# **Kernel Smoothing**

Lecture 23

Parametric vs non-parametric methods

Kernel density estimation / smoothing

Kernel density classification

Kernel regression

Kernel methods and the Kernel Trick (a preamble to support vector machines)

**Kernel** (first definition): a non-negative function that integrates to one

A **window** or weighting function for **local** computation / neighborhood methods

**Kernel** (second definition): A **similarity** function

Inner products in some feature space

#### **Parametric Methods**

Models governed by a vector  $\mathbf{w}$  of parameters

During training w is estimated

Training data are discarded and predictions made only with  $\boldsymbol{w}$ 

Number of parameters is fixed

Typically slower to train, faster to predict

Example: linear regression

$$\hat{f}(x) = \sum_{i=0}^{p} w_i x_i$$
Model weights learned from training data

#### **Non-Parametric Methods**

**All training data** (or a subset) are used in the prediction phase

These are **memory-based** or **instance-based** methods

Typically faster to train, slower to predict

Number of parameters (may) vary with the size of the training data

Example: k-nearest neighbor regression

$$\hat{f}(\mathbf{x}) = \frac{1}{k} \sum_{i \in \mathcal{N}_k(\mathbf{x})} \mathbf{y}_i$$

where

 $\mathcal{N}_k(x) \triangleq \{ \text{k nearest neighbors} \}$ 

Training data

Bishop, Pattern Recognition and Machine Learning, 2006

#### Supervised Learning Techniques

- Linear Regression
- K-Nearest Neighbors
  - Perceptron
  - Logistic Regression
  - Fisher's Linear Discriminant
  - Linear Discriminant Analysis
  - Quadratic Discriminant Analysis
  - Naïve Bayes
- Decision Trees and Random Forests
- Ensemble methods (bagging, boosting, stacking)
- Neural Networks

Appropriate for:

Classification

Regression

Can be used with many machine learning techniques

#### Non-Parametric Supervised Learning Techniques

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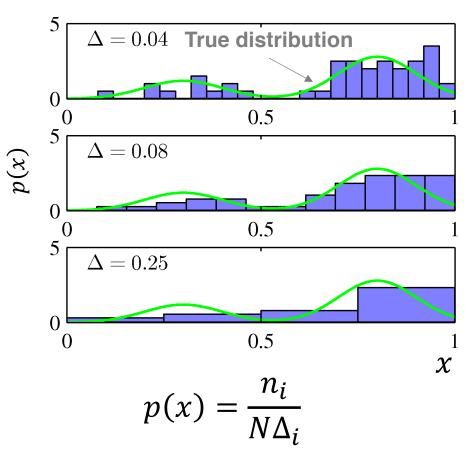
Appropriate for:

- Classification
- Regression

Can be used with many machine learning techniques

#### **Histogram Density Estimation**

Histogram



 $n_i$  = # observations of x falling in bin i

N = total # observations

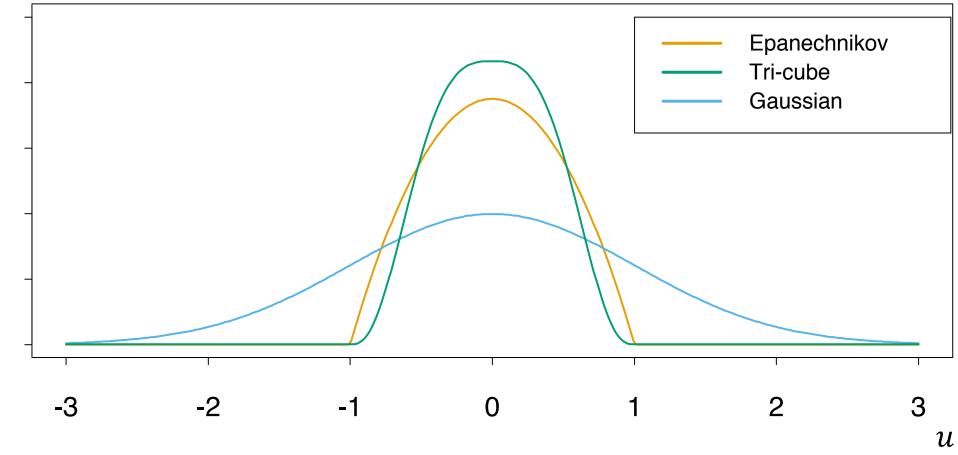
 $\Delta_i$  = width of bin i

Highly dependent on the choice of bin width,  $\Delta_i$ 

Has discontinuities at the bin edges

Local neighborhoods do appear to be helpful

#### Kernel **Functions** (window kernels)



Satisfy two properties:

$$k(u) \ge 0$$
$$\int k(u)du = 1$$

**Epanechnikov** 

$$k(u) = \frac{3}{4}(1 - u^2)$$
$$|u| \le 1$$

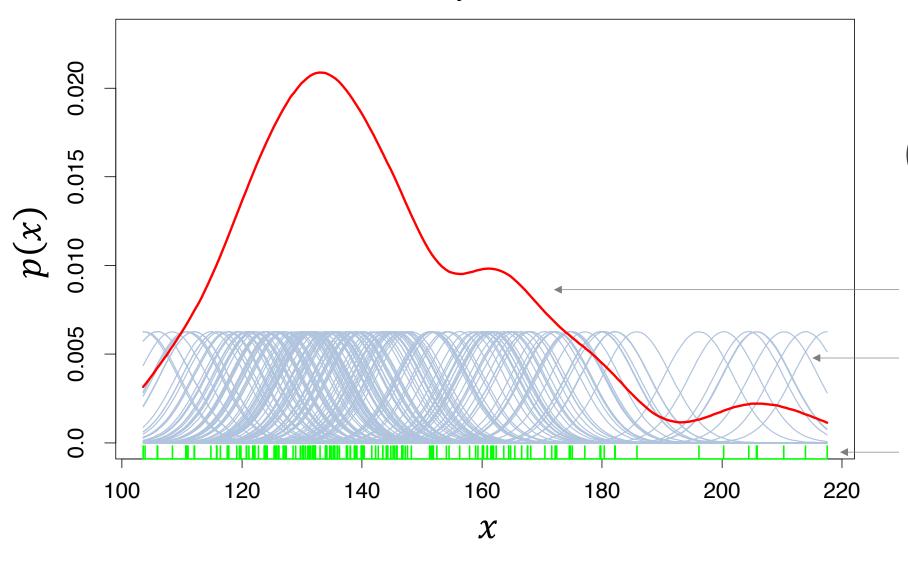
Tri-cube

$$k(u) = \frac{3}{4}(1 - u^2) \qquad k(u) = \frac{70}{81}(1 - |u^3|)^3 \qquad k(u) = \frac{1}{\sqrt{2\pi}}e^{-\frac{1}{2}u^2}$$
$$|u| \le 1 \qquad \qquad |u| \le 1 \qquad \qquad -\infty < u < \infty$$

Gaussian

$$k(u) = \frac{1}{\sqrt{2\pi}} e^{-\frac{1}{2}u^2}$$
$$-\infty < u < \infty$$

# Kernel Density Estimation a. k. a. Parzen Window Density Estimation



- Center a kernel function at each observation (x-value)
- Estimate p(x) by averaging across all the kernels at x

Density estimate

Kernel functions (Gaussian kernel shown)

Observations (x-values)

### **Kernel Density Estimation**

Center the kernel function at each x-value in the dataset:

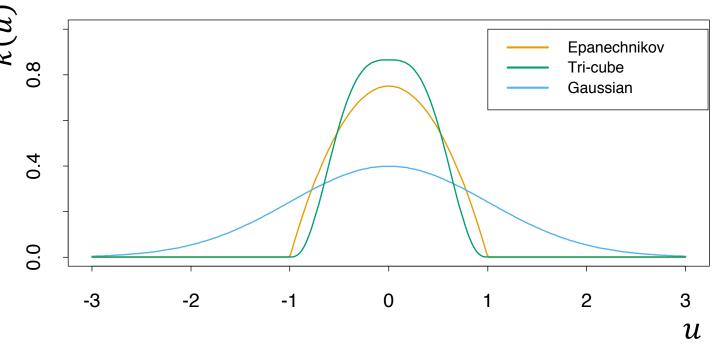
$$k(x - x_n) \qquad n = 1, 2, \dots, N$$

