

Unsupervised Learning

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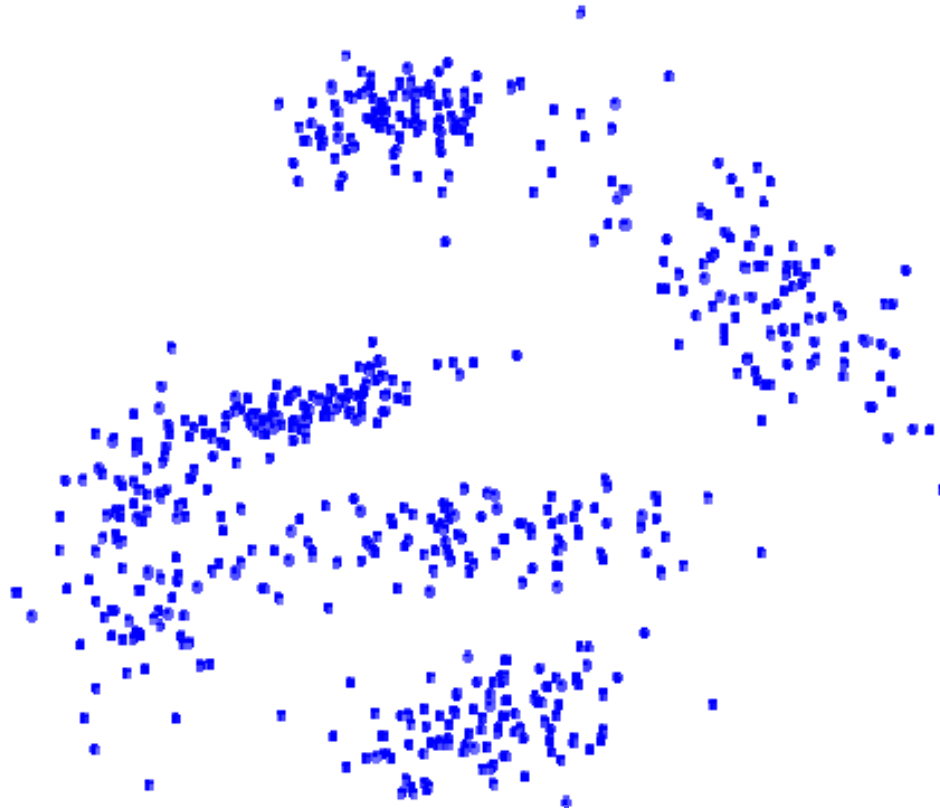
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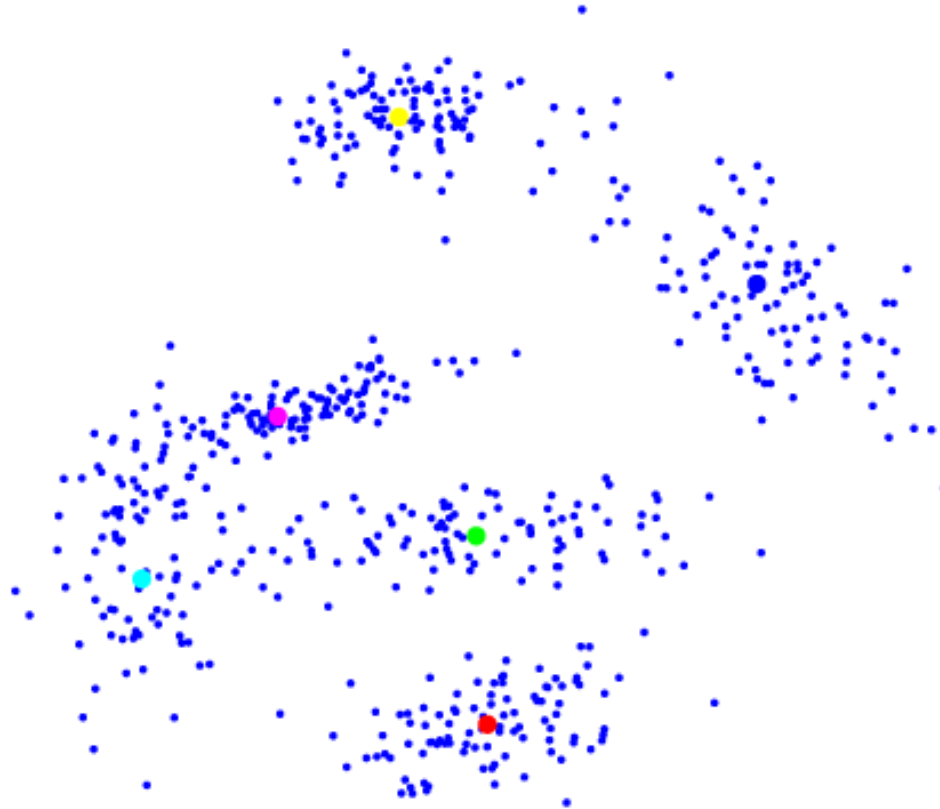
Content

- Fundamentals of Unsupervised Learning
 - K-means clustering
 - Principal component analysis
- Probabilistic Unsupervised Learning
 - Mixture Gaussians
 - EM Methods

K-Means Clustering



K-Means Clustering



K-Means Clustering

- Provide the number of desired clusters k
- Randomly choose k instances as seeds, one per each cluster, i.e. the centroid for each cluster
- Iterate
 - Assign each instance to the cluster with the closest centroid
 - Re-estimate the centroid of each cluster
- Stop when clustering converges
 - Or after a fixed number of iterations

K-Means Clustering: Centriod

- Assume instances are real-valued vectors

$$x \in \mathbb{R}^d$$

- Clusters based on centroids, center of gravity, or mean of points in a cluster C_k

$$\mu^k = \frac{1}{C_k} \sum_{x \in C_k} x$$

K-Means Clustering: Distance

- Distance to a centroid $L(x, \mu^k)$
- Euclidian distance (L2 norm)

$$L_2(x, \mu^k) = \|x - \mu^k\| = \sqrt{\sum_{m=1}^d (x_i - \mu_m^k)^2}$$

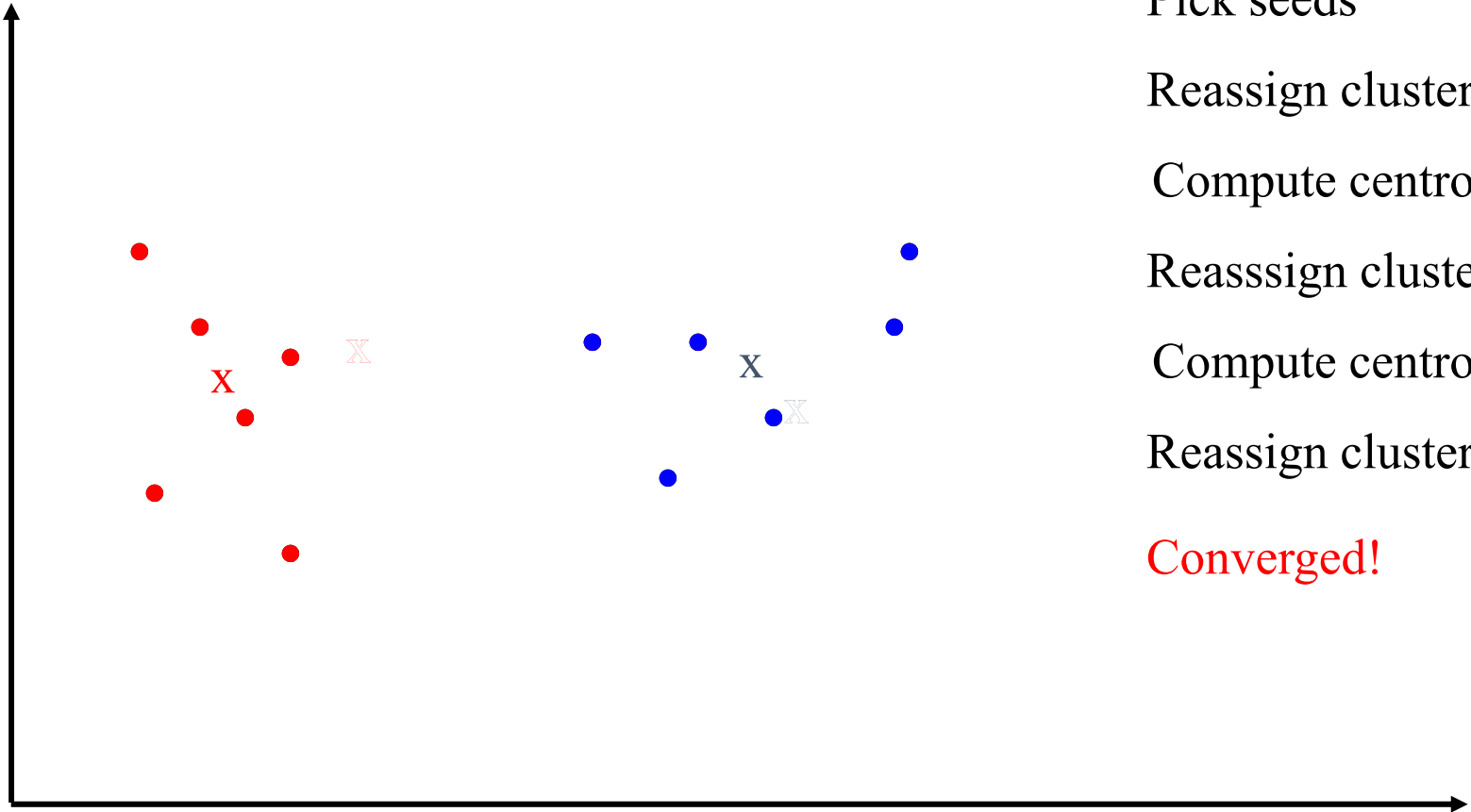
- Euclidian distance (L1 norm)

