Machine Learning for Data Science (CS4786)
Lecture 16

Probabilistic Modeling and EM Algorithm

Probabilistic Modeling

Probabilistic Model

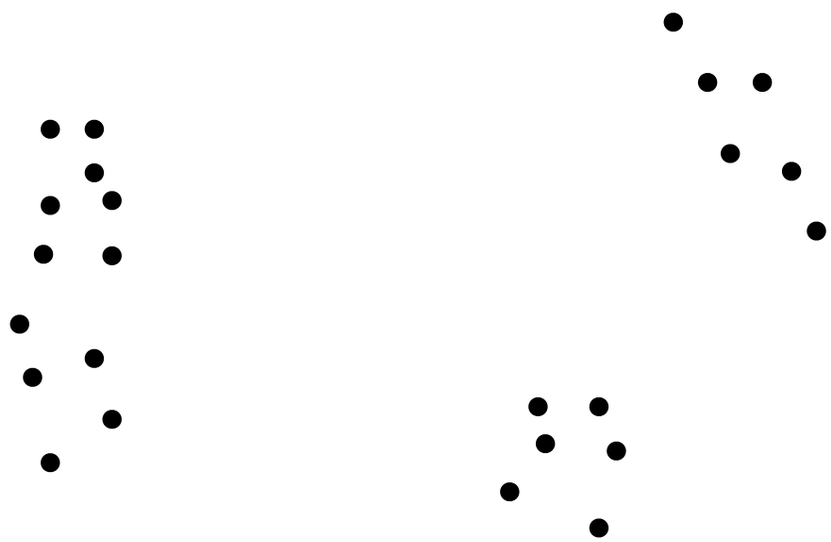
Data: $\mathbf{x}_1 \ldots, \mathbf{x}_n$

Probabilistic Model

$\theta \in \Theta$
P_θ explains data

Data: $\mathbf{x}_1 \ldots, \mathbf{x}_n$

Probabilistic Model



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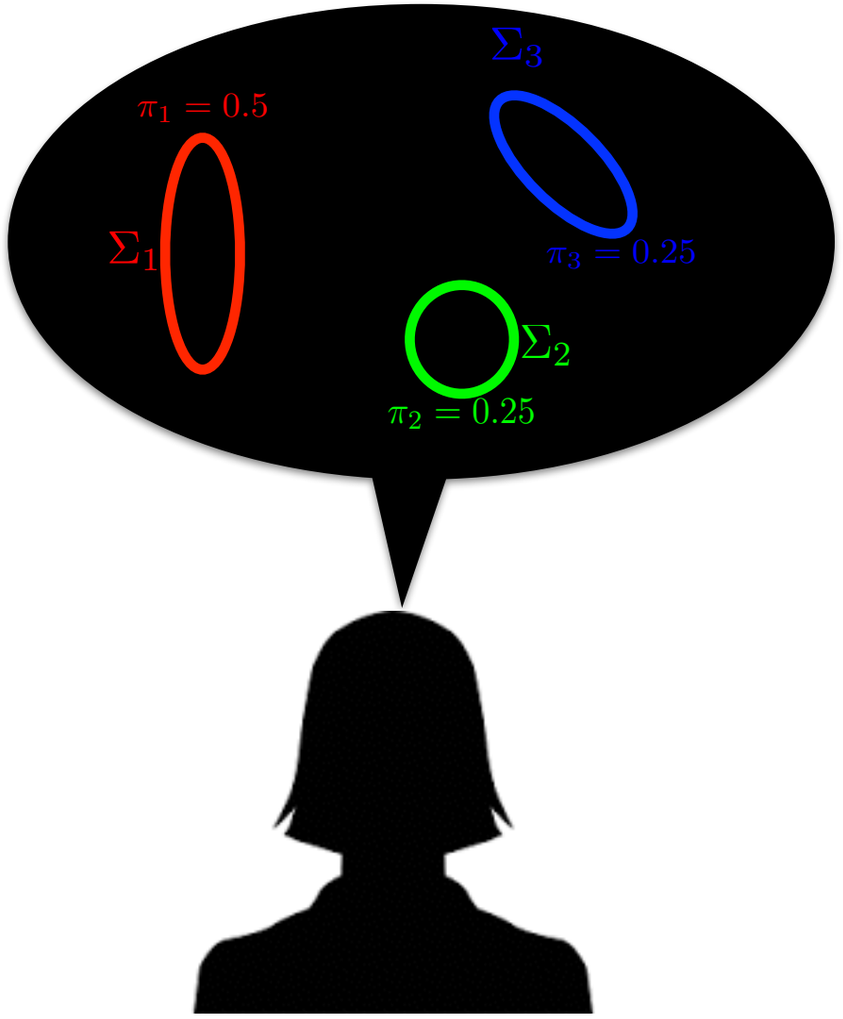