Average over all of the kernel functions to get the density estimate:

$$p(x) = \frac{1}{N} \sum_{n=1}^{N} k(x - x_n)$$

Note: we can scale the width of the kernel function with a scale factor, *h*:

$$k\left(\frac{x-x_n}{h}\right)$$

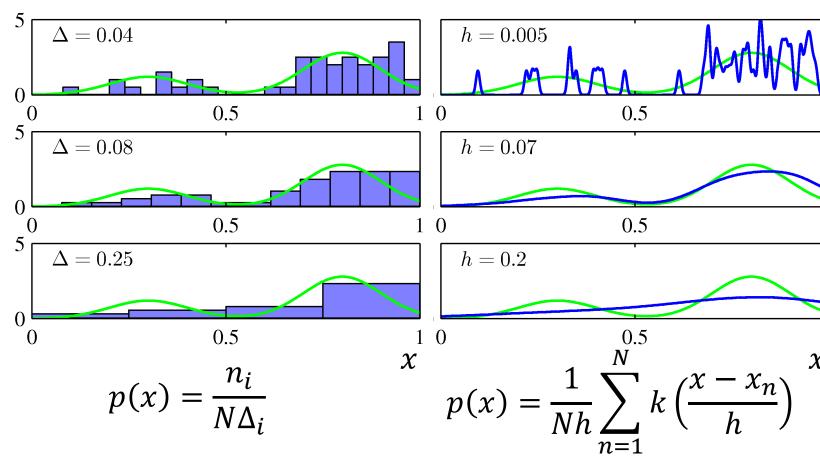


For kernel functions with **finite domains**, this means that each observation, x, will only affect the density estimate in a **neighborhood** close to the center of the kernel

#### **Kernel Density Estimation**



Kernel Density Estimation



Requires tuning h, the kernel width parameter

Computational cost of evaluating this density grows linearly with the size of the data

 $n_i$  = # observations of x falling in bin i

N = total # observations

 $\Delta_i$  = width of bin i

 $x_n$  = The n<sup>th</sup> observation of x

k = kernel function

h =width of the kernel

Bishop, Pattern Recognition and Machine Learning, 2006

p(x)

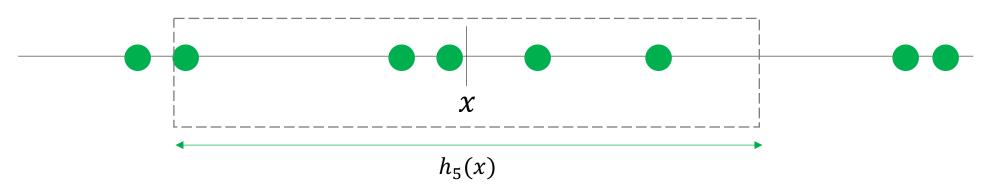
### **Nearest Neighbor Density Estimation**

Instead of fixing the window width and averaging the points within, we fix the number of points and vary the window width

1

For a point, x, determine the width of a window,  $h_K(x)$ , centered at x, that is just large enough to contain K points

Example for k = 5:



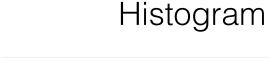
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The density estimate is computed as: p

$$p(x) = \frac{K}{Nh_K(x)}$$

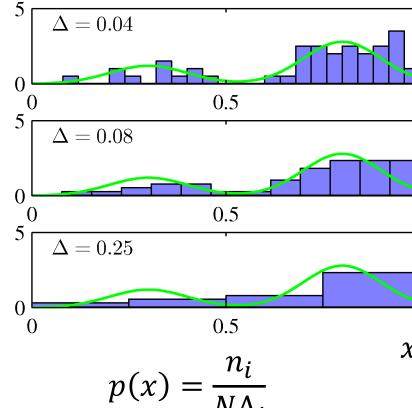
N = total number of observations

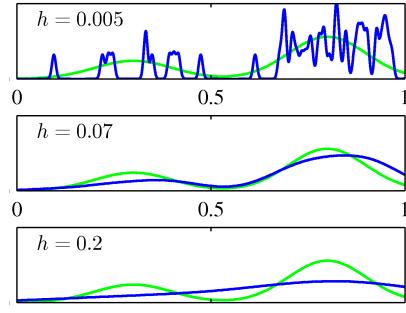
### **Kernel Density Estimation**

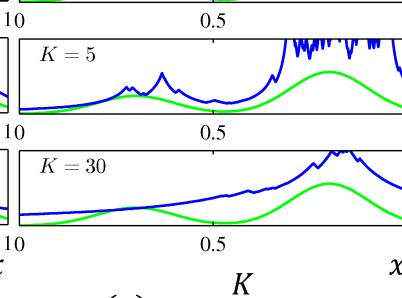


Kernel Density Estimation

Nearest Neighbor **Density Estimation** 







$$(x) = \frac{n_i}{N\Delta_i}$$

$$p(x) = \frac{1}{N} \sum_{n=1}^{N} \frac{1}{h} k \left( \frac{x - x_n}{h} \right)$$

