

# Principal component analysis

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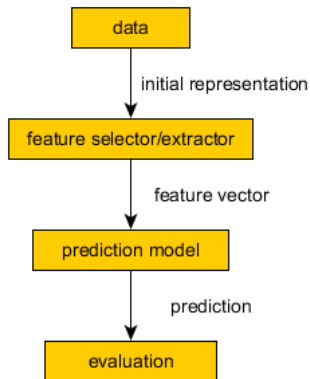
Yandex School of Data Analysis



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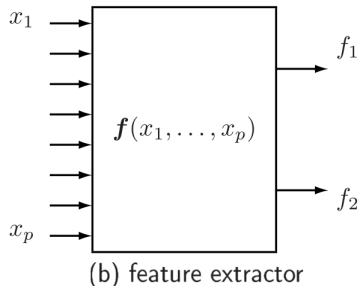
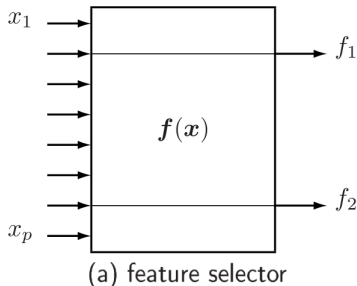
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# General modelling pipeline



# Dimensionality reduction

## Feature selection / Feature extraction



**Feature extraction:** find transformation of original data which extracts most relevant information for machine learning task.

We will consider unsupervised dimensionality reduction methods, which try to preserve geometrical properties of the data.

# Applications of dimensionality reduction

## Applications:

- visualization in 2D or 3D
- reduce operational costs (less memory, disk, CPU usage on data transfer)
- remove multi-collinearity to improve performance of machine-learning models

# Categorization

Supervision in dimensionality reduction:

- supervised (such as Fisher's direction)
- unsupervised

Mapping to reduced space:

- linear
- non-linear

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## 2 Principal component analysis

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# Best hyperplane fit

- For point  $x$  and subspace  $L$  denote:
  - $p$ -the projection of  $x$  on  $L$
  - $h$ -orthogonal complement
- $x = p + h$ ,  $\langle p, h \rangle = 0$ .

## Proposition 1

*For  $x$ , its projection  $p$  and orthogonal complement  $h$*

$$\|x\|^2 = \|p\|^2 + \|h\|^2.$$

- Prove proposition 1.
- For training set  $x_1, x_2, \dots, x_N$  and subspace  $L$  we can also find:
  - projections:  $p_1, p_2, \dots, p_N$
  - orthogonal complements:  $h_1, h_2, \dots, h_N$ .

# Best subspace fit

## Definition 1

Best-fit  $k$ -dimensional subspace for a set of points  $x_1, x_2, \dots, x_N$  is a subspace, spanned by  $k$  vectors  $v_1, v_2, \dots, v_k$ , solving

$$\sum_{n=1}^N \|h_n\|^2 \rightarrow \min_{v_1, v_2, \dots, v_k}$$

## Proposition 2

Vectors  $v_1, v_2, \dots, v_k$ , solving

$$\sum_{n=1}^N \|p_n\|^2 \rightarrow \max_{v_1, v_2, \dots, v_k}$$

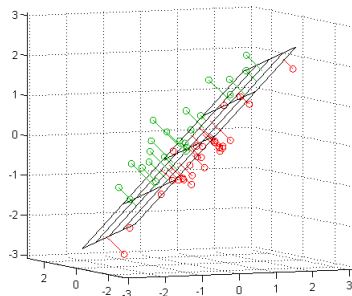
also define best-fit  $k$ -dimensional subspace.

- Prove 2 using proposition 1.

## Definition 2

Principal components  $a_1, a_2, \dots, a_k$  are vectors, forming orthonormal basis in the  $k$ -dimensional subspace of best fit.

# Best hyperplane fit



Subspace  $L_k$  or rank  $k$  best fits points  $x_1, x_2, \dots, x_D$ .

