MACHINE LEARNING

MEI/1

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Machine Learning

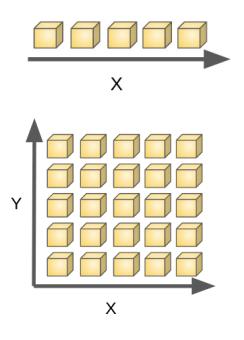
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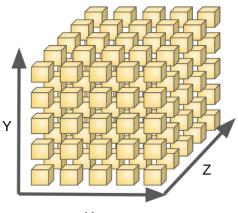
Syllabus

- Feature Representation
- Dimensionality Reduction
- PCA
 - Eigenvectors and Eigenvalues

Dimensionality Reduction

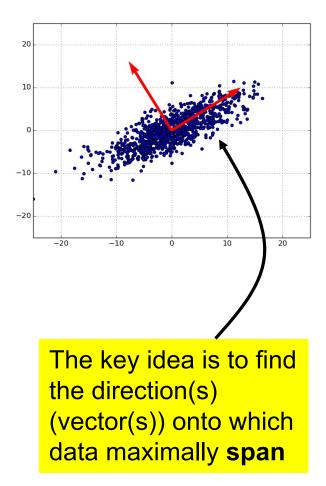
- The curse of dimensionality is one of the most classical phenomena in the development of Machine Learning systems.
 - In short, when the dimensionality increases, the volume of the space increases so fast that the data become sparse.
 - Sparsity is problematic for any method that requires statistical significance, i.e., densely populated spaces.
- For example, consider 100 evenly spaced sample points (instances) inside a unit interval.
 - On average, points will be separated around 10⁻²=0.01
- An equivalent sampling that will yield similar density in a 10-dimensional unit hypercube would require 10²⁰[=(10²)¹⁰] sample points





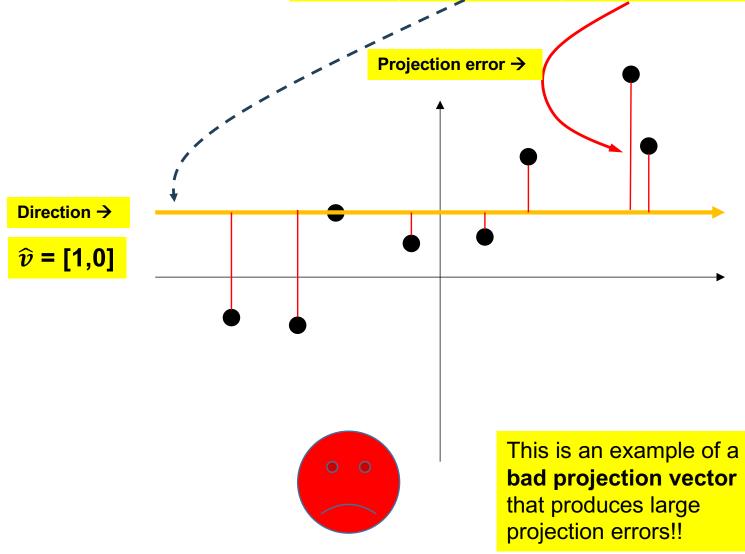
Dimensionality Reduction

- In statistics, machine learning, and information theory, dimensionality reduction is the process of reducing the number of random variables under consideration by obtaining a set of principal variables.
- In general, there are two families of methods to reduce the dimensionality of a data set:
 - Feature Selection. The idea is to find a subset of the original features that better represent the problem, i.e., that minimally decrease the amount of available information, when compared to the original dataset.
 - Most approaches are based in filters (based in information gain), wrappers (based in accuracy) and embedded (features iteratively selected/removed according to prediction errors)
 - Feature Extraction. It is often also designated as "Feature Projection" and the idea is to transform the original feature space into a space of fewer dimensions, while keeping as much of the original information as we can.
 - **Principal Component Analysis (PCA)** is the main technique in this family.



