

# Introduction to Machine Learning

## Principal Component Analysis

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## Principal Components Analysis

- Introduction to PCA

- Principle of Maximal Variance

- Defining Principal Components

- Dimensionality Reduction Using PCA

- PCA Algorithm

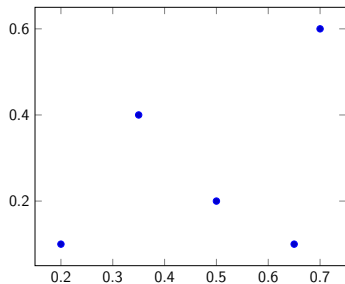
- Recovering Original Data

- Eigen Faces

## Kernel PCA

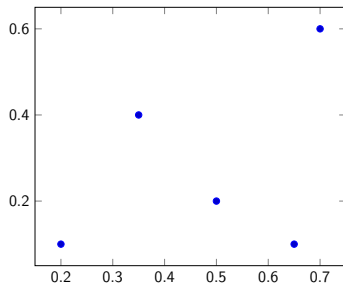
# Introduction to PCA

- ▶ Consider the following data points



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- ▶ *Embed* these points in 1 dimension
- ▶ What is the best way?
  - ▶ **Along the direction of the maximum variance**
  - ▶ Why?

# Why Maximal Variance?

- ▶ Least loss of information
- ▶ Best capture the “spread”

# Why Maximal Variance?

- ▶ Least loss of information
- ▶ Best capture the “spread”
- ▶ What is the direction of maximal variance?
- ▶ Given any direction ( $\hat{\mathbf{u}}$ ), the projection of  $\mathbf{x}$  on  $\hat{\mathbf{u}}$  is given by:

$$\mathbf{x}_i^\top \hat{\mathbf{u}}$$

- ▶ Direction of maximal variance can be obtained by maximizing

$$\begin{aligned} \frac{1}{N} \sum_{i=1}^N (\mathbf{x}_i^\top \hat{\mathbf{u}})^2 &= \frac{1}{N} \sum_{i=1}^N \hat{\mathbf{u}}^\top \mathbf{x}_i \mathbf{x}_i^\top \hat{\mathbf{u}} \\ &= \hat{\mathbf{u}}^\top \left( \frac{1}{N} \sum_{i=1}^N \mathbf{x}_i \mathbf{x}_i^\top \right) \hat{\mathbf{u}} \end{aligned}$$

# Finding Direction of Maximal Variance

- ▶ Find:

$$\max_{\hat{\mathbf{u}}: \hat{\mathbf{u}}^T \hat{\mathbf{u}} = 1} \hat{\mathbf{u}}^T \mathbf{S} \hat{\mathbf{u}}$$

where:

$$\mathbf{S} = \frac{1}{N} \sum_{i=1}^N \mathbf{x}_i \mathbf{x}_i^T$$

- ▶  $\mathbf{S}$  is the covariance matrix of the mean-centered data

# Defining Principal Components

- ▶ First PC: Eigen-vector of the (sample) covariance matrix with largest eigen-value
- ▶ Second PC?



# Defining Principal Components

- ▶ First PC: Eigen-vector of the (sample) covariance matrix with largest eigen-value
- ▶ Second PC?
- ▶ Eigen-vector with next largest value
- ▶ Variance of each PC is given by  $\lambda_i$
- ▶ Variance captured by first  $L$  PC ( $1 \leq L \leq D$ )

$$\frac{\sum_{i=1}^L \lambda_i}{\sum_{i=1}^D \lambda_i} \times 100$$

- ▶ What are eigen vectors and values?

$$\mathbf{A}\mathbf{v} = \lambda\mathbf{v}$$

$\mathbf{v}$  is eigen vector and  $\lambda$  is eigen-value for the **square matrix  $\mathbf{A}$**

- ▶ Geometric interpretation?

