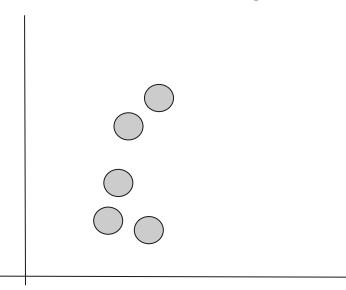
CIS 635 Knowledge Discovery & Data Mining

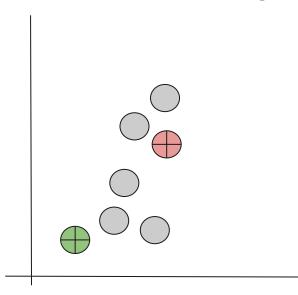
- k-means: Centroid Based
- Hierarchical clustering: Distance connectivity based
- GMM: Distribution based
- **DBSCAN**: Density Based

- k-means: Centroid Based
- **Hierarchical clustering**: Distance connectivity based
- GMM: Distribution based
- **DBSCAN**: Density Based

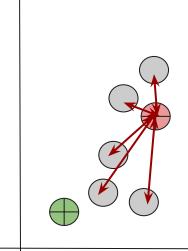
- Centroid based
- Only works for numeric data only
- Explicit k-centroid inputs (initialization)
- Iterative algorithm



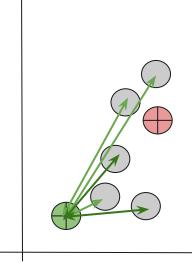
- 5 given data points of forms (x, y,): numeric



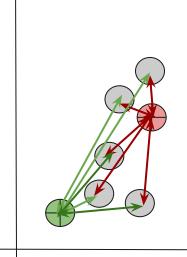
- 5 given data points of forms (x_i, y_i) : numeric
- For k = 2, the algorithm starts with two random (users can also provide based on their analysis/intuition) initials means $(m_x, m_y)_j$ where j = 1, ... k



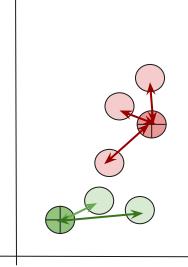
- 5 given data points of forms (x, y,): numeric
- For k = 2, the algorithm starts with two random (users can also provide based on their analysis/intuition) initials means $(m_x, m_y)_j$ where j = 1, ... k
- Compares distance between the means $(m_x, m_y)_i$ and all given (5) data points and for all j



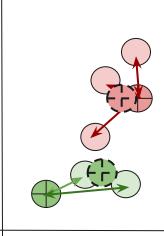
- 5 given data points of forms (x, y,): numeric
- For k = 2, the algorithm starts with two random (users can also provide based on their analysis/intuition) initials means $(m_x, m_y)_j$ where j= 1, .. k
- Compares distance between the means (m_x, m_y) and all given (5) data points and for all j



- 5 given data points of forms (x_i, y_i) : numeric
- For k = 2, the algorithm starts with two random (users can also provide based on their analysis/intuition) initials means $(m_x, m_y)_j$ where j = 1, ... k
- Compares distance between the means $(m_x, m_y)_i$ and all given (5) data points and for all j

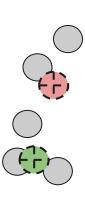


- 5 given data points of forms (x, y,): numeric
- For k = 2, the algorithm starts with two random (users can also provide based on their analysis/intuition) initials means $(m_x, m_y)_j$ where j= 1, .. k
- Compares distance between the means $(m_x, m_y)_j$ and all given (5) data points and for all j
- Assign cluster labels based on the distances measured.

