V2A2_Regression

December 14, 2018

```
In [1]: #!/usr/bin/env python
       # V2A2_Regression.py
        # Programmgeruest zu Versuch 2, Aufgabe 2
       import numpy as np
       import scipy.spatial
       from random import randint
        # base class for regressifiers
        # -----
       class Regressifier:
           11 11 11
           Abstract base class for regressifiers
           Inherit from this class to implement a concrete regression algorithm
           n n n
           def fit(self,X,T):
                                     # train/compute regression with lists of feature vectors
               11 11 11
               Train regressifier by training data X, T, should be overwritten by any derived
               :param X: Data matrix of size NxD, contains in each row a data vector of size .
               :param T: Target vector matrix of size NxK, contains in each row a target vect
               :returns: -
               n n n
               pass
                                   # predict a target vector given the data vector x
           def predict(self,x):
               Implementation of the regression algorithm; should be overwritten by any deriv
               :param x: test data vector of size D
               :returns: predicted target vector
               return None
           def crossvalidate(self,S,X,T,dist=lambda t: np.linalg.norm(t)): # do a S-fold cro
               11 11 11
               Do a S-fold cross validation
```

```
:param S: Number of parts the data set is divided into
:param X: Data matrix (one data vector per row)
:param T: Matrix of target vectors; T[n] is target vector of X[n]
:param dist: a fuction dist(t) returning the length of vector t (default=Eukli
:returns (E_dist,sd_dist,E_min,E_max) : mean, standard deviation, minimum, and
X,T=np.array(X),np.array(T)
                                                   # ensure array type
                                                   # N=number of data vectors
N=len(X)
perm = np.random.permutation(N)
                                                   # do a random permutation
X1,T1=[X[i] for i in perm], [T[i] for i in perm]
                                                  # ... to get random partit
idxS = [range(i*N//S,(i+1)*N//S) \text{ for } i \text{ in } range(S)] \# divide data set into } S p
E_dist, E_dist2, E_max, E_min=0, 0, -1, -1
                                                   # initialize first two mom
Erel_dist, Erel_dist2, Erel_max, Erel_min=0,0,-1,-1
                                                  # initialize first two mom
for idxTest in idxS:
                                                   # loop over all possible t
    # (i) generate training and testing data sets and train classifier
    idxLearn = [i for i in range(N) if i not in idxTest]
    if(S<=1): idxLearn=idxTest</pre>
   X_learn, T_learn = np.array([X1[i] for i in idxLearn]), np.array([T1[i] for
   X_test , T_test = np.array([X1[i] for i in idxTest ]), np.array([T1[i] for
                                                   # train regressifier
   self.fit(X_learn,T_learn)
    # (ii) test regressifier
   for i in range(len(X_test)): # loop over all data vectors to be tested
        # (ii.a) regress for i-th test vector
       xn_test = X_test[i].T
                                                       # data vector for test
                                                       # predict target value
       t_test = self.predict(xn_test)
        # (ii.b) check for regression errors
       t_true = T_test[i].T
                                                       # true target value
                                                       # (Euklidean) distance
       d=dist(t_test-t_true)
       dttrue=dist(t_true)
                                                       # length of t_true
       E_dist = E_dist+d
                                                       # sum up distances (fo
       E_dist2 = E_dist2+d*d
                                                       # sum up squared dista
       if(E_max<0)or(d>E_max): E_max=d
                                                       # collect maximal erro
       if(E_min<0)or(d<E_min): E_min=d</pre>
                                                       # collect minimal erro
       drel=d/dttrue
       Erel_dist = Erel_dist+drel
                                                       # sum up relative dist
       Erel_dist2 = Erel_dist2+(drel*drel)
                                                       # sum up squared relat
       if(Erel_max<0)or(drel>Erel_max): Erel_max=drel # collect maximal rela
       if(Erel_min<0)or(drel<Erel_min): Erel_min=drel # collect minimal rela
E_dist
           = E_dist/float(N)
                                                       # estimate of first mod
           = E_dist2/float(N)
E_dist2
                                                       # estimate of second m
                                                       # variance of error
Var_dist
           = E_dist2-E_dist*E_dist