# Dimensionality Reduction Using PCA

- ▶ Consider first  $L$  eigen values and eigen vectors
- ▶ Let  $\mathbf{W}$  denote the  $D \times L$  matrix with first  $L$  eigen vectors in the columns (sorted by  $\lambda$ 's)
- ▶ PC score matrix

$$\mathbf{Z} = \mathbf{XW}$$

- ▶ Each input vector ( $D \times 1$ ) is replaced by a shorter  $L \times 1$  vector

1. Center  $\mathbf{X}$

$$\mathbf{X} = \mathbf{X} - \hat{\boldsymbol{\mu}}$$

2. Compute sample covariance matrix:

$$\mathbf{S} = \frac{1}{N-1} \mathbf{X}^T \mathbf{X}$$

3. Find eigen vectors and eigen values for  $\mathbf{S}$
4.  $\mathbf{W}$  consists of first  $L$  eigen vectors as columns
  - ▶ Ordered by decreasing eigen-values
  - ▶  $\mathbf{W}$  is  $D \times L$
5. Let  $\mathbf{Z} = \mathbf{XW}$
6. Each row in  $\mathbf{Z}$  (or  $\mathbf{z}_i^T$ ) is the lower dimensional embedding of  $\mathbf{x}_i$

# Recovering Original Data

- ▶ Using  $\mathbf{W}$  and  $\mathbf{z}_i$

$$\hat{\mathbf{x}}_i = \mathbf{W}\mathbf{z}_i$$

- ▶ **Average Reconstruction Error**

$$J(\mathbf{W}, \mathbf{Z}) = \frac{1}{N} \sum_{i=1}^N \|\mathbf{x}_i - \hat{\mathbf{x}}_i\|^2$$

## Theorem (Classical PCA Theorem)

*Among all possible orthonormal sets of  $L$  basis vectors, PCA gives the solution which has the minimum reconstruction error.*

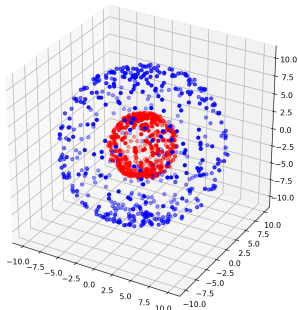
- ▶ Optimal “embedding” in  $L$  dimensional space is given by  $\mathbf{z}_i = \mathbf{W}^T \mathbf{x}_i$

## EigenFaces [2]

- ▶ **Input:** A set of images (of faces)
- ▶ **Task:** Identify if a new image is a face or not.

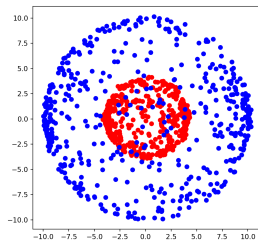
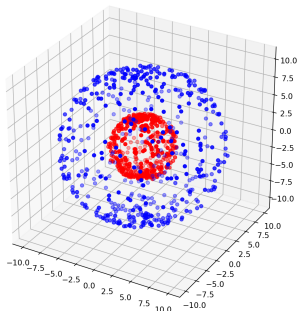
# Issues with PCA

- ▶ A linear transformation of the data
- ▶ Might not work everytime



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# Kernel PCA - Applying the kernel trick [1] I

- ▶ Assume a non-linear transformation of input:

$$\mathbf{x}_i \Rightarrow \Phi(\mathbf{x}_i), \Phi(\mathbf{x}_i) \in \mathbb{R}^M$$

- ▶ We will assume that the new data is mean centered:

$$\frac{1}{N} \sum_{i=1}^N \Phi(\mathbf{x}_i) = 0$$

- ▶ The covariance matrix of the projected data,  $\mathbf{C}$ :

$$\mathbf{C} = \frac{1}{N} \sum_{i=1}^N \Phi(\mathbf{x}_i) \Phi(\mathbf{x}_i)^\top$$

- ▶ The eigen vectors and eigen values of  $\mathbf{C}$  are given by:

$$\mathbf{C}\mathbf{v}_k = \lambda_k \mathbf{v}_k$$

where  $k = 1, 2, \dots, M$



# Kernel PCA - Applying the kernel trick [1] II

- ▶ Substituting the expression for  $\mathbf{C}$ :

$$\frac{1}{N} \sum_{i=1}^N \Phi(\mathbf{x}_i) \Phi(\mathbf{x}_i)^\top \mathbf{v}_k = \lambda_k \mathbf{v}_k$$

- ▶ Rearranging and assuming  $a_{ki} = \frac{\Phi(\mathbf{x}_i)^\top \mathbf{v}_k}{N \lambda_k}$

$$\mathbf{v}_k = \sum_{i=1}^N a_{ki} \Phi(\mathbf{x}_i)$$

- ▶ Substituting this back in the above equation

$$\frac{1}{N} \sum_{i=1}^N \Phi(\mathbf{x}_i) \Phi(\mathbf{x}_i)^\top \sum_{j=1}^N a_{kj} \Phi(\mathbf{x}_j) = \lambda_k \sum_{i=1}^N a_{ki} \Phi(\mathbf{x}_i)$$

Note the subscript  $j$  in the second summation on the left hand side.

