Machine Learning CS 165B

Prof. Matthew Turk

Wednesday, May 18, 2016

- Features (Ch. 10)

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Notes

Four Eyes Lab Open House Friday, May 27, 5-8pm 3rd floor Elings Hall

Along with the Media Arts and Technology End of Year Show Exhibition 5-9pm, throughout Elings Hall

http://show.mat.ucsb.edu

Summary: Distance methods and clustering

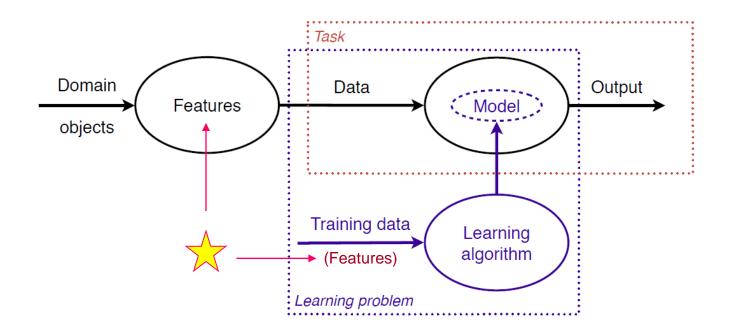
- Similarity is a function of distance
- Euclidian distance may not always be the right choice
- Nearest neighbor methods assign classes/clusters based on distances to points or exemplars, not based on computed boundaries
- For good clustering, we want high within-class (intra-class) similarity and low between-class (inter-class) similarity
- The scatter matrix is an important structure in clustering
- The K-means algorithm (and variations) is widely used

Features

Chapter 10 in the textbook (mostly 10.3)

Schedule change: skipping Chapter 9!!

A machine learning system



A **task** requires an appropriate mapping – a **model** – from data described by **features** to outputs. Obtaining such a model from training data is what constitutes a **learning problem**.

Tasks are addressed by **models**.

Learning problems are solved by **learning algorithms** that produce **models**.

Features

- Let's think about features (aka attributes) a bit more than we have....
- What's a feature? A mapping from the instance space \mathscr{L} to the feature domain \mathscr{F}

$$f_i: \mathcal{X} \to \mathcal{F}_i$$

- The model can be thought of as just a new feature a
 particular combination of the input features, constructed to
 solve the task at hand
 - E.g., the $\mathbf{w}^T \mathbf{x}$ value in a linear classifier or SVM
- There are different categories of features and of permissible operations on features

Quantitative features

- Measured on a meaningful numeric scale
- Domain is often real values
- E.g.: weight, height, age, angle, match to template, ...

Ordinal features

- The relevant information is the ordering, not the scale
- Domain is an ordered set
- E.g., rank, street addresses, preference, ratings, ...

Categorical features

- No scale or order information
- Domain is an unordered set
- E.g., colors, names, parts of speech, binary attributes, ...

Kinds of features, their properties, and allowable statistics:

Kind	Order	Scale	Tendency	Dispersion	Shape
Categorical Ordinal Quantitative	× √ √	× × √	mode median mean	n/a quantiles range, interquartile range, variance, standard devia- tion	n/a n/a skewness, kurtosis

Statistics – calculations on the features

Mode – the value that occurs most frequently

Median – the middle value in an ordered list

Mean (expected value) – the arithmetic average of the values

$$\{0, 0, 0, 1, 2, 9, 1000\} \implies Mode = 0$$

$$Median = 1$$

$$Mean = 144.57$$

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Statistics – calculations on the features

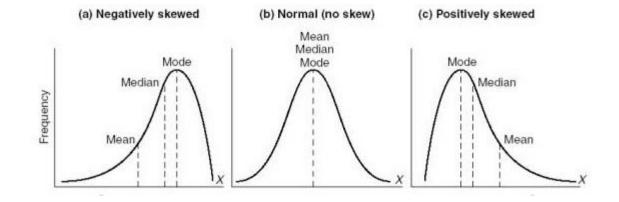
Range = (max. value - min. value)

Standard deviation (σ) = Square root of the variance (σ^2)

Skewness $\propto 3^{\text{rd}}$ moment (lack of symmetry: right or left skewed)