$$L_1(x, \mu^k) = |x - \mu^k| = \sum_{m=1}^d |x_i - \mu_m^k|$$

- Cosine distance

$$L_{\cos}(x, \mu^k) = 1 - \frac{x^\top \mu^k}{|x| \cdot |\mu^k|}$$

K-Means Example ($K=2$)



Pick seeds

Reassign clusters

Compute centroids

Reassign clusters

Compute centroids

Reassign clusters

Converged!

K-Means Time Complexity

- Assume computing distance between two instances is $O(d)$ where d is the dimensionality of the vectors
- Reassigning clusters: $O(knd)$ distance computations
- Computing centroids: Each instance vector gets added once to some centroid: $O(nd)$
- Assume these two steps are each done once for l iterations: $O(lknd)$

K-Means Clustering Objective

- The objective of K -means is to minimize the total sum of the squared distance of every point to its corresponding cluster centroid

$$\min_{\{\mu^k\}_{k=1}^K} \sum_{k=1}^K \sum_{x \in C_k} L(x - \mu^k) \qquad \mu^k = \frac{1}{C_k} \sum_{x \in C_k} x$$

- Finding the global optimum is NP-hard.
- The K -means algorithm is guaranteed to converge a local optimum.

Seed Choice

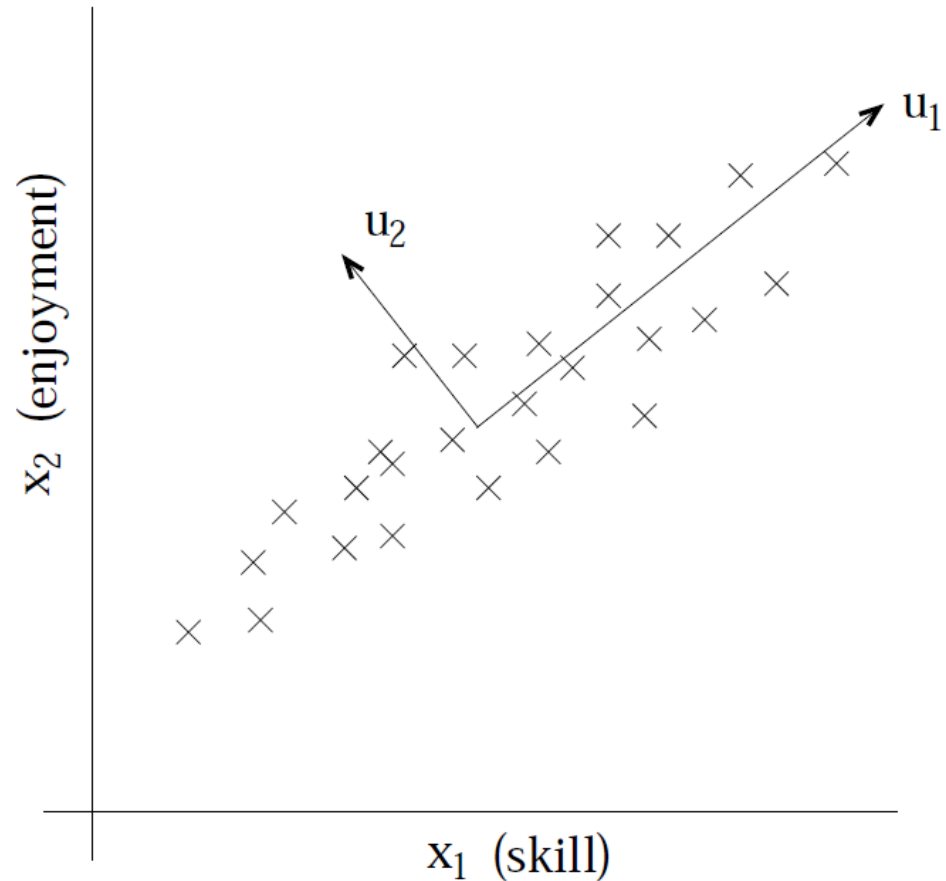
- Results can vary based on random seed selection.
- Some seeds can result in poor convergence rate, or convergence to sub-optimal clusterings.
- Select good seeds using a heuristic or the results of another method.

Clustering Applications

- Text mining
 - Cluster documents for related search
 - Cluster words for query suggestion
- Recommender systems and advertising
 - Cluster users for item/ad recommendation
 - Cluster items for related item suggestion
- Image search
 - Cluster images for similar image search and duplication detection
- Speech recognition or separation
 - Cluster phonetical features

Principal Component Analysis (PCA)

- An example of 2-dimensional data
 - x_1 : the piloting skill of pilot
 - x_2 : how much he/she enjoys flying
- Main components
 - u_1 : intrinsic piloting “karma” of a person
 - u_2 : some noise



Principal Component Analysis (PCA)

- PCA tries to identify the subspace in which the data approximately lies
- PCA uses an **orthogonal transformation** to convert a set of observations of possibly correlated variables into a set of values of **linearly uncorrelated variables** called principal components.
 - The number of principal components is less than or equal to the smaller of the number of original variables or the number of observations.

$$\mathbb{R}^d \rightarrow \mathbb{R}^k \quad k \ll d$$

PCA Data Preprocessing

- Given the dataset

$$D = \{x^{(i)}\}_{i=1}^m$$

- Typically we first pre-process the data to normalize its mean and variance

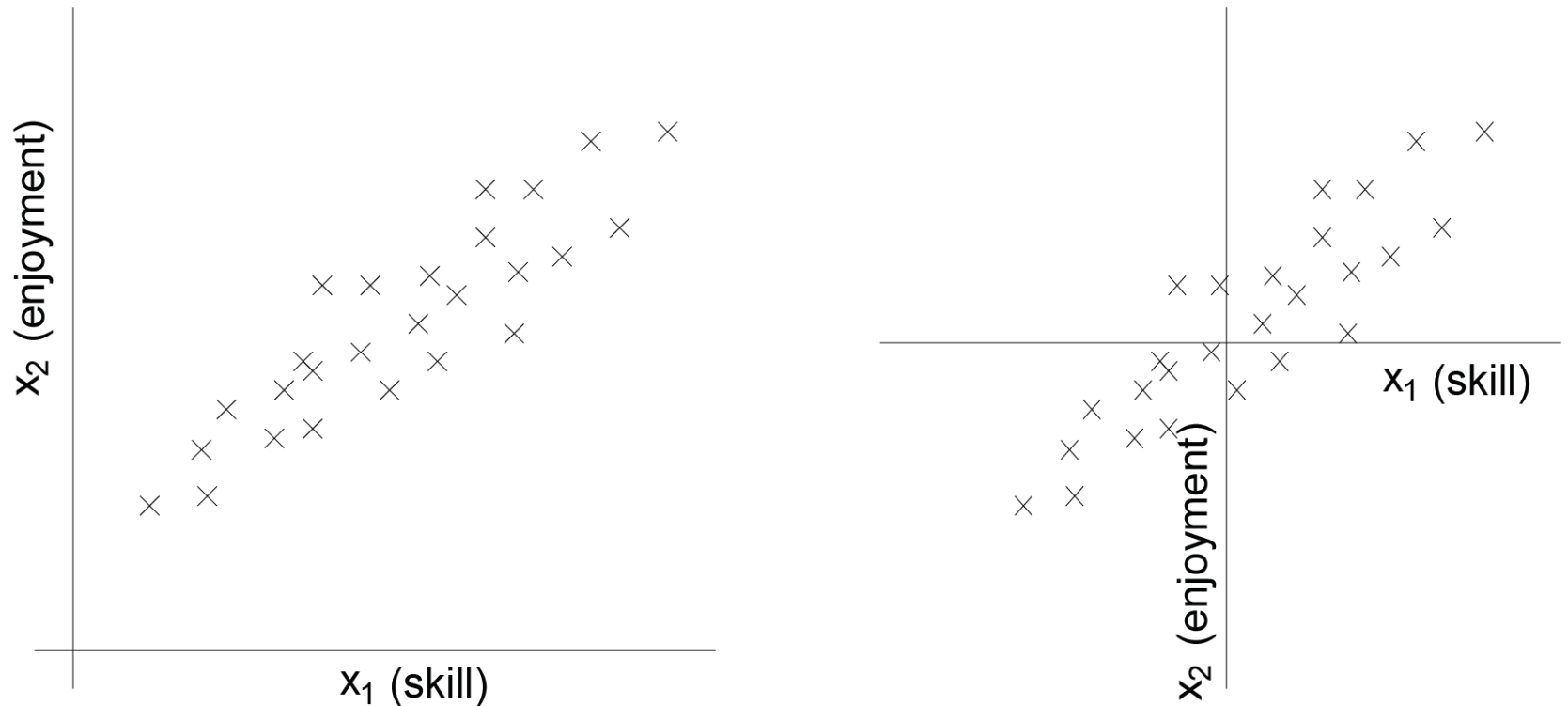
1. Move the central of the data set to 0

