IFT 6390 Fundamentals of Machine *Learning*

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Homework 1 - Practical Part

Parzen with soft windows (kernels)

In this Homework we will use *banknote authentication Data Set* as a toy dataset. It contains 1372 samples (one for each row), each with 4 features 1(the 4 first columns) and one label in {0,1} (the 5th column). It is recommended you download it here and then test your code by importing it like this:

```
import numpy as np
banknote = np.gen fromtxt('databanknote a uthentication.txt', delimiter=
',')
```

When the answer template in <u>solution.py</u> has "banknote" as an argument, you may assume that this argument is the dataset in numpy format. Your function should use this argument to perform computations, not a version of the dataset that you loaded by yourself.

```
In [4]: %matplotlib inline
        import numpy as np
        import random
        import matplotlib
        import matplotlib.pyplot as plt
        import time
        import sys
        IN_COLAB = 'google.colab' in sys.modules
        from numpy import argmax
        from keras.utils import to_categorical
        # FOR VISUALIZATION
        from ipywidgets import interact, SelectMultiple, fixed, Checkbox, IntRang
        eSlider, IntSlider, FloatSlider
        import ipywidgets as widgets
        matplotlib.rcParams['figure.figsize'] = [10,5]
        plt.style.use('fivethirtyeight')
```

```
In [5]: from google.colab import drive
    drive.mount('/content/drive', force_remount=True)

data = np.genfromtxt('/content/drive/My Drive/data_banknote_authenticatio
    n.txt', delimiter=',')

# Number of classes
label_list = np.unique(data[:,-1])
    n_classes = len(label_list)
```

Mounted at /content/drive

```
In [6]: print(data.shape)
```

```
[0. 1.]
In [7]: import numpy as np
        def split_dataset(banknote):
            #function split_dataset that splits the banknote dataset as follows:
            #•A training set consisting of the samples of the dataset with indice
        S
                #which have a remainder of either 0 or 1, or 2 when divided by 5.
            #•A validation set consisting of the samples of the dataset with indi
        ces
                #which have a remainder of 3 when divided by 5.
            #•A test set consisting of the samples of the dataset with indices
                #which have a remainder of 4 when divided by 5.
            data = banknote
            label_list = np.unique(data[:, -1])
            n_classes = len(np.unique(data[:,-1]))
            #sperate the indexes into three different sets
            train_indexes = [i for i in range(data.shape[0]) if i%5==0 or i%5==1
         or i%5==2]
            valid_indexes = [i for i in range(data.shape[0]) if i%5==3]
            test_indexes = [i for i in range(data.shape[0]) if i%5==4]
            #by the indexes of three sets, copy data into their sets
            train_set = data[train_indexes, :]
            valid_set = data[valid_indexes, :]
            test_set = data[test_indexes, :]
            #draw out the labels from train_set, valid_set and test_set.
            train_labels = train_set[:, -1].astype('int32')
            valid_labels = valid_set[:, -1].astype('int32')
            test_labels = test_set[:, -1].astype('int32')
            #draw out the feature from train_set, valid_set and test_set
            train_data = train_set[:, :-1]
            valid_data = valid_set[:, :-1]
            test_data = test_set[:, :-1]
            return train_data, train_labels, valid_data, valid_labels, test_data,
        test_labels, label_list, n_classes
In [8]: import numpy as np
        ####### DO NOT MODIFY THIS FUNCTION #######
        def draw_rand_label(x, label_list):
            seed = abs(np.sum(x))
            while seed < 1:</pre>
                seed = 10 * seed
            seed = int(10000000 * seed)
            np.random.seed(seed)
            return np.random.choice(label_list)
```

print(np.unique(data[:,-1]))

(1372, 5)

1.[4 points]

In [9]: def minkowski_mat(x, Y, p=2):

Question. Write functions that take that dataset as input and returnthe following statistics:

return (np.sum((np.abs(x - Y)) ** p, axis=1)) ** (1.0 / p)