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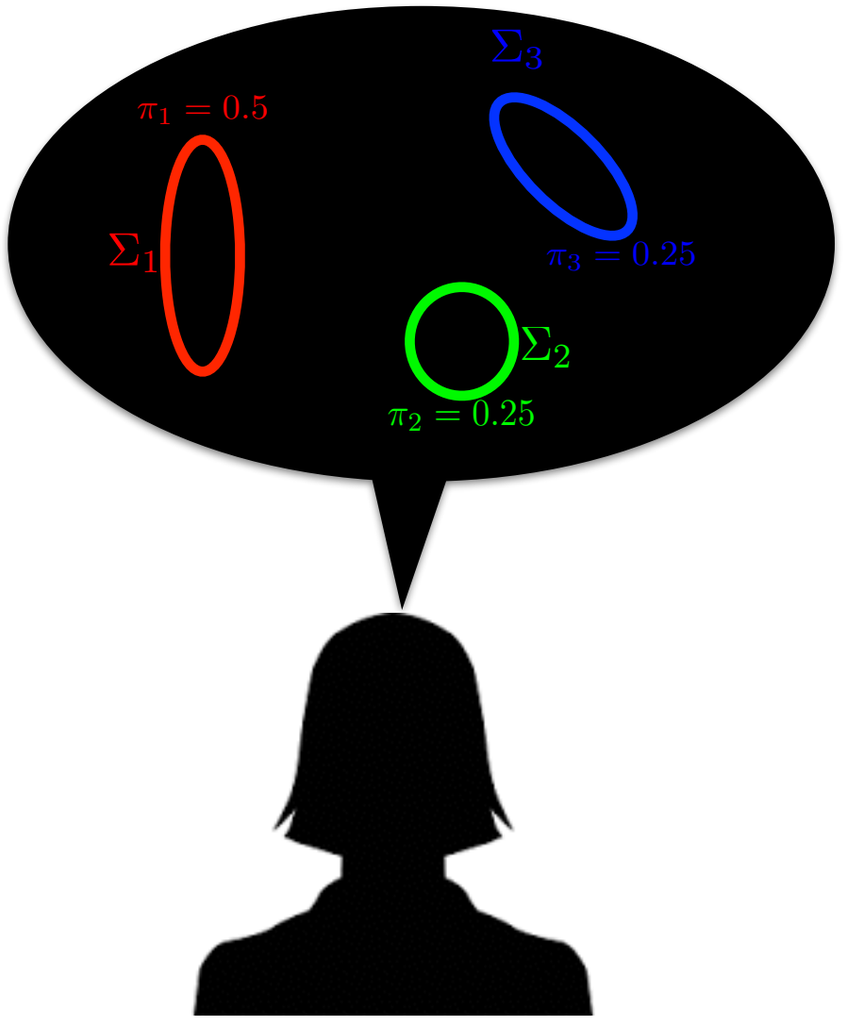

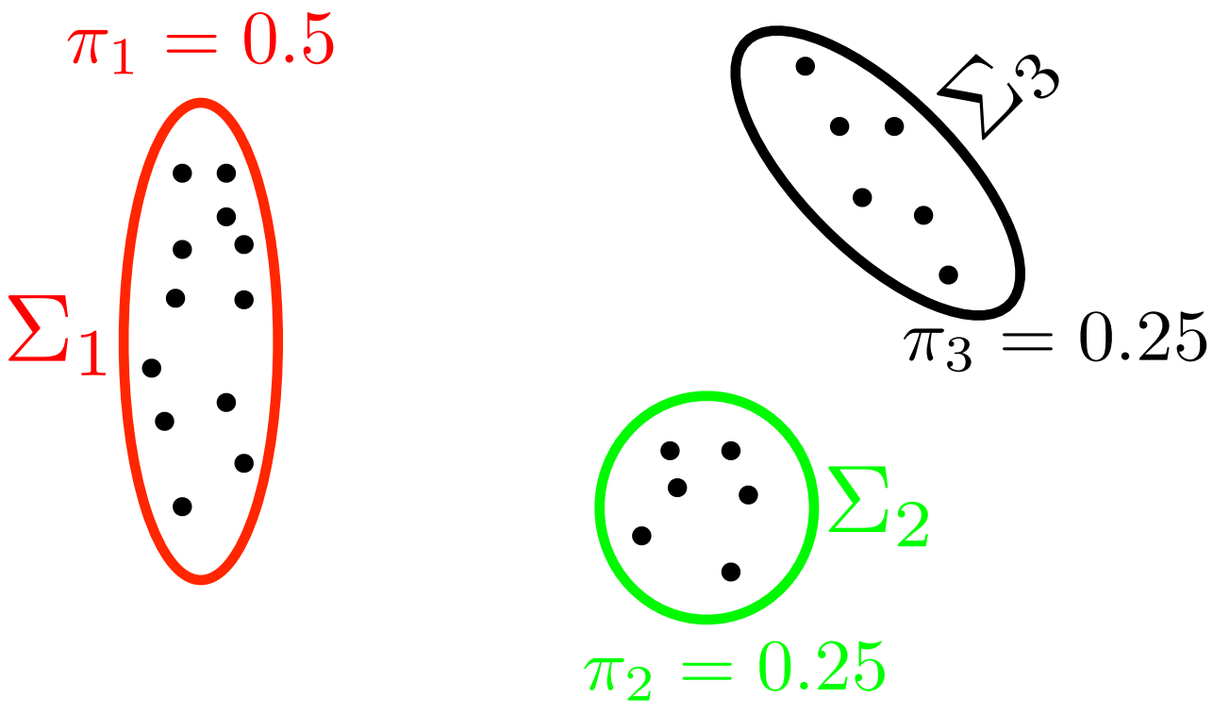

Probabilistic Models

- Set of models Θ consists of parameters s.t. P_θ for each $\theta \in \Theta$ is a distribution over data.

