SDS 383D, Exercises 3: Linear smoothing and Gaussian processes

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Curve fitting by linear smoothing

(A) Linear smoothing Suppose we want to estimate the value of the regression function y^* at some new point x^* , denoted $\hat{f}(x^*)$. Assume for the moment that f(x) is linear, and that y and x have already had their means subtracted, in which case $y_i = \beta x_i + \epsilon_i$.

Show that for the one-predictor case, your prediction $\hat{y}^* = \hat{f}(x^*) = \hat{\beta}x^*$ may be expressed as:

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

We can express $\hat{\beta}$ in matrix notation as:

$$\hat{\beta} = (X^T X)^{-1} X^T Y.$$

Plugging this in to find our predictor, \hat{y}^* we get,

$$\hat{y}^* = x^* (X^T X)^{-1} X^T Y$$

$$= \frac{X^T Y}{X^T X}$$

$$= \frac{\sum_{i=1}^n x_i x^* y_i}{\sum_{i=1}^n x_i^2}$$

The result is of the form $\hat{f}(x^*) = \sum_{i=1}^n w(x_i, x^*) y_i$ with the weights expressed as,

$$w(x_i, x^*) = \frac{x_i x^*}{\sum_{i=1}^n x_i^2}.$$

This resultant *smoother* has constant weights and produces predictions that lie on the line with slope $\hat{\beta}$. If the data is binned and locality preserved, this estimate is similar to a running line average.

$$w_K(x_i, x^*) = \left\{ \begin{array}{ll} 1/K, & x_i \text{ one of the } K \text{ closest sample points to } x^*\,, \\ 0, & \text{otherwise.} \end{array} \right.$$

With K-nearest-neighbor smoothing \hat{y}^* is essentially the arithmetic mean of the K y_i 's nearest x^* .

(B) A kernel function K(x) is a smooth function satisfying

$$\int_{\mathbb{R}} K(x)dx = 1 , \quad \int_{\mathbb{R}} xK(x)dx = 0 , \quad \int_{\mathbb{R}} x^2K(x)dx > 0 .$$

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A very simple example is the uniform kernel,

$$K(x) = \frac{1}{2}I(x) \quad \text{where} \quad I(x) = \left\{ \begin{array}{ll} 1, & |x| \leq 1 \\ 0, & \text{otherwise} \,. \end{array} \right.$$

Another common example is the Gaussian kernel:

$$K(x) = \frac{1}{\sqrt{2\pi}}e^{-x^2/2}$$
.

Kernels are used as weighting functions for taking local averages. Specifically, define the weighting function

$$w(x_i, x^*) = \frac{1}{h} K\left(\frac{x_j - x^*}{h}\right),$$

where h is the bandwidth. Using this weighting function in a linear smoother is called *kernel regression*. (The weighting function gives the unnormalized weights; you should normalize the weights so that they sum to 1.)

```
from __future__ import division
import numpy as np
import matplotlib.pyplot as plt
import sys
sys.path.append('../../scripts/')
from smoothers import kernel_smoother
```

```
# Create noisy data, y_i = sin(x_i) + e_i
x = np.linspace(0, 2*np.pi, num = 20)
y = np.sin(x) + np.random.normal(scale=0.2, size=x.shape)

# Create vector for fitting purposes
x_star = np.linspace(0, 2*np.pi)

# Instantiate array of bandwidths
h = np.array([0.25, 1.0])
```

```
# Plots data
plt.figure()
plt.plot(x, y, '.k', label='Noisy response')
plt.plot(x, np.sin(x), '--k', label='True function')
# Iterates over the smoother objects and plots functions for the uniform and gaussian kernels
for i in range(len(h)):
    plt.plot(x_star, H[i].predictor(kernel='uniform'), label='Uniform Kernel, h='+str(h[i]))
    plt.plot(x_star, H[i].predictor(kernel='gaussian'), label='Gaussian Kernel, h='+str(h[i]))
plt.legend(loc=0)
# plt.show()
plt.savefig('figures/kernel_smoother.pdf')
plt.close()
```