$$p(x) = \frac{K}{Nh_K(x)}$$

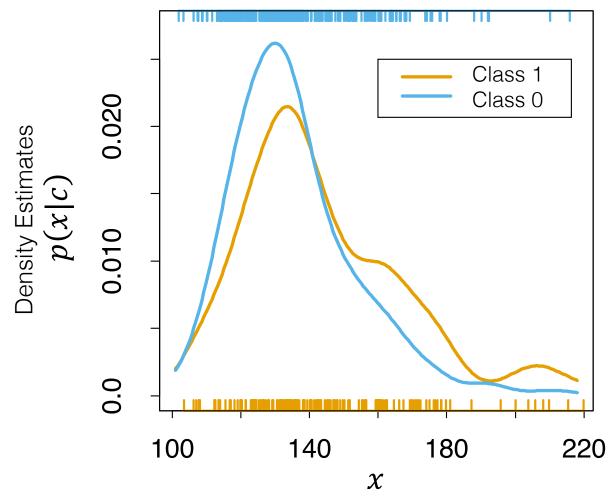
 $n_i = \#$  observations of x falling in bin i N = total # observations $\Delta_i$  = width of bin i

 $x_n$  = The n<sup>th</sup> observation of xk = kernel functionh =width of the kernel

K = # points in neighborhood  $h_K(x) = \text{minimum width containing}$ K observations

Bishop, Pattern Recognition and Machine Learning, 2006

p(x)



p(x|c) = class conditional distribution where c represents the class

Bayes' Rule 
$$p(c|x) = \frac{p(x|c)p(c)}{p(x)}$$

Since there are two classes:

$$p(x) = p(x|c_1)p(c_1) + p(x|c_0)p(c_0)$$

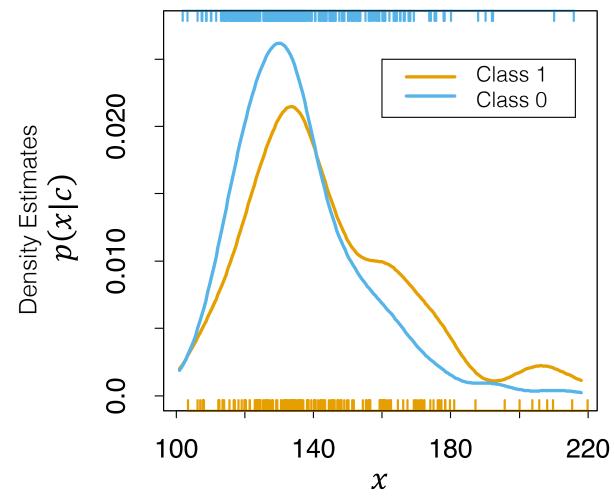
Therefore, we have:

$$p(c|x) = \frac{p(x|c)p(c)}{p(x|c_1)p(c_1) + p(x|c_0)p(c_0)}$$

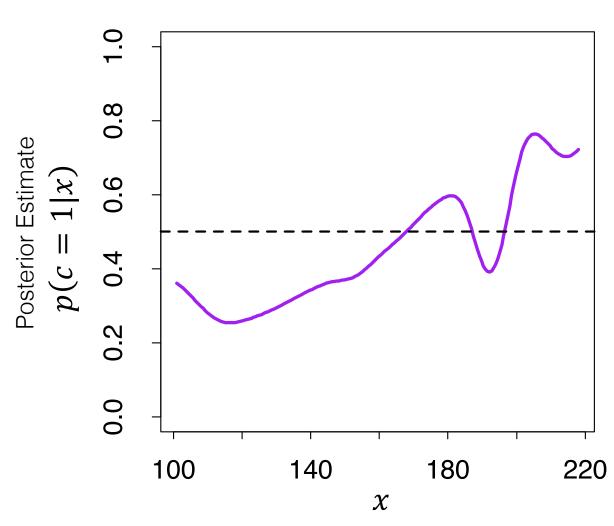
Class priors, p(c), could simply be the fraction of observations from each class