$$\mu = \frac{1}{m} \sum_{i=1}^m x^{(i)} \qquad x^{(i)} \leftarrow x^{(i)} - \mu$$

2. Unify the variance of each variable

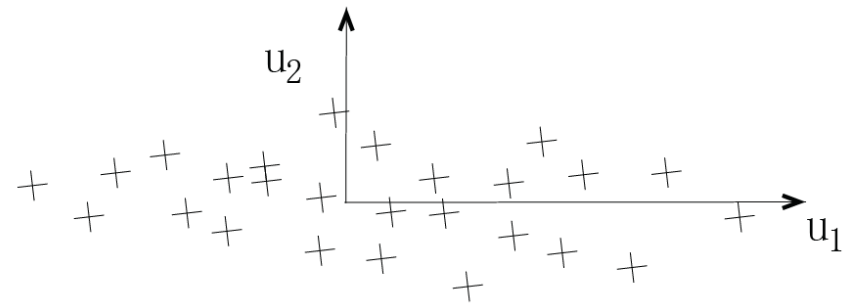
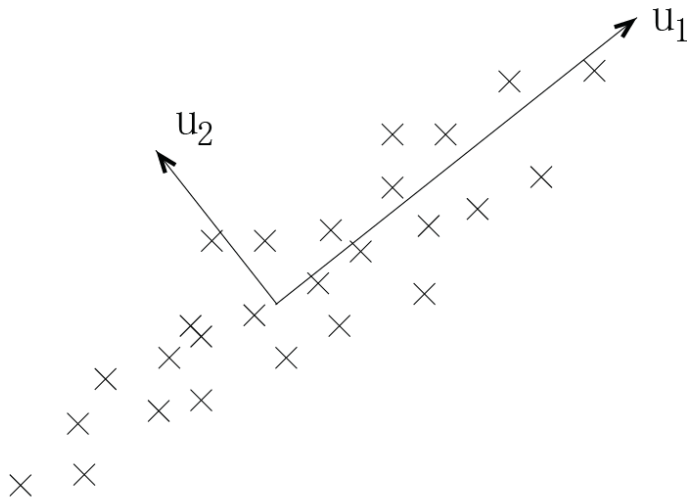
$$\sigma_j^2 = \frac{1}{m} \sum_{i=1}^m (x_j^{(i)})^2 \qquad x^{(i)} \leftarrow x^{(i)} / \sigma_j$$

PCA Data Preprocessing



- Zero out the mean of the data
- Rescale each coordinate to have unit variance, which ensures that different attributes are all treated on the same “scale”.

PCA Solution



- PCA finds the directions with the largest variable variance
 - which correspond to the eigenvectors of the matrix $X^T X$ with the largest eigenvalues

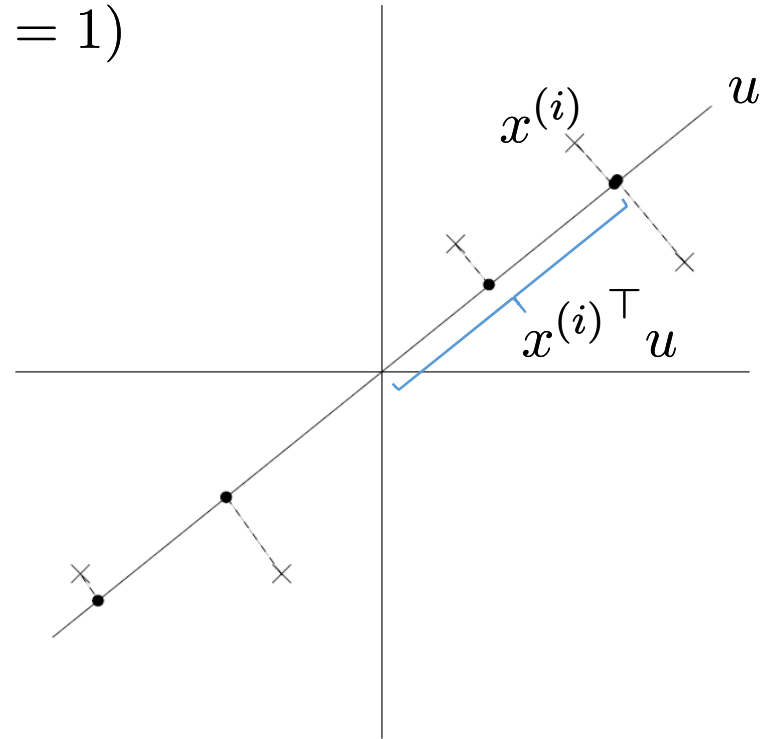
PCA Solution: Data Projection

- The projection of each point $x^{(i)}$ to a direction u ($\|u\| = 1$)

$$x^{(i)\top} u$$

- The variance of the projection

$$\begin{aligned}\frac{1}{m} \sum_{i=1}^m (x^{(i)\top} u)^2 &= \frac{1}{m} \sum_{i=1}^m u^\top x^{(i)} x^{(i)\top} u \\ &= u^\top \left(\frac{1}{m} \sum_{i=1}^m x^{(i)} x^{(i)\top} \right) u \\ &\equiv u^\top \Sigma u\end{aligned}$$



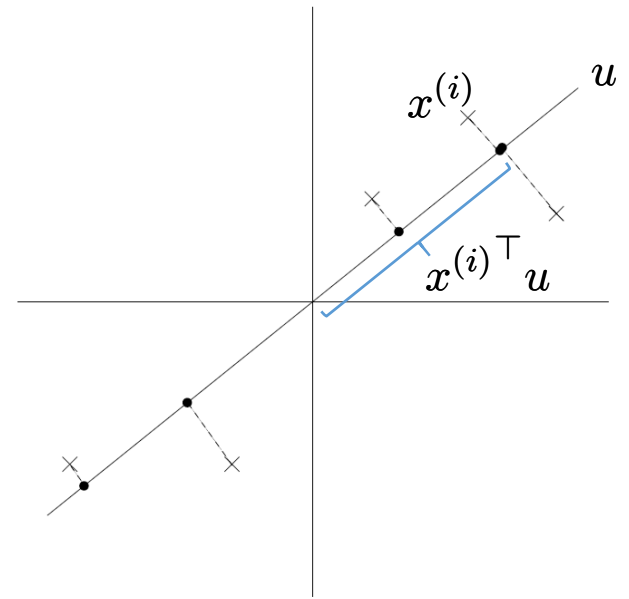
PCA Solution: Largest Eigenvalues

$$\begin{aligned} \max_u \quad & u^\top \Sigma u \\ \text{s.t.} \quad & \|u\| = 1 \end{aligned}$$

$$\Sigma = \frac{1}{m} \sum_{i=1}^m x^{(i)} x^{(i)\top}$$

- Find k principal components of the data is to find the k principal eigenvectors of Σ
 - i.e. the top- k eigenvectors with the largest eigenvalues
- Projected vector for $x^{(i)}$

$$y^{(i)} = \begin{bmatrix} u_1^\top x^{(i)} \\ u_2^\top x^{(i)} \\ \vdots \\ u_k^\top x^{(i)} \end{bmatrix} \in \mathbb{R}^k$$