- (a) Q1.feature_means : An array containing the empirical means of each feature, from all examples present in the dataset. Make sure to maintain the original order of features. e.g. : Q1.feature_means (banknote) = $[\mu_1, \mu_2, \mu_3, \mu_4]$
- (b) Q1.covariance_matrix : $A\ 4\times 4$ matrix that represents the em-pirical covariance matrix of features on the whole dataset.
- (c) Q1.feature_means_class_1 : An array containing the empiricalmeans of each feature, but only from examples in class 1. Thepossible classes in the banknote dataset are 0 and 1.e.g. : Q1.feature_means_class_1(banknote) = $[\mu_1, \mu_2, \mu_3, \mu_4]$.
- (d) Q1.covariance_matrix_class_1 : $A \ 4 \times 4$ matrix that represents the empirical covariance matrix of features, but only from exam-ples in class 1.

```
In [10]: import numpy as np
         class Q1:
             def feature_means(self, banknote):
                  # (a) Q1.feature_means : An array containing the empirical means
          of each feature,
                  # from all examples present in the dataset.
                  # Make sure to maintain the original order of features.
                  # e.g.: Q1.feature_means(banknote) = [\mu 1, \mu 2, \mu 3, \mu 4]
                  # banknote n*5
                  data = banknote[:, :-1]
                  \mu = \text{np.mean(data, axis=0)}
                  return μ
             def covariance_matrix(self, banknote):
                  # Q1.covariance_matrix : A 4×4 matrix that represents
                  # the em-pirical covariance matrix of features on the whole datas
         et.
                  data = banknote[:, :-1]
                  covMatrix = np.cov(data, rowvar=False)
                  return covMatrix
             def feature_means_class_1(self, banknote):
                  #Q1.feature_means_class_1 : An array containing the empiricalmean
         s of each feature, but only from examples in class 1.
                  #The possible classes in the banknote dataset are 0 and 1.e.g. :
          Q1. feature_means_class_1(banknote) = [\mu 1, \mu 2, \mu 3, \mu 4].
                  ind_class1 = [i for i in range(banknote.shape[0]) if banknote[i,
         -1]==1]
                             = np.mean(banknote[ind_class1, :-1], axis=0)
                  Сμ
                  return cµ
             def covariance_matrix_class_1(self, banknote):
                  #Q1.covariance_matrix_class_1 : A 4×4 matrix that represents t
         he empirical covariance matrix of features, but only from exam-ples in cl
         ass 1.
                  ind_class1 = [i for i in range(banknote.shape[0]) if banknote[i,
         -1]==1]
                  cCovMatrix = np.cov(banknote[ind_class1, :-1], rowvar=False)
                  return cCovMatrix
```

```
In [11]: q = Q1()
    print('print(q.feature_means(data)) = ')
    print(q.feature_means(data))
    print('print(q.covariance_matrix(data)) =')
    print(q.covariance_matrix(data))
```

```
print('print(q.feature_means_class_1(data)) = ')
print(q.feature_means_class_1(data))
print('q.covariance_matrix_class_1(data) = ')
print(q.covariance_matrix_class_1(data))
print(q.feature_means(data)) =
[ 0.43373526  1.92235312  1.39762712 -1.19165652]
print(q.covariance_matrix(data)) =
   8.08129912 4.40508287 -4.66632326
                                          1.65333797]
   4.40508287 34.44570968 -19.90511909 -6.490033
 [ -4.66632326 -19.90511909 18.57635938
                                         2.88724129]
   1.65333797 -6.490033
                           2.88724129
                                         4.4142562 ]]
print(q.feature_means_class_1(data)) =
[-1.86844256 -0.99357612 2.14827101 -1.24664075]
q.covariance_matrix_class_1(data) =
[[ 3.53884798  0.74923443  -4.69053721  1.26243851]
 [ 0.74923443 29.21276835 -25.24469813 -5.69675942]
 [ -4.69053721 -25.24469813 27.68665431 3.00778701]
   1.26243851 -5.69675942 3.00778701 4.28897428]]
```

1. [1 points]

Question. Implement Parzen with hard window parameter h. Use the standard Euclidean distance on the original features of the dataset. Your answer should have the following behavior:

f = HardParzen(h) initiates the algorithm with parameter h.

f.train(X, Y) trains the algorithm, where X is a $n \times m$ matrix of n training samples with m features, and Y is an array containing the n labels. The labels are denoted by integers, but the number of classes in Y can vary.

