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CS540 Introduction to Artificial Intelligence Lecture 16

Young Wu

Based on lecture slides by Jerry Zhu, Yingyu Liang, and Charles Dyer

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High Dimensional Data Motivation

- High dimensional data are training set with a lot of features.
- Document classification.
- Ø MEG brain imaging.
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Low Dimension Representation

- Unsupervised learning techniques are used to find low dimensional representation.
- Visualization.
- efficient storage.
- Better generalization.
- Olise removal.

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Dimension Reduction

- Rotate the axes so that they capture the directions of the greatest variability of data.
- The new axes (orthogonal directions) are principal components.

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Principal Component Analysis Description

- Find the direction of the greatest variability in data, call it $u_{1.}$
- Find the next direction orthogonal to *u*₁ of the greatest variability, call it *u*₂.
- Repeat until there are $u_1, u_2, ..., u_K$.

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Orthogonal Directions Definition

- In Euclidean space (L₂ norm), a unit vector u_k has length 1. $\|u_k\|_2 = u_k^T u_k = 1$
- Two vectors u_k, u_{k'} are orthogonal (or uncorrelated) if the dot product is 0.

$$u_k \cdot u_{k'} = u_k^T u_{k'} = 0$$

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Projection Definition

• The projection of x_i onto a unit vector u_k is the vector in the direction of u_k that is the closest to x_i .

proj
$$_{u_k}x_i = \left(\frac{u_k^T x_i}{u_k^T u_k}\right) u_k = u_k^T x_i u_k$$

• The length of the projection of x_i onto a unit vector u_k is $u_k^T x_i$.

$$\left\| \operatorname{proj}_{u_k} x_i \right\|_2 = u_k^T x_i$$

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Variance Definition

• The sample variance of a data set {*x*₁, *x*₂, ..., *x_n*} is the sum of the squared distance from the mean.

$$\begin{split} X &= \begin{bmatrix} x_1 \\ x_2 \\ \dots \\ x_n \end{bmatrix} \\ \hat{\mu} &= \frac{1}{n} \sum_{i=1}^n x_i \\ \hat{\Sigma} &= \frac{1}{n-1} \sum_{i=1}^n \left(x_i - \hat{\mu} \right) \left(x_i - \hat{\mu} \right)^T \end{split}$$

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Normalization Definition

• Normalize the data by subtracting the mean, then the variance expression can be simplified.

$$\hat{\Sigma} = \frac{1}{n-1} \sum_{i=1}^{n} x_i x_i^T = \frac{1}{n-1} X^T X$$

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Covariance Matrix

• $\hat{\Sigma}$ is an $m \times m$ matrix and it is usually called the sample covariance matrix. The diagonal elements are variances in each dimension.

$$\hat{\sigma}_j^2 = \hat{\Sigma}_{jj} = rac{1}{n-1}\sum_{i=1}^n x_{ij}^2$$

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Projected Variance

Note that x_{ij} = e_j^Tx_i, where e_j is the vector of 0 except it is 1 in coordinate j.

$$\hat{\sigma}_j^2 = e_j^T \hat{\Sigma} e_j = \frac{1}{n-1} e_j^T X^T X e_j$$
$$= \frac{1}{n-1} \sum_{i=1}^n \left(e_j^T x_i \right)^2$$

• The variance of the normalized x_i projected onto direction u_k has a similar expression.

$$u_k^T \hat{\Sigma} u_k = \frac{1}{n-1} u_k^T X^T X u_k$$
$$= \frac{1}{n-1} \sum_{i=1}^n \left(u_k^T x_i \right)^2$$

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Maximum Variance Directions Definition

• The goal is to find the direction that maximizes the projected variance.

$$\max_{u_k} u_k^T \hat{\Sigma} u_k \text{ such that } u_k^T u_k = 1$$
$$\Rightarrow \max_{u_k} u_k^T \hat{\Sigma} u_k - \lambda u_k^T u_k$$
$$\Rightarrow \hat{\Sigma} u_k = \lambda u_k$$