Eigendecomposition Revisit

- For a semi-positive square matrix $\Sigma_{d \times d}$
 - suppose u to be its eigenvector ($\|u\| = 1$)
 - with the scalar eigenvalue w $\Sigma u = wu$
 - There are d eigenvectors-eigenvalue pairs (u_i, w_i)
 - These d eigenvectors are orthogonal, thus they form an orthonormal basis

$$\sum_{i=1}^d u_i u_i^\top = I$$

- Thus any vector v can be written as

$$v = \left(\sum_{i=1}^d u_i u_i^\top \right) v = \sum_{i=1}^d (u_i^\top v) u_i = \sum_{i=1}^d v_{(i)} u_i$$

$$U = [u_1, u_2, \dots, u_d]$$

- $\Sigma_{d \times d}$ can be written as

$$\Sigma = \sum_{i=1}^d u_i u_i^\top \Sigma = \sum_{i=1}^d w_i u_i u_i^\top = U W U^\top$$

$$W = \begin{bmatrix} w_1 & 0 & \cdots & 0 \\ 0 & w_2 & \cdots & 0 \\ \vdots & \vdots & \ddots & 0 \\ 0 & 0 & \cdots & w_d \end{bmatrix}$$

Eigendecomposition Revisit

- Given the data $X = \begin{bmatrix} x_1^\top \\ x_2^\top \\ \vdots \\ x_n^\top \end{bmatrix}$ and its covariance matrix $\Sigma = X^\top X$
(here we may drop m for simplicity)

- The variance in direction u_i is

$$\|Xu_i\|^2 = u_i^\top X^\top Xu_i = u_i^\top \Sigma u_i = u_i^\top w_i u_i = w_i$$

- The variance in any direction v is

$$\|Xv\|^2 = \left\| X \left(\sum_{i=1}^d v_{(i)} u_i \right) \right\|^2 = \sum_{ij} v_{(i)} u_i^\top \Sigma u_i v_{(j)} = \sum_{i=1}^d v_{(i)}^2 w_i$$

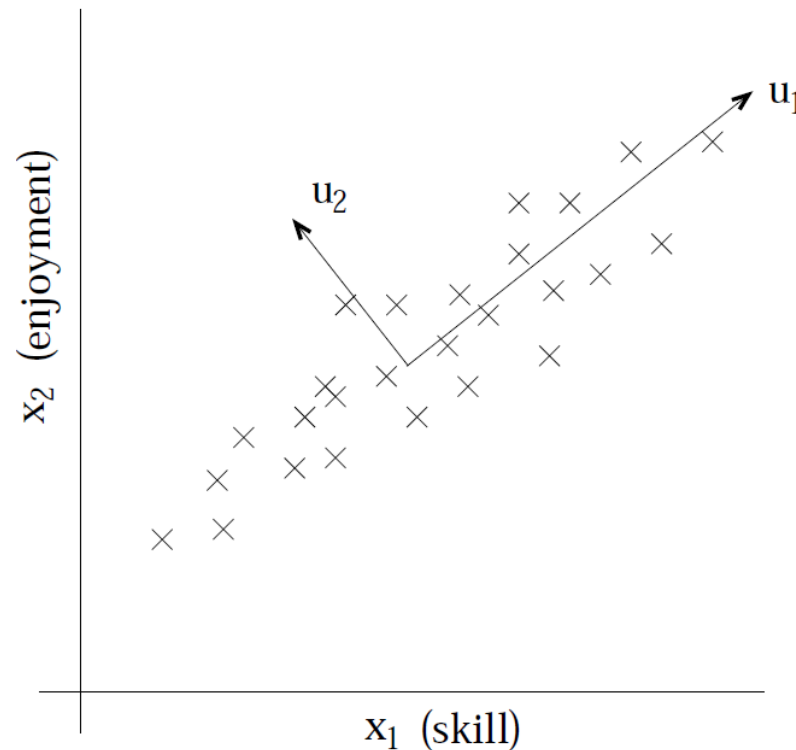
where $v_{(i)}$ is the projection length of v on u_i

- If $v^\top v = 1$, then $\arg \max_{\|v\|=1} \|Xv\|^2 = u_{(\max)}$

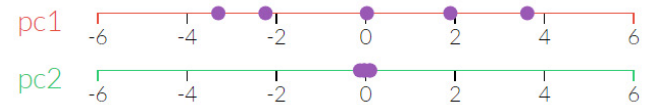
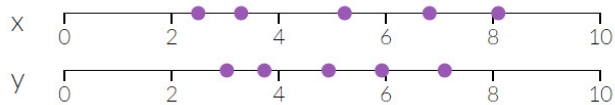
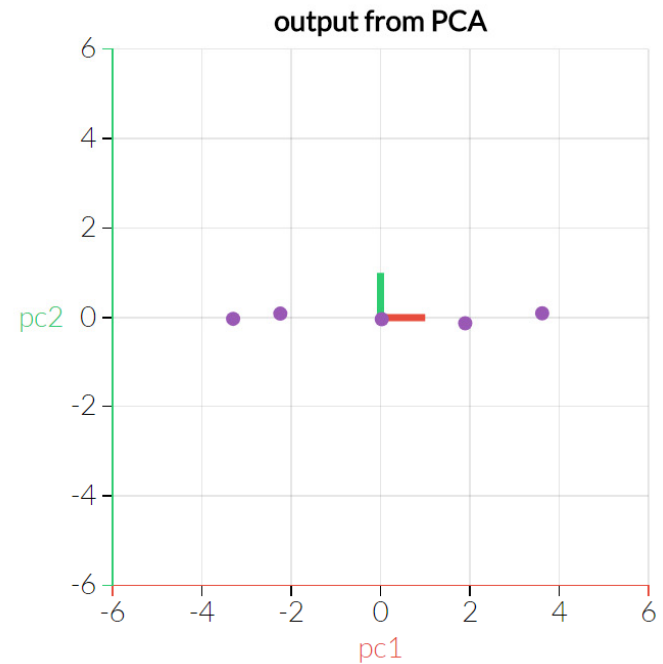
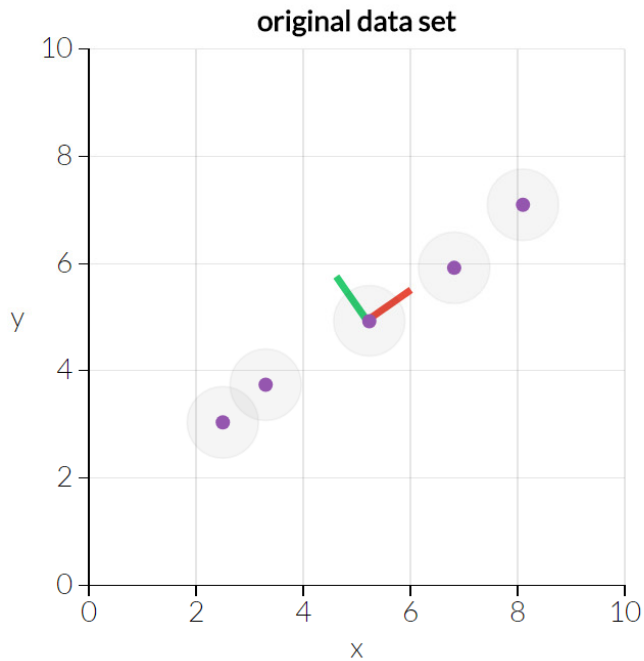
The direction of greatest variance is the eigenvector with the largest eigenvalue

PCA Discussion

- PCA can also be derived by picking the basis that minimizes the approximation error arising from projecting the data onto the k -dimensional subspace spanned by them.