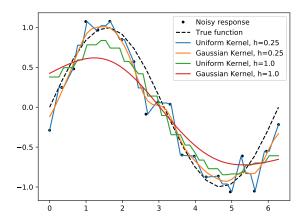


Figure 1: Uniform and Gaussian Kernel smoothing for different bandwidths

Cross validation

(A) Presumably a good choice of h would be one that led to smaller predictive errors on fresh data. Write a function or script that will: (1) accept an old ("training") data set and a new ("testing") data set as inputs; (2) fit the kernel-regression estimator to the training data for specified choices of h; and (3) return the estimated functions and the realized prediction error on the testing data for each value of h. This should involve a fairly straightforward "wrapper" of the function you've already written.

```
from __future__ import division
import numpy as np
import sys
sys.path.append('../../scripts/')
from smoothers import kernel_smoother
```

```
# Function for testing smoothers
def func(x):
    return np.sin(x)

# Initialize random seed
np.random.seed(3)

# Initialize x-vector
x = np.linspace(0, 2*np.pi, num=20)

# Training data of the form, y_i = f(x_i) + \epsilon_i
y_training = func(x) + np.random.normal(scale=0.2, size=x.shape)

# Testing data of the form, y_i = f(x_i) + \epsilon_i
y_test = func(x) + np.random.normal(scale=0.2, size=x.shape)
```

```
# Initialize list for storing kernel_smoother objects
H =[]

# Array of differing bandwidths to evaluate
h = np.array([0.2, 0.3, 0.4, 0.5, 0.6])

# Initialize MSE vector
mse = np.zeros(len(h))
```

```
# Iterates over bandwidth array and finds approximate MSE for each
for i in range(len(h)):
    H.append(kernel_smoother(x, y_training, x, h=h[i]))
    H[i].predictor()
    mse[i] = H[i].MSE(y_test)
print(mse)
```

```
jancabrera@phoenix$ python cross_validation_a.py
[0.07662882 0.0815833 0.0892882 0.10014996 0.11362704]
```

(B) Imagine a conceptual two-by-two table for the unknown, true state of affairs. The rows of the table are "wiggly function" and "smooth function," and the columns are "highly noisy observations" and "not so noisy observations." Simulate one data set (say, 500 points) for each of the four cells of this table, where the x's take values in the unit interval. Then split each data set into training and testing subsets. You choose the functions. Apply your method to each case, using the testing data to select a bandwidth parameter. Choose the estimate that minimizes the average squared error in prediction, which estimates the mean-squared error:

$$L_n(\hat{f}) = \frac{1}{n} \sum_{i=1}^{n^*} (y_i^* - \hat{y}_i^*)^2,$$

where (y_i^*, x_i^*) are the points in the test set, and \hat{y}_i^* is your predicted value arising from the model you fit using only the training data. Does your out-of-sample predictive validation method lead to reasonable choices of h for each case?

```
from __future__ import division
import numpy as np
import matplotlib.pyplot as plt
import sys
sys.path.append('../../scripts/')
from smoothers import kernel_smoother
```

```
# Function for testing smoothers
def func(x, period = 1):
    return np.sin(x*2*np.pi*period)
# Initialize random seed
np.random.seed(3)
# Initialize x-vector
x = np.linspace(0, 1, num=100)
# Noise level array, low and high
noise = np.array([0.05, 0.25])
noise_label = ['Low', 'High']
# Period of function, high period (wiggly), low period (smooth)
period = np.array([3, 0.5])
period_label = ['Wiggly', 'Smooth']
### Initialize lists for storing the different responses
# Training responses
Y_training = []
# Testing responses
Y_test = []
# True response
T = []
```

```
# Generates plots
i = 0
for n in range(len(noise)):
    for p in range(len(period)):

        # Instantiate kernel_smoother object
        y_smooth = kernel_smoother(x, Y_training[i], x)
        # Perform initial curve fit
        y_smooth.predictor()

    # Optimize bandwidth given the test data
        y_smooth.optimize_h(Y_test[i])

    # Perform curve fit with optimized bandwidth
        y_smooth.predictor()

# Predicted values for opimized bandwidth
        y_pred = y_smooth.y_star
```