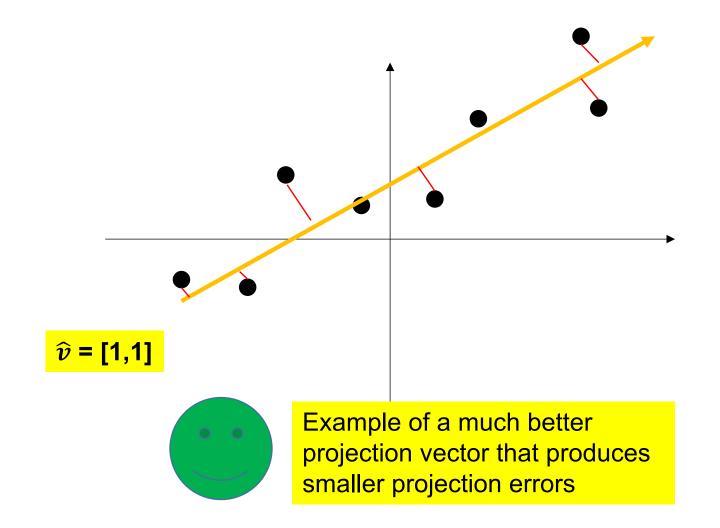
PCA

 Graphically, we are interested in finding the direction (vector in the original space) onto which the projected data provides the minimal projection error:



PCA

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PCA: Covariance

 The covariance can be obtained for any two dimensions (features) of a n-dimensional feature space

Y

Χ

X

- It is a measure of the joint variability of two features
 - If both variables vary in a direct way, the / covariance is positive
 - On the contrary, if both variables vary inversely, the variance values will be negative.
- The sign of the covariance shows the tendency in the linear relationship between the variables.
- The magnitude of the covariance is not easy to interpret because it is not normalized and hence depends on the magnitudes of the variables.
- The normalized version of the covariance, the correlation coefficient, however, shows by its magnitude the strength of the linear relation.

PCA: Covariance

 The distance between sample points and their mean is multiplied. Then, the result is divided by the number of data points minus 1:

$$cov(X,Y) = \frac{\sum_{i=1}^{n} (X_i - X^*)(Y_i - Y^*)}{n-1}$$

where X_i , Y_i are the ith data points, X^* , Y^* are the sample means and "n" is the number of data points.

- The results is meaningful essentially by analysing it's sign:
 - Positive: Both dimensions vary directly.
 - Negative: Both dimensions vary inversely.
 - Zero: Dimensions are independent.

PCA: Covariance Matrix

The Covariance Matrix C contains all covariance pair values between every possible dimensions of a feature space :

$$\boldsymbol{C} = [cij | cij = cov(Xi, Xj)]$$

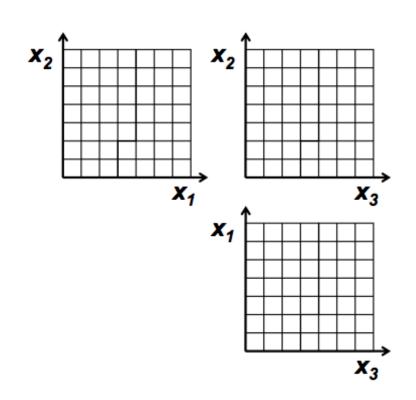
• For exemple, considering a three dimensional space {X, Y, Z}, the covariance matrix will correspond to:

$ \begin{vmatrix} cov(Y,X) & cov(Y,Y) & cov(Y,Z) \\ cov(Z,X) & cov(Z,Y) & cov(Z,Z) \end{vmatrix} $
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- Values along the main diagonal describe the variance of the corresponding dimension.
- Based on its definition, it is obvious that cov(X,Y)=cov(Y,X), i.e., the covariance matrix is symetric with respect to its main diagonal.