f.compute_predictions(X_test) computes the predicted labels and return them as an array of same size as X_test. X_test is a k × m matrix of k test samples with same number of features as X. This function is called only after training on (X,Y). If a test sample x has no neighbor within window h, the algorithm should choose a label at random by using **draw_rand_label(x, label_list)**, a function that is provided in the **solution.py** file, where label_list is the list of different classes present in Y , and x is the array of features of the corresponding point.

```
In [13]:
         class HardParzen:
             def __init__(self, h=0.4, dist_func=minkowski_mat):
                 #f = HardParzen(h) initiates the algorithm with parameter h
                 self.h = h #h is the threshold distance, h is a positive real.
                 self.dist_func = dist_func
             def train(self, train_inputs, train_labels):
                 #f.train(X, Y)^* trains the algorithm, where X is a n × m matrix
          of n training samples with m features,
                 #and Y is an array containing the n labels.
                 #The labels are denoted by integers, but the number of classes in
         Y can vary.
                 self.train_inputs = train_inputs
                 self.train_labels = train_labels
                 self.label_list = np.unique(data[:,-1]) #the elements of self.l
         abel_list are monotonically increasing
                 self.n_classes = len(self.label_list)
             def compute_predictions(self, test_data):
                 # f.compute_predictions(X_test) computes the predicted labels and
         return them
                 # as an array of same size as X_test.
                 \# X test is a k \times m matrix of k test samples with same number of
          features as X.
                 # This function is called only after training on (X, Y).
                 # If a test sample x has no neighbor within window h,
```

```
# the algorithm should choose a label at random by using draw_ran
        d_{abel}(x, label_{list}),
               # a function that is provided in the solution.py file, where labe
        l_list is the list of different classes present in Y ,
               \# and x is the array of features of the corresponding point.
               #counts n \times n_{classes} matrix to save the voting number of differe
        nt classes.
               #in counts, every elment initially set 0
               pred_test_labels = np.zeros((test_data.shape[0]))
               # For each test datapoint
               for (i, ex) in enumerate(test_data):
                   # i is the row index
                   # ex is the i'th row
                   # Find the distances between ex to training inputs point usin
        g dist_func
                                 = minkowski_mat(ex, self.train_inputs, p=2)
                   distances
                   #find the neighbour index with distance less than self.h
                   neighbour_index = [j for j in range(len(distances)) if distan
        ces[j] < self.h]</pre>
                   if len(neighbour_index)== 0:
                    k = draw_rand_label(ex, self.label_list) # k = a random cla
        ss for ex
                    pred_test_labels[i] = k
                   else:
                    neighb_train_labels = self.train_labels[neighbour_index]
                    # create an array count_class len(self.label_list) to save
         the count classes.
                    pre_class = self.label_list[0]
                    max\_count = 0
                    for (m,a) in enumerate(self.label_list):
                      # for each ex, count number of the different classes.
                      index_neigh_set = [j for j in range(len(neighb_train_labe)
        ls)) if neighb_train_labels[j]==a]
                      if len(index_neigh_set) > max_count:
                        max_count = len(index_neigh_set)
                        pre_class = self.label_list[m]
                    pred_test_labels[i] = pre_class
               pred_test_labels_int = pred_test_labels.astype(int)
               return pred_test_labels_int
In [14]: hd_parzen = HardParzen(h=3)
        train_data, train_labels, valid_data, valid_labels, test_data, test_label
        s, label_list, n_classes = split_dataset(data)
        hd_parzen.train(train_data, train_labels)
        hd_pred = hd_parzen.compute_predictions(test_data)
        print('test_labels = ')
        print(test_labels)
        print('hd_pred = ')
        print(hd_pred)
        test_labels =
```

```
1
1
1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1
hd_pred =
1
1
1
111111111111111
```

1. [5 points] Question. Implement Parzen with a soft window. We will use a radial basis function (RBF) kernel (also known as *Gaussian* kernel) with parameter σ. Use the standard Euclidean distance on the original features of the dataset. Please refer to the slides from the second week for the definition. The structure of your solution should be the same as in the previous question, but you will never need to draw a label at random with draw_rand_label(x, label_list). The class name is SoftRBFParzen.

```
In [15]:
         import math
         import numpy as np
         from numpy import argmax
         from keras.utils import to_categorical
         class SoftRBFParzen:
             def __init__(self, sigma_sq=1.0, n_dims=4):
                 self.n_dims = n_dims
                 self.mu = np.zeros(n_dims)
                 self.sigma_sq = sigma_sq #sigma_sq is the square of sigma.
             def train(self, train_inputs, train_labels):
                 # self.label_list = np.unique(train_labels)
                 self.train_inputs = train_inputs
                 self.train_labels = train_labels
                 self.label_list = np.unique(data[:, -1])
                 self.n_classes = len(self.label_list)
                 #onehot_train_lables is for predicting the class of test
                 #the function to_categorical() to code onehot
                 self.onehot_train_labels = to_categorical(self.train_labels)
                 # Here, calculate the mean and variance of the train_data data an
         d put it in self.mu
                 # self.sigma_sq is given
                 self.mu = np.mean(self.train_inputs, axis=0)
                 #self.sigma_sq = np.sum((self.train_inputs - self.mu) ** 2.0) /
          (self.n_dims * train_inputs.shape[0])
```

```
def GaussianKernel(self, dist=0):
                 #calculate Gaussian Kernel
                 #dist is the distance between two points, dim is the dimension nu
         mber
                 part1 = 1/ ((2*math.pi)**(self.n_dims/2)*(self.sigma_sq**self.n_d
         ims))
                 part2 = math.exp(-1/2*(dist**2)/self.sigma_sq)
                 return part1*part2
             def compute_predictions(self, test_data):
                 # Implement Parzen with a soft window.
                 # We will use a radial basis function (RBF) kernel (also known as
         Gaussian kernel) with parameter \sigma.
                 # Given parameter \sigma
                 classes_pred = np.zeros(test_data.shape[0], dtype=int)-1 #classes
         _pred initialize -1
                 self.onehot_train_lables = to_categorical(self.train_labels)
                 # For each test datapoint
                 for (i, ex) in enumerate(test_data):
                     # i is the row index
                     # ex is the i'th row test data
                     #count train_inputs.shape[0] by n_classes,
                     #it is used to record the multiplication between GaussianKern
         el and onehot-coding
                     counts = np.zeros_like(self.onehot_train_lables, dtype=float)
         #
                     # Find the distances to each training set point using dist_fu
         nc
                     #distances = self.dist_func(ex, self.train_inputs)
                     distances = minkowski_mat(ex, self.train_inputs, p=2)
                     total_kernel = 0.0
                     for (j, dist) in enumerate(distances):
                        # Go through the training set to calculate GaussianKernel
          * onehot_train_labels[i]
                        # Implement SoftRBFParzen with hard window parameter self.
         mu and self.sigma_sq here.
                        kernel
                                 = self.GaussianKernel(dist)
                        total kernel += kernel
                        #the two way to calculate counts
                        counts[j] = kernel * self.onehot_train_labels[j]
                     sum_counts = np.sum(counts, axis=0) / total_kernel
                     classes_pred[i] = self.label_list[np.argmax(sum_counts)] #n
         p.argmax return the index of max element in sum_counts, it is the index o
         f class.
                 return classes_pred
In [16]: softRBF = SoftRBFParzen(0.2)
         train_data, train_labels, valid_data, valid_labels, test_data, test_label
         s, label_list, n_classes = split_dataset(data)
         softRBF.train(train_data, train_labels)
         classes_pred = softRBF.compute_predictions(test_data)
         print('test_labels = ')
         print(test_labels)
         print('classes_pred = ')
         print(classes_pred)
```

