

# PRINCIPAL COMPONENT ANALYSIS

## Unsupervised Learning

**Lê Hồng Phương**

Data Science Laboratory

Vietnam National University, Hanoi

*<phuonglh@hus.edu.vn>*

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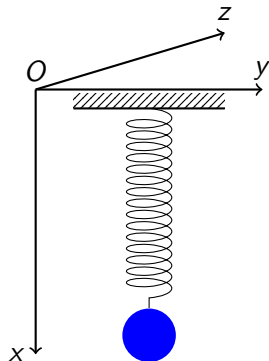
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# Overview

- Principal Component Analysis (PCA) is a simple, non-parametric method of *extracting relevant information* from noisy datasets.
- PCA provides a method to reduce a complex dataset to a lower dimension to reveal hidden properties/structures of the dataset.
- PCA is widely used in many forms of analysis: neuroscience, computer graphics, natural language processing, *etc.*

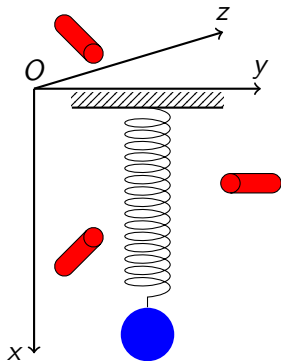
# Motivation: A Toy Example

- We are studying the motion of an ideal spring.
- This system consists of a ball of mass  $m$  attached to a massless, frictionless spring.
- The ball is released a small distance away from equilibrium (the spring is stretched).
- The spring oscillates indefinitely along the  $x$ -axis about its equilibrium at some frequency.



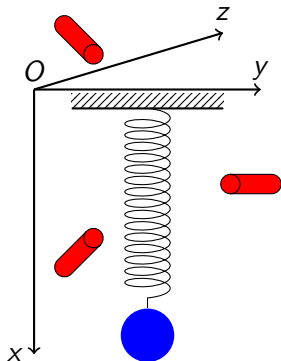
# Motivation: A Toy Example

- This is a standard problem in physics, the motion along the  $x$ -axis is solved by an explicit function of time.
  - The underlying dynamics can be expressed as a function of a single variable  $x$ .
- However, suppose that we do not know which axes and dimensions are important to measure.
- Thus, we decide to measure the ball's position in a three-dimensional space.
  - We place 3 cameras around our system of interest.



# Motivation: A Toy Example

- At 200 Hz, each camera records an image indicating a 2-dimensional position of the ball (a projection).
- Unfortunately, we do not even know what are the real “x”, “y” and “z”, so we choose 3 camera axes  $\{\vec{a}, \vec{b}, \vec{c}\}$  at some arbitrary angles w.r.t. the system.
- The angles between our measurements might not even be  $90^\circ$ !
- Now, we record the cameras for 2 minutes.
- How do we get from this dataset to a simple equation of  $x$ ?



# Motivation: A Toy Example

## Some common problems:

- We sometimes record more dimensions than we actually need.
- We have to deal with noise (e.g. air, imperfect cameras, friction. . .)

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**Goal:** PCA computes the most meaningful *basis* to re-express a noisy, garbled dataset.

- The new basis will filter out the noise and reveal hidden dynamics (e.g. the dynamics are along the x-axis).



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# A Naive Basis

A naive and simple choice of a basis is the identity matrix:

$$\mathbf{I} = \begin{pmatrix} \mathbf{e}_1 \\ \mathbf{e}_2 \\ \vdots \\ \mathbf{e}_D \end{pmatrix} = \begin{pmatrix} 1 & 0 & \cdots & 0 \\ 0 & 1 & \cdots & 0 \\ \vdots & \vdots & \ddots & \vdots \\ 0 & 0 & \cdots & 1 \end{pmatrix}$$

- Each row is a basis vector  $\mathbf{e}_i$  with  $D$  components.
- Every data point is a vector that lies in a  $D$ -dimensional vector space spanned by an orthonormal basis.
- *All vectors in this space are a linear combination of this set of unit length basis vectors.*

# Change of Basis

**PCA question:** *Is there another basis, which is a linear combination of the original basis, that best re-expresses our dataset?*

# Change of Basis

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**Note:** PCA makes a powerful assumption: *linearity*.

- The data characterizes/provides an ability to interpolate between the individual data points.