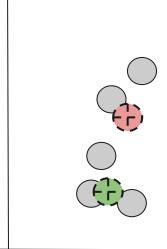


- 5 given data points of forms (x_i, y_j) : numeric For k = 2, the algorithm starts with two random (users can also provide based on their analysis/intuition) initials means $(m_x, m_y)_i$ where j =
- Compares distance between the means $(m_x, m_y)_i$ and all given (5) data points and for all j
- Assign cluster labels based on the distances measured.
- Update the two means:

 - $m_x = 1/|x_i|\sum x_i$ $m_y = 1/|y_i|\sum y_i$



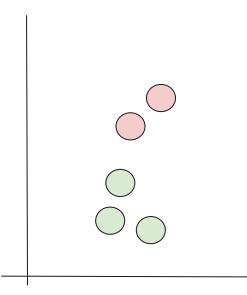
- 5 given data points of forms (x_i, y_j) : numeric For k = 2, the algorithm starts with two random (users can also provide based on their analysis/intuition) initials means $(m_x, m_y)_i$ where j =
- Compares distance between the means $(m_x, m_y)_i$ and all given (5) data points and for all j
- Assign cluster labels based on the distances measured.
- Update the two means:
 - $m_x = 1/|x_i|\sum x_i$ $m_y = 1/|y_i|\sum y_i$



- 5 given data points of forms (x_i, y_i) : numeric For k = 2, the algorithm starts with two random (users can also provide based on their analysis/intuition) initials means $(m_x, m_y)_i$ where j =1, .. k
- Compares distance between the means $(m_x, m_y)_i$ and all given (5) data points and for all *j*
- Assign cluster labels based on the distances measured.
- Update the two means:

 - $m_x = 1/|x_i|\sum x_i$ $m_y = 1/|y_i|\sum y_i$

- Iterate until the means $(m_x, m_y)_i$ don't move; or a predefined number of iterations are run



- 5 given data points of forms (x_i, y_i) : numeric For k = 2, the algorithm starts with two random (users can also provide based on their analysis/intuition) initials means $(m_x, m_y)_i$ where j =1, .. k
- Compares distance between the means $(m_x, m_y)_i$ and all given (5) data points and for all j
- Assign cluster labels based on the distances measured.
- Update the two means:
 - $m_x = 1/|x_i|\sum x_i$ $m_y = 1/|y_i|\sum y_i$
- Iterate until the means $(m_x, m_y)_i$ don't move; or a predefined number of iterations are run
- Likely the final cluster configuration

Given an initial set of k means $m_1^{(1)}$, ..., $m_k^{(1)}$ (see below), the algorithm proceeds by alternating between two steps:^[7]

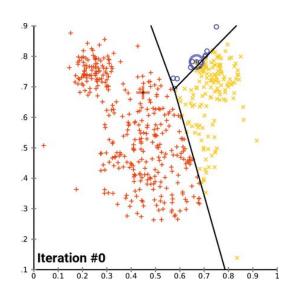
1. **Assignment step**: Assign each observation to the cluster with the nearest mean: that with the least squared Euclidean distance.^[8] (Mathematically, this means partitioning the observations according to the Voronoi diagram generated by the means.)

$$S_i^{(t)} = \left\{ x_p : \left\| x_p - m_i^{(t)}
ight\|^2 \leq \left\| x_p - m_j^{(t)}
ight\|^2 \, orall j, 1 \leq j \leq k
ight\},$$

where each x_p is assigned to exactly one $S^{(t)}$, even if it could be assigned to two or more of them.

2. Update step: Recalculate means (centroids) for observations assigned to each cluster.

$$m_i^{(t+1)} = rac{1}{\left|S_i^{(t)}
ight|} \sum_{x_j \in S_i^{(t)}} x_j$$

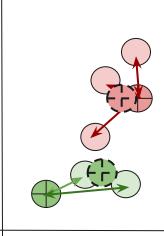


k-modes is the categorical equivalent!