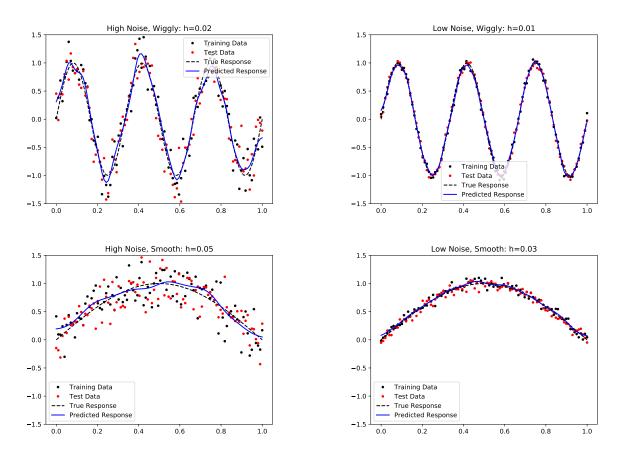


Figure 2: Bandwidths for various functions and noise levels

```
h_star = y_smooth.h

plt.figure()
plt.title(noise_label[n]+' Noise, '+ period_label[p] + ': h={:.2f}'.format(h_star[0]))
plt.plot(x, Y_training[i], '.k', label='Training Data')
plt.plot(x, Y_test[i], '.r', label='Test Data')
plt.plot(x, T[i], '--k', label='True Response')
plt.plot(x, y_pred, '-b', label='Predicted Response')
plt.legend(loc=0)
# plt.show()
plt.savefig('figures/cross_validation_'+noise_label[n]+'_'+period_label[p]+'.pdf')
plt.close()
i += 1
```

Local polynomial regression

Kernel regression has a nice interpretation as a "locally constant" estimator, obtained from locally weighted least squares. To see this, suppose we observe pairs (x_i, y_i) for i = 1, ..., n from our new favorite model, $y_i = f(x_i) + \epsilon_i$ and wish to estimate the value of the underlying function f(x) at some point x by weighted least squares. Our estimate is the scalar quantity

$$\hat{f}(x) = a = \arg\min_{\mathbb{R}} \sum_{i=1}^{n} w_i (y_i - a)^2,$$

where the w_i are the normalized weights (i.e. they have been rescaled to sum to 1 for fixed x). Clearly if $w_i = 1/n$, the estimate is simply \bar{y} , the sample mean, which is the "best" globally constant estimator. Using elementary

calculus, it is easy to see that if the unnormalized weights are

$$w_i \equiv w(x, x_i) = \frac{1}{h} K\left(\frac{x_i - x}{h}\right),$$

then the solution is exactly the kernel-regression estimator.

(A) A natural generalization of locally constant regression is local polynomial regression. For points u in a neighborhood of the target point x, define the polynomial

$$g_x(u; a) = a_0 + \sum_{k=1}^{D} a_j (u - x)^k$$

for some vector of coefficients $a = (a_0, \ldots, a_D)$. As above, we will estimate the coefficients a in $g_x(u; a)$ at some target point x using weighted least squares:

$$\hat{a} = \arg\min_{R^{D+1}} \sum_{i=1}^{n} w_i \{y_i - g_x(x_i; a)\}^2,$$

where $w_i \equiv w(x_i, x)$ are the kernel weights defined just above, normalized to sum to one. Derive a concise (matrix) form of the weight vector \hat{a} , and by extension, the local function estimate $\hat{f}(x)$ at the target value x. Life will be easier if you define the matrix X_x whose (i, j) entry is $(x_i - x)^{j-1}$, and remember that (weighted) polynomial regression is the same thing as (weighted) linear regression with a polynomial basis.