test_labels =

```
0
1
1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1
classes_pred =
0
1
1
1
1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1
```

1. [5 points]

ט טן

Question. Implement a function **split_dataset** that splits the banknote dataset as follows:

- A training set consisting of the samples of the dataset with indices which have a remainder of either 0 or 1, or 2 when divided by 5.
- A validation set consisting of the samples of the dataset with indices which have a remainder of 3 when divided by 5.
- A test set consisting of the samples of the dataset with indices which have a remainder of 4 when divided by 5.

For instance the element of index 14 (in the original dataset) should be part of the test set because the remainder of 14 divided by 5 is 4. Do not use random splitting for this exercise (even though it is generally a very good idea). The function should take as input the dataset and return the three sets as a tuple (train, validation, test), where each element of the tuple is a matrix with 5 columns (the 4 features and the labels, kept in the same order).

```
In [17]: import numpy as np

def split_dataset(banknote):
    #function split_dataset that splits the banknote dataset as follows:
    #*A training set consisting of the samples of the dataset with indice
s
    #which have a remainder of either 0 or 1, or 2 when divided by 5.
    #*A validation set consisting of the samples of the dataset with indices
    #which have a remainder of 3 when divided by 5.
#*A test set consisting of the samples of the dataset with indices
    #which have a remainder of 4 when divided by 5.
data = banknote
label_list = np.unique(data[:, -1])
n_classes = len(np.unique(data[:, -1]))
```

```
#sperate the indexes into three different sets
   train_indexes = [i for i in range(data.shape[0]) if i%5==0 or i%5==1
or i%5==2]
   valid_indexes = [i for i in range(data.shape[0]) if i%5==3]
    test_indexes = [i for i in range(data.shape[0]) if i%5==4]
   #by the indexes of three sets, copy data into their sets
   train_set = data[train_indexes, :]
   valid_set = data[valid_indexes, :]
   test_set = data[test_indexes, :]
   #draw out the labels from train_set, valid_set and test_set.
    train_labels = train_set[:, -1].astype('int32')
   valid_labels = valid_set[:, -1].astype('int32')
    test_labels = test_set[:, -1].astype('int32')
   #draw out the feature from train_set, valid_set and test_set
   train_data = train_set[:, :-1]
   valid_data = valid_set[:, :-1]
   test_data = test_set[:, :-1]
   return train_data, train_labels, valid_data, valid_labels, test_data,
test_labels, label_list, n_classes
```

```
In [18]: train_data, train_labels, valid_data, valid_labels, test_data, test_label
    s, label_list, n_classes = split_dataset(data)
    print('data.shape = ', data.shape)
    print('train_data.shape = ', train_data.shape)
    print('valid_data.shape = ', valid_data.shape)
    print('test_data.shape = ', test_data.shape)

data.shape = (1372, 5)
    train_data.shape = (824, 4)
    valid_data.shape = (274, 4)
    test_data.shape = (274, 4)
```

1. [10 points] Question. Implement two functions ErrorRate.hard_parzen and ErrorRate.soft_parzen that compute the error rate (i.e. the proportion of missclassifications) of the HardParzen and SoftRBFParzen algorithms. The expected behavior is as follows: test_error = ErrorRate(x_train, y_train, x_val, y_val) initiates the class and stores the

test_error = ErrorRate(x_train, y_train, x_val, y_val) initiates the class and stores the training and validation sets, where x_t are matrices with 4 feature columns, and y_t are arrays containing the labels.

test_error.hard_parzen(h) takes as input the window parameter h and returns as a float error rate on x_val and y_val of the Hard- Parzen algorithm that has been trained on x_train and y_train. **test_error.soft_parzen(\sigma)** works just like with Hard Parzen, but with the SoftRBFParzen algorithm.