- Centroid based
- Only works for categorical data
- Explicit k-centroid inputs (initialization)
- Iterative algorithm
- Only

- 1. Pick K observations at random and use them as leaders/clusters
- 2. Calculate the dissimilarities and assign each observation to its closest cluster
- 3. Define new modes for the clusters
- 4. Repeat 2-3 steps until there are is no re-assignment required

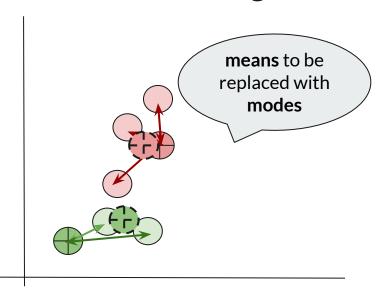
- One simple metric: number of categorical value match
- Whiteboarding



- 5 given data points of forms (x_i, y_j) : numeric For k = 2, the algorithm starts with two random (users can also provide based on their analysis/intuition) initials means $(m_x, m_y)_i$ where j =
- Compares distance between the means $(m_x, m_y)_i$ and all given (5) data points and for all j
- Assign cluster labels based on the distances measured.
- Update the two means:

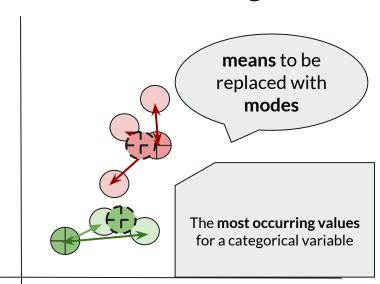
 - $m_x = 1/|x_i|\sum x_i$ $m_y = 1/|y_i|\sum y_i$

k-modes Clustering



- 5 given data points of forms (x_i, y_i) : **categorical** For k = 2, the algorithm starts with two random (users can also provide based on their analysis/intuition) initials modes $(m_x, m_y)_i$ where j =
- Compares distance between the modes $(m_x, m_y)_i$ and all given (5) data points and for all i
- Assign cluster labels based on the distances measured.
- Update the two modes:
 - m_x = 1/|x_i| mode(x_i) m_y = 1/|y_i| mode(x_i)

k-modes Clustering



- 5 given data points of forms (x_i, y_i) : categorical For k = 2, the algorithm starts with two random (users can also provide based on their analysis/intuition) initials modes $(m_v, m_v)_i$ where j =
- Compares distance between the modes (m_v, m_v) , and all given (5) data points and for all *i*
- Assign cluster labels based on the distances measured.
- Update the two modes:
 - m_x = 1/|x_i| mode(x_i) m_y = 1/|y_i| mode(x_i)

How about when you have mixed data types?

- Either you have to convert data in one type
 - What could be the challenges?

How about when you have mixed data types?

- Either you have to convert data in one type
 - What could be the challenges?
- Or, you have to have an algorithm that updates centroids

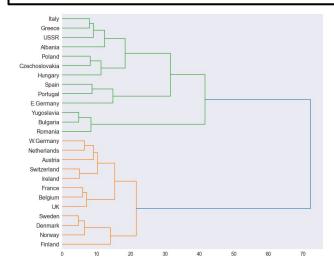
- K-means: Centroid Based
- Hierarchical clustering: Distance connectivity based
- GMM: Distribution based
- DBSCAN: Density Based

Hierarchical clustering

Agglomerative: This is a "bottom up" approach. Each observation starts as a new cluster, and pairs of clusters are merged as one moves up the hierarchy.

Divisive: This is a "top down" approach. All data starts as on cluster, and recursively splits into two/multiple clusters.