# Properties of PCA

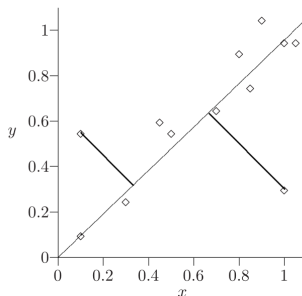
- Properties:
  - Not invariant to translation:
    - Before applying PCA, it is recommended to center objects:

$$x \leftarrow x - \mu \text{ where } \mu = \frac{1}{N} \sum_{n=1}^N x_n$$

- Not invariant to scaling:
  - scale features to have unit variance

## Example: line of best fit

- In PCA the sum of squared perpendicular distances to line is minimized:

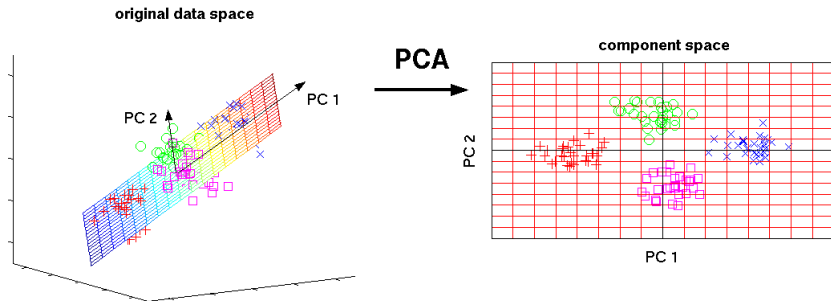


- *What is the difference with least squares minimization in regression?*

## 2 Principal component analysis

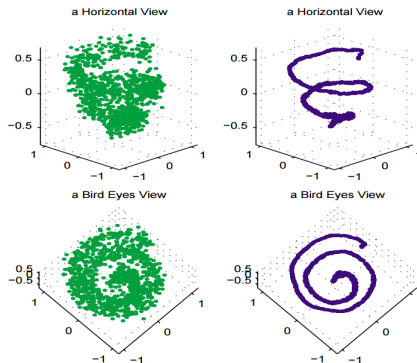
- Definition
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# Visualization



# Data filtering

Remove noise to get a cleaner picture of data distribution:



X. Huo and Jihong Chen (2002). Local linear projection (LLP). First IEEE Workshop on Genomic Signal Processing and Statistics (GENSIPS), Raleigh, NC, October.  
<http://www.gensips.gatech.edu/proceedings/>.



## Economic description of data

Faces database:

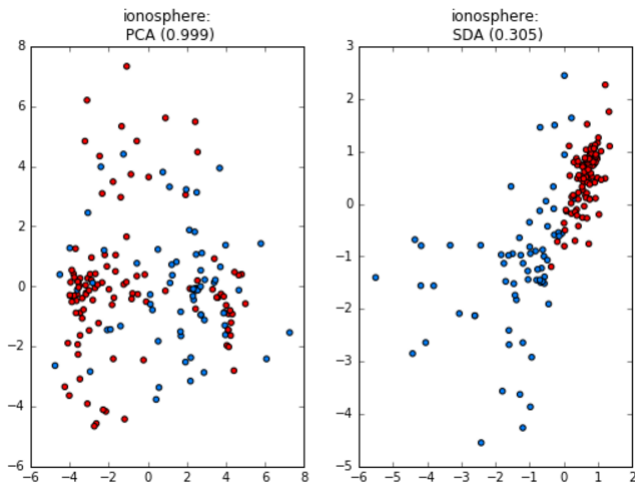


# Eigenfaces

Eigenvectors are called eigenfaces. Projections on first several eigenfaces describe most of face variability.

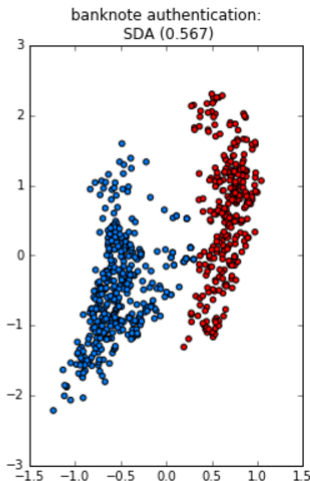
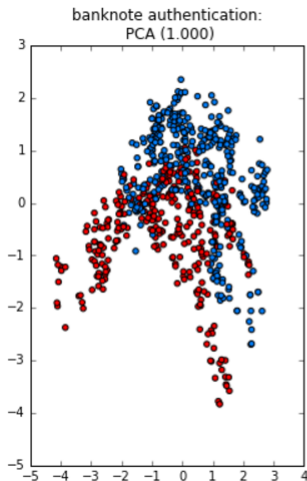


# PCA vs. SDA (not covered here)



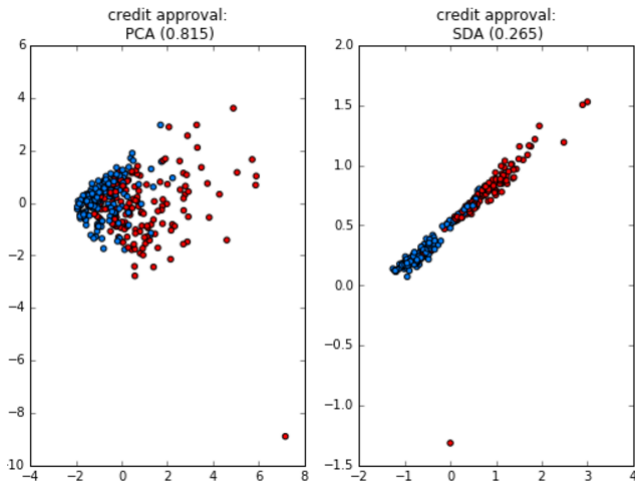
Title format: dataset, method (quality of approximation (2)).