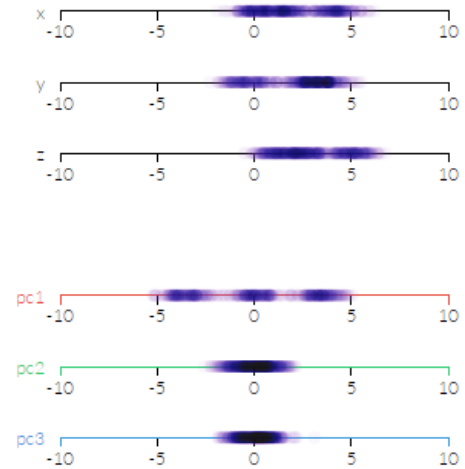
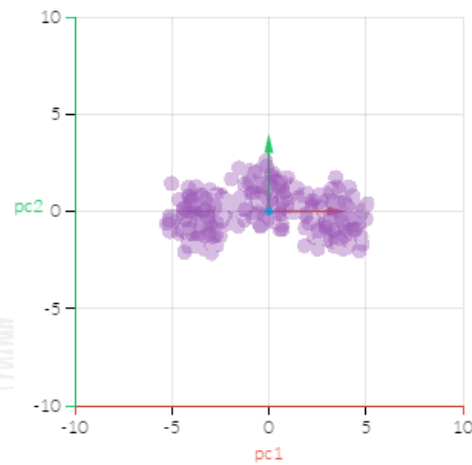
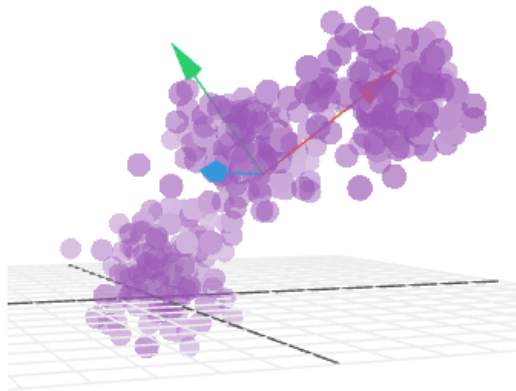


PCA Visualization



<http://setosa.io/ev/principal-component-analysis/>

PCA Visualization

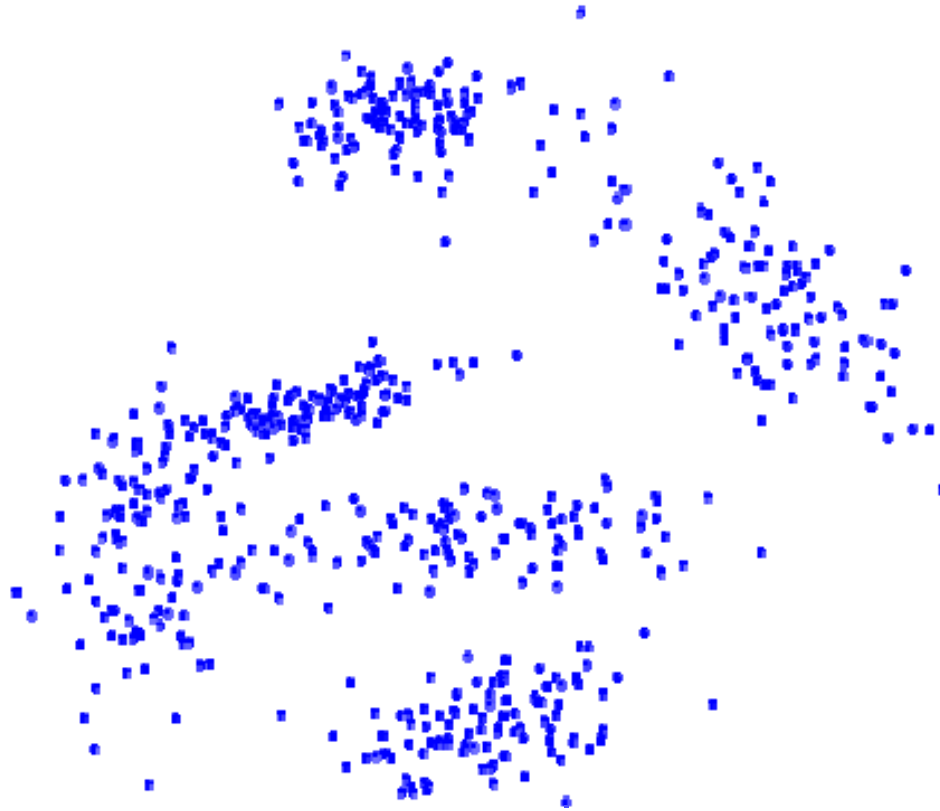


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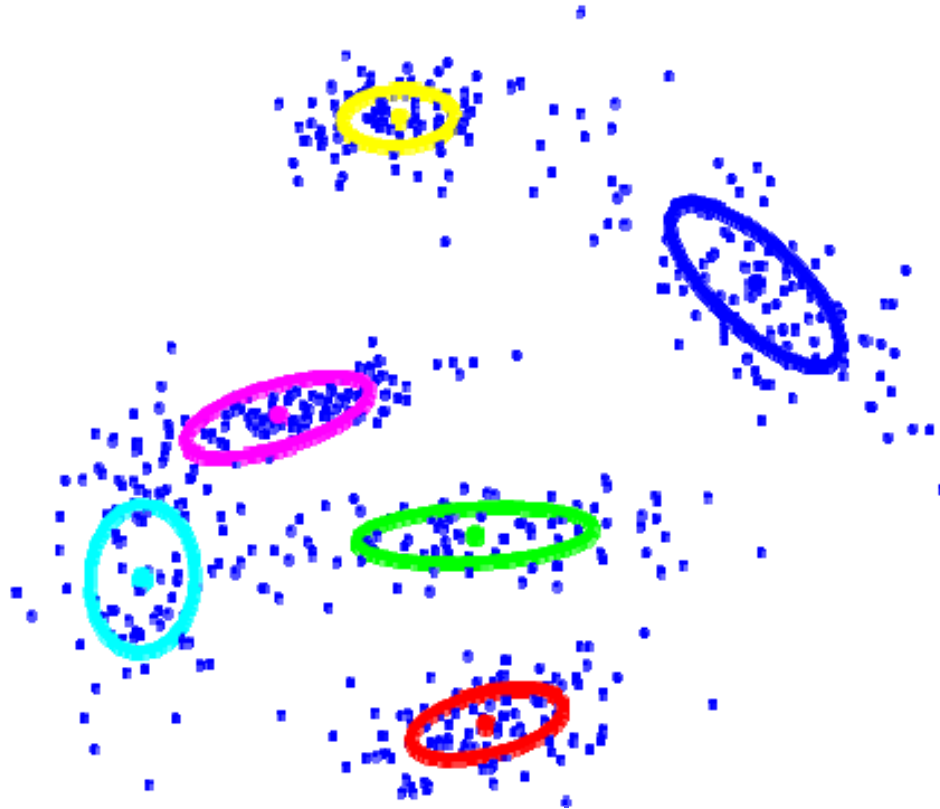
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Mixture Gaussian



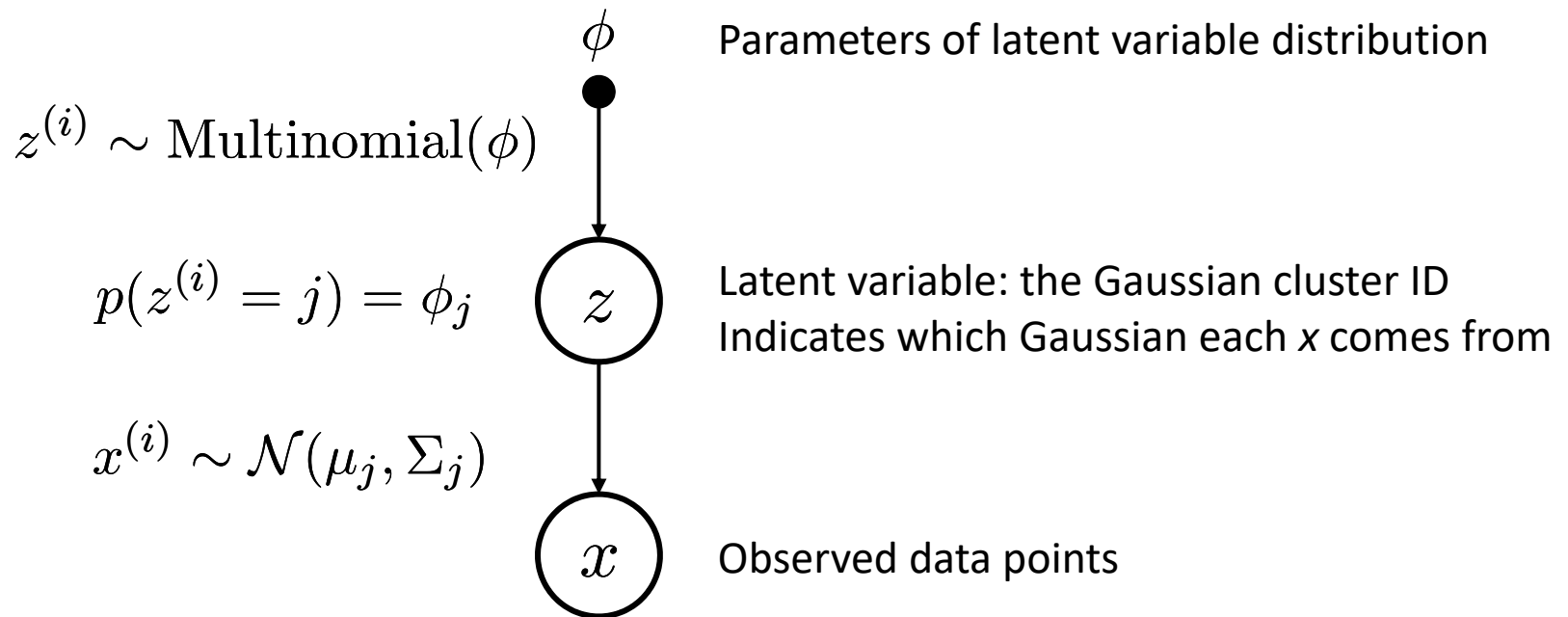
Mixture Gaussian



Graphic Model for Mixture Gaussian

- Given a training set $\{x^{(1)}, x^{(2)}, \dots, x^{(m)}\}$
- Model the data by specifying a joint distribution