Grouping countries according to their protein consumption. (dendrogram graph below)



- K-means: Centroid Based
- **Hierarchical clustering**: Distance connectivity based
- GMM: Distribution based
- DBSCAN: Density Based

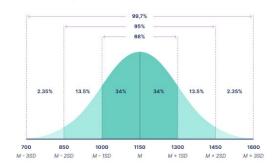
Basic idea from the model title itself:

Gaussian/Normal distribution: distribution modeling some continuous variables such as: population height/weight in a certain region, yearly sales of a business etc.

$$X \sim \mathcal{N}(\mu, \sigma^2)$$

$$X \sim \mathcal{N}(\mu, \sigma^2)$$
$$p(x) = \frac{1}{\sigma\sqrt{2\pi}}e^{-\frac{1}{2}(\frac{x-\mu}{\sigma})^2}$$

Using the empirical rule in a normal distribution



Basic idea from the model title itself:

Gaussian/Normal distribution: distribution modeling some continuous variables such as: population height/weight in a certain region, yearly sales of a business etc.

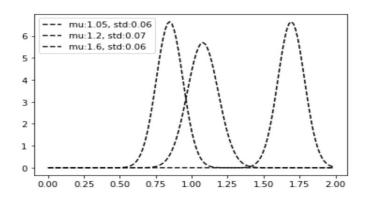
An example problem

- Display the weight distribution of grade 5,6 and 10 students
- Choose an x (confusing between g 5 and 6) and explain through words

$$X \sim \mathcal{N}(\mu, \sigma^2)$$

$$p(x) = \frac{1}{\sigma\sqrt{2\pi}}e^{-\frac{1}{2}(\frac{x-\mu}{\sigma})^2}$$

$$p(x) = \frac{1}{\sigma\sqrt{2\pi}}e^{-\frac{1}{2}(\frac{x-\mu}{\sigma})^2}$$



Basic idea from the model title itself:

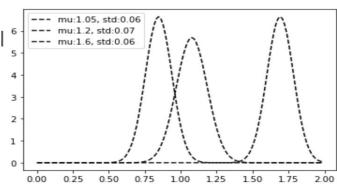
- Gaussian/Normal distribution: distribution modeling some continuous variables such as: population height/weight in a certain region, yearl sales of a business etc.
- *Mixture*: more than one object/component

$$p(x) = \sum_{k=1}^{K} \pi_k \mathcal{N}(x|\mu_k, \sigma_k)$$
 $p(\mathbf{x}) = \sum_{k=1}^{K} \pi_k \mathcal{N}(\mathbf{x}|\boldsymbol{\mu}_k, \boldsymbol{\Sigma}_k).$

$$X \sim \mathcal{N}(\mu, \sigma^2)$$

$$X \sim \mathcal{N}(\mu, \, \sigma^2)$$

$$p(x) = \frac{1}{\sigma \sqrt{2\pi}} e^{-\frac{1}{2}(\frac{x-\mu}{\sigma})^2}$$



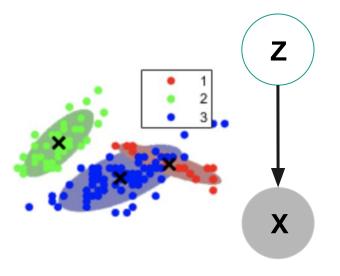
- Fig at the right shows the graphical model (B Net) of GMM
- X: feature vector; in our example case a vector with apples (size, color)
- Z: encoding of clusters (1-of-K is 1, rest are 0s), K is the number of clusters.
- Essentially GMM models(learns) the joint distribution (an example of what we call a generative model)

$$p(\mathbf{x}, \mathbf{z}) = p(\mathbf{z})p(\mathbf{x}|\mathbf{z})$$
$$p(\mathbf{x}) = \sum_{\mathbf{z}} p(\mathbf{z})p(\mathbf{x}|\mathbf{z}) = \sum_{k=1}^{K} \pi_k \mathcal{N}(\mathbf{x}|\boldsymbol{\mu}_k, \boldsymbol{\Sigma}_k)$$

$$0 \leqslant \pi_k \leqslant 1$$

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

$$p(\mathbf{z}) = \prod_{k=1}^K \pi_k^{z_k}.$$



Model parameters(all k s)