We can construct a classifier:

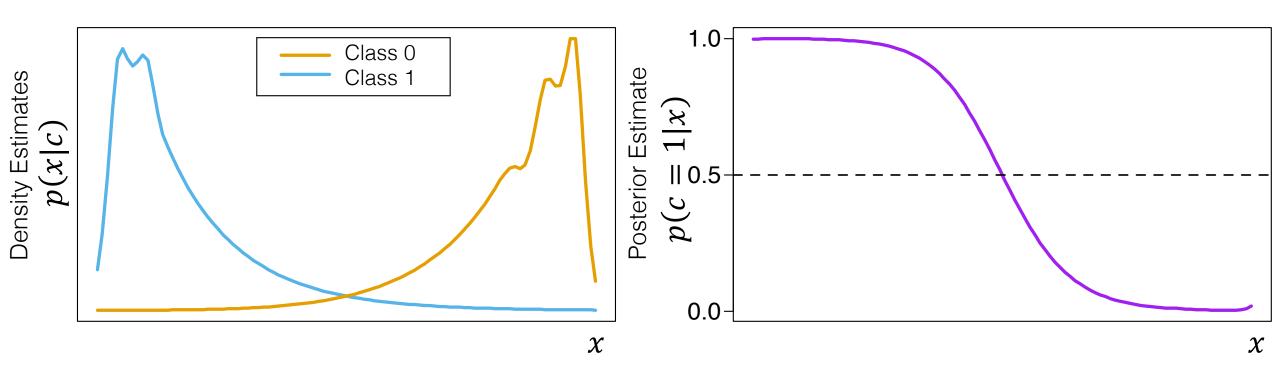
Predict class 
$$c = \hat{f}(x) = \arg \max_{c} p(c|x)$$



p(x|c) = class conditional distribution where c represents the class



p(c) = class prior (# of samples from that class in training set)
Hastie, Tibshirani, and Friedman, The Elements of Statistical learning, 2001



We may not need all of the interesting structure of each class density, but will only need to estimate the posterior near the decision boundary (or boundaries)

Hastie, Tibshirani, and Friedman, The Elements of Statistical learning, 2001

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...and recall Naïve Bayes

#### Assumption: Given the class, the features are independent

$$P(y = i | x_1, x_2, ..., x_D) \propto P(y = i) \prod_{j=1}^{D} P(x_j | y = i)$$

This implies we estimate the density of each feature **separately** 

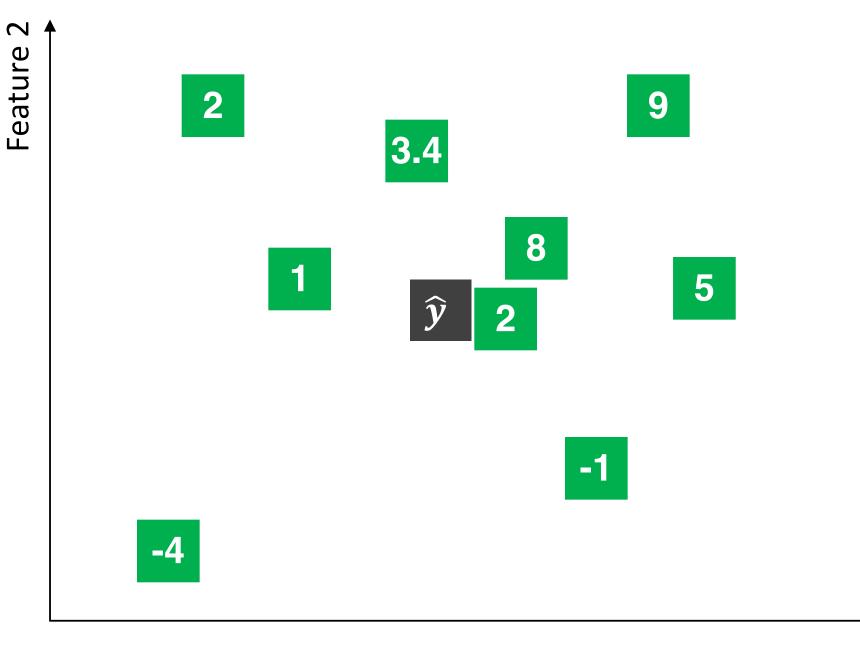
We can easily apply Naïve Bayes using kernel density estimates for each feature, then multiply them together