- Learning: Estimate $\theta^* \in \Theta$ that best models given data

Maximum Likelihood Principal

Pick $\theta \in \Theta$ that maximizes probability of observation

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Often referred to as frequentist view

Maximum Likelihood Principal

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$$\theta_{MLE} = \text{argmax}_{\theta \in \Theta} \underbrace{\log P_\theta(x_1, \ldots, x_n)}_{\text{Likelihood}}$$

- A priori all models are equally good, data could have been generated by any one of them

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Say you had a prior belief about models provided by $P(\theta)$

Pick $\theta \in \Theta$ that is most likely given data

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There are Bayesian and there Bayesians

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Maximize a posteriori probability of model given data

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The Bayesian Choice

Don't pick any $\theta^* \in \Theta$

- Model is simply an abstraction
- We have a prosteriori distribution over models, why pick one θ?

$$P(X|\text{data}) = \sum_{\theta \in \Theta} P(X, \theta|\text{data}) = \sum_{\theta \in \Theta} P(X|\theta) P(\theta|\text{data})$$

Lets get back to GMM

Hard Gaussian Mixture Model

- For all $j \in [K]$, initialize cluster centroids $\hat{\mathbf{r}}_j^0$, ellipsoids $\hat{\Sigma}_j^0$ and initial proportions π^0 randomly and set $m = 1$
- Repeat until convergence (or until patience runs out)
  1. For each $t \in \{1, \ldots, n\}$, set cluster identity of the point

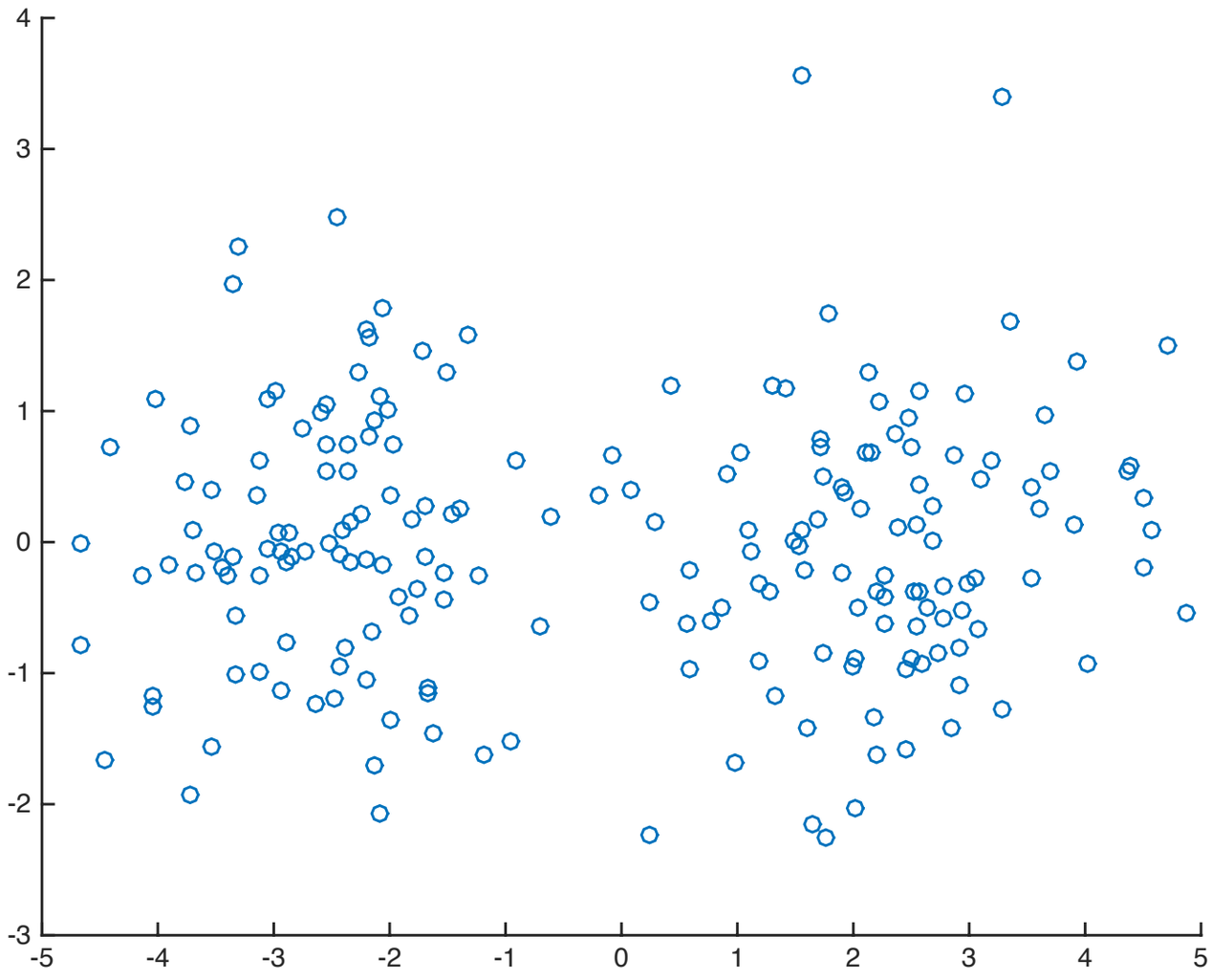
$$\hat{c}^m(\mathbf{x}_t) = \arg\max_{j \in [K]} \; p(\mathbf{x}_t, \hat{\mathbf{r}}_j^{m-1}, \hat{\Sigma}^{m-1}) \times \pi^m(j)$$