# Change of Basis

Let  $\mathbf{X}$  and  $\mathbf{Z}$  be  $N \times D$  matrices related by a linear transformation  $\theta$ :

$$\mathbf{X}\theta = \mathbf{Z}$$

- $\mathbf{X}$  is the original recorded dataset;
- $\mathbf{Z}$  is a re-representation of that dataset.

# Change of Basis

$$\mathbf{X}\theta = \mathbf{Z}$$

This change of basis has some interpretations:

- $\theta$  is a matrix that transforms  $\mathbf{X}$  to  $\mathbf{Z}$ .
- Geometrically,  $\theta$  is a rotation and a stretch which transforms  $\mathbf{X}$  into  $\mathbf{Z}$ .
- The columns of  $\theta$  are a set of new basis vectors for expressing the rows of  $\mathbf{X}$ :

$$\begin{pmatrix} - & \mathbf{x}_1 & - \\ - & \mathbf{x}_2 & - \\ \dots & \dots & \dots \\ - & \mathbf{x}_N & - \end{pmatrix} \begin{pmatrix} | & | & \vdots & | \\ \theta_1 & \theta_2 & \vdots & \theta_D \\ | & | & \vdots & | \end{pmatrix} = \mathbf{Z}$$

# Change of Basis

- Each row of  $\mathbf{Z}$  is

$$\mathbf{z}_i = (\mathbf{x}_i \cdot \theta_1, \mathbf{x}_i \cdot \theta_2, \dots, \mathbf{x}_i \cdot \theta_D).$$

- Each element of  $\mathbf{z}_i$  is a dot product of  $\mathbf{x}_i$  with the corresponding column in  $\theta$ .
- That is, the  $j$ -th element of  $\mathbf{z}_i$  is a projection of  $\mathbf{x}_i$  onto the  $j$ -th column of  $\theta$ .

# Change of Basis

- By assuming linearity, the problem reduces to finding the appropriate change of basis.
- The column vectors  $\theta_1, \theta_2, \dots, \theta_D$  will become the **principal components** of  $\mathbf{X}$ .



# Change of Basis

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## Questions:

- What is the best way to re-express  $\mathbf{X}$ ?
- What is a good choice of basis  $\theta$ ?

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# Noise

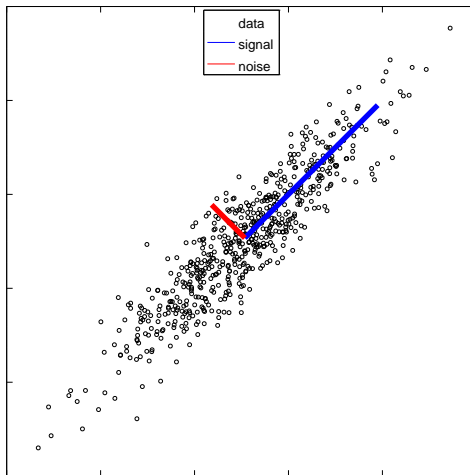
- Noise in any dataset must be low, otherwise, no useful information of a system can be extracted.
- All noise is measure relative to the measurement. A common measure is the *signal-to-noise ratio* (SNR), or a ratio of variance  $\sigma^2$ :

$$\text{SNR} = \frac{\sigma_{\text{signal}}^2}{\sigma_{\text{noise}}^2}$$

- A high SNR ( $\text{SNR} \gg 1$ ) indicates high precision data, while a low SNR indicates noise contaminated data.

# Noise

The SNR measures how “fat” the oval is.



# Noise

- In our example, any individual camera should record motion in a straight line.
- Therefore, any spread deviating from straight-line motion must be noise.

# Redundancy

- Redundancy is more tricky issue. Multiple sensors record the same dynamics information.
- A simple way to quantify the redundancy between measurements is to calculate the their **covariance**.
- Covariance is a measure of how much two random variables change together:
  - If the variables tend to show similar behavior (greater/greater, smaller/smaller), the covariance is positive.
  - If the variables tend to show opposite behavior (greater/smaller, smaller/greater), the covariance is negative.
- The sign of the covariance therefore shows the tendency in the *linear relationship* between the variables.