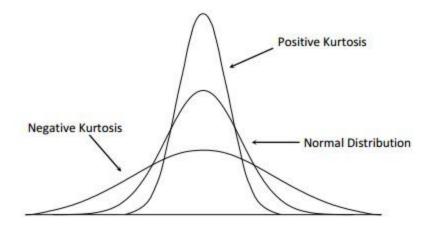
Kurtosis $\propto 4^{\text{th}}$ moment (peakedness relative to a normal distribution)

Skewness and kurtosis





Skewness:



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Statistics – calculations on the features

Percentiles: p^{th} percentile = the value such that p percent of the instances fall below it

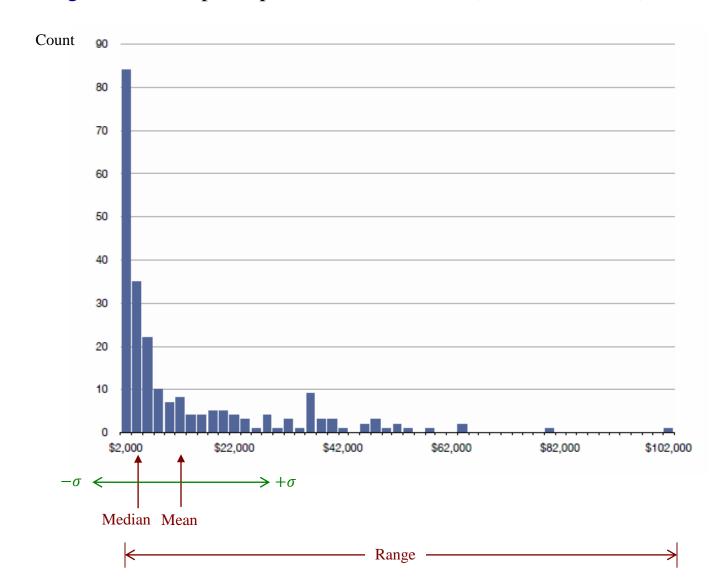
Deciles – multiples of 10 percentiles (10th, 20th, etc.)

Quartiles – multiples of 25 percentiles (25th, 50th, etc.)

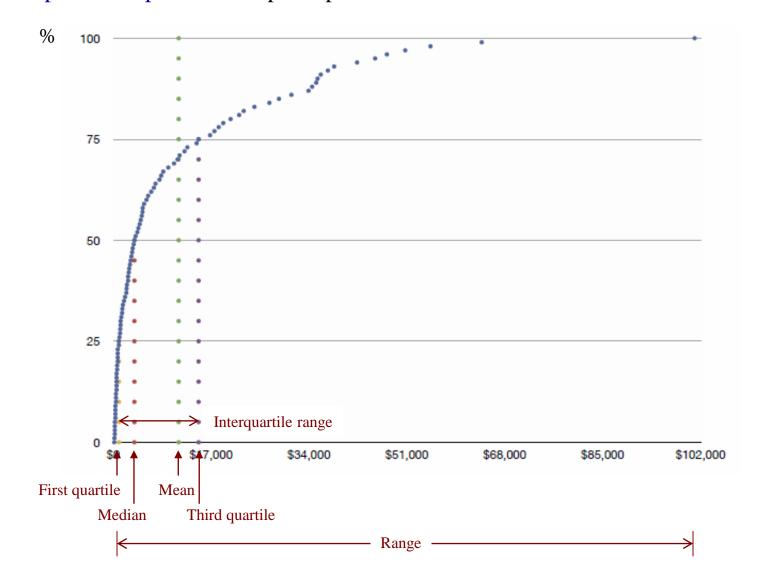
Interquartile range – the difference between the first and third quartiles (75th and 25th percentiles)

Histogram

A histogram of GDP per capita for 231 countries (bin size = \$2000):



A percentile plot of GDP per capita for 231 countries:

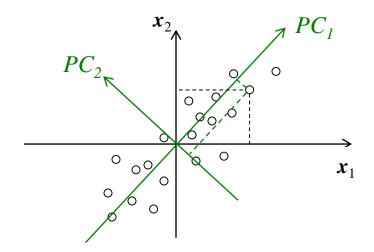


Feature construction and selection

- New features can be constructed from the given feature set
 - E.g., kernel methods do this, sometimes for great benefit
- Features may be combined in various ways, e.g.:
 - N-grams groups of N consecutive elements
 - E.g., trigrams: "my first teacher," "my first job," "my first the"
 - Cartesian products of categorical features
 - E.g., Shapes × Colors = { blue circle, red square, green line, ... }
 - w vector for a perceptron (combining training data point features)
 - Kernel function κ (e.g., 2D to 3D transformation)
- Typical approach: Generate new features, then select a suitable subset prior to learning
 - How to determine/evaluate suitability? How many to keep?
 - Many different approaches to this problem....

PCA and eigendecomposition

- One of the most widely used feature construction/selection techniques is Principal Component Analysis (PCA)
 - PCA constructs new features that are linear combinations of given features
- Computed eigenvectors and eigenvalues hold useful information
- Often used for dimensionality reduction, finding the intrinsic linear structure in the data

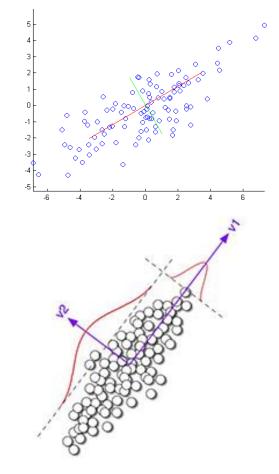


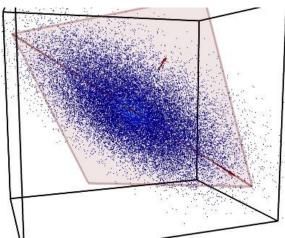
Given features 1 and 2 (x_1, x_2) <u>Computed</u> features 1 and 2 (green axes)

$$PC_1 = \underset{y}{\operatorname{argmax}}(y^T X)(X^T y)$$
(maximize variance of points projected onto unit vector y)

PCA and eigendecomposition

- The first principal component is the direction of maximum (1D) variance in the data
- The second principal component is the direction, perpendicular to the 1st PC, of maximum variance in the data
 - Etc. for additional PCs
- For N-dimensional data, there are N principal components
 - But perhaps only k of them are useful (k < N) dimensionality reduction!
- PCA typically assumes zero-mean data
 - First subtract the mean from each data point;
 centroid is thus (0, 0)





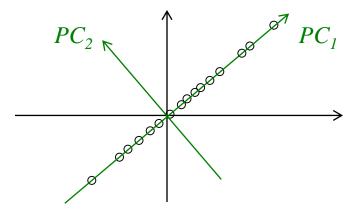
PCA and eigendecomposition

- For *n* points of dimension *d*, let $X = [x_1 \ x_2 \ ... \ x_n] \ (d \times n)$
- The eigenvectors of $S = X_z X_z^T (d \times d)$ are the principal components of the data, ordered by decreasing eigenvalues

$$Su_i = \lambda_i u_i \rightarrow SU = U\Lambda$$

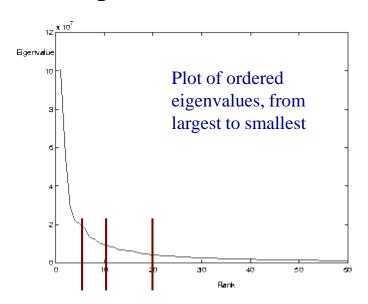
where U is the matrix of eigenvectors (columns of U) and Λ is a diagonal matrix of eigenvalues diag $(\lambda_i, \lambda_2, ..., \lambda_d)$

- The eigenvector u_i associated with the largest eigenvalue λ_i is the first principal component
- If rank(S) < d, then some eigenvalues will be zero



PCA for dimensionality reduction

- So the eigenvalues can give clues to the inherent dimensionality of the data, or at least provide a way to more efficiently approximate high-dimensional with lowerdimensional feature vectors
- For example:



60-dimensional data (60 eigenvectors and eigenvalues)