  2. For each $j \in [K]$, set new representative as

$$\hat{\mathbf{r}}_j^m = \frac{1}{|\hat{C}_j^m|} \sum_{\mathbf{x}_t \in \hat{C}_j^m} \mathbf{x}_t \qquad \hat{\Sigma}^m = \frac{1}{|C_j|} \sum_{t \in C_j} (\mathbf{x}_t - \hat{\mathbf{r}}_j^m)(\mathbf{x}_t - \hat{\mathbf{r}}_j^m)^\top \qquad \pi_j^m = \frac{|C_j^m|}{n}$$

  3. $m \leftarrow m + 1$

Pitfall of Hard Assignment



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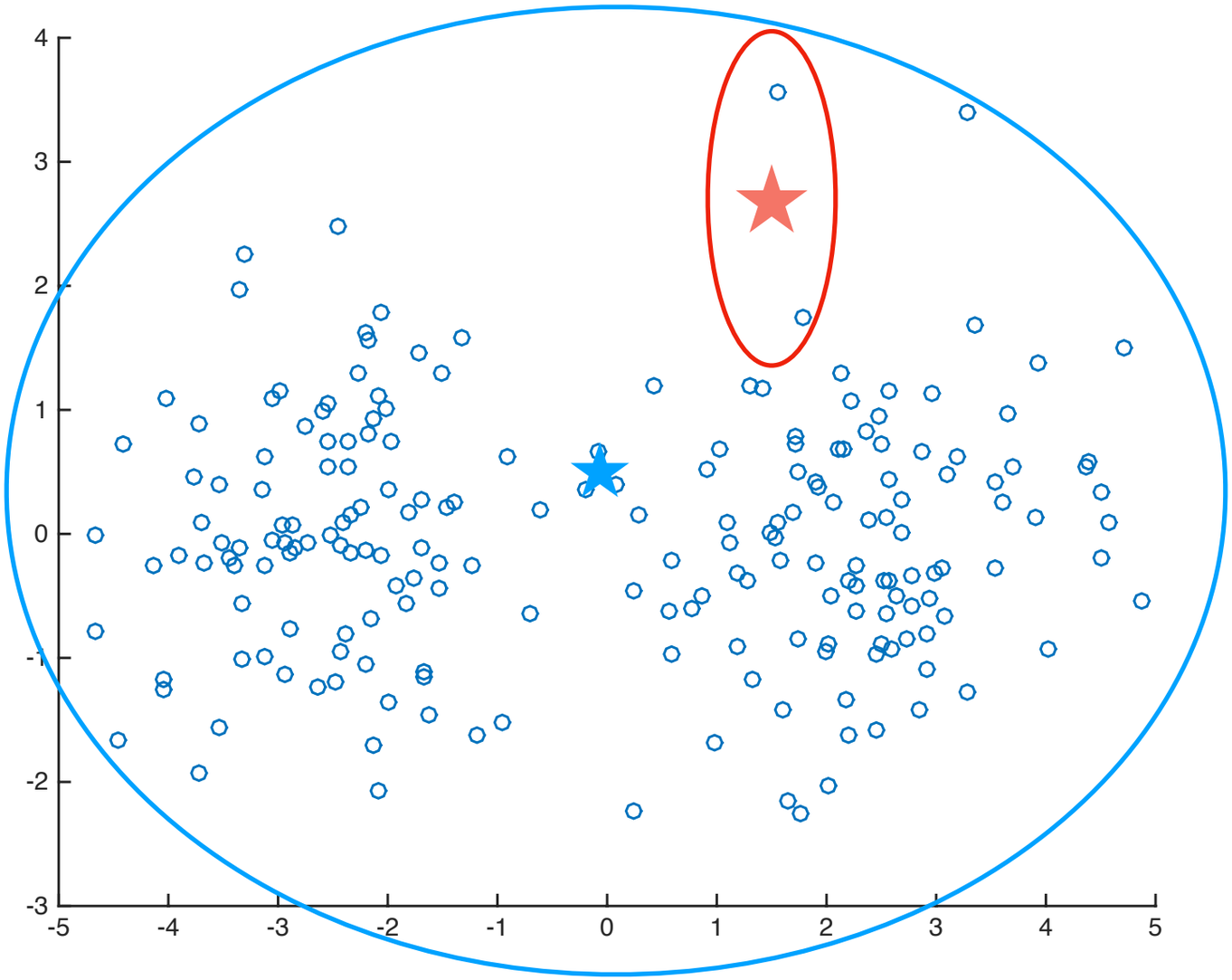