Hastie, Tibshirani, and Friedman, The Elements of Statistical learning, 2001

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(and smoothing)

## K Nearest Neighbor Regression



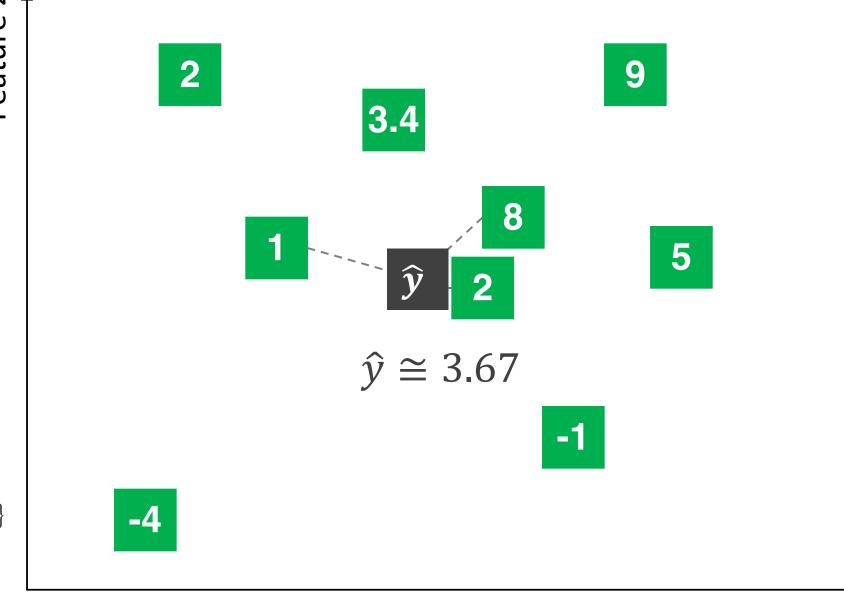
Feature 1

## K Nearest Neighbor Regression

Example for k = 3

$$\hat{y} = \frac{1}{k} \sum_{i \in \mathcal{N}_k(x)} y_i$$
Training data

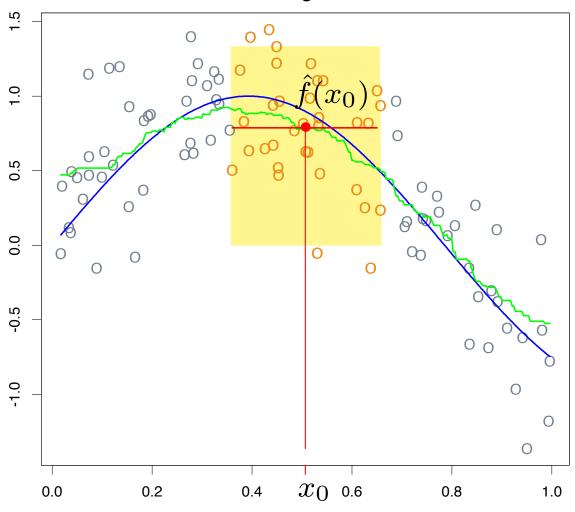
where  $\mathcal{N}_k(x) \triangleq \{k \text{ nearest neighbors}\}$ 



#### Feature 1

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#### Nearest-Neighbor Kernel



This produces a bumpy, discontinuous estimate,  $\hat{f}(x)$ 

This is also a **local estimate** of the average of this function

This discontinuity can be avoided by using a **smoothing kernel** 

Target function (blue); prediction (green); observations (circles); observations included in estimate of  $\hat{f}(x_0)$  (orange circles)