$$p(x^{(i)}, z^{(i)}) = p(x^{(i)} | z^{(i)}) p(z^{(i)})$$



Data Likelihood

- We want to maximize

$$\begin{aligned}l(\phi, \mu, \Sigma) &= \sum_{i=1}^m \log p(x^{(i)}; \phi, \mu, \Sigma) \\&= \sum_{i=1}^m \log \sum_{z^{(i)}=1}^k p(x^{(i)}|z^{(i)}; \mu, \Sigma)p(z^{(i)}; \phi) \\&= \sum_{i=1}^m \log \sum_{j=1}^k \mathcal{N}(x^{(i)}|\mu_j, \Sigma_j)\phi_j\end{aligned}$$

- No closed form solution by simply setting

$$\frac{\partial l(\phi, \mu, \Sigma)}{\partial \phi} = 0 \quad \frac{\partial l(\phi, \mu, \Sigma)}{\partial \mu} = 0 \quad \frac{\partial l(\phi, \mu, \Sigma)}{\partial \Sigma} = 0$$

Data Likelihood Maximization

- For each data point $x^{(i)}$, latent variable $z^{(i)}$ indicates which Gaussian it comes from
- If we knew $z^{(i)}$, the data likelihood

$$\begin{aligned} l(\phi, \mu, \Sigma) &= \sum_{i=1}^m \log p(x^{(i)}; \phi, \mu, \Sigma) \\ &= \sum_{i=1}^m \log p(x^{(i)} | z^{(i)}; \mu, \Sigma) p(z^{(i)}; \phi) \\ &= \sum_{i=1}^m \log \mathcal{N}(x^{(i)} | \mu_{z^{(i)}}, \Sigma_{z^{(i)}}) + \log p(z^{(i)}; \phi) \end{aligned}$$

Data Likelihood Maximization

- Given $z^{(i)}$, maximize the data likelihood

$$\max_{\phi, \mu, \Sigma} l(\phi, \mu, \Sigma) = \max_{\phi, \mu, \Sigma} \sum_{i=1}^m \log \mathcal{N}(x^{(i)} | \mu_{z^{(i)}}, \Sigma_{z^{(i)}}) + \log p(z^{(i)}; \phi)$$

- It is easy to get the solution

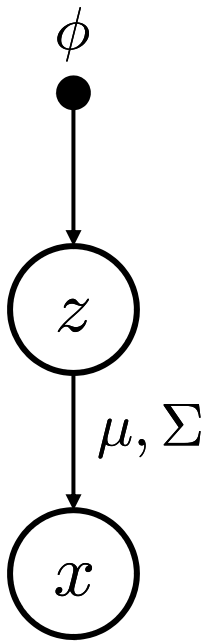
$$\phi_j = \frac{1}{m} \sum_{i=1}^m 1\{z^{(i)} = j\}$$

$$\mu_j = \frac{\sum_{i=1}^m 1\{z^{(i)} = j\} x^{(i)}}{\sum_{i=1}^m 1\{z^{(i)} = j\}}$$

$$\Sigma_j = \frac{\sum_{i=1}^m 1\{z^{(i)} = j\} (x^{(i)} - \mu_j)(x^{(i)} - \mu_j)^\top}{\sum_{i=1}^m 1\{z^{(i)} = j\}}$$

Latent Variable Inference

- Given the parameters μ, Σ, ϕ , it is not hard to infer the posterior of the latent variable $z^{(i)}$ for each instance



$$\begin{aligned} p(z^{(i)} = j | x^{(i)}; \phi, \mu, \Sigma) &= \frac{p(z^{(i)} = j, x^{(i)}; \phi, \mu, \Sigma)}{p(x^{(i)}; \phi, \mu, \Sigma)} \\ &= \frac{p(x^{(i)} | z^{(i)} = j; \mu, \Sigma) p(z^{(i)} = j; \phi)}{\sum_{l=1}^k p(x^{(i)} | z^{(i)} = l; \mu, \Sigma) p(z^{(i)} = l; \phi)} \end{aligned}$$

where

- The prior of $z^{(i)}$ is $p(z^{(i)} = j; \phi)$
- The likelihood is $p(x^{(i)} | z^{(i)} = j; \mu, \Sigma)$

Expectation Maximization Methods

- E-step: infer the posterior distribution of the latent variables given the model parameters
- M-step: tune parameters to maximize the data likelihood given the latent variable distribution
- EM methods
 - Iteratively execute E-step and M-step until convergence

EM Methods for Mixture Gaussians

- Mixture Gaussian example

Repeat until convergence: {

(E-step) For each i, j , set

$$w_j^{(i)} = p(z^{(i)} = j, x^{(i)}; \phi, \mu, \Sigma)$$

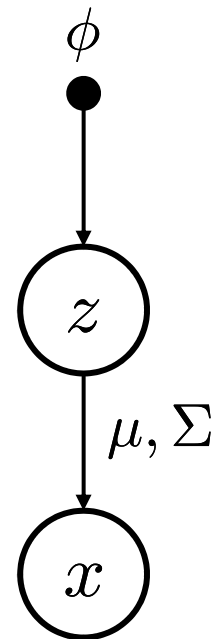
(M-step) Update the parameters

$$\phi_j = \frac{1}{m} \sum_{i=1}^m w_j^{(i)}$$

$$\mu_j = \frac{\sum_{i=1}^m w_j^{(i)} x^{(i)}}{\sum_{i=1}^m w_j^{(i)}}$$

$$\Sigma_j = \frac{\sum_{i=1}^m w_j^{(i)} (x^{(i)} - \mu_j)(x^{(i)} - \mu_j)^\top}{\sum_{i=1}^m w_j^{(i)}}$$

}



General EM Methods

- Claims:
 1. After each E-M step, the data likelihood will not decrease.
 2. The EM algorithm finds a (local) maximum of a latent variable model likelihood
- Now let's discuss the general EM methods and verify its effectiveness of improving data likelihood and its convergence

Jensen's Inequality

Theorem. Let f be a convex function, and let X be a random variable.

Then:

$$\mathbb{E}[f(X)] \geq f(\mathbb{E}[X])$$

- Moreover, if f is strictly convex, then

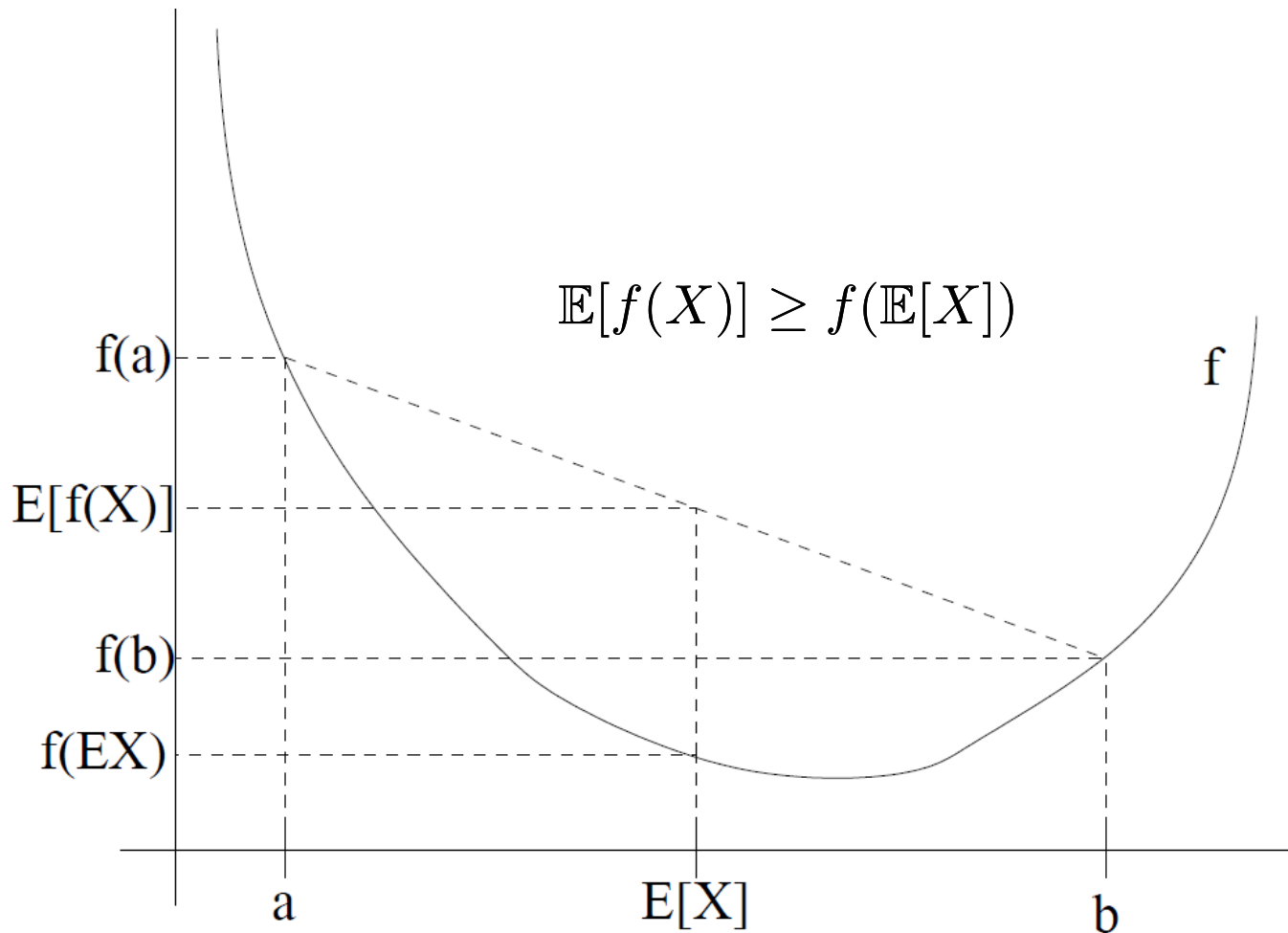
$$\mathbb{E}[f(X)] = f(\mathbb{E}[X])$$

holds true if and only if

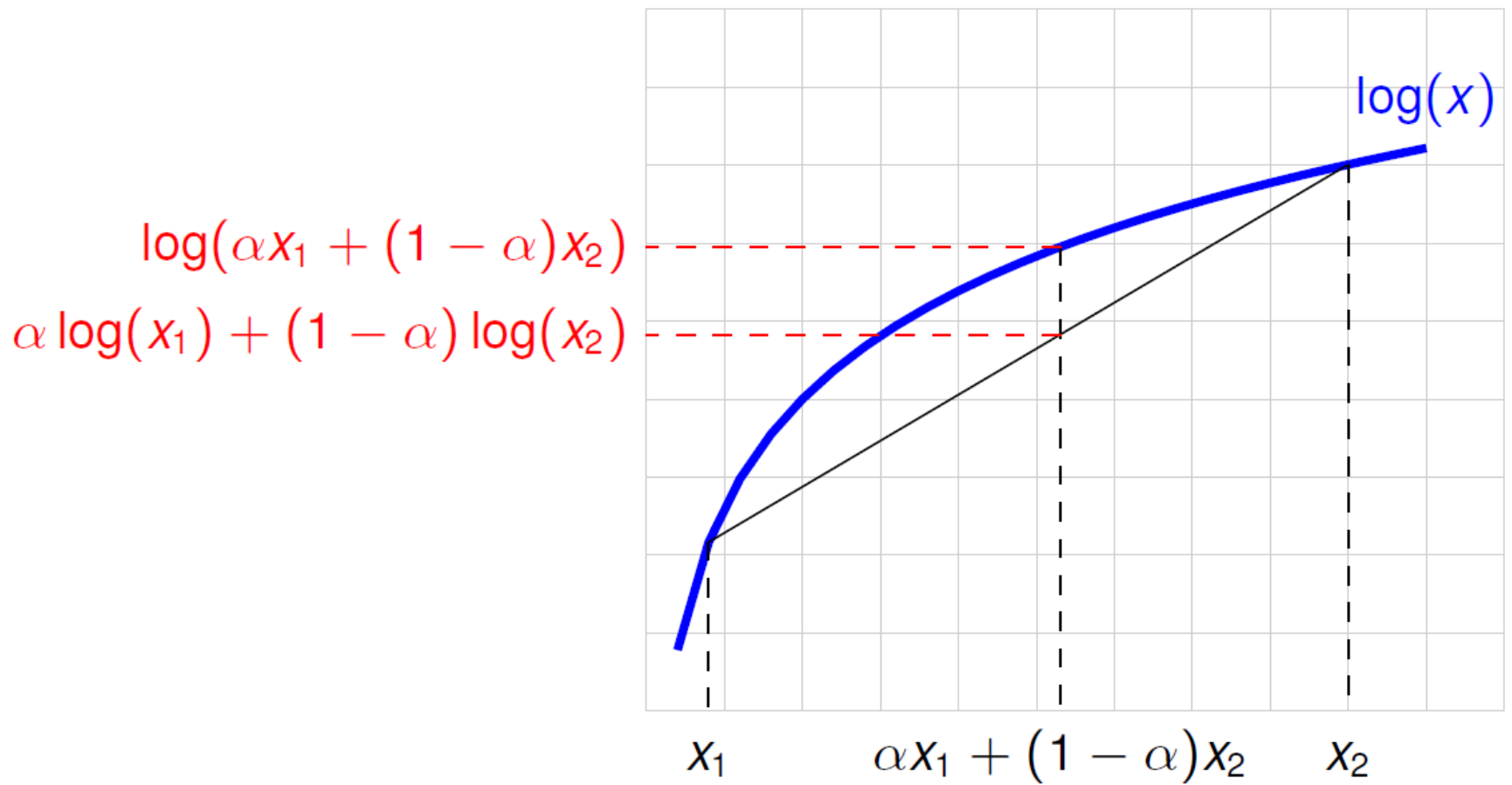
$$X = \mathbb{E}[X]$$

with probability 1 (i.e., if X is a constant).

Jensen's Inequality



Jensen's Inequality



General EM Methods: Problem

- Given the training dataset

$$D = \{x_i\}_{i=1,2,\dots,N}$$

let the machine learn the data underlying patterns

- Assume latent variables

$$z \rightarrow x$$

- We wish to fit the parameters of a model $p(x,z)$ to the data, where the log-likelihood is

$$\begin{aligned} l(\theta) &= \sum_{i=1}^N \log p(x; \theta) \\ &= \sum_{i=1}^N \log \sum_z p(x, z; \theta) \end{aligned}$$

General EM Methods: Problems

- EM methods solve the problems where
 - Explicitly find the maximum likelihood estimation (MLE) is hard

$$\theta^* = \arg \max_{\theta} \sum_{i=1}^N \log \sum_z p(x^{(i)}, z^{(i)}; \theta)$$

- But given $z^{(i)}$ observed, the MLE is easy

$$\theta^* = \arg \max_{\theta} \sum_{i=1}^N \log p(x^{(i)} | z^{(i)}; \theta)$$

- EM methods give an efficient solution for MLE, by iteratively doing
 - E-step: construct a (good) lower-bound of log-likelihood
 - M-step: optimize that lower-bound

General EM Methods: Lower Bound

- For each instance i , let q_i be some distribution of $z^{(i)}$

$$\sum_z q_i(z) = 1, \quad q_i(z) \geq 0$$

- Thus the data log-likelihood

$$\begin{aligned} l(\theta) &= \sum_{i=1}^N \log p(x^{(i)}; \theta) = \sum_{i=1}^N \log \sum_{z^{(i)}} p(x^{(i)}, z^{(i)}; \theta) \\ &= \sum_{i=1}^N \log \sum_{z^{(i)}} q_i(z^{(i)}) \frac{p(x^{(i)}, z^{(i)}; \theta)}{q_i(z^{(i)})} \\ &\geq \sum_{i=1}^N \sum_{z^{(i)}} q_i(z^{(i)}) \log \frac{p(x^{(i)}, z^{(i)}; \theta)}{q_i(z^{(i)})} \quad \text{Lower bound of } l(\vartheta) \end{aligned}$$

Jensen's inequality
-log(x) is a convex function

General EM Methods: Lower Bound

$$l(\theta) = \sum_{i=1}^N \log p(x^{(i)}; \theta) \geq \sum_{i=1}^N \sum_{z^{(i)}} q_i(z^{(i)}) \log \frac{p(x^{(i)}, z^{(i)}; \theta)}{q_i(z^{(i)})}$$

- Then what $q_i(z)$ should we choose?