# Redundancy

The covariance between two jointly distributed real-valued random variables  $X$  and  $Y$  is defined as:

$$\begin{aligned}\sigma_{XY}^2 &= \mathbb{E}[(X - \mathbb{E}[X])(Y - \mathbb{E}[Y])] \\ &= \mathbb{E}[XY] - \mathbb{E}[X]\mathbb{E}[Y].\end{aligned}$$

Two important facts about covariance:

- $\sigma_{XY}^2 = 0$  iff  $X$  and  $Y$  are entirely uncorrelated.
- $\sigma_{XY}^2 = \sigma_X^2$  iff  $X = Y$ .

# Redundancy

- For random vectors  $\mathbf{X}$  and  $\mathbf{Y}$ , both of dimension  $D$ , their  $D \times D$  *covariance matrix* is

$$\sigma_{\mathbf{X}\mathbf{Y}}^2 = \mathbb{E}[\mathbf{X}\mathbf{Y}^T] - \mathbb{E}[\mathbf{X}]\mathbb{E}[\mathbf{Y}^T].$$

- For a vector  $X$  of  $D$  jointly distributed real-valued random variables, its covariance matrix is

$$\Sigma(\mathbf{X}) = \sigma_{\mathbf{X}\mathbf{X}}^2.$$



# Redundancy

The Iris dataset:

$$\Sigma(\mathbf{X}) = \begin{pmatrix} 0.665822 & -0.026056 & 1.235005 & 0.500998 \\ -0.026056 & 0.190509 & -0.308566 & -0.111119 \\ 1.235005 & -0.308566 & 3.071335 & 1.279612 \\ 0.500998 & -0.111119 & 1.279612 & 0.576284 \end{pmatrix}$$

Note that the diagonal elements of  $\Sigma(\mathbf{X})$  are the variances of particular features.

# Redundancy

If  $\mathbf{X}$  is a dataset containing  $N$  examples, each example  $\mathbf{x}_i$  has  $D$  features with *zero mean*. Then:

$$\Sigma(\mathbf{X}) = \frac{1}{N-1} \mathbf{X}^T \mathbf{X}.$$

$\Sigma(\mathbf{X})$  is a square symmetric  $D \times D$  matrix.

# Diagonalize the Covariance Matrix

- Our goal is to reduce redundancy, then we want each feature to co-vary as little as possible with other features.
- In order to remove redundancy, we want that all the covariances between separate features to be zero.
- That is, we want to transform from  $\mathbf{X}$  to  $\mathbf{Z}$  such that  $\Sigma(\mathbf{Z})$  is a diagonal matrix.

# Diagonalize the Covariance Matrix

- There are many methods for diagonalizing  $\Sigma(\mathbf{Z})$ . PCA uses the easiest method.
- First, PCA assumes that all basis vectors are orthonormal, that is

$$\theta_i \cdot \theta_j \equiv \delta(i = j).$$

In other words,  $\theta$  is an orthonormal matrix.

- Second, PCA assumes the directions with the largest variances are the most “*important*”, or most “*principal*”.

# Diagonalize the Covariance Matrix

How PCA works:

- First, it selects a normalized direction in  $D$ -dimensional space along which the variance in  $\mathbf{Z}$  is maximized. It saves this as  $\theta_1$ .
- Then, it find another direction  $\theta_2$  along which the variance is maximized. Because of the orthonormality condition, it restricts the search to all directions perpindicular to all previous selected directions ( $\theta_2 \cdot \theta_1 = 0$ ).
- This continues until  $D$  directions are selected.
- The resulting ordered set of  $\theta_j$  are the **principal components**.

# PCA Problem

## Problem

*Find some orthonormal matrix  $\theta$  where  $\mathbf{Z} = \mathbf{X}\theta$  such that  $\Sigma(\mathbf{Z}) = \frac{1}{N-1} \mathbf{Z}^T \mathbf{Z}$  is diagonalized.*

The columns of  $\theta$  are the principal components of  $\mathbf{X}$ .

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# Solving PCA: Eigenvectors of Covariance

We have

$$\begin{aligned}\Sigma(\mathbf{Z}) &= \frac{1}{N-1} \mathbf{Z}^T \mathbf{Z} \\ &= \frac{1}{N-1} (\mathbf{X} \boldsymbol{\theta})^T (\mathbf{X} \boldsymbol{\theta}) \\ &= \frac{1}{N-1} \boldsymbol{\theta}^T \mathbf{X}^T \mathbf{X} \boldsymbol{\theta} \\ &= \frac{1}{N-1} \boldsymbol{\theta}^T \mathbf{A} \boldsymbol{\theta},\end{aligned}$$

where we define  $\mathbf{A} = \mathbf{X}^T \mathbf{X}$ , which is a *symmetric* matrix.