                                                       # standard deviation o
\mathtt{sd\_dist}
          = np.sqrt(Var_dist)
                                                       # estimate of first mod
Erel_dist = Erel_dist/float(N)
Erel_dist2 = Erel_dist2/float(N)
                                                       # estimate of second m
Varrel_dist = Erel_dist2-Erel_dist*Erel_dist
                                                       # variance of error
sdrel_dist = np.sqrt(Varrel_dist)
                                                       # standard deviation o
return (E_dist,sd_dist,E_min,E_max), (Erel_dist,sdrel_dist,Erel_min,Erel_max)
```

```
# DataScaler: scale data to standardize data distribution (for mean=0, standard deviat
       # ------
       class DataScaler:
           11 11 11
           Class for standardizing data vectors
           Some regression methods require standardizing of data before training to avoid num
           def __init__(self,X):
                                              # X is data matrix, where rows are data vector
               Constructor: Set parameters (mean, std,...) to standardize data matrix X
               : param X: Data matrix of size NxD the standardization parameters (mean, std, .
               :returns: object of class DataScaler
               self.meanX = np.mean(X,0)
                                             # mean values for each feature column
               self.stdX = np.std(X,0)
                                             # standard deviation for each feature column
               if isinstance(self.stdX,(list,tuple,np.ndarray)):
                   self.stdX[self.stdX==0]=1.0 # do not scale data with zero std (that is, co
               else:
                   if(self.stdX==0): self.stdX=1.0 # in case stdX is a scalar
               self.stdXinv = 1.0/self.stdX # inverse standard deviation
           def scale(self,x):
                                              # scales data vector x to mean=0 and std=1
               scale data vector (or data matrix) x to mean=0 and s.d.=1
               :param x: data vector or data matrix
               :returns: scaled (standardized) data vector or data matrix
               n n n
               return np.multiply(x-self.meanX,self.stdXinv)
           def unscale(self,x):
                                              # unscale data vector x to original distributi
               unscale data vector (or data matrix) x to original data ranges
               :param x: standardized data vector or data matrix
               :returns: unscaled data vector or data matrix
               return np.multiply(x,self.stdX)+self.meanX
           def printState(self):
               print standardization parameters (mean value, standard deviation (std), and in
               print("mean=",self.meanX, " std=",self.stdX, " std_inv=",self.stdXinv)
In [3]: from itertools import *
```

```
_____
       # -----
       # function to compute polynomial basis functions
       def phi_polynomial(x,deg=1):
                                         # x should be list or np.array or 1xD matrix; r
          polynomial basis function vector; may be used to transform a data vector x into a
          :param x: data vector to be transformed into a feature vector
          :param deg: degree of polynomial
          :returns phi: feature vector
          Example: phi polynomial (x,3) returns for one-dimensional x the vector [1, x, x*x,
                                          # ensure that x is a 1D array (first row of x)
          x=np.array(np.mat(x))[0]
          D=len(x)
          \#assert\ (D==1)\ or\ ((D>1)\ and\ (deg<=3)),\ "phi_polynomial(x,deg)\ not\ implemented\ for
              phi = np.array([x[0]**i for i in range(deg+1)])
          else:
              phi = np.array(phi_helper(x.tolist(),deg))
              #phi = np.array([])
              #if(deg>=0):
                  phi = np.concatenate((phi,[1]))  # include degree 0 terms
                   if(deg>=1):
                      phi = np.concatenate((phi,x)) # includes degree 1 terms
                      if(deg>=2):
                          for i in range(D):
                             phi = np.concatenate((phi, [x[i]*x[j] for j in range(i+1)])
                          if(deq>=3):
                             for i in range(D):
                                 for j in range(i+1):
                                    phi = np.concatenate((phi, [x[i]*x[j]*x[k] for k in 
                             # EXTEND CODE HERE FOR deq>3!!!!
          return phi.T # return basis function vector (=feature vector corresponding to dat
       def phi_helper(x, deg):
          if deg <= 0:
              return [1]
          else:
              return phi_helper(x, deg-1) + [reduce(lambda a,b:a*b,combi) for combi in combi:
In [4]: # -----
       # Least Squares (ML) linear regression with sum of squares Regularization,
       # ------
       class LSRRegressifier(Regressifier):
          Class for Least Squares (or Maximum Likelihood) Linear Regressifier with sum of sq
           11 11 11
```