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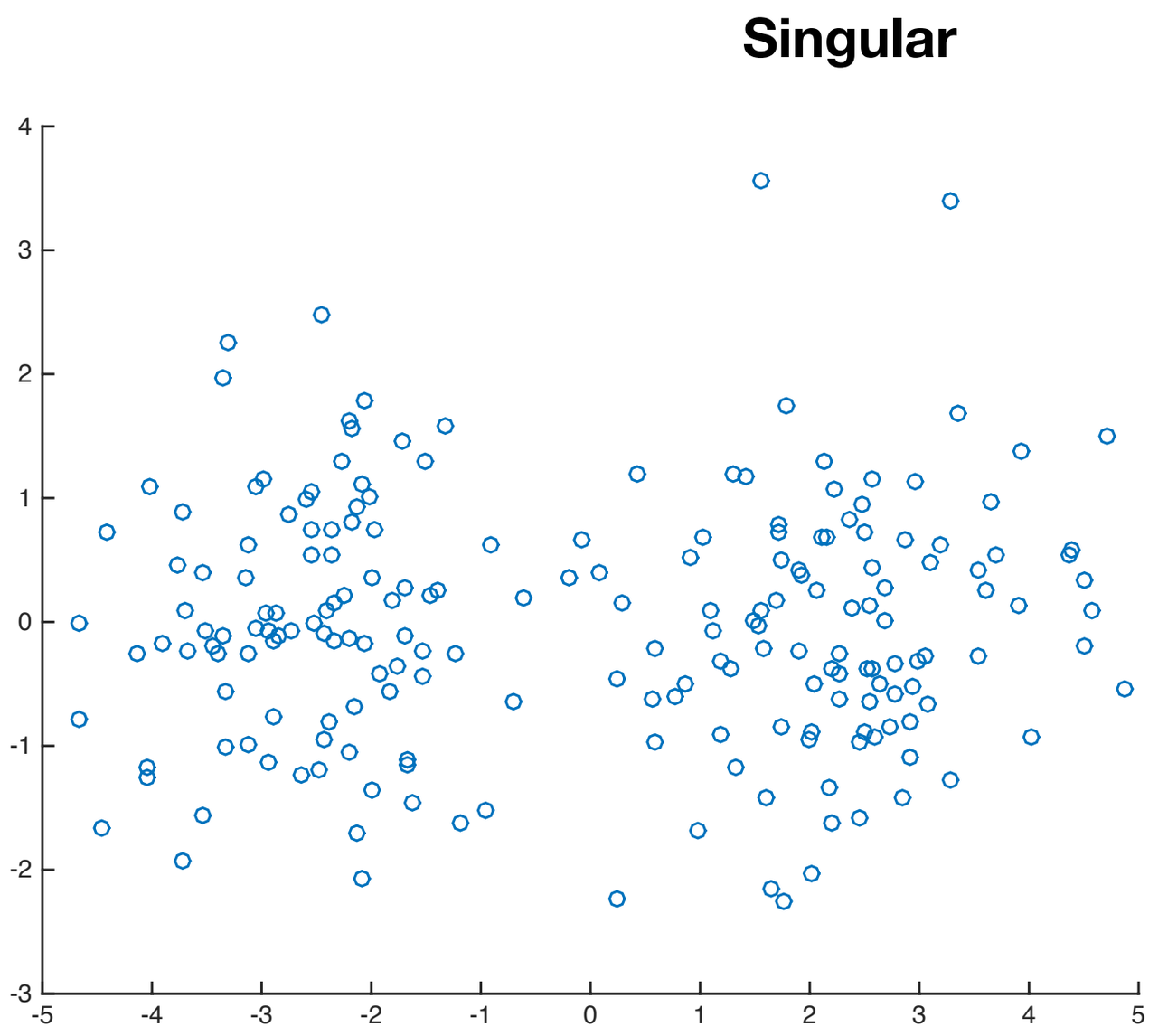

