# Machine Learning

Non-parametric Classification: k-NN Method and its variants. Common Metrics. Classification Mean Error.

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ML Research







Classification Mean Test Error





- O Classification Mean Test Error
- 2 Euclidean and Manhattan Distance





- O Classification Mean Test Error
- 2 Euclidean and Manhattan Distance
- **3** 1-NN





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- **3** 1-NN
- 0 k-NN





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- 2 Euclidean and Manhattan Distance
- **3** 1-NN
- 4 k-NN
- Weighted k-NN





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- Weighted k-NN
- Template Selection





#### Mean test error

- X set of objects descriptions, Y set of objects labels
- Unknown target dependency: mapping  $y: X \to Y$
- Finite training set:  $X^m = \{(x_1, y_1), \dots, (x_m, y_m)\}$ , so as  $y_i = y(x_i)$
- Finite test set:  $X_t^q = \{(x_1, y_1), \dots, (x_q, y_q)\}$ , so as  $y_i = y(x_i)$





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#### Mean test error for classification

The main goal is to train an algorithm  $a: X \to Y$  on the train set  $X^m$  so as the **mean error** on the test set is minimal:  $R(a, X_t^q) = \frac{1}{q} \sum_{i=1}^q [a(x_i) \neq y(x_i)] \to min_a$ 





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#### Common metrics

- Object  $x \in X$  is represented in the  $\mathbb{R}^n$  space:  $x = (x^1, \dots, x^n) n$ -dimensional vector
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#### Euclidean metric

Euclidean, or  $L_2$ -distance, between 2 points x and y from  $\mathbb{R}^n$  is:

$$d_2(x,y) = ||x-y||_2 = \sqrt{\sum_{i=1}^n (x_i - y_i)^2}$$





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#### Manhattan metric

Manhattan, or  $L_1$ -distance, between 2 points x and y from  $\mathbb{R}^n$  is:

$$d_1(x,y) = ||x-y||_1 = \sum_{i=1}^n |x_i - y_i|$$





# Parametric and non-parametric machine learning methods

#### Parametric methods

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- Parameters are found as a solution of the optimization problem





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### Non-parametric methods

Nonparametric methods are methods that are not parametric methods.

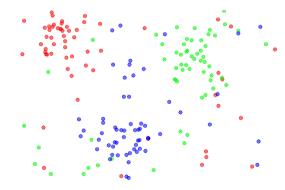
• Examples: Metric algorithms, kernel methods

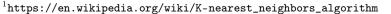




# Basic Assumption

- "Close" objects usually lie in the same class
- Proximity is specified by the metric
- Typical example <sup>1</sup>





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- Method parameter: metric
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- Ease of implementation (there is no training procedure as such in the naive implementation)
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### Disadvantages

- Instability to outliers
- Ambiguity of classification at equal distances to two objects
- The need to store the entire training set
- The search algorithm is computationally complex (if the training sample is quite large)
- Distance value is not taken into account

WAP

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# k-nearest neighbors (k-NN, kNN) method

- ullet Method parameter: metric,  ${f k}$
- Algorithm: using a given metric, we search for the **k** closest objects in the training set and classify the object as the **majority class** of the **k** objects

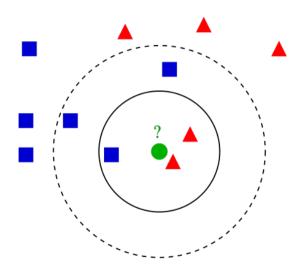
### Advantages

- Ease of implementation
- Good interpretability
- Parameter k can be optimized using cross-validation

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# k-NN method







# Weighted k-NN method

- Method parameters: metric, k, weights
- Algorithm: using a given metric, we look for the k closest objects in the training set and classify the object by weighted voting

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- Fixed object weights





# Weighted k-NN among a set of templates

- Method parameters: metric, k, weights, template selection method
- Algorithm: using a given metric, we look for the k closest objects among the templates selected from the training set and classify the object by weighted voting

### Advantages

- Ease of implementation
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- Parameter k can be optimized using cross-validation

### Disadvantages

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- Classification ambiguity at equal distances to two objects
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# Template selection

#### Task

Get approximately the same quality of the algorithm with less stored data.

It is even possible to obtain an improvement in quality because outliers will be removed during the template selection process.

#### Ideas

- Object clustering
- Greedy algorithm





# Template selection by k-means clustering method

#### Task

$$V = \sum_{i=1}^{k} \sum_{x \in S_i} (x - \mu_i)^2 \to \min_{S_i},$$

where k is the number of clusters,  $S_i$  is the resulting clusters,  $\mu_i$  is the center of mass of the  $S_i$  cluster.

### Algorithm

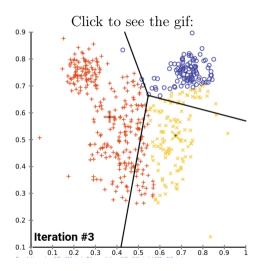
- $\bullet$  k elements are randomly selected from the sample and declared as centroids
- ② For fixed **centroids**, each sample element belongs to one of the clusters
- 3 For fixed **clusters**, centroids are calculated
- Steps 2 and 3 are repeated until convergence (or exhausting the computation/time budget)

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# Visualization of k-means







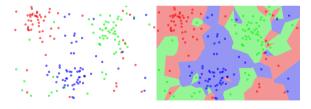
- 1st image: dataset
- 2nd image: 1-NN w/o template selection
- 3rd image: templates
- 4th image: 1-NN w/ template selection







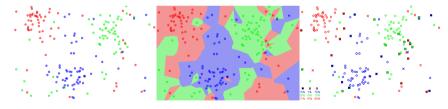
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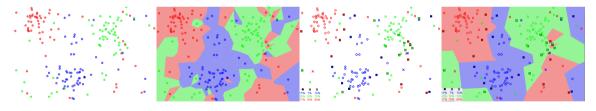
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### Additional modification: RadiusNN

#### Idea

Sometimes it makes sense to look for neighbors at a distance no greater than some radius r

#### Parameter r

Instead of the input parameter for the number of neighbors, the radius is used





# Mandatory external links to read

- Read the Introduction to K-Nearest Neighbor (kNN)
  - ▶ Main <u>source</u> ("K Nearest Neighbor (KNN) in R"section is fully optional)
  - ▶ Additional <u>one</u> (mostly section "*The KNN Algorithm*"), including the section 3.5 from "The Hundred-Page Machine Learning Book" (see "*References*" course page).





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- The method has a large number of variations for customization





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  - ▶ Number of nearest neighbors
  - ▶ Weights in the weighted version of the method





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  - ▶ Algorithm for selecting templates





# Thank you!