We first begin by plugging in $g_x(x_i; a)$ into the formula for \hat{a} , expanding terms and collecting them in matrix form.

$$\hat{a} = \arg\min_{R^{D+1}} \sum_{i=1}^{n} w_i \left\{ y_i - \left[a_0 + \sum_{k=1}^{D} a_j (x_i - x)^k \right] \right\}^2$$

$$= \arg\min_{R^{D+1}} \sum_{i=1}^{n} w_i \left\{ y_i - a_0 - a_1 (x_i - x) - a_2 (x_i - x)^2 - \dots - a_D (x_i - x)^D \right\}^2$$

$$= \arg\min_{R^{D+1}} (y - Xa)^T W (y - Xa)$$

We are then interested in minimizing the expression. This is done by setting the derivative with respect to a of the expression to zero. The term in the brackets are expanded.

$$0 = \frac{d}{da} \left[(y - Xa)^T W (y - Xa) \right]$$

$$= \frac{d}{da} \left[(y^T - a^T X^T) W (y - Xa) \right]$$

$$= \frac{d}{da} \left[y^T W y - y^T W X a - a^T X^T W y + a^T X^T W X a \right]$$

$$= \frac{d}{da} \left[y^T W y - 2y^T W X a + a^T X^T W X a \right]$$

Using the hint from exercise one that $\frac{\partial (z^T A z)}{\partial z} = (A + A^T)z$, and $\frac{\partial b^T z}{\partial z} = b$.

$$0 = -2y^T W X + 2X^T W X a$$
$$y^T W X = X^T W X a$$

We then evaluate the expression and solve for a.

$$\hat{a} = (X^T W X)^{-1} y^T W X$$

Notice that this has the same form as the linear case:

$$\hat{a} = Hy$$

With the hat matrix being, $H = (X^T W X)^{-1} X^T W$. The approximate estimate at our target value is then $\hat{f}(x) = \begin{bmatrix} 1 & 0 & \dots & 0 \end{bmatrix} \hat{a}$

(B) From this, conclude that for the special case of the local linear estimator (D=1), we can write $\hat{f}(x)$ as a linear smoother of the form

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

where the unnormalized weights are

$$w_i(x) = K\left(\frac{x-x_i}{h}\right) \left\{ s_2(x) - (x_i - x)s_1(x) \right\}$$

$$s_j(x) = \sum_{i=1}^n K\left(\frac{x-x_i}{h}\right) (x_i - x)^j.$$

From $\hat{f}(x) = \begin{bmatrix} 1 & 0 & \dots & 0 \end{bmatrix} \hat{a}$, we expand the matrices for the D = 1 case and multiply the matrices together.

$$\hat{f}(x) = \begin{bmatrix} 1 & 0 \end{bmatrix} \left\{ \begin{bmatrix} (x_1 - x)^0 & (x_1 - x) \\ \vdots & \vdots \\ (x_n - x)^0 & (x_n - x) \end{bmatrix}^T \begin{bmatrix} w_1 & \dots & 0 \\ \vdots & \ddots & \vdots \\ 0 & \dots & w_n \end{bmatrix} \begin{bmatrix} (x_1 - x)^0 & (x_1 - x) \\ \vdots & \vdots & \vdots \\ (x_n - x)^0 & (x_n - x) \end{bmatrix} \right\}^{-1} \begin{bmatrix} (x_1 - x)^0 & (x_1 - x) \\ \vdots & \vdots & \vdots \\ (x_n - x)^0 & (x_n - x) \end{bmatrix}^T \begin{bmatrix} w_1 & \dots & 0 \\ \vdots & \ddots & \vdots \\ 0 & \dots & w_n \end{bmatrix} \begin{bmatrix} y_1 \\ \vdots \\ y_n \end{bmatrix}$$