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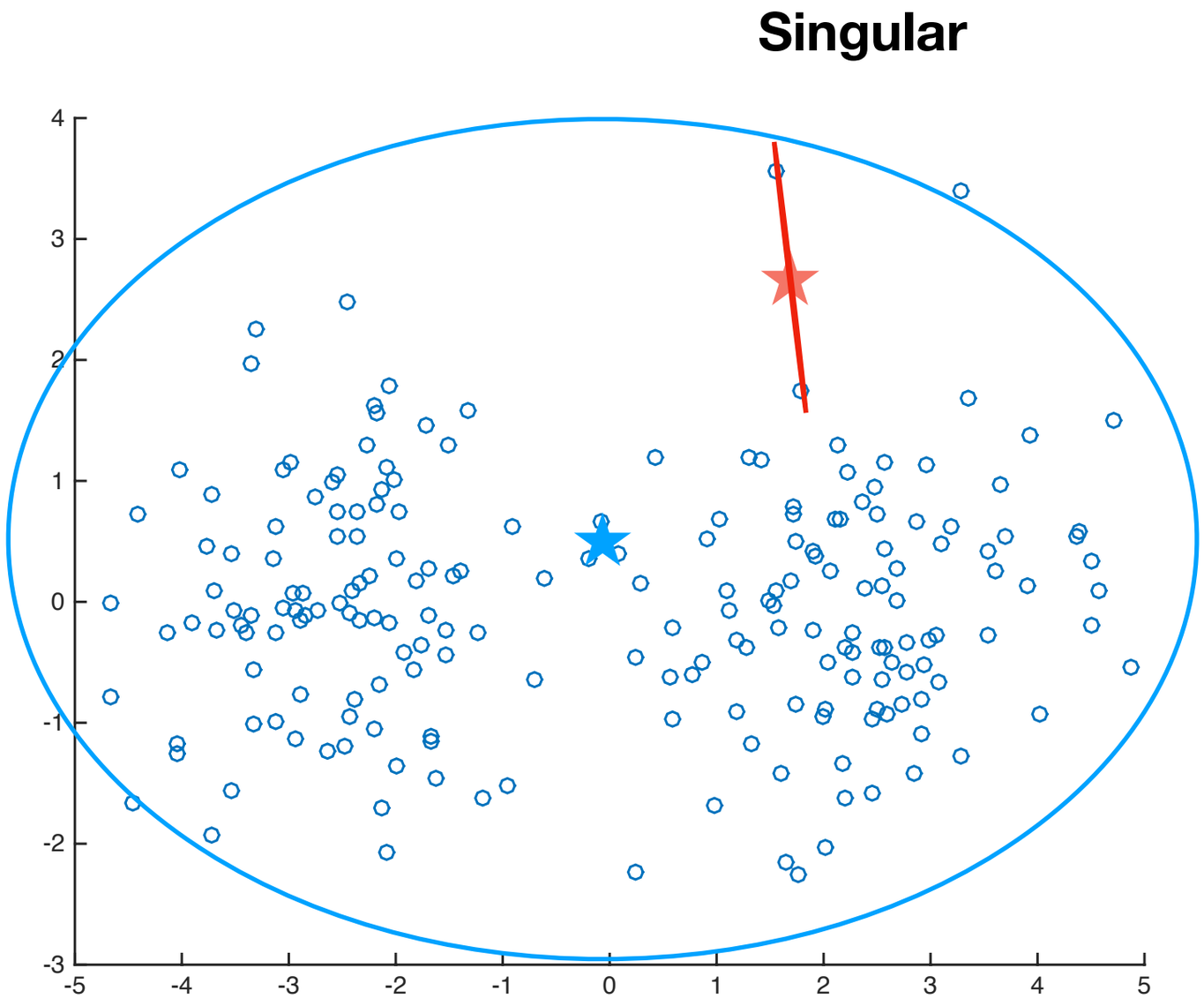
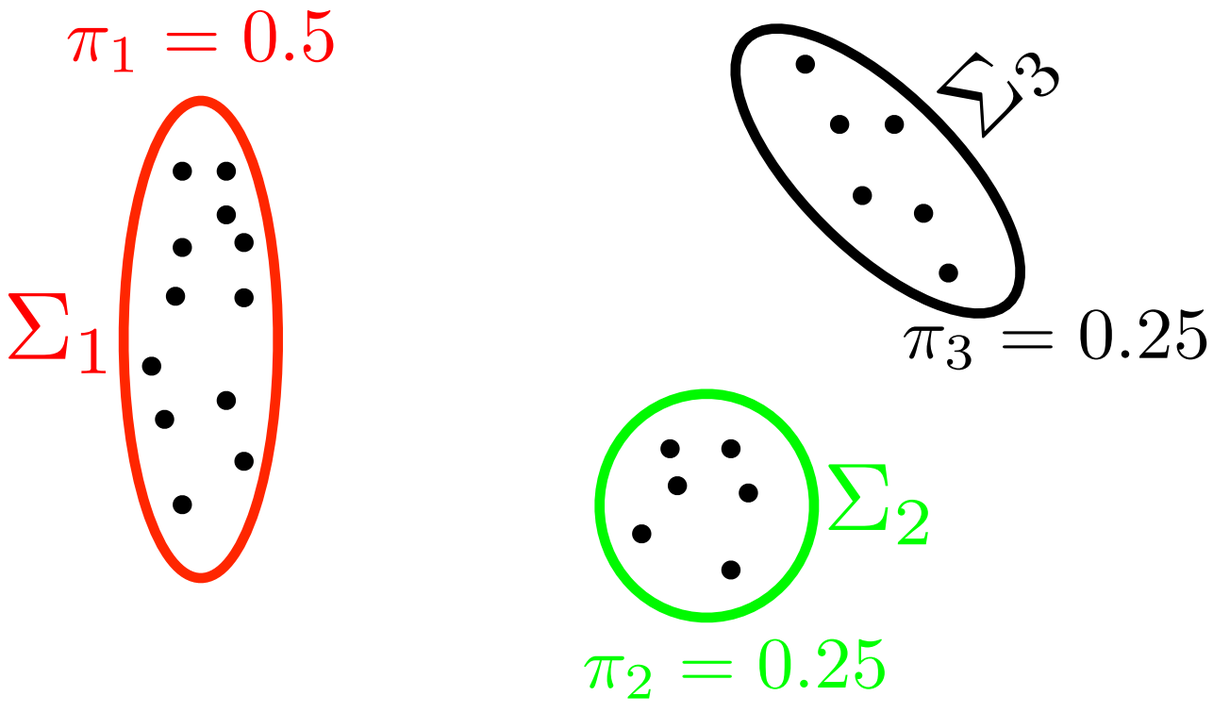