from functools import reduce

```
def __init__(self,lmbda=0,phi=lambda x: phi_polynomial(x,1),flagSTD=0,eps=0.000001
   Constructor of class LSRegressifier
    :param lmbda: Regularization coefficient lambda
    :param phi: Basis-functions used by the linear model (default linear polynomia
    :param flagSTD: If >0 then standardize data X and target values T (to mean 0 a
    :param eps: maximal residual value to tolerate (instead of zero) for numerical
    :returns: -
   self.lmbda=lmbda
                           # set regression parameter (default 0)
                           # set basis functions used for linear regression (defau
   self.phi=phi
   self.flagSTD=flagSTD; # if flaq >0 then data will be standardized, i.e., scal
                           # maximal residual value to tolerate (instead of zero)
   self.eps=eps;
def fit(self,X,T,lmbda=None,phi=None,flagSTD=None): # train/compute LS regression
    Train regressifier (see lecture manuscript, theorem 3.11, p33)
    :param X: Data matrix of size NxD, contains in each row a data vector of size .
    :param T: Target vector matrix of size NxK, contains in each row a target vect
    :param lmbda: Regularization coefficient lambda
    :param phi: Basis-functions used by the linear model (default linear polynomia
    :param flagSTD: If >0 then standardize data X and target values T (to mean 0 a
    :returns: flagOK: if >0 then all is ok, otherwise matrix inversion was bad con
    # (i) set parameters
   if lmbda==None: lmbda=self.lmbda
                                           # reset regularization coefficient?
   if phi==None: phi=self.phi
                                           # reset basis functions?
   if flagSTD==None: flagSTD=self.flagSTD # standardize data vectors?
    # (ii) scale data for mean=0 and s.d.=0 ?
   if flagSTD>0:
                                           # if yes, then...
       self.datascalerX=DataScaler(X)
                                           # create datascaler for data matrix X
        self.datascalerT=DataScaler(T)
                                           # create datascaler for target matrix T
       X=self.datascalerX.scale(X)
                                           # scale all features (=columns) of data
       T=self.datascalerT.scale(T)
                                           # ditto for target matrix T
    # (iii) compute weight matrix and check numerical condition
   flagOK,maxZ=1,0;
                                           # if <1 then matrix inversion is numeri
   try:
       self.N,self.D = X.shape
                                           \# data matrix X has size N x D (N is nu
       self.M = self.phi(self.D*[0]).size # get number of basis functions
        \#self.K = T.shape[1]
                                            # DELTE dummy code (just required for
       PHI = np.array([np.transpose(phi(x)) for x in X])
       PHIT_PHI_lmbdaI = np.add(np.dot(PHI.T,PHI),np.dot(self.lmbda,np.identity(self.lmbda)
        #print(PHIT_PHI_lmbdaI)
        # REPLACE dummy code: compute PHI_T*PHI+lambda*I
       PHIT_PHI_lmbdaI_inv = np.linalg.inv(PHIT_PHI_lmbdaI)
                                                                      # REPLACE dum
        self.W_LSR = np.dot(np.dot(PHIT_PHI_lmbdaI_inv,PHI.T),T)# REPLACE dummy co
        # (iv) check numerical condition
```

```
Z=np.dot(PHIT_PHI_lmbdaI,PHIT_PHI_lmbdaI_inv)-np.identity(self.M) # REPLA
                                        \max Z = np.max(Z)
                                                                               # REPLACE dummy code: Compute maximum (absolute) compo
                                        assert maxZ<=self.eps</pre>
                                                                                                                 # maxZ should be <eps for good condition
                                except:
                                        flagOK=0;
                                        print("EXCEPTION DUE TO BAD CONDITION:flagOK=", flagOK, " maxZ=", maxZ, " or print("EXCEPTION DUE TO BAD CONDITION:flagOK=", flagOK, " maxZ=", maxZ, " or print("EXCEPTION DUE TO BAD CONDITION:flagOK=", flagOK, " maxZ=", maxZ, " or print("EXCEPTION DUE TO BAD CONDITION:flagOK=", flagOK, " maxZ=", maxZ, " or print("EXCEPTION DUE TO BAD CONDITION:flagOK=", flagOK, " maxZ=", maxZ=", maxZ, " or print("EXCEPTION DUE TO BAD CONDITION:flagOK=", flagOK, " maxZ=", maxZ, " or print("EXCEPTION DUE TO BAD CONDITION:flagOK=", flagOK, " maxZ=", maxZ="
                                return flagOK
                        def predict(self,x,flagSTD=None):
                                                                                                        # predict a target value given data vector
                                predicts the target value y(x) for a test vector x
                                 :param x: test data vector of size D
                                 :param flagSTD: If >0 then standardize data X and target values T (to mean 0 a
                                 :returns: predicted target vector y of size K
                                 n n n
                                if flagSTD==None: flagSTD=self.flagSTD
                                                                                                                           # standardazion?