Then, include in your report a single plot with two lines:

(a) Hard Parzen window's classification error on the validation set of banknote, when trained on the training set (see question 4) for the following values of h:

```
h \in \{0.01, 0.1, 0.2, 0.3, 0.4, 0.5, 1.0, 3.0, 10.0, 20.0\}
```

(b) RBF Parzen's classification error on the validation set of ban- knote, when trained on the training set (see question 4) for the following values of σ :

```
\sigma \in \{0.01, 0.1, 0.2, 0.3, 0.4, 0.5, 1.0, 3.0, 10.0, 20.0\}
```

The common x-axis will represent either h or σ . Always label your axes and lines in the plot! Give a detailed discussion of your observations.

```
def confusion_matrix(true_labels, pred_labels):
    matrix = np.zeros((n_classes, n_classes))
    for (true, pred) in zip(true_labels, pred_labels):
        matrix[int(true - 1), int(pred - 1)] += 1
    return matrix

def comput_test_error(conf_mat):
    #compute test_error from n by n comfusion_matrix
    sum_preds = np.sum(conf_mat)
    sum_correct = np.sum(np.diag(conf_mat))
    return 1.0-float(sum_correct)/float(sum_preds)

#the following is for testing the above functions confusion matrix() and
```

```
In [20]: #the following is for testing the above functions confusion_matrix() and
    comput_test_error()
    import numpy as np

a = np.array([1, 2, 1, 2, 1])
    b = np.array([2, 1, 1, 1, 2])
    matrix = confusion_matrix(a, b)

error = comput_test_error(matrix)
    print(error)
```

0.8

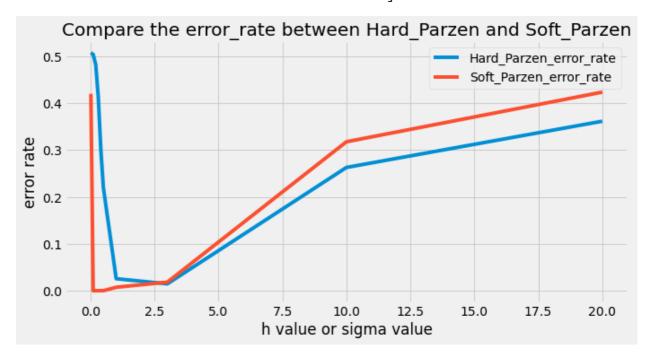
```
In [21]:
         class ErrorRate:
             def __init__(self, x_train, y_train, x_val, y_val):
                 # initiates the class and stores the training and validation set
         S,
                 # where x_train and y_train are matrices with 4 feature columns,
                 # and x_{val} and y_{val} are arrays containing the labels.
                 self.x_train = x_train
                 self.y\_train = y\_train
                 self.x_val = x_val
                 self.y_val
                              = y_val
                 self.h
                          = 1.0
                 self.sigma_sq = 1.0
             def hard_parzen(self, h):
                 #takes as input the window parameter h and
                 #returns as a float the error rate on x_val and y_val of the Hard
         Parzen algorithm
                 #that has been trained on x_train and y_train.
                 self.h = h
                 pass
                 x_hard_parzen
                                    = HardParzen(self.h)
                 x_hard_parzen.train(self.x_train, self.x_val)
                 y_pred_test_labels = x_hard_parzen.compute_predictions(self.y_tra
         in)
                 y_confusion_matrix = confusion_matrix(self.y_val, y_pred_test_lab
         els)
                 y_error_rate
                                    = comput_test_error(y_confusion_matrix)
                                    = HardParzen(self.h)
                 y_hard_parzen
                 y_hard_parzen.train(self.y_train, self.y_val)
                 x_pred_test_labels = y_hard_parzen.compute_predictions(self.x_tra
         in)
                 x_confusion_matrix = confusion_matrix(self.x_val, x_pred_test_lab
         els)
```

```
x_error_rate
                          = comput_test_error(x_confusion_matrix)
        return y_error_rate, x_error_rate
    def soft_parzen(self, sigma):
        #takes as input the parameter sigma and
        #returns as a float the error rate on x_val and y_val of the Soft
RBFParzen algorithm
        #that has been trained on x_train and y_train.
        self.sigma_sq = sigma**2
        pass
        x_soft_RBFParzen = SoftRBFParzen(self.sigma_sq)
        x_soft_RBFParzen.train(self.x_train, self.x_val)
        y_pred_test_labels = x_soft_RBFParzen.compute_predictions(self.y_
train)
       y_confusion_matrix = confusion_matrix(self.y_val, y_pred_test_lab
els)
       y_error_rate
                          = comput_test_error(y_confusion_matrix)
        y_soft_RBFParzen = SoftRBFParzen(self.sigma_sq)
        y_soft_RBFParzen.train(self.y_train, self.y_val)
        x_pred_test_labels = y_soft_RBFParzen.compute_predictions(self.x_
train)
       x_confusion_matrix = confusion_matrix(self.x_val, x_pred_test_lab
els)
                           = comput_test_error(x_confusion_matrix)
       x_error_rate
        return y_error_rate, x_error_rate
s, label_list, n_classes = split_dataset(data)
arr_h
           = np.array([0.01, 0.1, 0.2, 0.3, 0.4, 0.5, 1.0, 3.0, 10.0, 20.
```

```
In [22]: train_data, train_labels, valid_data, valid_labels, test_data, test_label
         0])
         arr\_sigma = np.array([0.01, 0.1, 0.2, 0.3, 0.4, 0.5, 1.0, 3.0, 10.0, 20.
         0])
         # create a ErrorRate class and initialize it
         error_rate = ErrorRate(train_data, valid_data, train_labels, valid_labels
         )
         hp_error_rate = np.zeros_like(arr_h, dtype=float)
         sp_error_rate = np.zeros_like(arr_sigma, dtype=float)
         for i in range(len(arr_h)):
           hp_val_error_rate, hp_train_error_rate = error_rate.hard_parzen(arr_h[i
         ])
           hp_error_rate[i] = hp_val_error_rate
         for j in range(len(arr_sigma)):
           sp_val_error_rate, sp_train_error_rate = error_rate.soft_parzen(arr_sig
         ma[j])
           sp_error_rate[j] = sp_val_error_rate
         print(hp_error_rate)
         print(sp_error_rate)
         #draw a single plot with two lines
         plt.plot(arr_h, hp_error_rate, label='Hard_Parzen_error_rate')
         plt.plot(arr_h, sp_error_rate, label='Soft_Parzen_error_rate')
         plt.xlabel("h value or sigma value")
         plt.ylabel('error rate')
         plt.title("Compare the error_rate between Hard_Parzen and Soft_Parzen")
         plt.legend()
         plt.show()
```