MLE for GMM

Say we knew model parameters, how do we assign clusters?

Given probability of each point belonging to each of the clusters, how do we compute model parameters?

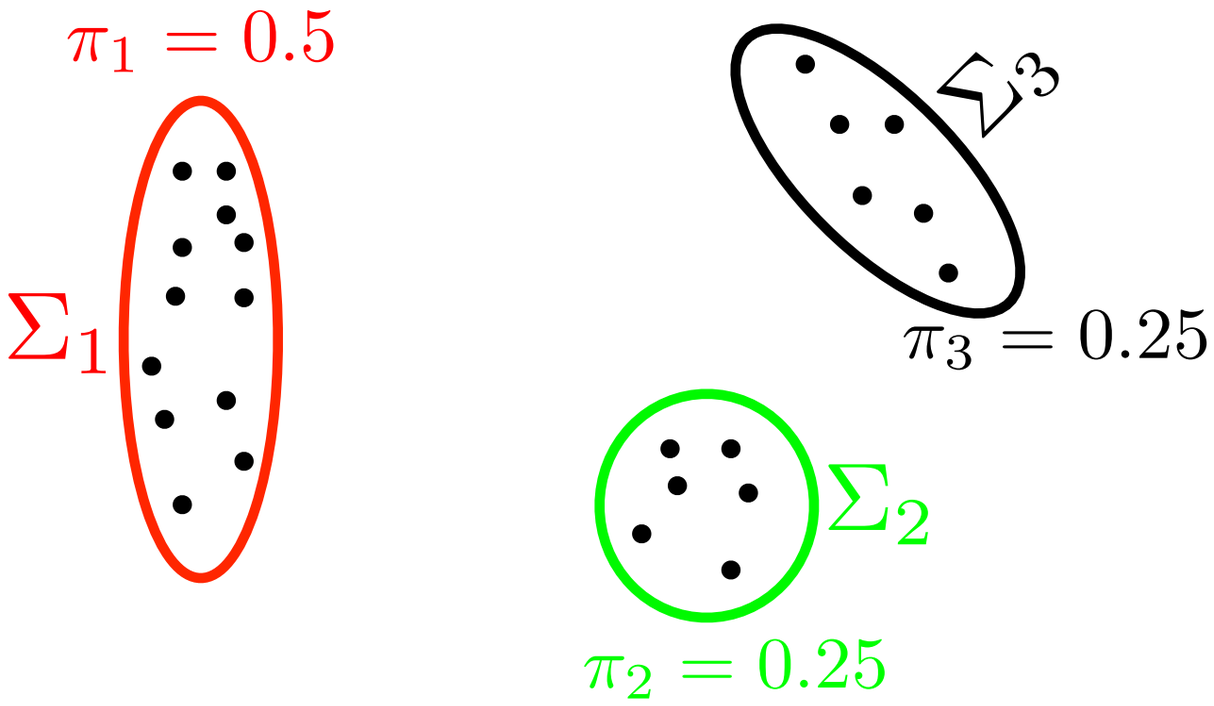


MLE for GMM

Say we knew model parameters, ~~how do we assign clusters?~~

what are the probabilities of points falling in each of the clusters?

Given probability of each point belonging to each of the clusters, how do we compute model parameters?