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Eigenvalue Definition

• The λ represents the projected variance.

$$u_k^T \hat{\Sigma} u_k = u_k^T \lambda u_k = \lambda$$

The larger the variance, the larger the variability in direction u_k. There are m eigenvalues for a symmetric positive semidefinite matrix (for example, X^TX is always symmetric PSD). Order the eigenvectors u_k by the size of their corresponding eigenvalues λ_k.

$$\lambda_1 \geqslant \lambda_2 \geqslant \dots \geqslant \lambda_m$$

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Eigenvalue Algorithm

• Solving eigenvalue using the definition (characteristic polynomial) is computationally inefficient.

$$\left(\hat{\Sigma} - \lambda_k I\right) u_k = 0 \Rightarrow \det \left(\hat{\Sigma} - \lambda_k I\right) = 0$$

• There are many fast eigenvalue algorithms that computes the spectral (eigen) decomposition for real symmetric matrices. Columns of *Q* are unit eigenvectors and diagonal elements of *D* are eigenvalues.

$$\hat{\Sigma} = PDP^{-1}, D$$
 is diagonal
= QDQ^T , if Q is orthogonal, i.e. $Q^TQ = I$

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Principal Component Analysis Algorithm

- Input: instances: {x_i}ⁿ_{i=1}, the number of dimensions after reduction K < m.
- Output: K principal components.
- Find the largest K eigenvalues $\lambda_1 \ge \lambda_2 \ge ... \ge \lambda_K$.
- Return the corresponding unit orthogonal eigenvectors $u_1, u_2...u_K$.

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Number of Dimensions

- There are a few ways to choose the number of principal components *K*.
- K can be selected given prior knowledge or requirement.
- K can be the number of non-zero eigenvalues.
- *K* can be the number of eigenvalues that are large (larger than some threshold).

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Reduced Feature Space

- The original feature space is m dimensional. $(x_{i1}, x_{i2}, ..., x_{im})^T$
- The new feature space is K dimensional. $\left(u_1^T x_i, u_2^T x_i, ..., u_K^T x_i\right)^T$
- Other supervised learning algorithms can be applied on the new features.

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Reconstruction Error

• Reconstruction error is the squared error (distance) between the original data and its projection onto u_k .

$$\left\|x_{i}-\left(u_{k}^{T}x_{i}\right)u_{k}\right\|^{2}$$

• Finding the variance maximizing directions is the same as finding the reconstruction error minimizing directions.

$$\frac{1}{n}\sum_{i=1}^{n}\left\|x_{i}-\left(u_{k}^{T}x_{i}\right)u_{k}\right\|^{2}$$

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Reconstruction Error Diagram

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Eigenface Discussion

- Eigenfaces are eigenvectors of face images (pixel intensities or HOG features).
- Every face can be written as a linear combination of eigenfaces. The coefficients determine specific faces.

$$x_i = \sum_{k=1}^m \left(u_k^T x_i \right) u_k \approx \sum_{k=1}^K \left(u_k^T x_i \right) u_k$$

• Eigenfaces and SVM can be combined to detect or recognize faces.

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Autoencoder Discussion

- A multi-layer neural network with the same input and output $y_i = x_i$ is called an autoencoder.
- The hidden layers have fewer units than the dimension of the input *m*.
- The hidden units form an encoding of the input with reduced dimensionality.

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Kernel PCA Discussion

• A kernel can be applied before finding the principal components.

$$\hat{\Sigma} = \frac{1}{n-1} \sum_{i=1}^{n} \varphi(x_i) \varphi(x_i)^{T}$$

- The principal components can be found without explicitly computing $\varphi(x_i)$, similar to the kernel trick for support vector machines.
- Kernel PCA is a non-linear dimensionality reduction method.

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T-Distributed Stochastic Neighbor Embedding

- t-distributed stochastic neighbor embedding is another non-linear dimensionality reduction method used mainly for visualization.
- Points in high dimensional spaces are embedded in 2 or 3-dimensional spaces to preserve the distance (neighbor) relationship between points.

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Embedding Diagram