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holds true if and only if

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with probability 1 (i.e., if X is a constant).

General EM Methods: Lower Bound

$$l(\theta) = \sum_{i=1}^N \log p(x^{(i)}; \theta) \geq \sum_{i=1}^N \sum_{z^{(i)}} q_i(z^{(i)}) \log \frac{p(x^{(i)}, z^{(i)}; \theta)}{q_i(z^{(i)})}$$

- Then what $q_i(z)$ should we choose?
- In order to make above inequality tight (to hold with equality), it is sufficient that

$$p(x^{(i)}, z^{(i)}; \theta) = q_i(z^{(i)}) \cdot c$$

- We can derive

$$\log p(x^{(i)}; \theta) = \log \sum_{z^{(i)}} p(x^{(i)}, z^{(i)}; \theta) = \log \sum_{z^{(i)}} q(z^{(i)})c = \sum_{z^{(i)}} q_i(z^{(i)}) \log \frac{p(x^{(i)}, z^{(i)}; \theta)}{q_i(z^{(i)})}$$

- As such $q_i(z)$ is written as the posterior distribution

$$q_i(z^{(i)}) = \frac{p(x^{(i)}, z^{(i)}; \theta)}{\sum_z p(x^{(i)}, z; \theta)} = \frac{p(x^{(i)}, z^{(i)}; \theta)}{p(x^{(i)}; \theta)} = p(z^{(i)} | x^{(i)}; \theta)$$

General EM Methods

Repeat until convergence: {

(E-step) For each i , set

$$q_i(z^{(i)}) = p(z^{(i)} | x^{(i)}; \theta)$$

(M-step) Update the parameters

$$\theta = \arg \max_{\theta} \sum_{i=1}^N \sum_{z^{(i)}} q_i(z^{(i)}) \log \frac{p(x^{(i)}, z^{(i)}; \theta)}{q_i(z^{(i)})}$$

}

Convergence of EM

- Denote $\vartheta^{(t)}$ and $\vartheta^{(t+1)}$ as the parameters of two successive iterations of EM, we prove that

$$l(\theta^{(t)}) \leq l(\theta^{(t+1)})$$

which shows EM always monotonically improves the log-likelihood, thus ensures EM will at least converge to a local optimum.

Proof of EM Convergence

- Start from $\vartheta^{(t)}$, we choose the posterior of latent variable

$$q_i^{(t)}(z^{(i)}) = p(z^{(i)} | x^{(i)}; \theta^{(t)})$$

- This choice ensures the Jensen's inequality holds with equality

$$l(\theta^{(t)}) = \sum_{i=1}^N \log \sum_{z^{(i)}} q_i^{(t)}(z^{(i)}) \frac{p(x^{(i)}, z^{(i)}; \theta^{(t)})}{q_i^{(t)}(z^{(i)})} = \sum_{i=1}^N \sum_{z^{(i)}} q_i(z^{(i)}) \log \frac{p(x^{(i)}, z^{(i)}; \theta^{(t)})}{q_i^{(t)}(z^{(i)})}$$

- Then the parameters $\vartheta^{(t+1)}$ are then obtained by maximizing the right hand side of above equation

$$\begin{aligned} \text{• Thus } l(\theta^{(t+1)}) &\geq \sum_{i=1}^N \sum_{z^{(i)}} q_i^{(t)}(z^{(i)}) \log \frac{p(x^{(i)}, z^{(i)}; \theta^{(t+1)})}{q_i^{(t)}(z^{(i)})} && \text{[lower bound]} \\ &\geq \sum_{i=1}^N \sum_{z^{(i)}} q_i^{(t)}(z^{(i)}) \log \frac{p(x^{(i)}, z^{(i)}; \theta^{(t)})}{q_i^{(t)}(z^{(i)})} && \text{[parameter optimization]} \\ &= l(\theta^{(t)}) \end{aligned}$$

Remark of EM Convergence

- If we define

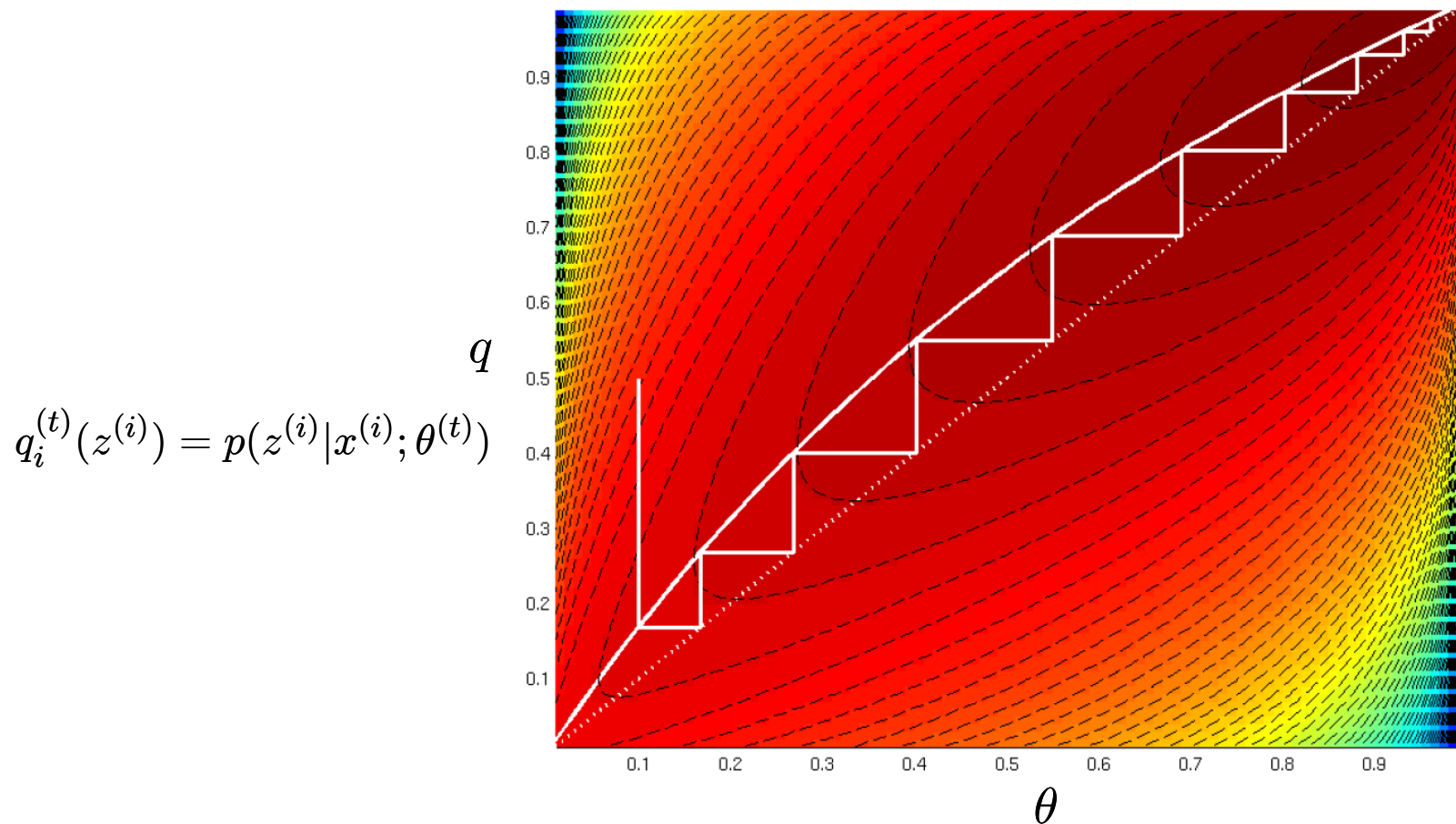
$$J(q, \theta) = \sum_{i=1}^N \sum_{z^{(i)}} q_i(z^{(i)}) \log \frac{p(x^{(i)}, z^{(i)}; \theta)}{q_i(z^{(i)})}$$

Then we know

$$l(\theta) \geq J(q, \theta)$$

- EM can also be viewed as a coordinate ascent on J
 - E-step maximizes it w.r.t. q
 - M-step maximizes it w.r.t. ϑ

Coordinate Ascent in EM



$$\theta^{(t+1)} = \arg \max_{\theta} \sum_{i=1}^N \sum_{z^{(i)}} q_i^{(t)}(z^{(i)}) \log \frac{p(x^{(i)}, z^{(i)}; \theta)}{q_i^{(t)}(z^{(i)})}$$