## MLCC 2017 Local Methods and Bias Variance Trade-Off

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#### **About this class**

- 1. Introduce a basic class of learning methods, namely local methods.
- 2. Discuss the fundamental concept of **bias-variance** trade-off to understand parameter tuning (a.k.a. model selection)

### **Outline**

Learning with Local Methods

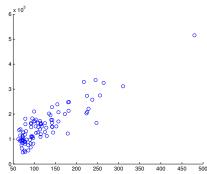
From Bias-Variance to Cross-Validation

What is the price of one house given its area?

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Area $\left(m^2 ight)$	Price (€)
$x_1 = 62$	$y_1 = 99,200$
$x_2 = 64$	$y_2 = 135,700$
$x_3 = 65$	$y_3 = 93,300$
$x_4 = 66$	$y_4 = 114,000$
	:

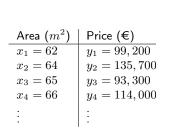


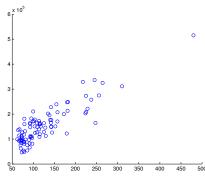
Let S the houses example dataset (n = 100)

$$S = \{(x_1, y_1), \dots, (x_n, y_n)\}\$$

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Let S the houses example dataset (n = 100)

$$S = \{(x_1, y_1), \dots, (x_n, y_n)\}\$$

Given a new point  $x^*$  we want to predict  $y^*$  by means of S.

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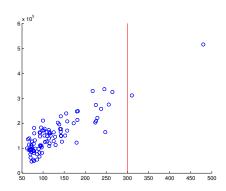
# **Example**

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$x_{95} = 310$ $x_{96} = 480$	$y_{95} = 311,200$ $y_{96} = 515,400$
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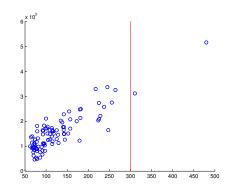


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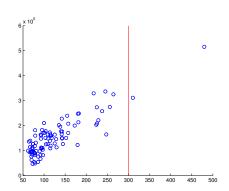


What is its price?

**Nearest Neighbor**:  $y^*$  is the same of the closest point to  $x^*$  in S.

$$y^* = 311,200$$

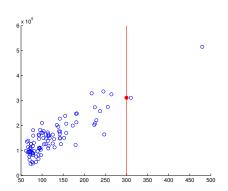
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$$y_* = y_j$$
  $j = \arg\min_{i=1,...,n} ||x - x_i||$ 

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16

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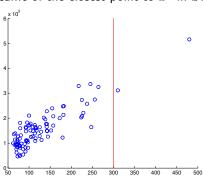
In general let  $d: \mathbb{R}^D \times \mathbb{R}^D$  a distance on the input space, then

$$f(x) = y_j$$
  $j = \arg\min_{i=1,\dots,n} d(x, x_i)$ 

### **Extensions**

Nearest Neighbor takes  $y^*$  is the same of the closest point to  $x^*$  in S.

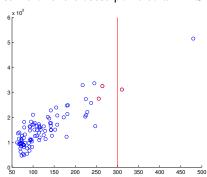
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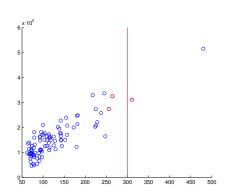


Can we do better? (for example using more points)

**K-Nearest Neighbor**:  $y^*$  is the mean of the values of the K closest point to  $x^*$  in S. If K=3 we have

$$y^* = \frac{274,600 + 324,900 + 311,200}{3} = 303,600$$

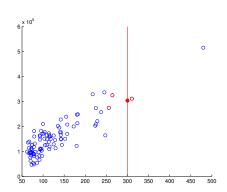
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## K-Nearest Neighbors (cont.)

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▶ Computational cost  $O(nD + n \log n)$ : compute the n distances  $||x - x_i||$  for  $i = \{1, ..., n\}$  (each costs O(D)). Order them  $O(n \log n)$ .