/usr/local/lib/python3.6/dist-packages/ipykernel_launcher.py:68: RuntimeW arning: invalid value encountered in true_divide

```
[0.50729927 0.50364964 0.48175182 0.41240876 0.29927007 0.2189781 0.02554745 0.01459854 0.26277372 0.36131387] [0.41970803 0. 0. 0. 0. 0. 0. 0. 0. 0. 0.00729927 0.01824818 0.31751825 0.42335766]
```



Give a detailed discussion of your observations.

Answer: For hard parzen algorithm, when h=2.5,3.0, it will obtain the lowest error rate. When h<2.5, the less h is, the more error rate is. When h>3.0, the bigger h is, the more error rate is. That is to say, distance parameter h is too small and too big, that cause error rate to increase.

For soft parzen algorithm, when $\sigma=0.1,0.2,0.3,0.4,0.5$, error rate is near 0. When sigma=0.01, error rate is very big.Furthermore, when sigma>0.5, the bigger sigma is, the larger error rate is. That is to say, the proper sigma parameter $\sigma\in[0.1,0.2,0.3,0.4,0.5]$

1. [5 points] **Question**. Implement a function **get_test_errors** that uses the evaluated validation errors from the previous question to select h^* and σ^* , then computes the error rates on the test set. The value h^* is the one (among the proposed set in question 5) that results in the smallest validation error for Parzen with hard window, and σ^* is the parameter (among the proposed set in question 5) that results in the smallest validation error for Parzen with RBF. The function should take as input the dataset and split it using question 4. The expected output is an array of size 2, the first value being the error rate on the test set of Hard Parzen with parameter h^* (trained on the training set), and the second value being the error rate on the test set of Soft RBF Parzen with parameter σ^* (trained on the training set).

```
In [23]: def get_test_errors(banknote, star_h=3.0, star_sigma=0.3):
    pass

# function get_test_errors that uses the evaluated validation errors
# from the previous question to select h* and \(\sigma^*\),
# then computes the error rates on the test set.
# The value h* is the one (among the proposed set in question 5)
# that results in the smallest validation error for Parzen with hard
window.
# \(\sigma^*\) is the parameter (among the proposed set in question 5)
# that results in the smallest validation error for Parzen with RBF.

train_data, train_labels, valid_data, valid_labels, test_data, test_l
abels, label_list, n_classes = split_dataset(data)
```

the value star_h is the one (among the proposed set in question 5)

```
# that results in the smallest validation error for Parzen with hard
window
   x_hard_parzen = HardParzen(star_h)
    x_hard_parzen.train(train_data, train_labels)
    y_hp_pred_test_lab = x_hard_parzen.compute_predictions(test_data)
   y_hp_conf_matrix = confusion_matrix(test_labels, y_hp_pred_test_lab
)
   y_hp_error_rate = comput_test_error(y_hp_conf_matrix)
    star_sigma_sq = star_sigma**2
    # \sigma* is the parameter (among the proposed set in question 5)
    # that results in the smallest validation error for Parzen with RBF.
   x_soft_RBFParzen = SoftRBFParzen(star_sigma_sq)
    x_soft_RBFParzen.train(train_data, train_labels)
   y_soft_pred_test_lab = x_soft_RBFParzen.compute_predictions(test_data
)
   y_soft_conf_matrix = confusion_matrix(test_labels, y_soft_pred_test
_lab)
   y_soft_error_rate = comput_test_error(y_soft_conf_matrix)
    # expected output is an array of size 2,
   # the first value being the error rate on the test set of Hard Parzen
with parameter h*
    # the second value being the error rate on the test set of Soft RBF P
arzen with parameter \sigma*
   list_error_rate = [y_hp_error_rate, y_soft_error_rate]
    hp_sp_error_rate = np.array(list_error_rate, dtype=float)
    return hp_sp_error_rate
```

```
In [26]: get_test_errors(data, star_h=3.0, star_sigma=0.1)
Out[26]: array([0.00729927, 0. ])
```