# Kernel PCA - Applying the kernel trick [1] III

- ▶ Now multiplying both sides with  $\Phi(\mathbf{x}_i)^\top$ :

$$\frac{1}{N} \Phi(\mathbf{x}_i)^\top \sum_{i=1}^N \Phi(\mathbf{x}_i) \Phi(\mathbf{x}_i)^\top \sum_{j=1}^N a_{kj} \Phi(\mathbf{x}_j) = \lambda_k \Phi(\mathbf{x}_i)^\top \sum_{i=1}^N a_{ki} \Phi(\mathbf{x}_i)$$

which is the same as:

$$\sum_{i=1}^N \underbrace{\Phi(\mathbf{x}_i)^\top \Phi(\mathbf{x}_i)} \sum_{j=1}^N a_{kj} \underbrace{\Phi(\mathbf{x}_i)^\top \Phi(\mathbf{x}_j)} = N \lambda_k \sum_{i=1}^N \underbrace{a_{ki} \Phi(\mathbf{x}_i)^\top \Phi(\mathbf{x}_i)}$$

- ▶ Let  $k()$  be a function, such that:  $k(\mathbf{x}_i, \mathbf{x}_j) = \Phi(\mathbf{x}_i)^\top \Phi(\mathbf{x}_j)$
- ▶ The above expression can be written as:

$$\sum_{i=1}^N k(\mathbf{x}_i, \mathbf{x}_i) \sum_{j=1}^N a_{kj} k(\mathbf{x}_i, \mathbf{x}_j) = N \lambda_k \sum_{i=1}^N k(\mathbf{x}_i, \mathbf{x}_i)$$

# Kernel PCA - Applying the kernel trick [1] IV

- ▶ Consider the  $N \times 1$  vector,  $\mathbf{a}_k$  and  $N \times N$  matrix,  $\mathbf{K}$ , such that

$$\mathbf{a}_k = \begin{bmatrix} a_{k1} \\ a_{k2} \\ \vdots \\ a_{kN} \end{bmatrix}$$

and,

$$\mathbf{K}[i][j] = k(\mathbf{x}_i, \mathbf{x}_j)$$

- ▶ The above expression can be written, using matrix notation, as:

$$\mathbf{K}^2 \mathbf{a}_k = \lambda_k N \mathbf{K} \mathbf{a}_k$$

- ▶ To solve for  $\mathbf{a}_k$ , we can solve the following:

$$\mathbf{K} \mathbf{a}_k = \lambda_k N \mathbf{a}_k$$

# Projecting data using Kernel PCA

- ▶ For a new data instance,  $\mathbf{x}^*$ , the  $k^{\text{th}}$  entry of the corresponding  $\mathbf{z}^*$  will be:

$$\begin{aligned}z_k^* &= \Phi(\mathbf{x}^*)^\top \mathbf{v}_k \\&= \Phi(\mathbf{x}^*)^\top \sum_{i=1}^N a_{ki} \Phi(\mathbf{x}_i) \\&= \sum_{i=1}^N a_{ki} \Phi(\mathbf{x}^*)^\top \Phi(\mathbf{x}_i) \\&= \sum_{i=1}^N a_{ki} k(\mathbf{x}^*, \mathbf{x}_i)\end{aligned}$$

# Centering the projected data

- ▶ How do we ensure that the projected new features have a zero mean, without doing the actual projection?
- ▶ Use the **Gram matrix**:

$$\tilde{\mathbf{K}} = \mathbf{K} - \mathbf{1}_N \mathbf{K} - \mathbf{K} \mathbf{1}_N + \mathbf{1}_N \mathbf{K} \mathbf{1}_N$$

where  $\mathbf{1}_N$  is a  $N \times N$  matrix with all entries equal to  $\frac{1}{N}$

# Kernel PCA Algorithm

1. Given the data set  $\mathbf{X}$  and a kernel function  $k()$ , construct the kernel matrix  $\mathbf{K}$
2. Compute the Gram matrix  $\tilde{\mathbf{K}}$
3. Find eigen-vectors of  $\tilde{\mathbf{K}}$
4. Use top  $M$  eigen-vectors to project a new data instance,  $\mathbf{x}^*$  to the corresponding  $\mathbf{z}^*$

# Kernel PCA - Final Thoughts

- ▶ Very sensitive to the kernel choice and kernel parameters
- ▶ Slow ( $O(N^3)$ )
- ▶ Recovering original data is not straightforward as linear PCA



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