(Soft) Gaussian Mixture Model

- For all $j \in [K]$, initialize cluster centroids $\hat{\mathbf{r}}_j^0$ and ellipsoids $\hat{\Sigma}_j^0$ randomly and set $m = 1$
- Repeat until convergence (or until patience runs out)
  1. For each $t \in \{1, \ldots, n\}$, set cluster identity of the point

$$Q_t^m(j) = p(\mathbf{x}_t, \hat{\mathbf{r}}_j^{m-1}, \hat{\Sigma}^{m-1}) \times \pi^m(j)$$

  2. For each $j \in [K]$, set new representative as

$$\hat{\mathbf{r}}_j^m = \frac{\sum_{t=1}^n Q_t(j)\mathbf{x}_t}{\sum_{t=1}^n Q_t(j)} \qquad \hat{\Sigma}^m = \frac{\sum_{t=1}^n Q_t(j)(\mathbf{x}_t - \hat{\mathbf{r}}_j^m)(\mathbf{x}_t - \hat{\mathbf{r}}_j^m)^\top}{\sum_{t=1}^n Q_t(j)}$$

$$\pi_j^m = \frac{\sum_{t=1}^n Q_t(j)}{n}$$

  3. $m \leftarrow m + 1$

Expectation Maximization Algorithm

- For demonstration we shall consider the problem of finding MLE (MAP version is very similar)
- Initialize $\theta^{(0)}$ arbitrarily, repeat unit convergence:

(E step) For every t, define distribution Q_t over the latent variable c_t as:

$$Q_t^{(i)}(c_t) = P(c_t|x_t, \theta^{(i-1)})$$

(M step)

$$\theta^{(i)} = \text{argmax}_{\theta \in \Theta} \sum_{t=1}^{n} \sum_{c_t} Q_t^{(i)}(c_t) \log P(x_t, c_t|\theta)$$

Example: EM for GMM

- E step: For every $k \in [K]$,

$$Q_t^{(i)}(c_t = k) = P\left(c_t = k | x_t, \theta^{(i-1)}\right) = P\left(x_t | c_t = k, \theta^{(i-1)}\right) \times P\left(c_t = k | \theta^{(i-1)}\right)$$
$$\propto \underbrace{\phi\left(x_t; \mu_k^{(i-1)}, \Sigma_k^{(i-1)}\right)}_{\text{gaussian p.d.f.}} \times \pi_k^{(i-1)}$$

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- M step: Given $Q_1, \ldots, Q_n$, we need to find

$$\theta^{(i)} = \operatorname*{argmax}_{\theta \in \Theta} \sum_{t=1}^{n} \sum_{k=1}^{K} Q_t^{(i)}(k) \log P(x_t, c_t = k | \theta)$$
$$= \operatorname*{argmax}_{\theta} \sum_{t=1}^{n} \sum_{k=1}^{K} Q_t^{(i)}(k) \left(\log P(x_t | c_t = k, \theta) + \log P(c_t = k | \theta)\right)$$
$$= \operatorname*{argmax}_{\pi, \mu_{1,\ldots,K}, \Sigma_{1,\ldots,K}} \sum_{t=1}^{n} \sum_{c_t=1}^{K} Q_t^{(i)}(k) \left(\log \phi\left(x_t; \mu_k, \Sigma_k\right) + \log \pi_k\right)$$

Example: EM for GMM

For every $k \in [K]$, the maximization step yields,

$$\mu_k^{(i)} = \frac{\sum_{t=1}^n Q_t^{(i)}(k) x_t}{\sum_{t=1}^n Q_t(k)}, \quad \Sigma_k^{(i)} = \frac{\sum_{t=1}^n Q_t^{(i)}(k) \left(x_t - \mu_k^{(i)}\right)\left(x_t - \mu_k^{(i)}\right)^\top}{\sum_{t=1}^n Q_t(k)}$$

$$\pi_k^{(i)} = \frac{\sum_{t=1}^n Q_t^{(i)}(k)}{n}$$

Let us derive this!

Why should EM work?

A very high level view:

- Performing E-step will never decrease log-likelihood (or log a posteriori)

Why should EM work?

A very high level view:

- Performing E-step will never decrease log-likelihood (or log a posteriori)
- Performing M-step will never decrease log-likelihood (or log a posteriori)

Why should EM work?

Steps to show that $\log \mathrm{Lik}(\theta^{(i)}) \geq \log \mathrm{Lik}(\theta^{(i-1)})$:

$$\log P_{\theta^{(i)}}(x_1, \ldots, x_n)$$

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Log(average) > average of Log

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Why should EM work?

- Likelihood never decreases
- So whenever we converge we converge to a local optima
- However problem is non-convex and can have many local optimal
- In general no guarantee on rate of convergence
- In practice, do multiple random initializations and pick the best one!

EM Algorithm Generally

- More generally, EM can be used to learn any probabilistic model with some Latent (unseen) variables and some observed variables whenever
  - Its is easy to find parameters given distribution/ observation for all variables
  - Given all parameters finding distribution for latent variables is easy

How to choose K (no. of clusters)

- Elbow method:
  - plot Objective versus K, typically it monotonically decreases.
  - Pick point where there is a kink
    - Intuition: look at rate of change
- Add to objective penalty (+ pen(K)) and minimize, pen increases with K
  - intuition we prefer smaller number of clusters
  - Use prior knowledge to pick p
  - (AIC, BIC etc can been seen to be specific cases)
- We can leave the burden of choosing K to the probabilistic model