

# Machine Learning and Intelligent Systems

K-Nearest Neighbors

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# Definition

#### Intuition

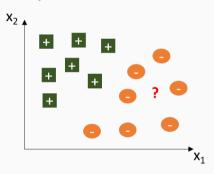
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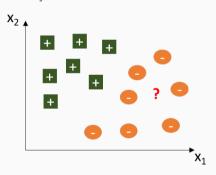


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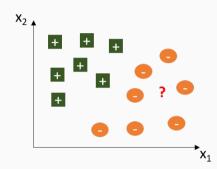


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**Assumption:** Similar inputs have similar outputs

# **Data Assumptions:**

Similar x have similar y

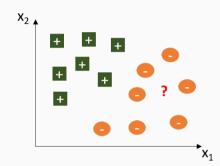


## **Data Assumptions:**

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 $y \in \mathbb{R}$  - regression

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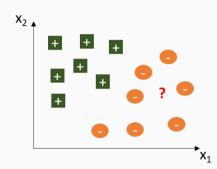
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#### **Definition of the** *k* **nearest neighbors:**

Given a test point  $\mathbf{x}^*$ , let us denote  $\mathcal{S}_{\mathbf{x}^*}$  as its set of k nearest neighbors. Formally:



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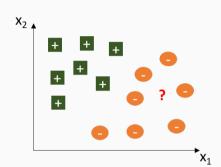
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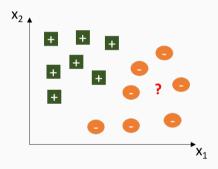
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$$\mathcal{S}_{\mathbf{x}^*} \subseteq \mathcal{D}$$
 .s.t.  $|\mathcal{S}_{\mathbf{x}^*}| = k$ 

$$\forall (\mathbf{x}', \mathbf{y}') \in \mathcal{D} \setminus \mathcal{S}_{\mathbf{x}^*} \operatorname{dist}(\mathbf{x}^*, \mathbf{x}') \ge \max_{(\mathbf{x}'', \mathbf{y}'') \in \mathcal{S}_{\mathbf{x}^*}} \operatorname{dist}(\mathbf{x}^*, \mathbf{x}'')$$

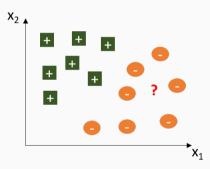
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**Regression:** The output is the average of the values of k nearest neighbors

$$\hat{\mathbf{y}} = h(\mathbf{x}^*) = \frac{1}{k} \sum_{(\mathbf{x}_i'', \mathbf{y}_i'') \in \mathcal{S}_{\mathbf{x}^*}} \mathbf{y}_i'' \tag{1}$$



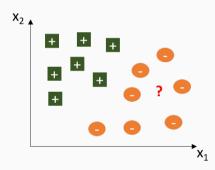
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 (1)

**Classification:** An unseen point  $x^*$  is classified by a majority vote of its k nearest neighbors:

$$\hat{\mathbf{y}} = h(\mathbf{x}^*) = \mathsf{mode}(\{\mathbf{y}_i'' : (\mathbf{x}_i'', \mathbf{y}_i'') \in \mathcal{S}_{\mathbf{x}^*}\}) \tag{2}$$



#### **Distance Metric**

The definition of the set  $\mathcal{S}_{x^*}$  is strongly dependent on the distance metric (dist) that is used. The better the chosen metric reflects the similarity among points the better the resulting model will be.

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One of the most common metrics used is the Minkowski distance:

$$\operatorname{dist}(\mathbf{x}, \mathbf{z}) = \left(\sum_{j=1}^{D} |x_j - z_j|^p\right)^{1/p} \tag{3}$$

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The main advantage of this metric is its generality. It contain many well-known distances as special cases.

**Question:** Can you identify what case is p = 2?

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It does not attempt to construct a general internal model, but simply stores instances of the training data.

This type of approach is referred as instance-based learning or non-generalizing learning.

Question: Can you spot the differences?

**Placement Exam** 



# **Summary: Notation**

Symbol	Reads as
X	Input variable $(\mathbb{R}^D)$
$\mathbf{x}_{i}$	$i^{th}$ feature vector. Observed value of $X$ .
$\mathbf{X} = (\mathbf{x}_1, \mathbf{x}_2, \dots, \mathbf{x}_N)^T$	Matrix of $N$ input $D$ -dimensional vectors $\mathbf{x}_i$
×j	$j^{th}$ element of the $i^{th}$ input vector $\mathbf{x}_i$ , i.e. $\mathbf{x}_i^j$
Y	Output variable $(\mathcal{C})$
y <sub>i</sub>	i <sup>th</sup> output label
$\mathbf{y} = (\mathbf{y}_1, \dots, \mathbf{y}_N)^T$	Observed vector of outputs $y_i$
x*	Test point (unseen data)
ŷ	Prediction for x*

Table 1: Different notation for the input and output variables

# Further Reading and Useful Material

Source	Notes
The Elements of Statistical Learning	Ch. 2