$$oldsymbol{\pi}, oldsymbol{\mu}, oldsymbol{\Sigma}$$

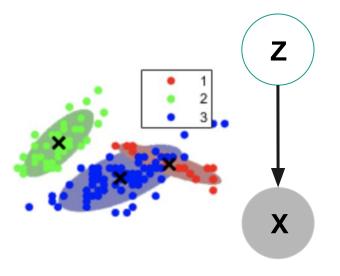
- Fig at the right shows the graphical model (B Net) of GMM
- X: feature vector; in our example case a vector with apples (size, color)
- Z: encoding of clusters (1-of-K is 1, rest are 0s), K is the number of clusters.
- Essentially GMM models(learns) the joint distribution (an example of what we call a generative model)

$$p(\mathbf{x}, \mathbf{z}) = p(\mathbf{z})p(\mathbf{x}|\mathbf{z})$$
$$p(\mathbf{x}) = \sum_{\mathbf{z}} p(\mathbf{z})p(\mathbf{x}|\mathbf{z}) = \sum_{k=1}^{K} \pi_k \mathcal{N}(\mathbf{x}|\boldsymbol{\mu}_k, \boldsymbol{\Sigma}_k)$$

$$0 \leqslant \pi_k \leqslant 1$$

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

$$p(\mathbf{z}) = \prod_{k=1}^K \pi_k^{z_k}.$$



Model parameters(all k s)

$$oldsymbol{\pi}, oldsymbol{\mu}, oldsymbol{\Sigma}$$

 If given model parameters, using Bayes rule, we can estimate the class conditional probabilities for a given query X

$$p(z_k = 1|\mathbf{x}) = \frac{p(z_k = 1)p(\mathbf{x}|z_k = 1)}{\sum_{j=1}^{K} p(z_j = 1)p(\mathbf{x}|z_j = 1)}$$
$$= \frac{\pi_k \mathcal{N}(\mathbf{x}|\boldsymbol{\mu}_k, \boldsymbol{\Sigma}_k)}{\sum_{j=1}^{K} \pi_j \mathcal{N}(\mathbf{x}|\boldsymbol{\mu}_j, \boldsymbol{\Sigma}_j)}.$$



Maximum Likelihood Learning

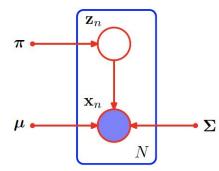
- Start with random parameter initializations
- Optimize the following likelihood function for N data points

$$\ln p(\mathbf{X}|\boldsymbol{\pi}, \boldsymbol{\mu}, \boldsymbol{\Sigma}) = \sum_{n=1}^{N} \ln \left\{ \sum_{k=1}^{K} \pi_k \mathcal{N}(\mathbf{x}_n | \boldsymbol{\mu}_k, \boldsymbol{\Sigma}_k) \right\}$$

- Some popular techniques:
 - Expectation Maximization algorithm
 - We can also use gradient based optimization techniques

$$\log p(X) = \log(p(Z)p(X|Z)) =$$

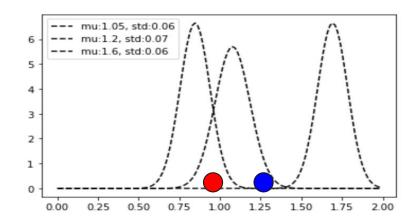
$$= \log p(Z) + \log p(X|Z)$$



Why do we need GMM?

 Hard cluster assignments (k-means, hierarchical clustering)

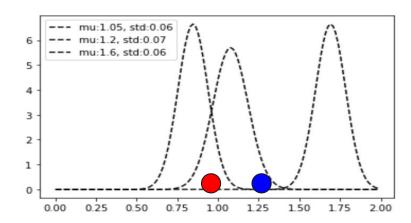
Gmm 3 components



Why do we need GMM?

