find that the MLE parameters are

$$\mu_{\text{MLE}} = \frac{1}{n} \sum_{i=1}^{n} x_{i}$$

$$\Sigma_{\text{MLE}} = \frac{1}{n} \sum_{i=1}^{n} (x_{i} - \mu_{\text{MLE}}) (x_{i} - \mu_{\text{MLE}})^{T}$$

- For GMM, If we knew the cluster assignment z_i for each x_i ,
 - we could compute the MLEs for each cluster.

Cluster Responsibilities: Some New Notation

• Denote the probability that observed value x_i comes from cluster j by

$$\gamma_i^j = \mathbb{P}(Z = j \mid X = x_i).$$

- The **responsibility** that cluster j takes for observation x_i .
- Computationally,

$$\gamma_i^j = \mathbb{P}(Z = j \mid X = x_i).
= p(Z = j, X = x_i)/p(x)
= \frac{\pi_j \mathcal{N}(x_i \mid \mu_j, \Sigma_j)}{\sum_{c=1}^k \pi_c \mathcal{N}(x_i \mid \mu_c, \Sigma_c)}$$

- The vector $(\gamma_i^1, \dots, \gamma_i^k)$ is exactly the **soft assignment** for x_i .
- Let $n_c = \sum_{i=1}^n \gamma_i^c$ be the number of points "soft assigned" to cluster c.

EM Algorithm for GMM: Overview

- **1** Initialize parameters μ , Σ , π .
- "E step". Evaluate the responsibilities using current parameters:

$$\gamma_i^j = \frac{\pi_j \mathcal{N}(x_i \mid \mu_j, \Sigma_j)}{\sum_{c=1}^k \pi_c \mathcal{N}(x_i \mid \mu_c, \Sigma_c)},$$

for i = 1, ..., n and j = 1, ..., k.

3 "M step". Re-estimate the parameters using responsibilities:

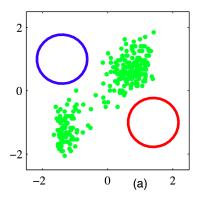
$$\mu_c^{\text{new}} = \frac{1}{n_c} \sum_{i=1}^n \gamma_i^c x_i$$

$$\Sigma_c^{\text{new}} = \frac{1}{n_c} \sum_{i=1}^n \gamma_i^c (x_i - \mu_{\text{MLE}}) (x_i - \mu_{\text{MLE}})^T$$

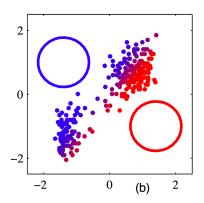
$$\pi_c^{\text{new}} = \frac{n_c}{n_c},$$

Repeat from Step 2, until log-likelihood converges.

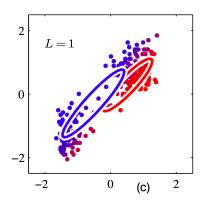
Initialization



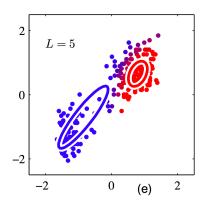
• First soft assignment:



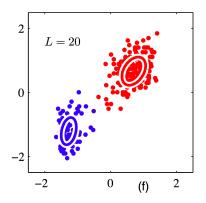
• First soft assignment:



After 5 rounds of EM:



After 20 rounds of EM:



Relation to K-Means

- EM for GMM seems a little like k-means.
- In fact, there is a precise correspondence.
- First, fix each cluster covariance matrix to be $\sigma^2 I$.
- As we take $\sigma^2 \to 0$, the update equations converge to doing *k*-means.
- If you do a quick experiment yourself, you'll find
 - Soft assignments converge to hard assignments.
 - Has to do with the tail behavior (exponential decay) of Gaussian.