1. [5 points] **Question**. Include in your report a discussion on the running time complexity of these two methods. How does it vary for each method when the hyperparameter h or σ changes? Why?

```
In [32]: import time
         from datetime import timedelta
         train_data, train_labels, valid_data, valid_labels, test_data, test_label
         s, label_list, n_classes = split_dataset(data)
         repeat_error_rate = ErrorRate(train_data, valid_data, train_labels, valid
         _labels)
         #print(repeat_error_rate.hard_parzen(arr_sigma[i]))
         arr_h = np.array([0.01, 0.1, 0.2, 0.3, 0.4, 0.5, 1.0, 3.0, 10.0, 20.0])
         hp_time = np.zeros_like(arr_h, dtype=float)
         for i in range(len(arr_h)):
           hp_start_time = time.time()
           repeat_error_rate.hard_parzen(arr_h[i])
           hp_time[i] = time.time() - hp_start_time
         arr_sigma = np.array([0.01, 0.1, 0.2, 0.3, 0.4, 0.5, 1.0, 3.0, 10.0, 20.0)
         sp_time = np.zeros_like(arr_sigma, dtype=float)
         for j in range(len(arr_sigma)):
           sp_start_time = time.time()
           repeat_error_rate.soft_parzen(arr_sigma[j])
                      = time.time() - sp_start_time
           sp_time[j]
```

```
print("hp_time = ")
print(hp_time)
print("sp_time = ")
print(sp_time)
```

/usr/local/lib/python3.6/dist-packages/ipykernel_launcher.py:68: RuntimeW arning: invalid value encountered in true_divide

```
hp_time =
[0.18668962 0.18660641 0.18985415 0.17572403 0.17302465 0.17333174
0.17466164 0.27933335 1.27485037 1.98644376]
sp_time =
[2.7162919 2.70941401 2.73479557 2.67663622 2.68910503 2.67660809
2.62141061 2.62936401 2.62370086 2.60773039]
```

How does it vary for each method when the hyperparameter h or σ changes? Why?

Answer: For hard Parzen, With the distance h increasing, time-cost increases. Because when h becomes large, the number of the involved points rises up. So the algorithm complexity shoot up. But for **soft Parzen**, the cost time of soft parzen doesn't change much, with the increasement of σ . Because every time, it computes all points, which doesn't change. So the algorithm complexity keeps nearly stable.

1. [5 points] Question.

Implement a random projection (Gaussian sketch) map to be used on the input data: Your function $project_data$ should accept as input a feature matrix X of dimension $n \times 4$, as well as a 4×2 matrix A encoding our projection.

Define $p: x \mapsto \frac{1}{\sqrt{2}}A^Tx$ and use this random projection map to reduce the dimension of the inputs (feature vectors of the dataset) from 4 to 2.

Your function should return the output of the map p when applied to X, in the form of a $n \times 2$ matrix.

```
e.g. project\_data(X_{n,4},A_{4,2}) = X_{n,2}^{proj}
```

```
import numpy as np
import math

def random_projections(X, A):
    pass
    proj_X = np.dot(X, A)/math.sqrt(2)
    return proj_X
```

```
In [34]: # the following is for testing function random_projections()
    x = np.array([[1, 2, 3, 4], [5, 6, 7, 8], [9, 10, 11, 12]])
    A = np.array([[0.1, 0.2], [0.3, 0.4], [0.5, 0.6], [0.7, 0.8]])
    proj = random_projections(x, A)

    print(x)
    print(type(x))
    print(d)
    print(proj)
```

[12.586500/1 15.55634919]]

1. [10 points] **Question**. Similar to Question 5, compute the validation errors of Hard Parzen classifiers trained on 500 random projections of the training set, for

```
h \in \{0.01, 0.1, 0.2, 0.3, 0.4, 0.5, 1.0, 3.0, 10.0, 20.0\}
```

The validation errors should be computed on the projected validation set, using the same matrix A. To obtain random projections, you may draw A as 8 independent variables drawn uniformly from a gaussian distribution of mean 0 and variance 1.

You can for example store these validation errors in a 500×9 matrix, with a row for each random projection and a column for each value of h.

Do the same thing for RBF Parzen classifiers, for

```
\sigma \in 0.01, 0.1, 0.2, 0.3, 0.4, 0.5, 1.0, 3.0, 10.0, 20.0
```

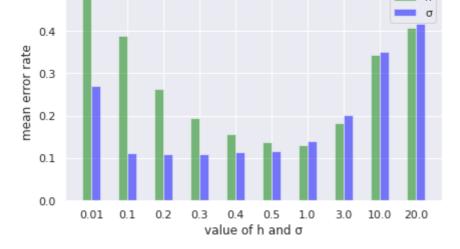
Plot and include in your report in the same graph the average values of the validation errors (over all random projections) for each value of h and σ , along with error bars of length equal to 0.2× the standard deviations.

