Gaussian Mixture Model

~~lets took at GMM~~

lets consider the data to be in 1-dimension

Suppose 'N' number of points in 1-d is given and we are asked to divide it into 'K' clusters.

So we estimate μ (mean) and σ^2 (variance) for each k

Using E-M Algorithm (Expectation-Maximization) we estimate the model parameter ~~iteratively~~ iteratively.

E - estimation step
M - Maximization step

Implementation of em algorithm for GMM in 1D is given below:

initialization:

Initial means (μ_k) → ~~using~~ choose random point ~~given~~ from given for each k

Initial variances (σ_k^2) → variance of respective points.

Initial mixing coefficient (π_k) → $1/k$

mixing coefficient defines how much points among given points belong to a respective cluster

$$\sum_{k}^{K} \pi_k = 1$$

pdf for univariate GD:-

$$p(x \mid \mu, \sigma) = \frac{1}{\sqrt{2\pi\sigma^2}} \exp\left[\frac{-(x-\mu)^2}{2\sigma^2}\right]$$

~~Variance~~

log :-

$$-\frac{1}{2} \times \log(2\pi) - \log(\sigma) - \left(\frac{(x-\mu)^2}{2\sigma^2}\right)$$

E-step

compute γ_{nk} (i.e probability that data point x_n belongs to K)

$$\gamma_{nk} = \frac{\pi_k \, N(x_n \mid \mu_k, \sigma_k^2)}{\sum_{j=1}^{K} \pi_j \, N(x_n, \mu_j, \sigma_j^2)}$$

Using log-sum-exp trick to handle small probability (when x_n is far from μ_k)

$$\log(\gamma_{nk}) = \log(\pi_k) + \log(N(x_n \mid \mu_k, \sigma_k^2))$$

$$-\log\left(\sum_{j=1}^{K} \pi_j \, N(x_n \mid \mu_j, \sigma_j^2)\right)$$

Also

$$\log\left(\sum_{j=1}^{K} \pi_j \, N(x_n \mid \mu_j, \sigma_j^2)\right)$$

$$= \log\left(\sum_{j=1}^{K} \exp\left(\log(\pi_j) + \log(N(\dots))\right)\right)$$

Apply it in find equation of $\log(\gamma_{nk})$

M-step

update parameter based on γ_{nk}

$$\pi_k = \frac{\sum \gamma_{nk}}{N}$$

these are parameters for each k cluster

$$\mu_k = \frac{\sum_{n=1}^{N} \gamma_{nk} x_n}{\sum_{n=1}^{N} \gamma_{nk}}$$

$$\sigma_k^2 = \frac{\sum_{n=1}^{N} \gamma_{nk} (x_n - \mu_k)^2}{\sum_{n=1}^{N} \gamma_{nk}}$$

Using the log-sum trick:

$$\pi_k = \frac{\sum_{n=1}^{N} \exp(\log(\gamma_{nk}))}{N}$$

$$\mu_k = \frac{\sum_{n=1}^{N} (\exp(\log(r_{nk})) \cdot x_n)}{\sum_{n=1}^{N} \exp(\log(r_{nk}))}$$

$$\sigma_k^2 = \frac{\sum_{n=1}^{N} (\exp(\log(r_{nk})) \cdot (x_n - \mu_k)^2)}{\sum_{n=1}^{N} \exp(\log(r_{nk}))}$$

Compare the values of parameters with previous iteration, if the difference is below a threshold, we can say it converging (so we can stop)

In my code, I didn't consider changing π_k as the initial mean is calculated randomly and not by K-means.

Other converging criteria is the use of log likelihood

$$\Rightarrow \sum_{n=1}^{N} \log \left(\sum_{k=1}^{K} \pi_k \mathcal{N}(x_n | \mu_k, \sigma_k^2) \right)$$

using log-sum-exp trick

If its change in adjacent iteration are below a threshold we can stop.

finally, for a point x_1

if $r_{1k_1} > r_{1k_2}$, then we can say

that x_1 belongs to k_1 (probability is high)

μ_k and σ_k gives us the final params of k gaussian distributions

Note: In my code all the parameters must converge and also log likelihood must be stable to end the iterations.