$$= \begin{bmatrix} 1 & 0 \end{bmatrix} \begin{bmatrix} \sum_{i=1}^n w_i \\ \sum_{i=1}^n w_i \\ w_i(x_i - x) \end{bmatrix}^{-1} \begin{bmatrix} \sum_{i=1}^n w_i y_i \\ \sum_{i=1}^n (x_i - x) \end{bmatrix}^{-1} \begin{bmatrix} \sum_{i=1}^n w_i y_i \\ \sum_{i=1}^n (x_i - x) \end{bmatrix}^{-1} \begin{bmatrix} \sum_{i=1}^n w_i y_i \\ \sum_{i=1}^n (x_i - x) \end{bmatrix}^{-1} \begin{bmatrix} \sum_{i=1}^n w_i y_i \\ \sum_{i=1}^n (x_i - x) \end{bmatrix}^{-1} \begin{bmatrix} \sum_{i=1}^n w_i y_i \\ y_i \end{bmatrix}$$

We note that for a two-by-two matrix the inverse can be found as

$$\begin{bmatrix} a & b \\ c & d \end{bmatrix}^{-1} = \frac{1}{ad - bc} \begin{bmatrix} d & -b \\ -c & a \end{bmatrix}$$

Evaluating the inverse and multiplying the resultant matrices together we obtain the following.

$$\hat{f}(x) = \left[\sum_{i=1}^{n} w_{i} \sum_{i=1}^{n} w_{i}(x_{i} - x)^{2} - \left(\sum_{i=1}^{n} w_{i}(x_{i} - x)\right)^{2}\right]^{-1} \left[1 \quad 0\right] \left[\sum_{i=1}^{n} w_{i}(x_{i} - x)^{2} - \sum_{i=1}^{n} w_{i}(x_{i} - x)\right] \left[\sum_{i=1}^{n} w_{i}y_{i} + \sum_{i=1}^{n} w_{i}(x_{i} - x)\right] \left[\sum_{i=1}^{n} w_{i}y_{i} + \sum_{i=1}^{n} w_{i}(x_{i} - x)\right] \left[\sum_{i=1}^{n} w_{i}(x_{i}$$

With $s_1 = \sum_{i=1}^n w_i(x_i - x)$ and $s_2 = \sum_{i=1}^n w_i(x_i - x)^2$ we obtain the form we were seeking:

$$\hat{f}(x) = \frac{\sum_{i=1}^{n} w_i y_i (s_2 - (x_i - x)s_1)}{\sum_{i=1}^{n} w_i (s_2 - (x_i - x)s_1)}.$$

(C) Suppose that the residuals have constant variance σ^2 (that is, the spread of the residuals does not depend on x). Derive the mean and variance of the sampling distribution for the local polynomial estimate $\hat{f}(x)$ at some arbitrary point x. Note: the random variable $\hat{f}(x)$ is just a scalar quantity at x, not the whole function.

$$E \left[\hat{f}(x) \right] = E \left[\begin{bmatrix} 1 & 0 & \dots & 0 \end{bmatrix} Hy \right]$$
$$= \begin{bmatrix} 1 & 0 & \dots & 0 \end{bmatrix} HE[y]$$
$$= \begin{bmatrix} 1 & 0 & \dots & 0 \end{bmatrix} Hf(x)$$

$$var[\hat{f}(x)] = var(\begin{bmatrix} 1 & 0 & \dots & 0 \end{bmatrix} H y)$$

$$= \{ \begin{bmatrix} 1 & 0 & \dots & 0 \end{bmatrix} H \}^T var(y) \{ \begin{bmatrix} 1 & 0 & \dots & 0 \end{bmatrix} H \}$$