                                if flagSTD>0: x=self.datascalerX.scale(x) # if yes, then scale x before comp
                                phi_of_x = self.phi(x)
                                                                                                                           # compute feature vector phi_of_x |
                                y=np.dot(self.W_LSR.T, phi_of_x)
                                                                                                                                            # REPLACE dummy code: com
                                if flagSTD>0: y=self.datascalerT.unscale(y) # scale prediction back to origina
                                return y
                                                                                      # return prediction y for data vector x
In [5]: # -----
                # KNN regression
                class KNNRegressifier(Regressifier):
                         Class for fast K-Nearest-Neighbor-Regression using KD-trees
                         11 11 11
                        def __init__(self,K,flagKLinReg=0):
                                Constructor of class KNNRegressifier
                                 :param K: number of nearest neighbors that are used to compute prediction
                                 :flagKLinReg: if >0 then the do a linear (least squares) regression on the the
                                                             otherwise just take the mean of the K nearest neighbors target v
                                 :returns: -
                                 11 11 11
                                self.K = K
                                                                                                                         # K is number of nearest-neighbors
                                self.X, self.T = [],[]
                                                                                                                         # initially no data is stored
                                self.flagKLinReg=flagKLinReg
                                                                                                                         # if flag is set then do a linear r
                        def fit(self,X,T): # train/compute regression with lists of data vectors X and tar
                                 Train regressifier by stroing X and T and by creating a KD-Tree based on X
                                 :param X: Data matrix of size NxD, contains in each row a data vector of size .
                                 :param T: Target vector matrix of size NxK, contains in each row a target vect
```

```
11 11 11
              self.X, self.T = np.array(X),np.array(T) # just store feature vectors X and
              self.N, self.D = self.X.shape
                                                      \# store data number N and dimension
              self.kdtree = scipy.spatial.KDTree(self.X) # do an indexing of the feature vec
           def predict(self,x,K=None,flagKLinReg=None): # predict a target value given data
              n n n
              predicts the target value y(x) for a test vector x
              :param x: test data vector of size D
              :param K: number of nearest neighbors that are used to compute prediction
              :flagKLinReq: if >0 then the do a linear (least squares) regression on the the
                           otherwise just take the mean of the K nearest neighbors target v
              :returns: predicted target vector of size K
              if(K==None): K=self.K
                                                      # do a K-NN search...
              if(flagKLinReg==None): flagKLinReg=self.flagKLinReg # if flag >0 then do a reg
              nn = self.kdtree.query(x,K)
                                                      # get indexes of K nearest neighbor
              if K==1: idxNN=[nn[1]]
                                                      # cast nearest neighbor indexes nn
              else: idxNN=nn[1]
              t out=0
              if(self.flagKLinReg==0):
                  # just take mean value of KNNs
                  t_out=np.mean([self.T[i] for i in idxNN])
              else:
                  # do a linear regression of the KNNs
                  lsr=LSRRegressifier(lmbda=0.0001,phi=lambda x:phi_polynomial(x,1),flagSTD=
                  lsr.fit(self.X[idxNN],self.T[idxNN])
                  t_out=lsr.predict(x)
              return t_out
# main
        # Module test
        # **********************************
        if __name__ == '__main__':
           print("\n----")
           print("Example: 1D-linear regression problem")
           print("----")
            # (i) generate data
           N = 100
                                   # parameters of line
           w0, w1=4,2
           X=np.zeros((N,1))
                                   # x data: allocate Nx1 matrix as numpy ndarray
           X[:,0]=np.arange(0,50.0,50.0/N) # equidistant sampling of the interval [0,50)
                                  # target values: allocate Nx1 matrix as numpy ndarray
           T=np.zeros((N,1))
            sd_noise = 1.0
                                   # noise power (=standard deviation)
           T=T+w1*X+w0 + np.random.normal(0,sd_noise,T.shape) # generate noisy target value
```