# Solving PCA: Eigenvectors of Covariance

## Theorem

*A symmetric matrix is diagonalized by a matrix of its orthonormal eigenvectors.*

Because of this theorem, there exists a diagonal matrix **D** such that

$$\mathbf{A} = \mathbf{E}\mathbf{D}\mathbf{E}^T,$$

where **E** is a matrix of eigenvectors of **A** arranged as columns.

# Solving PCA: Eigenvectors of Covariance

- The matrix  $\mathbf{A}$  has  $L \leq D$  orthonormal eigenvectors where  $L$  is the rank of the matrix.
- The rank of  $\mathbf{A}$  is less than  $D$  when  $\mathbf{A}$  degenerate, or all data occupy a subspace of dimension  $L < D$ .
- So, we select the matrix  $\theta$  to be a matrix where each column  $\theta_j$  is an eigenvector of  $\mathbf{X}^T \mathbf{X}$ .
- By this selection, we have  $\theta = \mathbf{E}$ . So,

$$\mathbf{A} = \theta \mathbf{D} \theta^T.$$

# Solving PCA: Eigenvectors of Covariance

Therefore,

$$\begin{aligned}\Sigma(\mathbf{Z}) &= \frac{1}{N-1} \theta^T \mathbf{A} \theta \\ &= \frac{1}{N-1} \theta^T (\theta \mathbf{D} \theta^T) \theta \\ &= \frac{1}{N-1} (\theta^T \theta) \mathbf{D} (\theta^T \theta) \\ &= \frac{1}{N-1} (\theta^{-1} \theta) \mathbf{D} (\theta^{-1} \theta) \\ &= \frac{1}{N-1} \mathbf{D}.\end{aligned}$$

That is, the choice of  $\theta$  diagonalizes  $\Sigma(\mathbf{Z})$ .

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# Theoretical Basis

## Theorem

*The inverse of an orthogonal matrix is its transpose.*

## Theorem

*If  $\mathbf{X}$  is any matrix, the matrices  $\mathbf{X}^T \mathbf{X}$  and  $\mathbf{X} \mathbf{X}^T$  are both symmetric.*

## Theorem

*A matrix is symmetric if and only if it is orthogonally diagonalizable.*

# Theoretical Basis

## Theorem

*A symmetric matrix is diagonalized by a matrix of its orthonormal eigenvectors.*

## Theorem

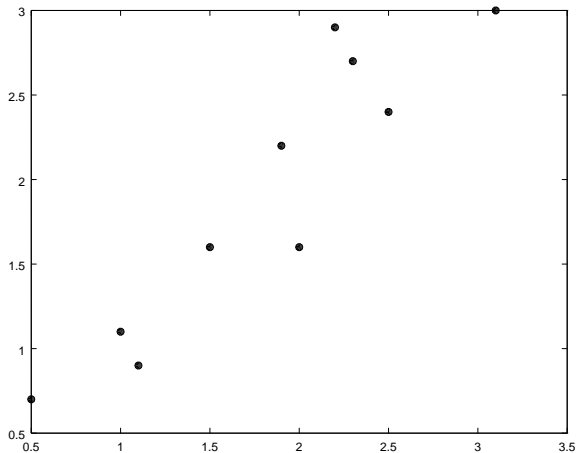
*For any arbitrary  $N \times D$  matrix  $\mathbf{X}$ , the symmetric matrix  $\mathbf{X}^T \mathbf{X}$  has a set of orthonormal eigenvectors  $\{\mathbf{v}_1, \mathbf{v}_2, \dots, \mathbf{v}_D\}$  and a set of associated eigenvalues  $\{\lambda_1, \lambda_2, \dots, \lambda_D\}$ . The set of vectors  $\{\mathbf{X} \mathbf{v}_1, \mathbf{X} \mathbf{v}_2, \dots, \mathbf{X} \mathbf{v}_D\}$  form an orthogonal basis, where each vector  $\mathbf{X} \mathbf{v}_j$  is of length  $\sqrt{\lambda_j}$ .*

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# Example 1: Toy Dataset

| $x_1$ | $x_2$ |
|-------|-------|
| 2.5   | 2.4   |
| 0.5   | 0.7   |
| 2.2   | 2.9   |
| 1.9   | 2.2   |
| 3.1   | 3.0   |
| 2.3   | 2.7   |
| 2.0   | 1.6   |
| 1.0   | 1.1   |
| 1.5   | 1.6   |
| 1.1   | 0.9   |

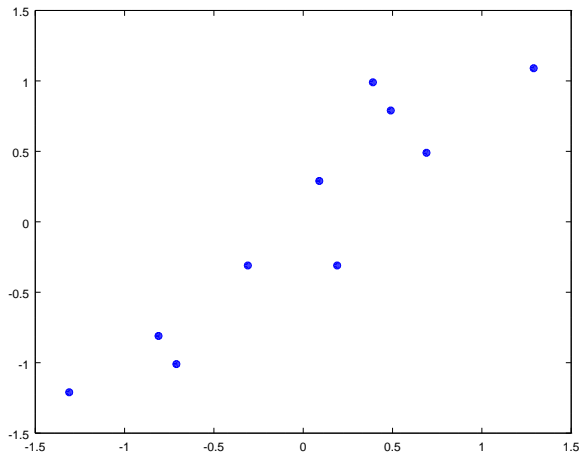




# Example 1: Toy Dataset

| $x_1$ | $x_2$ |
|-------|-------|
| 0.69  | 0.49  |
| -1.31 | -1.21 |
| 0.39  | 0.99  |
| 0.09  | 0.29  |
| 1.29  | 1.09  |
| 0.49  | 0.79  |
| 0.19  | -0.31 |
| -0.81 | -0.81 |
| -0.31 | -0.31 |
| -0.71 | -1.01 |

$$\mu = (1.81, 1.91)$$



# Example 1: Toy Dataset

$$\theta = \begin{pmatrix} 0.67787 & -0.73518 \\ 0.73518 & 0.67787 \end{pmatrix}$$

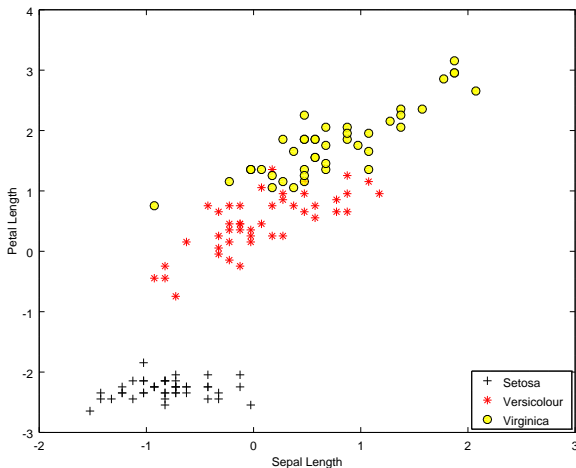
$$\mathbf{Z} = \begin{pmatrix} 0.827970 & -0.175115 \\ -1.777580 & 0.142857 \\ 0.992197 & 0.384375 \\ 0.274210 & 0.130417 \\ 1.675801 & -0.209498 \\ 0.912949 & 0.175282 \\ -0.099109 & -0.349825 \\ -1.144572 & 0.046417 \\ -0.438046 & 0.017765 \\ -1.223821 & -0.162675 \end{pmatrix}$$

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# Iris Dataset

Four features, reorder features as “Sepal Length”, “Petal Length”, “Sepal Width”, “Petal Width”.



# Iris Dataset

$$\theta = \begin{pmatrix} 0.356687 & 0.657221 & 0.578737 & 0.325419 \\ 0.858455 & -0.176179 & -0.060299 & -0.477891 \\ -0.079358 & 0.729440 & -0.589941 & -0.337032 \\ 0.359904 & -0.070280 & -0.559819 & 0.743056 \end{pmatrix}$$

First and second principal component:

$$\theta_1 = \begin{pmatrix} 0.356687 \\ 0.858455 \\ -0.079358 \\ 0.359904 \end{pmatrix}; \quad \theta_2 = \begin{pmatrix} 0.657221 \\ -0.176179 \\ 0.729440 \\ -0.070280 \end{pmatrix}$$