$$= \sigma^2 \{ \begin{bmatrix} 1 & 0 & \dots & 0 \end{bmatrix} H \}^T \{ \begin{bmatrix} 1 & 0 & \dots & 0 \end{bmatrix} H \}$$

$$y_i = f(x_i) + \epsilon_i$$

(D) We don't know the residual variance, but we can estimate it. A basic fact is that if x is a random vector with mean μ and covariance matrix Σ , then for any symmetric matrix Q of appropriate dimension, the quadratic form x^TQx has expectation

$$E(x^T Q x) = \operatorname{tr}(Q \Sigma) + \mu^T Q \mu.$$

Write the vector of residuals as $r = y - \hat{y} = y - Hy$, where H is the smoothing matrix. Compute the expected value of the estimator

$$\hat{\sigma}^2 = \frac{\|r\|_2^2}{n - 2\text{tr}(H) + \text{tr}(H^T H)},$$

and simplify things as much as possible. Roughly under what circumstances will this estimator be nearly unbiased for large n? Note: the quantity $2\text{tr}(H) - \text{tr}(H^T H)$ is often referred to as the "effective degrees of freedom" in such problems.

$$\begin{split} \mathbf{E}\left[\hat{\sigma}^2\right] &= \mathbf{E}\left[\frac{(y-Hy)^T(y-Hy)}{n-2\mathrm{tr}(H)+\mathrm{tr}(H^TH)}\right] \\ &= \frac{\mathbf{E}[(y-Hy)^T(y-Hy)]}{n-2\mathrm{tr}(H)+\mathrm{tr}(H^TH)} \\ &= \frac{\mathbf{E}[y^Ty]-2\mathbf{E}[y^THy]+\mathbf{E}[y^TH^THy]}{n-2\mathrm{tr}(H)+\mathrm{tr}(H^TH)} \end{split}$$

$$E[y^T y] = \operatorname{tr}(\Sigma) + f(x)^T f(x)$$

$$= n\sigma^2 + f(x)^T f(x)$$

$$E[y^T H y] = \operatorname{tr}(H\Sigma) + f(x)^T H f(x)$$

$$= \sigma^2 \operatorname{tr}(H) + f(x)^T H f(x)$$

$$E[y^T H^T H y] = \operatorname{tr}(H^T H \Sigma) + f(x)^T H^T H f(x)$$

$$= \sigma^2 \operatorname{tr}(H^T H) + f(x)^T H^T H f(x)$$

$$\begin{split} & \mathrm{E}\left[\hat{\sigma}^{2}\right] = \frac{n\sigma^{2} + f(x)^{T}f(x) - 2\sigma^{2}\mathrm{tr}(H) - 2f(x)^{T}Hf(x) + \sigma^{2}\mathrm{tr}(H^{T}H) + f(x)^{T}H^{T}Hf(x)}{n - 2\mathrm{tr}(H) + \mathrm{tr}(H^{T}H)} \\ & = \frac{\sigma^{2}[n - 2\mathrm{tr}(H) + 2\mathrm{tr}(H^{T}H)] + f(x)^{T}f(x) - 2f(x)^{T}Hf(x) + f(x)^{T}H^{T}Hf(x)}{n - 2\mathrm{tr}(H) + \mathrm{tr}(H^{T}H)} \\ & = \sigma^{2} + \frac{\|f(x) - Hf(x)\|_{2}^{2}}{n - 2\mathrm{tr}(H) + \mathrm{tr}(H^{T}H)} \end{split}$$

Unbiased when the difference, f(x) - Hf(x) is small

(E) Write a new R function that fits the local linear estimator using a Gaussian kernel for a specified choice of bandwidth h. Then load the data in "utilities.csv" into R. This data set shows the monthly gas bill (in dollars) for a single-family home in Minnesota, along with the average temperature in that month (in degrees F), and the number of billing days in that month. Let y be the average daily gas bill in a given month (i.e. dollars divided by billing days), and let x be the average temperature. Fit y versus x using local linear regression and some choice of kernel. Choose a bandwidth by leave-one-out cross-validation.