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- ▶ **General Metric** d f is the same, but  $j_1, \ldots, j_K$  are defined as  $j_1 = \arg\min_{i \in \{1, \ldots, n\}} d(x, x_i)$  and  $j_t = \arg\min_{i \in \{1, \ldots, n\} \setminus \{j_1, \ldots, j_{t-1}\}} d(x, x_i)$  for  $t \in \{2, \ldots, K\}$

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Closer points to  $x^*$  should influence more its value PARZEN WINDOWS:

$$\hat{f}(x) = \frac{\sum_{i=1}^{n} y_i k(x, x_i)}{\sum_{i=1}^{n} k(x, x_i)}$$

where k is a similarity function

- $k(x,x') \geq 0$  for all  $x,x' \in \mathbb{R}^D$
- $\blacktriangleright \ k(x,x') \to 1 \ \text{when} \ \|x-x'\| \to 0$
- $k(x, x') \to 0$  when  $||x x'|| \to \infty$

### Examples of k

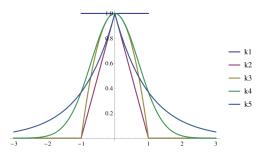
$$\blacktriangleright k_1(x,x') = \operatorname{sign}\left(1 - \frac{\|x - x'\|}{\sigma}\right)_{\perp}$$
 with a  $\sigma > 0$ 

$$\blacktriangleright \ k_2(x,x') = \left(1 - \frac{\|x - x'\|}{\sigma}\right)_\perp$$
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• 
$$k_3(x,x')=\left(1-\frac{\|x-x'\|^2}{\sigma^2}\right)_+$$
 with a  $\sigma>0$ 

• 
$$k_4(x,x') = e^{-\frac{\|x-x'\|^2}{2\sigma^2}}$$
 with a  $\sigma > 0$ 

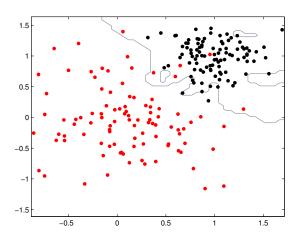
• 
$$k_5(x,x')=e^{-\frac{\|x-x'\|}{\sigma}}$$
 with a  $\sigma>0$ 



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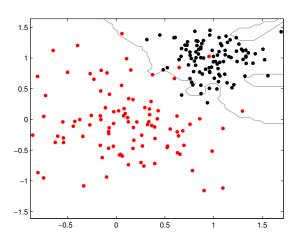
### K-NN example

 $K\operatorname{-Nearest}$  neighbor depends on K. When K=1



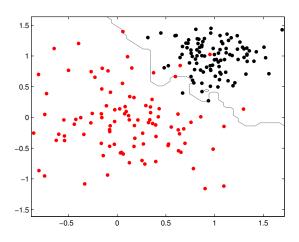
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 $K\operatorname{-Nearest}$  neighbor depends on K. When K=2

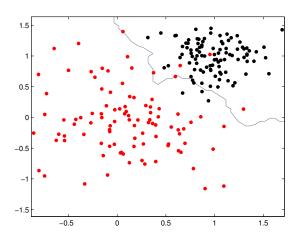


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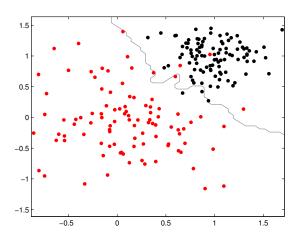
 $K\operatorname{-Nearest}$  neighbor depends on K. When K=3



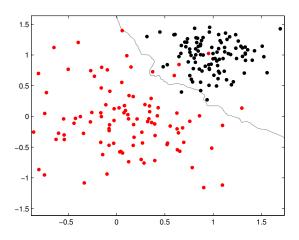
 $K\operatorname{-Nearest}$  neighbor depends on K. When K=4