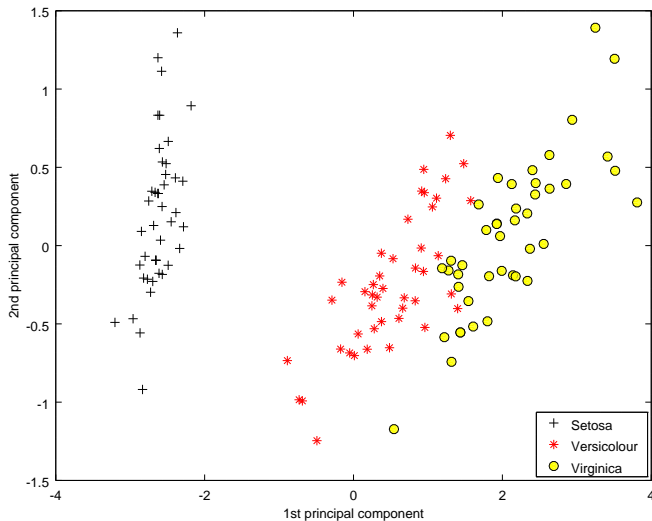
# Iris Dataset

Projection of  $\mathbf{X}$  into 2 two-dimensional space:

$$\mathbf{Z} = \mathbf{X} * [\theta_1, \theta_2]$$

This can be viewed as a “data compression” technique (dimensionality reduction).

# Iris Dataset



# Iris Dataset

- In practice, if we were using a learning algorithm (linear regression, neural networks, ...), we could now use the projected data instead of the original data.
- By using the projected data, we can train our model faster as there are less dimensions in the input.



# Data Reconstruction

- After projecting the data onto the lower dimensional space, we can approximately recover the data by projecting them back to the original high dimensional space:

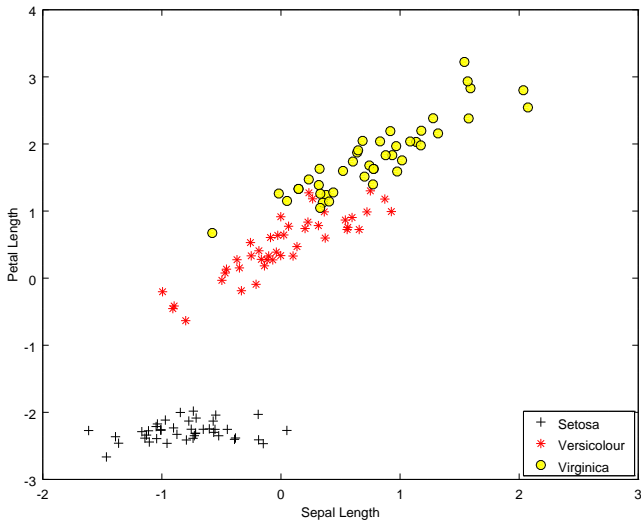
$$\mathbf{X}' = \mathbf{Z} \theta^T,$$

where  $\theta = [\theta_1, \theta_2, \dots, \theta_K]$  contains  $K$  principal components.

- The recovered data  $\mathbf{X}'$  is generally a coarsened-grained version of the original data  $\mathbf{X}$ :
  - Some information is lost, some hidden semantics/structures are retained.
- Reconstruction error:

$$J(\theta) = \frac{1}{N} \sum_{i=1}^N \|\mathbf{x}_i - \mathbf{x}'_i\|^2.$$

# Data Reconstruction – Iris Dataset



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# Face Image Dataset

- We run PCA on face images to see how it can be used in practice for dimension reduction.
- The face image dataset contains 5000 face images, each of size  $32 \times 32$  in grayscale.<sup>1</sup>
- Each row of  $\mathbf{X}$  corresponds to one face image (a row vector of length 1024).

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<sup>1</sup>A subset of the Labeled Face in the Wild Home.

# Face Image Dataset – 100 Original Faces



# Face Image Dataset – 36 Principal Components



# Face Image Dataset – 100 Principal Components

Original faces



Recovered faces



# Exercises

- 1 Implement the PCA algorithm.
- 2 Test the algorithm on different datasets.
- 3 Run a classification algorithm on the projected Iris dataset (using first two principal components) and report the classification accuracy.