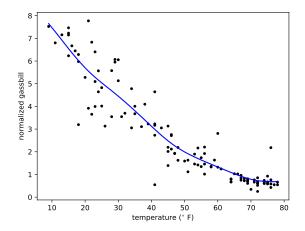


Figure 3: Fit using LOOCV, h = 6.87

- (F) Inspect the residuals from the model you just fit. Does the assumption of constant variance (homoscedasticity) look reasonable? If not, do you have any suggestion for fixing it?
- (G) Put everything together to construct an approximate point-wise 95% confidence interval for the local linear model (using your chosen bandwidth) for the value of the function at each of the observed points x_i for the utilities data. Plot these confidence bands, along with the estimated function, on top of a scatter plot of the data.

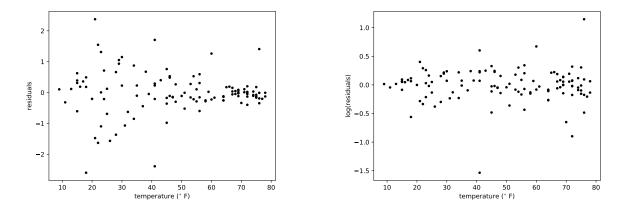


Figure 4: Residuals and log-residuals

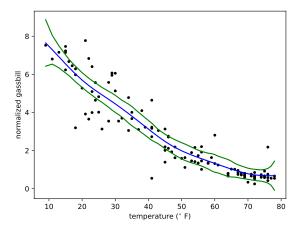


Figure 5: Fit using LOOCV, h = 6.87

A Kernel Smoothing Code

```
from __future__ import division
import numpy as np
from scipy.optimize import minimize
from numpy.linalg import inv
class kernels:
    This class contains kernel functions for kernel smoothing
   def __init__(self):
       pass
   def uniform(x):
       Parameters
          x: float
               argument to be evaluated
       Returns
           k: float
              uniform kernel evaluation at x
               .. math:: k(x) = 1/2 I(x), with I(x) = 1, if |x| \land |x|
       if np.abs(x) \le 1:
           k = 0.5
       else:
           k = 0
       return k
    def gaussian(x):
       Parameters
           x: float
              argument to be evaluated
       Returns
           k: float
               gaussian kernel evaluation at x
               .. math:: k(x) = \frac{1}{\sqrt{2}} e^{-x^2/2}
       return 1/np.sqrt(2*np.pi)*np.exp(-x**2/2)
class kernel_smoother:
    This class returns a vector of smoothed values given feature and response vectors
    def __init__(self, x, y, x_star, kernel='gaussian', h=1, D=1):
       Parameters
           x: float
               Feature vector
           y: float
               Response vector
            x star: float
               Scalar or vector to be evaluated
           kernel: str (optional)
            Kernel type to be used: Available kernels are gaussian,
```

```
h: float (optional)
           Bandwidth
       D: int (optional)
           Order of polynomial smoother
   self.x = x
   self.y = y
   self.x_star = x_star
   self.h = h
   self.D = D
   self.kernel=kernel
   if x_star.shape != x:
       raise Warning('Feature vector and evaluation vector are not the same length. Residuals \leftrightarrow
           will not be calculated.')
def local_general(self):
   Returns
       y_star: float, len(x_star)
           Predictor for x_star,
           .. math:: y_i^* = [1 \ 0 \ ...] \ H_i \ y
           .. math:: w(x_i, x^*) = \frac{1}{2}K( \frac{x_i - x^*}{h}), \sum_{i=1}^n w(x_i, \leftarrow x_i)
               x^* = 1
   Attributes
       sigma: float, len(x_star)
           Standard error at x_star
       residuals: float, len(x_star)
           residuals of prediction from data (only calculated if len(y_star) == (len(y)))
           .. math:: r_i = y_i - \hat{y}_i
       Hat_matrix: list
           first elements in hat matrix for a given x_star
    # Instantiate vectors and matrices for storing calculations
   self.y_star = np.zeros(self.x_star.shape)
   self.sigma = np.zeros(self.x_star.shape)