How do your results compare to the previous ones?

```
In [35]:
         import random
         import numpy
         def create_mat_A(mu=0.0, sigma=1.0):
           list = [random.normalvariate(mu, sigma) for i in range(8)]
           mat_A = np.array(list).reshape((4, 2))
           return mat_A
In [58]: import numpy as np
         import matplotlib.pyplot as plt
         import matplotlib as mpl
         import seaborn as sns
         mu, sigma = 0.0, 1.0
         train_data, train_labels, valid_data, valid_labels, test_data, test_label
         s, label_list, n_classes = split_dataset(data)
                   = np.array([0.01, 0.1, 0.2, 0.3, 0.4, 0.5, 1.0, 3.0, 10.0, 20.0)
         arr_h
         arr_sigma = np.array([0.01, 0.1, 0.2, 0.3, 0.4, 0.5, 1.0, 3.0, 10.0, 20.0)
         ])
         #store these validation errors in a 500 \times 10 matrix
         pj_hp_valid_errors = np.zeros((500,10), dtype=float)
         pj_sp_valid_errors = np.zeros((500,10), dtype=float)
         for j in range(500):
           # Implement a random projection (Gaussian sketch) map to be used on the
         input data
                           = create_mat_A(0.0, 1.0)
           mat_A
           proj_train_data = random_projections(train_data, mat_A)
           proj_valid_data = random_projections(valid_data, mat_A)
           proj_test_data = random_projections(test_data, mat_A)
           proj_error_rate = ErrorRate(proj_train_data, proj_valid_data,
                                        train_labels, valid_labels)
           for k in range(len(arr_h)):
             pj_hp_val_error_rate, pj_hp_train_error_rate = proj_error_rate.hard_p
         arzen(arr_h[k])
             pj_hp_valid_errors[j, k] = pj_hp_val_error_rate
```

```
for m in range(len(arr_sigma)):
    pj_sp_val_error_rate, pj_sp_train_error_rate = proj_error_rate.soft_p
arzen(arr_sigma[m])
    pj_sp_valid_errors[j, m] = pj_sp_val_error_rate
mean_error_rate_every_h = np.mean(pj_hp_valid_errors, axis=0)
mean_error_rate_every_sigma = np.mean(pj_sp_valid_errors, axis=0)
std_error_rate_every_h = np.std(pj_hp_valid_errors, axis=0)
std_error_rate_every_sigam = np.std(pj_sp_valid_errors, axis=0)
print("mean_error_rate_every_h = ")
print(mean_error_rate_every_h)
print("mean_error_rate_every_sigma = ")
print(mean_error_rate_every_sigma)
# ----- begin drawing plot-----
# Plot Plot and include in your report in the same graph the average valu
es of the validation errors
# (over all random projections) for each value of h and \sigma,
# along with error bars of length equal to 0.2× the standard deviations
sns.set(color_codes=True)
mpl.rcParams["font.sans-serif"] = ["SimHei"]
mpl.rcParams["axes.unicode_minus"] = False
# parameter of bar
Y1_mean_h = mean_error_rate_every_h
Y2_mean_sigma = mean_error_rate_every_sigma
        = np.arange(len(arr_h))
# arr_h & arr_sigma
bar_width = 0.25
tick_label = ['0.01', '0.1', '0.2', '0.3', '0.4', '0.5', '1.0', '3.0', '1
0.0', '20.0']
# draw the bar
plt.bar(X, Y1_mean_h, bar_width, align="center", color="green", label="h"
, alpha=0.5)
plt.bar(X+bar_width, Y2_mean_sigma, bar_width, color="blue", align="cente
r", \
        label="\sigma", alpha=0.5)
plt.xlabel("value of h and \sigma")
plt.ylabel("mean error rate")
plt.title('validation errors for each h and \sigma')
plt.xticks(X+bar_width/2, tick_label)
plt.legend()
plt.show()
plt.savefig('result.png',dpi = 400)
/usr/local/lib/python3.6/dist-packages/ipykernel_launcher.py:68: RuntimeW
arning: invalid value encountered in true_divide
mean_error_rate_every_h =
[0.48983212\ 0.38908029\ 0.26436496\ 0.19394891\ 0.15684672\ 0.13865693
0.13018978 0.18288321 0.34280292 0.4069854 ]
mean_error_rate_every_sigma =
\lceil 0.2709781 \quad 0.11218978 \quad 0.10920438 \quad 0.10972993 \quad 0.11321898 \quad 0.11757664
0.1399562 0.20151095 0.35027737 0.41785401
```

h h



<Figure size 432x288 with 0 Axes>

How do your results compare to the previous ones?

- (1) Comparing with the previous ones, the entire plot of this result is similar for hard parzen algorithm and soft parzen algorithm.
- (2) But the lowest points have a little change. For hard parzen algorith, h=0.5,1.0, their mean error rates obtain the lowest point. But in the previous, h=1.0,3.0, their mean error rates is lowest.
- (3) For soft parzen algorith, this results is same as the previous ones. When $\sigma=0.1,0.2,0.3,0.4,0.5$, their mean error rates is lowest. (4) From this example, we find that the way of random projections of the data set, can keep the relative distance between the points, reduce the algorithm complexity.