:returns: -

```
print("X=",X)
          print("T=",T)
          # (ii) define basis functions (phi should return list of basis functions; x shoul
                                          # degree of polynomial
          phi=lambda x: phi_polynomial(x,1) # define phi by polynomial basis-functions u
          print("phi(4)=", phi([4]))
                                         # print basis function vector [1, x, x*x ...
          print("phi([1,2])=", phi([1,2]))
                                          # print basis function vector for two-dim.
          # (iii) compute LSR regression
          print("\n----")
          print("Do a Least-Squares-Regression")
          print("----")
          lmbda=0;
          lsr = LSRRegressifier(lmbda,phi)
          lsr.fit(X,T)
          print("lsr.W_LSR=",lsr.W_LSR)
                                       # weight vector (should be approximately
          x=np.array([3.1415]).T
          print("prediction of x=",x,"is y=",lsr.predict(x))
          # do S-fold crossvalidation
          err_abs,err_rel = lsr.crossvalidate(S,X,T)
          print("LSRRegression cross-validation: absolute errors (E,sd,min,max)=", err_abs,
          # (iv) compute KNN-regression
          print("\n----")
          print("Do a KNN-Regression")
          print("----")
          K=2;
          knnr = KNNRegressifier(K)
          knnr.fit(X,T)
          print("prediction of x=",x,"is y=",knnr.predict(x))
          # do S-fold crossvalidation
          err_abs,err_rel = knnr.crossvalidate(S,X,T)
          print("KNNRegression cross-validation: absolute errors (E,sd,min,max)=", err_abs,
._____
```