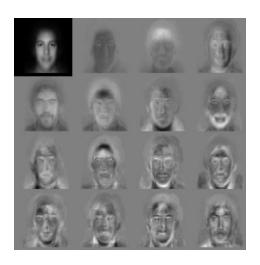
Many of the eigenvalues are small, meaning that their associated eigenvectors don't contribute much to the representation of the data

We can choose a cutoff - say, only use the first 20 eigenvectors (or 10, or 5)

 A well-known technique for face recognition based on computing eigenvectors of a training set of face images, i.e., Eigenfaces



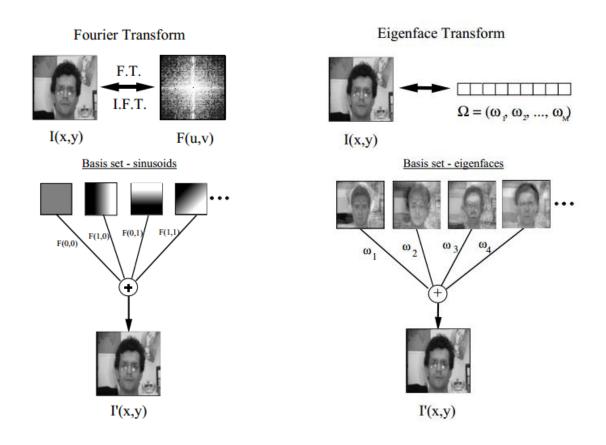
Eigenfaces 1



Eigenfaces 2

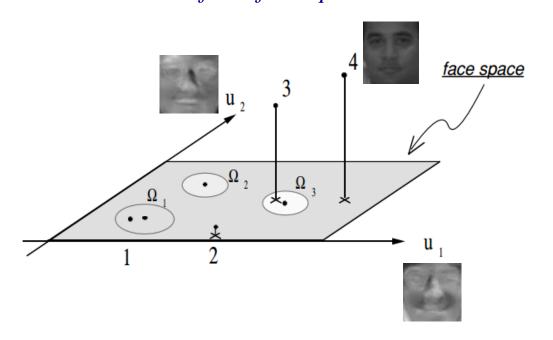
Keep in mind: an image is just an N-dimensional point or vector (where rows \times cols = N)

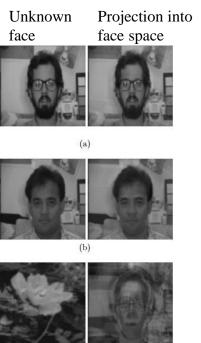
• Eigenvectors (eigenfaces) can be thought of as *basis vectors* for reconstructing data (face images)



- The Eigenfaces span a (relatively) low-dimensional *face space*, representing all possible face images
- A new (unknown) face image is projected into the face space (reconstructed by the Eigenfaces)

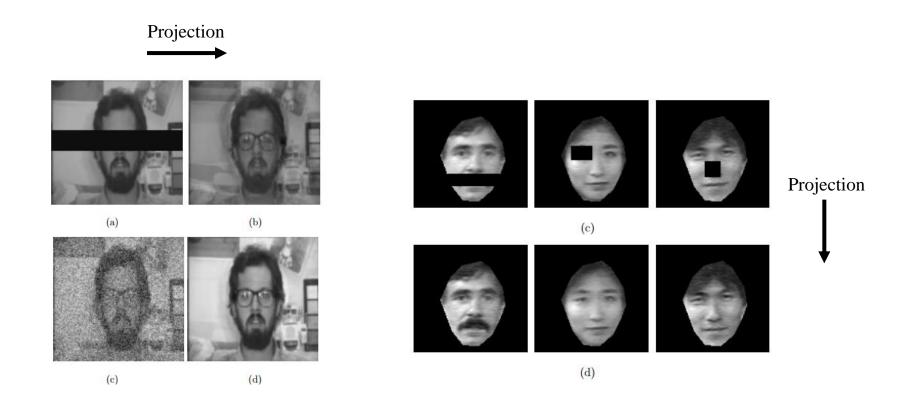
 The distance between the face image and its reconstruction is the distance from face space





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- The *distance from face space* measure could be used for face detection: Does this image (or part of an image) look like a face?
- If yes, then use the Eigenface weights (projections) as features for classification (classes Ω_1 , Ω_2 , etc.)
 - E.g., using k-NN