PCA: Covariance Matrix

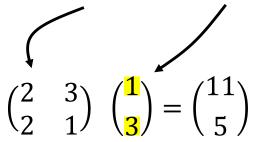
• Exercise. Obtain the covariance matrix for the given data set:



Obs.	X1	X2	Х3		
1	2	2	4		
2	3	4	6		
3	5	4	2		
4	6	6	4		

	X1	X ₂	X ₃	Х1-µ1	Х2-µ2	Х3-µ3	(x 1-μ1) ²	(x ₂ -μ ₂) ²	(x ₃ -μ ₃) ²	(x1-μ1)(x2-μ2)	(x ₁ -μ ₁)(x ₃ -μ ₃)	(x2-μ2)(x3-μ3)
1												
2												
3												
4												

• Consider the multiplication of a **matrix** by a **vector**:

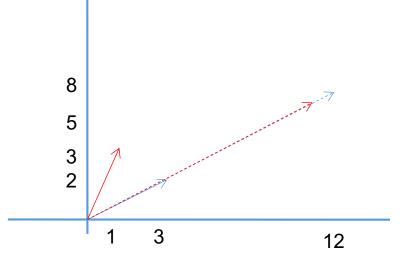


However, there are some particularly interesting vectors:

$$\begin{pmatrix} 2 & 3 \\ 2 & 1 \end{pmatrix} \begin{pmatrix} 3 \\ 2 \end{pmatrix} = \begin{pmatrix} 12 \\ 8 \end{pmatrix} = 4 \begin{pmatrix} 3 \\ 2 \end{pmatrix}$$

- In the first case, the resulting vector is not a multiple of the original vector.
- Oppositelly, in the second case, the resulting vector (12,8) is a multiple of the multiplier.
- As such, the latter is an eigenvector.
 - The correspondong eigenvalue is "4"

• By analysing the direction of the original and resultant vectors:



- Considering the matrix as a transformation, it can be concluded that in the second case, the direction was not changed. This is the key insight the notion of eigenvector.
 - The given matrix does not change the direction of its eigenvectors.

- As we've seen, the notion of **eigenvalue** is strongly related to the **eigenvector**.
 - It is the value that should by multiplied by the eigenvector to obtain the original vector.
 - In the above example, 4 was the eigenvalue that corresponds to the given eigenvector.

 As such, eigenvalues and eigenvectors come in pairs and are always inter-related.

- As a summary, the eigenvectors of a matrix correspond to the directions that are not changed by the (transformation) matrix.
- Not all matrices have eigenvectors.
- Matrices have to be square.
- A (n x n) matrix has at most "n" eigenvectors.
- The set of eigenvectors of a matrix (image) is widely used to describe the spatial content of that image (feature).
- In MATLAB, this eigenanalysis is made by the "eig()" function:
 - [V,D] = eig(A)
 - Returns the eigenvectors (D) and corresponding eigenvalues (V) of matrix A.
- In Python, this can simply be done by:
 - V, D = LA.eig(A)

- There is an important property to be stressed: the eigenvectors of a matrix are orthogonal. This is to say that they form an orthogonal basis of the matrix.
 - We are able to express every point of a data set by linear combinations of its basis-vectors.
 - This is specially usefull for the analysis of principal components (PCA).
 - It is usual to determine the eigenvectors/eigenvalues in their normalized version, i.e., with length normalized to 1.
 - As previously seen, the length of a vector does not affect its property of being (or not) an eigenvector.
 - Hence, having an eigenvector $(x_1, ..., x_n)$ it is usual to divide each component by the norm of this vector, in order to obtain length "1":
 - $||(x_1,...,x_n)|| = sqrt(x_1^2 + ... + x_n^2)$

• Exercise

• Determine, from the following vectors, which are eigenvectors of the matrix given below and, if positive, determine the corresponding eigenvalue.

• Matrix:

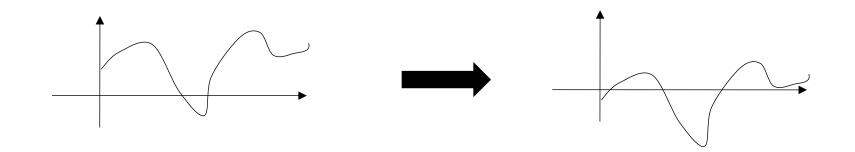
$$\begin{bmatrix} 3 & 0 & 1 \\ -4 & 1 & 2 \\ -6 & 0 & -2 \end{bmatrix}$$

• Vectors:

$$\begin{bmatrix} 2 & [-1 & [-1 & [0 & [3 \\ 2 & 0 & 1 & 1 & 2 \\ -1] & 2] & 3] & 0] & 1 \end{bmatrix}$$

- The Principal Component Analysis (PCA) it's a well known way to detect patterns on data, by expressing it on a way that enhances similarities or differences.
- Detecting patterns on high dimensional data is a hard task, either for humans or machines.
 - Requires huge amounts of data. An empirical rule says that at the minimum, d² instances are required to analyze a ddimensional data set.
- PCA is typically used to compress data (reduce dimensionality), without loosing substantial information.