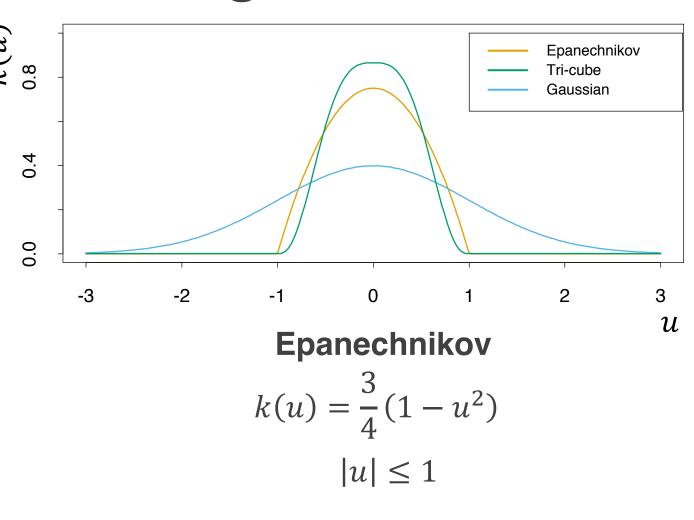
### Nadaraya-Watson Kernel Regression

Local kernel weighted average regression

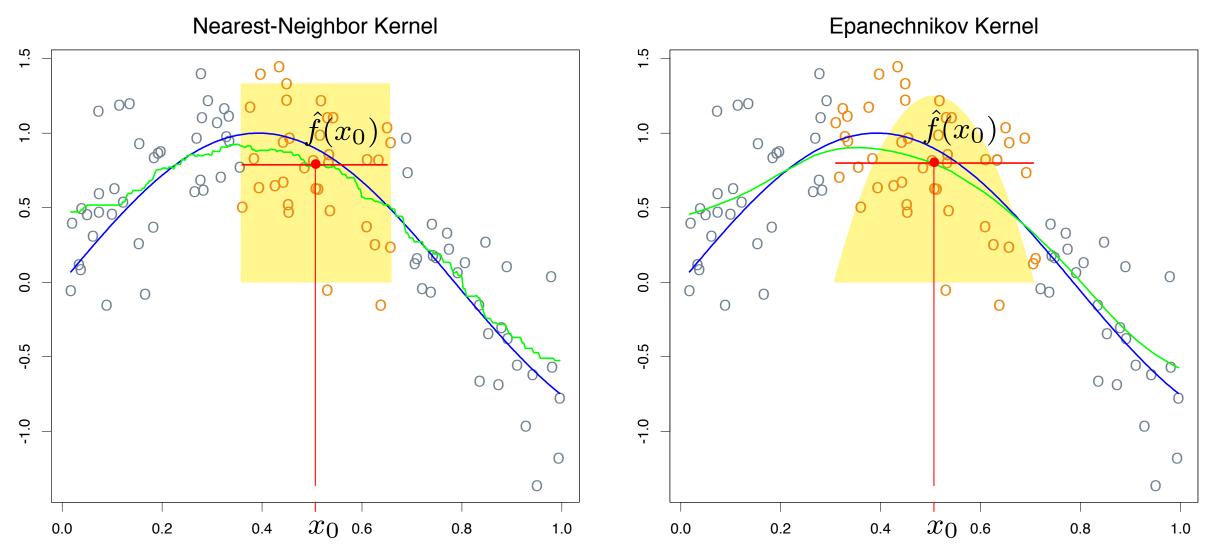
Weight the average for a given x value, by its neighbors, as defined by your choice of kernel

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

For kernels with a finite domain, this only uses a subset of the training data

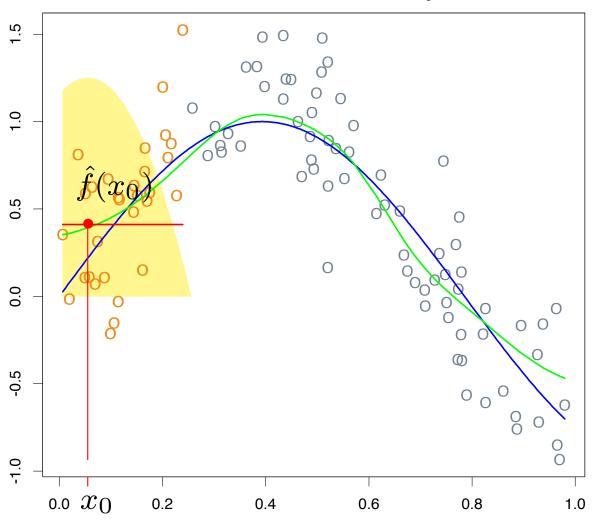


#### **Continuous, smooth function**



Target function (blue); prediction (green); observations (circles); observations included in estimate of  $\hat{f}(x_0)$  (orange circles)