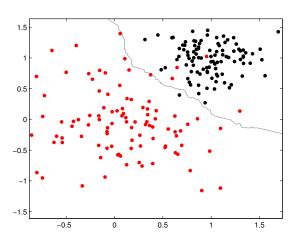
 $K\mbox{-Nearest}$  neighbor depends on  $K\mbox{.}$  When K=5



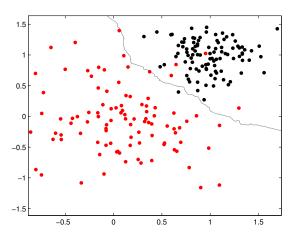
 $K ext{-Nearest neighbor depends on }K.$  When K=9



 $K\mbox{-Nearest}$  neighbor depends on K. When K=15



K-Nearest neighbor depends on K.



Changing K the result changes a lot! How to select K?

#### **Outline**

Learning with Local Methods

From Bias-Variance to Cross-Validation

- ▶  $S = (x_i, y_i)_{i=1}^n$  training set. Name  $Y = (y_1, \dots, y_n)$  and  $X = (x_1^\top, \dots, x_n^\top)$ .
- $lackbox{W} K \in \mathbb{N}$  hyperparameter of the learning algorithm
- $\hat{f}_{S,K}$  learned function (depends on S and K)

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The expected loss  $\mathcal{E}_K$  is

$$\mathcal{E}_K = \mathbb{E}_S \mathbb{E}_{x,y} (y - \hat{f}_{S,K}(x))^2$$

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Ideally! (In practice we don't have access to the distribution)

- ► We can still try to understand the above minimization problem: does a solution exists? What does it depend on?
- Yet, ultimately, we need something we can compute!

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$$\mathcal{E}_K(x) = \mathbb{E}_S(f_*(x) - \hat{f}_{S,K}(x))^2 + \sigma^2$$

. . .

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$$\mathcal{E}_K(x) = \underbrace{(f_*(x) - \mathbb{E}_X \tilde{f}_{S,K}(x))^2}_{\text{bias}} + \underbrace{\mathbb{E}_S(\tilde{f}_{S,K}(x) - \hat{f}_{S,K}(x))^2 + \sigma^2}_{\text{variance}}$$

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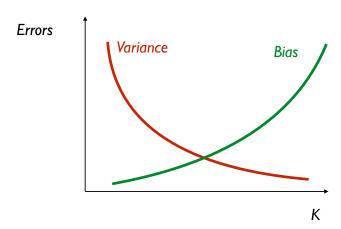
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#### Bias Variance trade-off



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- ▶ an optimal parameter exists and
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How to choose K in practice?

▶ Idea: train on some data and validate the parameter on new unseen data as a proxy for the ideal case.

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The above procedure can be repeated to augment stability and  ${\cal K}$  selected to minimize error over trials.

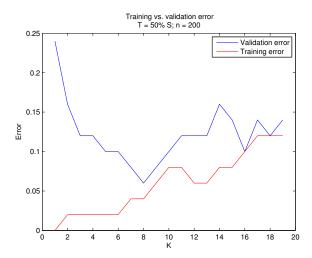
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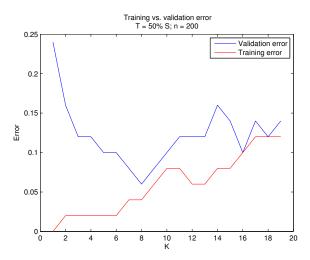
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There are other related parameter selection methods (k-fold cross validation, leave-one out...).

# **Training and Validation Error behavior**



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$$\hat{K} = 8$$
.

### Wrapping up

In this class we made our first encounter with learning algorithms (local methods) and the problem of tuning their parameters (via bias-variance trade-off and cross-validation) to avoid overfitting and achieve generalization.

#### **Next Class**

High Dimensions: Beyond local methods!