regularization parameter

par_lambda = 0

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- [4.5]
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T= [[3.96539552]

- [5.40900355]
- [6.80479406]
- [8.03458322]
- [8.3053615]
- [8.73249338]
- [8.77151182]
- [9.63643954]
- [12.01447684]
- [14.10503997]
- [14.10000997]
- [14.53639261]
- [16.56342681]
- [15.55803052]
- [17.54062696]
- [18.10718482]
- [20.19849118]
- [19.69947678]
- [21.30766924]
- [21.40690018]
- [22.14653271]
- [22.140002/1_
- [22.94023538]
- [25.36725429]
- [26.22817455]
- [27.78178612]
- [28.41925754]
- [27.71492035]
- [29.23188689]
- [29.9793288]
- [31.95354147]
- [32.47758882]
- [32.27269297]
- [36.15945205]
- [00:10010200]
- [35.22589939]
- [38.2674975]
- [39.33027519] [38.71212506]
- [40.19112542]
- [40.37822356]
- [42.08476023]
- [41.09623498]
- [43.07393777]
- [44.58435281]
- [46.23278059]
- [46.80294075]
- [48.494492]
- [47.73994133]
- [50.94662245]
- [50.73035166]
- [50.82361583]

- [54.41661601]
- [54.32981823]
- [55.41985461]
- [54.82862886]
- [58.04696979]
- [57.30488021]
- [59.91427266]
- [60.44129026]
- [62.53071307]
- [60.97598519]
- [62.26103967]
- [64.32530732]
- [64.36329645]
- [68.68025287]
- [66.79291027]
- [68.05828692]
- [69.51147636]
- [69.78195668]
- [72.01666756]
- [72.378812]
- [73.80618286]
- [73.32671575]
- [74.58556612]
- [75.35734468]
- [77.2098026]
- [79.00187307]
- [79.67579371]
- [79.86904299]
- [80.98003463]
- [82.03965717]
- [82.55033359]
- [81.91388516]
- [86.1249457]
- [84.73057236]
- [86.42908599]
- [86.79082706]
- [88.72028873]
- [90.40730705] [91.78703418]
- [91.15956868]
- [93.33433493]
- [93.12277541]
- [96.2559933]
- [95.55246418]
- [95.90050568]
- [97.07683317]
- [98.35082574]
- [99.51604217]

```
[100.21538238]
 [103.11867279]
 [102.75984867]]
phi(4) = [1 \ 4]
phi([1,2])= [1 1 2]
Do a Least-Squares-Regression
lsr.W_LSR= [[4.09052788]
 [1.99478017]]
prediction of x = [3.1415] is y = [10.35712978]
LSRRegression cross-validation: absolute errors (E,sd,min,max)= (0.7210491003592423, 0.5070165
  -----
Do a KNN-Regression
_____
prediction of x = [3.1415] is y = 9.203975679168057
KNNRegression cross-validation: absolute errors (E,sd,min,max)= (1.1743092886650586, 0.9170722
In [16]: temp = []
        for i,e in enumerate(KNNErrors):
            temp.append((e[2][3],i))
        temp.sort()
        #print(temp)
        smallest_mean= KNNErrors[1][2]
        smallest_sd= KNNErrors[1][2]
        smallest_min= KNNErrors[13][2]
        smallest_max= KNNErrors[1][2]
        print(smallest_mean)
        print(smallest_sd)
        print(smallest_min)
        print(smallest_max)
        print(KNNErrors[1])
(0.04137596473341779, 0.07148122842087293, 6.810069605730603e-05, 0.44728370446074633)
(0.04137596473341779, 0.07148122842087293, 6.810069605730603e-05, 0.44728370446074633)
(0.1297162287702424, 0.39638164915995594, 5.395317915671042e-05, 3.1265953272944778)
(0.04137596473341779, 0.07148122842087293, 6.810069605730603e-05, 0.44728370446074633)
(2, (1.0953802187892574, 0.8689819772601937, 0.003291802811396849, 3.884868724720441), (0.0413)
```

0.1 a) Versuchen Sie zunächst den Aufbau des Moduls V2A2_Regression.py zu verstehen:

Betrachten Sie den Aufbau des Moduls durch Eingabe von pydoc V2A2_Regressifier. Welche Klassen gehören zu dem Modul und welchen Zweck haben sie jeweils?

- DataScaler: Standatisiert Daten
- Regressifier: Abstrakte Klasse für Regressierer
- KNNRegressifier: ermöglicht Regression mithilfe von Fast-KNN
- LSRRegressifier: ermöglicht Regression mithilfe von Fehlerquadratsummen

Betrachten Sie nun die Basis-Klasse Regressifier im Quelltext: Wozu dienen jeweils die Methoden fit(self,X,T), predict(self,x) und crossvalidate(self,S,X,T) ?

- fit(self,X,T): Verlangt nach der implimentierung einer Methode die den Regressierer mit X und T trainiert.
- predict(self,x): Verlangt nach der implimentierung einer Methode die auf x eine Regression anwendet und eine Vorhersage zurückliefert
- crossvalidate(self,S,X,T): Macht S-fache Kreuzvaliedierung mit den Daten,Labeln (X,T) und liefert als Ergebniss Informationen über die relativen und absoluten Fehlerwerte

Worin unterscheidet sich crossvalidate(.) von der entsprechenden Methode für Klassifikation (siehe vorigen Versuch)?

- hier kann man noch angeben mit welcher Längenfunktion gearbeitet werden soll, und es werden andere Fehlerstatistiken zurückgegeben
- b) Betrachten Sie nun die Funktion phi_polynomial(x,deg):

Was berechnet die Funktion? Welches Ergebnis liefert phi_polynomial([3],5)? Welches Ergebnis liefert phi_polynomial([3,5],2)?

```
In [74]: phi_polynomial([3],5)
Out[74]: array([ 1,  3,  9,  27,  81, 243])
In [75]: phi_polynomial([3,5],2)
Out[75]: array([ 1,  3,  5,  9, 15, 25])
```

• phi_polynomial(x,deg): berechnet die Basisfunktion für die Werte x mit dem grad deg

Geben Sie eine allgemeine Formel an für das Ergebnis von phi_polynomial([x1,x2],2)?

• [1, x1, x2, (x1)š, x1*x2, (x2)š]

Wozu braucht man diese Funktion im Zusammenhang mit Regression?

• Das Kreuzprodukt aus dieser Funktion und dem Zielwertevektor geben die Prognose bei der Regression

Bis zu welchem Polynomgrad kann die Funktion Basisfunktionen berechnen? Erweitern Sie die Funktion mindestens bis Grad 5.