   self.residuals = np.zeros(self.x_star.shape)
   self.Hat_matrix = []
   # Vector for determining y_star
   e_1 = np.append(1, np.zeros(self.D))
    # Iterate through each value in y_star
   for i in range(self.y_star.shape[0]):
       # Instantiate a weights vector
       weights = np.zeros(self.x.shape)
        # Instantiates X matrix
       X = np.transpose(np.ones((self.x.shape[0], self.D+1)))
        # Populates X matrix given the order of the smoother, D
       for d in range(self.D):
           X[d+1] = (self.x - self.x_star[i]) **(d+1)
       X = np.transpose(X)
        # Iterates through feature/response vectors
       for j in range(self.x.shape[0]):
           self.h)
```

```
# Normalizes weights such that \sum_{i=1}^n w(x_i, x^*) = 1
        weights = weights/weights.sum()
        # Construct weights matrix
        W = np.diag(weights)
        # Calculate hat matrix; H = (X^T W X)^{-1} X^T W
        H = (inv(np.transpose(X) @ W @ X) @ np.transpose(X) @ W)
        # Append hat matrix with current H matrix (needed for LOOCV)
        self.Hat_matrix.append(e_1 @ H)
        \# \hat{f}(x^*) = e_1 H y
        self.y_star[i] = e_1 @ H @ self.y
        # Calculates residuals and confidence interval if y_star and y shapes are equivalent
        if self.y_star.shape == self.y.shape:
            self.residuals[i] = self.y[i] - self.y_star[i]
    # Residual sum of squared errors
   rss = (self.residuals**2).sum()
    # Approximate standard error of fit
    sigma_hat = rss/(len(self.y_star)-1)
    # Calculates approximate standard error
    for i in range(self.sigma.shape[0]):
        self.sigma[i] = np.sqrt(np.transpose(self.Hat_matrix[i]) @ (self.Hat_matrix[i]) ★↔
            sigma_hat)
   return self.y_star
def MSE(self, y_test):
    Parameters
       y_test: float
          Response vector
    Returns
       MSE: float
           Average squared error for y_test given y
           .. math:: MSE \\approx \\frac{1}{n}\\sum(y_test - y_star)**2
    self.y_test = y_test
   return np.mean((y_test - self.y_star)**2)
def MSE_optimization(self, y_test):
   Parameters
      y_test: float
           Response vector
   Returns
       h star: float
           Uses BFGS minimization method to minimize the MSE by varying h_star
    self.y_test = y_test
    def func(X):
       Y = kernel_smoother(self.x, self.y, self.x_star, kernel=self.kernel, h=X, D=self.D)
       Y.local_general()
       return Y.MSE(self.y_test)
   h_star = minimize(func, 1)
    self.h = h_star.x
   return h_star.x
```

```
def LOOCV(self):
    Returns
      loocv: float
          Leave one out cross validation statistic (LOOCV)
           .. math: loocv = \sum_{i=1}^n (\sqrt{y_i - \hat{y}_i}{1-H_{ii}})^2
    loocv = np.zeros(len(self.y))
    for i in range(len(loocv)):
       loocv[i] = ((self.y[i] - self.y_star[i])/(1 - self.Hat_matrix[i][i]))**2
    return loocv.sum()
def LOOCV_optimization(self):
    Returns
      h_star: float
          Uses BFGS minimization method to minimize the LOOCV by varying h_star
    def func(X):
       Y = kernel_smoother(self.x, self.y, self.x_star, kernel=self.kernel, h=X, D=self.D)
       Y.local_general()
       return Y.LOOCV()
    h_star = minimize(func, 0.7)
    self.h = h_star.x
    return self.h
```