- Step 1. The analysis of principal components requires a data set (with dimension n) and cardinality (k).
- Step 2. Removal of energy. For each dimension, the corresponding mean is subtrated to each component. As such, all dimensions of the data set have zero energy.



- Step 3. Calculus of the covariance matrix. Here, the relationships between independent components are detected, together with an assessment of the data dispersion in each dimension (by analysing the main diagonal components).
- Step 4. As the covariance matrix is square, it is possible to obtain the set of eigenvectors and corresp^oonding eigenvalues.
- **Step 4.1.** Eigenvectors normalization. All eigenvectors are normalized to have norm equal to 1.

- Step 5. Selection of components. The set of eigenvectors is sorted by decreasing order, considering the corresponding eigenvalues. From this set, the "k₁" principal components are selected.
 - This is the step that performs the reduction of dimensionality.
- Step 6. A transformation matrix is built, by concatenating the eigenvectors selected in the previous step.
 - This matrix will be used to represent all points in the reduced dimensionality feature space. MAT=[vect1, vect2, ... Vectk₁]

 Step 7. Data Transformation. As the transformation matrix has "d" lines (corresponding to the dimension of the original feature space and k₁ columns (corresponding to the dimension of the new feature space), when multipling each original data point by the transformation matrix, we obtain a vector of k₁ components. These are the new representation of the data points, in the principal components space.

$[1 x d] x [d x k_1] = [1 x k_1]$

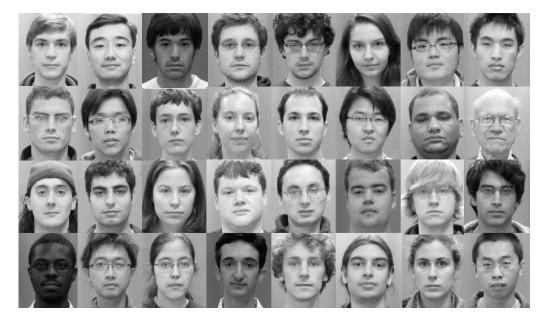
• How to choose the value of "k"?

- The previously described process does not give any information about a strategy to select the dimensionality of the principal components feature space.
- There is no formal rule. However, some heuristics about what is generally better exist.
- Usually, the variation in magnitude of consecutive eigenvalues (after sorting) is measured. When changes in magnitude are higher than a threshold, the selection process is stopped.
- But most frequently, the proportion of the data variability that is kept by the selected components is considered as the main criterium.
 - Typically, we are interested in keeping around 90, 95, 99% of the original data variability.
 - The analysis can be done by measuring the proportion of the sum of eigenvalues:

• Variability:
$$\frac{\sum_{i=1}^{k} \lambda_i}{\sum_{i=1}^{d} \lambda_i}$$
, "k": number of selected vectors, and "d": dimensionality

PCA: Example

- Consider a set of 128 face grayscale images (with dimensions 64 x 64).
 - Each image is represented by a 64 x 64 matrix = (4096), where each position represents the intensity at a point (0: black pixel, ... 255: White pixel)
- Each face can be regarded as a point represented in a 4096 dimensional feature space



PCA: Example

- We can use the PCA to select the principal components in this space (i.e., the directions in which the elements mostly span (vary)).
 - In pratice, the eigenvectors (each one with dimension 4096) with the largest corresponding eigenvalues will be selected.
- Next, each original face can be represented as a weighted combination of the top-k eigenvalues.
 - In such case, each face will actually be represented by weights α : α_1 , ... α_k
 - The PCA can be also regarded as a way to represent a face, with much less information than the originally used, while keeping the most importante information.
- Further, the facial recognition process can be done in the new feature space of (much more) reduced dimension, i.e., typically k << d (original space).

PCA: Example

• Example of the 16 principal components (eigenvectors with the largest eigenvalues) from the above data set:

