Lecture 23. Gaussian Mixture Models.

COMP90051 Statistical Machine Learning

Lecturer: Ben Rubinstein



This lecture

- Unsupervised learning
 - Diversity of problems
 - * k-means refresher
- Gaussian mixture model (GMM)
 - A probabilistic approach to clustering
 - * The GMM model
 - GMM clustering as an optimisation problem
- Starting Expectation-Maximisation (EM) algorithm

Unsupervised Learning

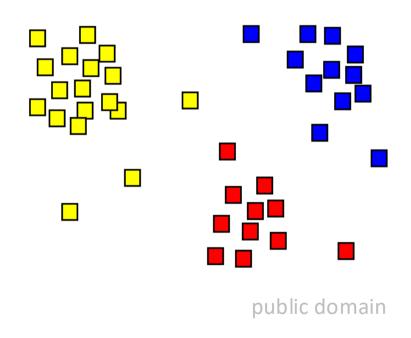
A large branch of ML that concerns with learning the structure of the data in the absence of labels

Main learning paradigms so far

- Supervised learning: Overarching aim is making predictions from data
- We studied methods in the context of this aim: e.g. linear/logistic regression, DNN, SVM
- We had instances $x_i \in \mathbb{R}^m$, i = 1, ..., n and corresponding labels y_i for model fitting, aiming to predict labels for new instances
- Can be viewed as a function approximation problem, but with a big caveat: ability to generalise is critical
- Bandits: a setting of partial supervision where subroutine in contextual bandits requires supervised learning

Now: Unsupervised learning

- In unsupervised learning, there is no dedicated variable called a "label"
- Instead, we just have a set of points $x_i \in \mathbb{R}^m$, i = 1, ..., n
- Aim of unsupervised learning is to explore the structure (patterns, regularities) of data



The aim of "exploring the structure" is vague

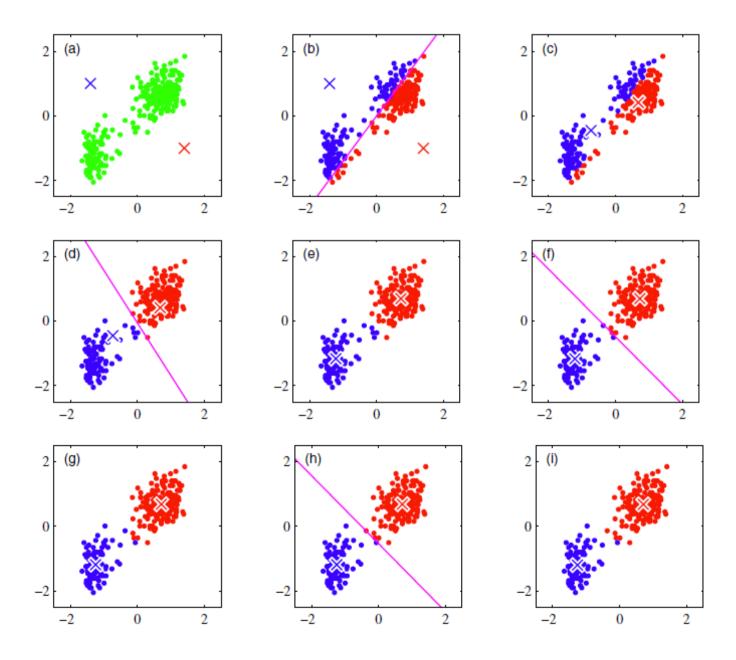
Unsupervised learning tasks

- Diversity of tasks fall into unsupervised learning category
 - Clustering (now)
 - Dimensionality reduction (autoencoders)
 - Learning parameters of probabilistic models (before/now)
- Applications and related tasks are numerous :
 - * Market basket analysis. E.g., use supermarket transaction logs to find items that are frequently purchased together
 - * Outlier detection. E.g., find potentially fraudulent credit card transactions
 - Often unsupervised tasks in (supervised) ML pipelines

(Refresher?) k-means clustering

- 1. Initialisation: choose k cluster centroids randomly
- 2. <u>Update</u>:
 - a) Assign points to the nearest* centroid
 - b) Compute centroids under the current assignment
- 3. Termination: if no change then stop
- 4. Go to Step 2
- *Distance represented by choice of metric typically L_2 Still one of the most popular data mining algorithms.

Refresher: k-means clustering



Requires specifying the number of clusters in advance

Measures
"dissimilarity" using
Euclidean distance

Finds "spherical" clusters

An iterative optimization procedure

Data: Old Faithful
Geyser Data: waiting
time between
eruptions and the
duration of eruptions

Figure: Bishop, Section 9.1

Mini Summary

- Unsupervised learning
 - * Face value: drop labels from training. That's it
 - Actually: catch-all for many many ML tasks, even as steps in supervised learning pipelines
- Refresher: k-means
 - Import next as we introduce GMMs

Next: The Gaussian mixture model

Gaussian Mixture Model

A probabilistic view of clustering. Simple example of a latent variable model.

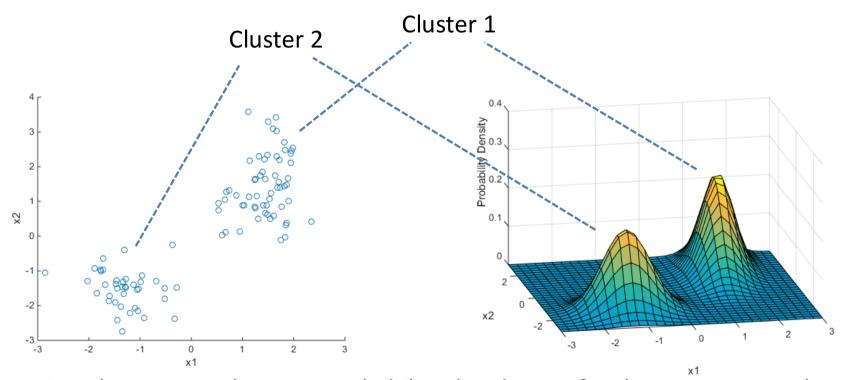
Modelling uncertainty in data clustering

- k-means clustering assigns each point to exactly one cluster
 - Does this make sense for points that are between two clusters?
 - * Clustering is often not well defined to begin with!
- Like k-means, a probabilistic mixture model requires the user to choose the number of clusters in advance
- Unlike k-means, the probabilistic model grants us power to express uncertainly about the origin of each point
 - * Each point originates from cluster c with probability w_c , c = 1, ..., k
- That is, each point still originates from one particular cluster (aka component), but we are not sure from which one
- Next
 - Clustering becomes model fitting in probabilistic sense. Philosophically satisfying.
 - Individual components modelled as Gaussians
 - Fitting illustrates general Expectation Maximization (EM) algorithm

Clustering: probabilistic model

Data points x_i are independent and identically distributed (i.i.d.) samples from a mixture of K distributions (components)

Each component in the mixture is what we call a cluster



In principle, we can adopt any probability distribution for the components, however, the normal distribution is a common modelling choice \rightarrow Gaussian Mixture Model

Normal (aka Gaussian) distribution

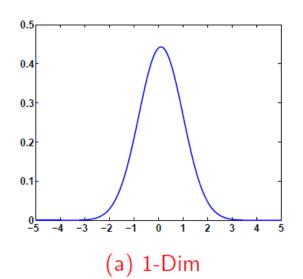
Recall that a 1D Gaussian is

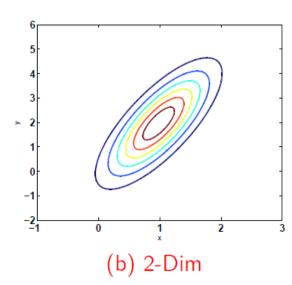
$$\mathcal{N}(x|\mu,\sigma) \equiv \frac{1}{\sqrt{2\pi\sigma^2}} \exp\left(-\frac{(x-\mu)^2}{2\sigma^2}\right)$$

And a d-dimensional Gaussian is

$$\mathcal{N}(\boldsymbol{x}|\boldsymbol{\mu},\boldsymbol{\Sigma}) \equiv (2\pi)^{-\frac{d}{2}}|\boldsymbol{\Sigma}|^{-\frac{1}{2}} \exp\left(-\frac{1}{2}(\boldsymbol{x}-\boldsymbol{\mu})^T \boldsymbol{\Sigma}^{-1}(\boldsymbol{x}-\boldsymbol{\mu})\right)$$

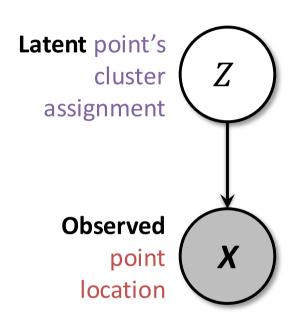
- * Σ is a PSD symmetric $d \times d$ matrix, the covariance matrix
- * |Σ| denotes determinant
- * No need to memorize the full formula.





Gaussian mixture model (GMM): One point

- Cluster assignment of point
 - Multinomial distribution on k outcomes
 - * P(Z = j) described by $P(C_j) = w_j \ge 0$ with $\sum_{j=1}^k w_j = 1$
- Location of point
 - Each cluster has its own Gaussian distribution
 - Location of point governed by its cluster assignment
 - * $P(X|Z=j) = \mathcal{N}(\mu_j, \Sigma_j)$ class conditional density
- Model's **parameters**: w_j , μ_j , Σ_j , j=1,...,k



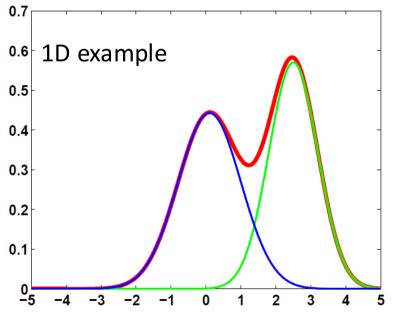
From marginalisation to mixture distribution

- When fitting the model to observations, we'll be maximising likelihood of observed portions of the data (the X's) not the latent parts (the Z's)
- Marginalising out the Z's derives the "familiar" mixture distribution
- Gaussian mixture distribution:

$$P(\mathbf{x}) \equiv \sum_{j=1}^{k} w_j \mathcal{N}(\mathbf{x} | \boldsymbol{\mu}_j, \boldsymbol{\Sigma}_j)$$

$$\equiv \sum_{j=1}^{k} P(C_j) P(\mathbf{x} | \boldsymbol{C}_j)$$

- A convex combination of Gaussians
- Simply marginalisation at work

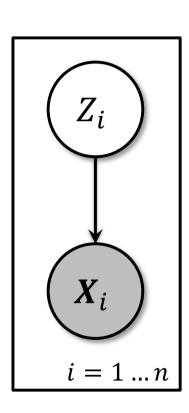


Mixture and individual component densities are re-scaled for visualisation purposes

Figure: Bishop

Clustering as model estimation

- Given a set of data points, we assume that data points are generated by a GMM
 - Each point in our dataset originates from our mixture distribution
 - * Shared parameters between points: independence assumption!!
- Clustering now amounts to finding parameters of the GMM that "best explains" observed data
- Call upon old friend MLE principle to find parameter values that maximise $p(x_1, ..., x_n)$



Mini Summary

- GMM is just another D-PGM
- Some variables are observed some latent
- Convenient to model location as generated by cluster assignment
- Shared clusters arise from independence b/w points
- Mixture distribution arises algebraically from marginalisation

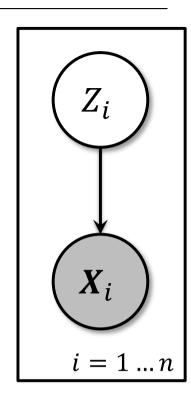
Next: MLE to fit the model, again motivating EM algorithm

Motivating (again) Expectation-Maximisation Algorithm

We want to implement MLE but we have unobserved r.v.'s that prevent clean decomposition as happens in fully observed settings

Fitting the GMM

• Modelling the data points as independent, aim is to find $P(C_j)$, μ_j , Σ_j , j=1,...,k that maximise $P(x_1,...,x_n) = \prod_{i=1}^n \sum_{j=1}^k P(C_j) P(x_i|C_j)$ where $P(x|C_j) \equiv \mathcal{N}(x|\mu_j,\Sigma_j)$ Can be solved analytically?



 Taking the derivative of this expression is pretty awkward, try the usual log trick

$$\log P(\mathbf{x}_1, \dots, \mathbf{x}_n) = \sum_{i=1}^n \log \left(\sum_{j=1}^k P(C_j) P(\mathbf{x}_i | \mathbf{C}_j) \right)$$

→ Expectation-Maximisation (EM)

Motivation of EM

- Consider a parametric probabilistic model $p(X|\theta)$, where X denotes data and θ denotes a vector of parameters
- According to MLE, we need to maximise $p(X|\theta)$ as a function of θ
 - * equivalently maximise $\log p(X|\theta)$



- There can be a couple of issues with this task
- Sometimes we don't observe some of the variables needed to compute the log likelihood
 - Example: GMM cluster membership Z is not known in advance
- Sometimes the form of the log likelihood is inconvenient to work with
 - Example: taking a derivative of GMM log likelihood results in a cumbersome equation

Expectation-Maximisation (EM) Algorithm

- Initialisation Step:
 - * Initialize K clusters: C_1 , ..., C_K (μ_i, Σ_i) and $P(C_i)$ for each cluster j.
- Iteration Step:
 - * Estimate the cluster of each datum $p(C_i | x_i)$



Re-estimate the cluster parameters



 $(\mu_j, \Sigma_j), p(C_j)$ for each cluster j

Summary

- Unsupervised learning
 - Diversity of problems
- Gaussian mixture model (GMM)
 - A probabilistic approach to clustering
 - * The GMM model
 - GMM clustering as an optimisation problem
- MLE: Motivating Expectation Maximization (EM)

Next lecture: Getting to the bottom of EM