## PCA vs. SDA (not covered here)



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## Quality of approximation

Consider vector  $x$ . Since all  $D$  principal components form a full orthonormal basis,  $x$  can be written as

$$x = \langle x, a_1 \rangle a_1 + \langle x, a_2 \rangle a_2 + \dots + \langle x, a_D \rangle a_D$$

Let  $p^K$  be the projection of  $x$  onto subspace spanned by first  $K$  principal components:

$$p^K = \langle x, a_1 \rangle a_1 + \langle x, a_2 \rangle a_2 + \dots + \langle x, a_K \rangle a_K$$

Error of this approximation is

$$h^K = x - p^K = \langle x, a_{K+1} \rangle a_{K+1} + \dots + \langle x, a_D \rangle a_D$$

## Quality of approximation

Using that  $a_1, \dots, a_D$  is an orthonormal set of vectors, we get

$$\begin{aligned}\|x\|^2 &= \langle x, x \rangle = \langle x, a_1 \rangle^2 + \dots + \langle x, a_D \rangle^2 \\ \|p^K\|^2 &= \langle p^K, p^K \rangle = \langle x, a_1 \rangle^2 + \dots + \langle x, a_K \rangle^2 \\ \|h^K\|^2 &= \langle h^K, h^K \rangle = \langle x, a_{K+1} \rangle^2 + \dots + \langle x, a_D \rangle^2\end{aligned}$$

We can measure how well first  $K$  components describe our dataset  $x_1, x_2, \dots, x_N$  using relative loss

$$L(K) = \frac{\sum_{n=1}^N \|h_n^K\|^2}{\sum_{n=1}^N \|x_n\|^2} \quad (1)$$

or relative score

$$S(K) = \frac{\sum_{n=1}^N \|p_n^K\|^2}{\sum_{n=1}^N \|x_n\|^2} \quad (2)$$

Evidently  $L(K) + S(K) = 1$ .



## Contribution of individual component

Contribution of  $a_k$  for explaining  $x$  is  $\langle x, a_k \rangle^2$ .

Contribution of  $a_k$  for explaining  $x_1, x_2, \dots, x_N$  is:

$$\sum_{n=1}^N \langle x_n, a_k \rangle^2$$

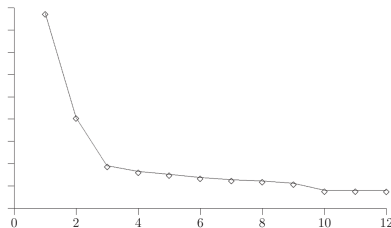
Explained variance ratio:

$$E(a_k) = \frac{\sum_{n=1}^N \langle x_n, a_k \rangle^2}{\sum_{d=1}^D \sum_{n=1}^N \langle x_n, a_d \rangle^2} = \frac{\sum_{n=1}^N \langle x_n, a_k \rangle^2}{\sum_{n=1}^N \|x_n\|^2}$$

- Explained variance ratio measures relative contribution of component  $a_k$  to explaining our dataset  $x_1, \dots, x_N$ .
- Note that  $\sum_{k=1}^K E(a_k) = S(K)$ .

## How many principal components to select?

- Data visualization: 2 or 3 components.
- Take most significant components until their variance falls sharply down:



- Or take minimum  $K$  such that  $L(K) \leq t$  or  $S(K) \geq 1 - t$ , where typically  $t = 0.95$ .

# Conclusion<sup>1</sup>

- For  $x \in \mathbb{R}^D$  there exist  $D$  principal components.
- Principal component  $a_i$  is the  $i$ -th eigenvector of  $X^T X$ , corresponding to  $i$ -th largest eigenvalue  $\lambda_i$ .
- Sum of squared projections onto  $a_i$  is  $\|Xa_i\|^2 = \lambda_i$ .
- *Explained variance ratio* by component  $a_i$  is equal to

$$\frac{\lambda_i}{\sum_{d=1}^D \lambda_d}$$

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<sup>1</sup>Compare dimensionality reduction with PCA and regularization as means of simplification of prediction model.