• momentan bis grad 3, siehe Verbesserung im code

```
In [83]: phi_polynomial([3,5],5)
```

c) Betrachten Sie die Klasse LSRRegressifier:

Welche Art von Regressions-Modell berechnet diese Klasse?

Fehlerquadratsummenregression mit Regularisierung

Wozu dienen jeweils die Parameter Imbda, phi, flagSTD und eps?

- lmbda: Regularisierungsparameter
- phi: Basisfunktion
- flagSTD: gibt an ob die Daten Standatisiert werden sollen
- eps: höchster erlaubter Fehlerrestwert

Welche Rolle spielt hier die Klasse DataScaler? In welchen Methoden und zu welchem Zweck werden die Daten ggf. umskaliert? Welches Problem kann auftreten wenn man dies nicht tut? Wozu braucht man die Variablen Z und maxZ in der Methode fit(.)?

- DataScaler wird in fit() und predict() verwendet um Daten zu Standartisieren
- nicht scalierte Daten werde können kleinste abweichungen in Gewichten zu großen und/oder unterschiedlichen Änderungen führen.
- Z ist die Matrix mit den Fehlerwerten die durch Invertierung entstehen.
- maxZ ist der gröSSte Wert aus Z und wird verwendet um zu bestimmen ob der Fehlerwert annehmbar ist

Vervollständigen Sie die Methoden fit(self,X,T,...) und predict(self,x,...) (vgl. vorige Aufgabe).

- siehe code
- d) Betrachten Sie die Klasse KNNRegressifier: Welche Art von Regressions-Modell berechnet diese Klasse?
- fast K-Nearest-Neighbor-Regression

Wozu dienen jeweils die Parameter K und flagKLinReg?

- K: Anzahl der NN die verwendet werden um eine Vorhersage zu treffen
- flagKLinReg: gibt an ob eine Regression verwendet werden soll oder nur der Mittelwert aus den NN genommen wird

Beschreiben Sie kurz in eigenen Worten (2-3 Sätze) auf welche Weise die Prädiktion y(x) berechnet wird.

- Es werden die KNNs für x ermittelt. Damit wird ein LSRRegressifier trainiert, der dann nach einer prediction gefragt wird.
- e) Betrachten Sie abschlieSSend den Modultest: Beschreiben Sie kurz was im Modultest passiert.
- es wird ein Linearer Datensatz erstellt

- die Funktion phi wird neu definiert und auf Grad 2 festgesetzt
- LSR Regression wird durchgeführt und Kreuzvalidiert
- KNN Regression wird durchgeführt und Kreuzvalidiert

Welche Gewichte W werden gelernt? Wie lautet also die gelernte Prädiktionsfunktion? Welche Funktion sollte sich idealerweise (für N) ergeben?

```
w0 = [3.78705442]
w1 = [2.01036841]
y = 3.787 + 2.010*x
y = 4 + 2*x \text{ für N}
```

Welche Ergebnisse liefert die Kreuzvalidierung? Was bedeuten die Werte?

- (E,sd,min,max)= (0.8335277318740052, 0.6423712068147653, 0.030450286307427632, 3.059745154084382)
- E ist der mittlere Fehlerwert
- sd ist die Standartabweichung der Fehler
- min ist der kleinste Fehlerwert
- max ist der gröSSte Fehlerwert

Vergleichen und Bewerten Sie die Ergebnisse von Least Squares Regression gegenüber der KNN-Regression (nach Optimierung der Hyper-Parameter , K, ...).

```
LSR:
                              (E,sd,min,max)=
                                                 (0.7210491003592423,
                                                                         0.5070165487566235,
          absolute
                     errors
0.0066007843635347285, 2.780939336314603)
KNN:
         absolute
                    errors
                             (E,sd,min,max)=
                                                 (1.1743092886650586,
                                                                         0.9170722237516987,
0.0038023455924474092, 4.2882465799825695)
                                                                       0.028216719179408916,
LSR:
       relative
                 errors
                           (E,sd,min,max)=
                                              (0.02240702968734219,
0.0001525998169486947, 0.14635559861181296)
```

KNN: relative errors (E,sd,min,max)= (0.05354319692710508, 0.13202983166413382, 5.0979643788569236e-05, 1.0603171407681475)

LSR scheint in fast allen Bereichen wesentlich bessere Ergebnisse als KNNR zu liefern.