#### N-W Kernel at Boundary



At the boundaries these estimates can be badly biased due to kernel asymmetry
This can also happen in the interior if the points aren't equally spaced

One solution: fit straight lines instead of constants locally

This leads to **local linear regression** 

Target function (blue); prediction (green); observations (circles); observations included in estimate of  $\hat{f}(x_0)$  (orange circles)

#### **Local linear regression** least squares:

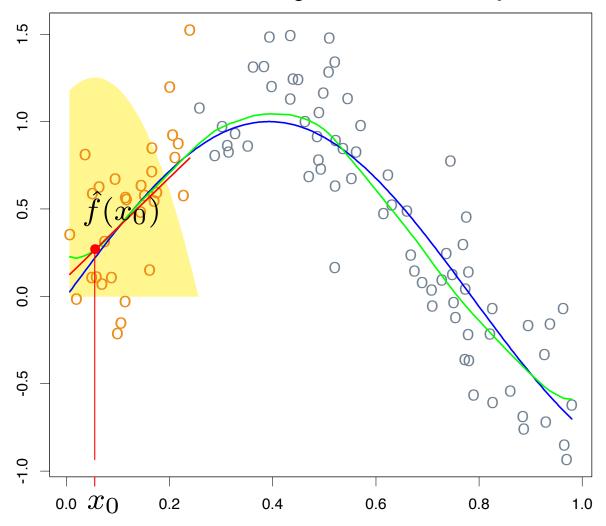
$$\min_{\boldsymbol{w}(\boldsymbol{x})} \sum_{i=1}^{N} k(\boldsymbol{x} - \boldsymbol{x}_i) [y_i - \boldsymbol{w}(\boldsymbol{x})^T \boldsymbol{x}_i]^2$$
Weighted Linear regression by kernel squared error

We fit this model on the local data to evaluate a **single point**:

$$\hat{f}(\boldsymbol{x}_0) = \boldsymbol{w}(\boldsymbol{x}_0)^T \boldsymbol{x}_0$$

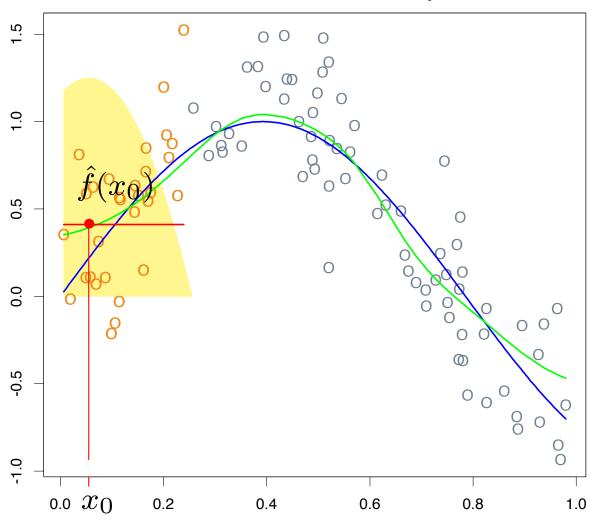
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#### Local Linear Regression at Boundary

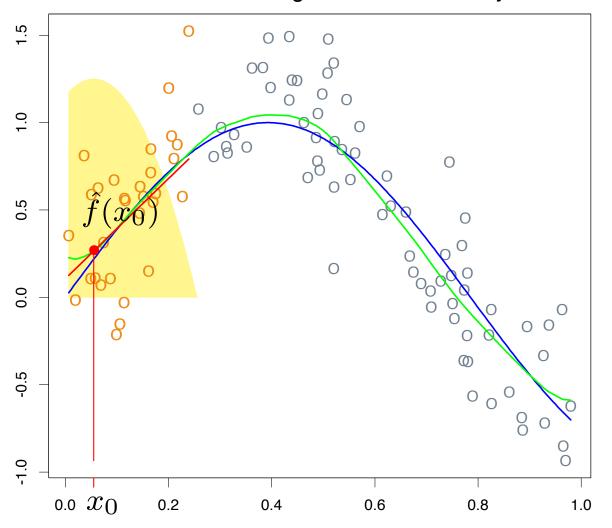


Hastie, Tibshirani, and Friedman, The Elements of Statistical learning, 2001

#### N-W Kernel at Boundary



#### Local Linear Regression at Boundary



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