- Hard cluster assignments (k-means, hierarchical clustering)
- Soft assignments: cluster assignment probability scores, p(Z|X)

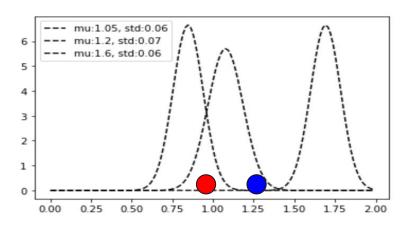
Gmm 3 components



Why do we need GMM?

- Hard cluster assignments (k-means, hierarchical clustering)
- Soft assignments: cluster assignment probability scores, p(Z|X)
 - GMM offers us a probability distribution over the (features & clusters) space, and this can be used as a part of a larger/complex modeling tasks, P(X, Z)

Gmm 3 components





Notebook presentation

- K-means: Centroid Based
- **Hierarchical clustering**: Distance connectivity based
- GMM: Distribution based
- DBSCAN: Density Based

DBSCAN

- Density-Based Spatial Clustering of Applications with Noise
- Two parameters:
 - **minPts**: The minimum number of points (a threshold) clustered together to be considered dense.
 - **eps** (ϵ): A distance measure that will be used to locate the points in the neighborhood of any point.
- The algorithm proceeds by arbitrarily picking up a point in the dataset.
- If there are at least 'minPoint' points within a radius of 'ε' to the point then we consider all these points to be part of the same cluster.
- The clusters are then expanded by recursively repeating the neighborhood calculation for each neighboring point

Noise point

Core point

Border point

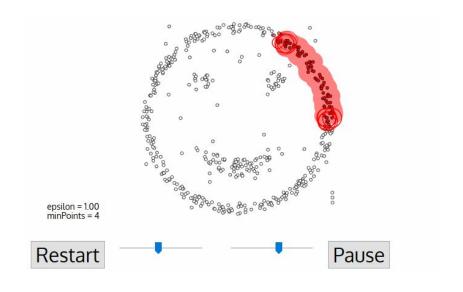
MinPts = 3

"The idea that a cluster in data space is a **contiguous region** of **high point density**, separated from other such clusters by contiguous regions of **low point density**."

Model details with visual demonstration

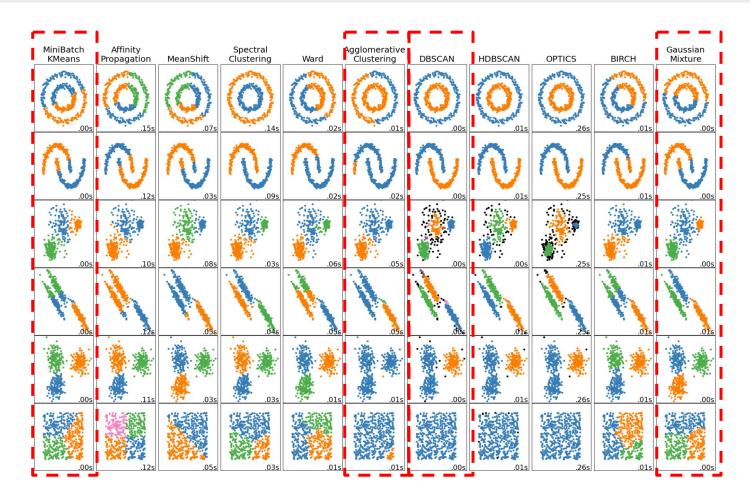
DBSCAN

- Two parameters:
 - *minPts*: The minimum number of points (a threshold) clustered together to be considered dense.
 - *eps* (ε): A distance measure that will be used to locate the points in the neighborhood of any point.
- The algorithm proceeds by arbitrarily picking up a point in the dataset.
- If there are at least 'minPoint' points within a radius of 'ε' to the point then we consider all these points to be part of the same cluster.
- The clusters are then expanded by recursively repeating the neighborhood calculation for each neighboring point



Reference

sklearn



sklearn packages

Clustering: sklearn

Clustering: sklearn (API reference)