Mustafa Yesilyurt, CS 178, Prof. Alexander Ihler, HW 1, 10/9/19

Problem 1: Python & Data Exploration

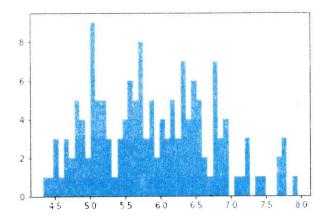
```
In [2]: import numpy as np
    import matplotlib.pyplot as plt

iris = np.genfromtxt("data/iris.txt",delimiter=None) # load the text file
    Y = iris[:,-1] # target value (iris species) is the last column
    X = iris[:,0:-1] # features are the other columns
```

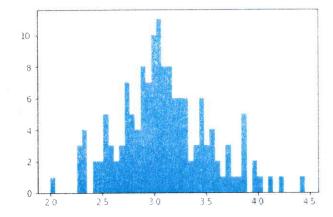
First question: Use X.shape to get the number of features and the number of data points. Report both numbers, mentioning which number is which. (5 points)

Second question: For each feature, plot a histogram (plt.hist) of the data values. (5 points)

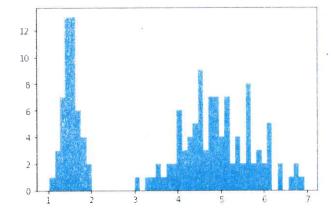
histogram for feature 1



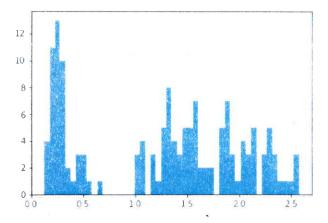
histogram for feature 2



histogram for feature 3



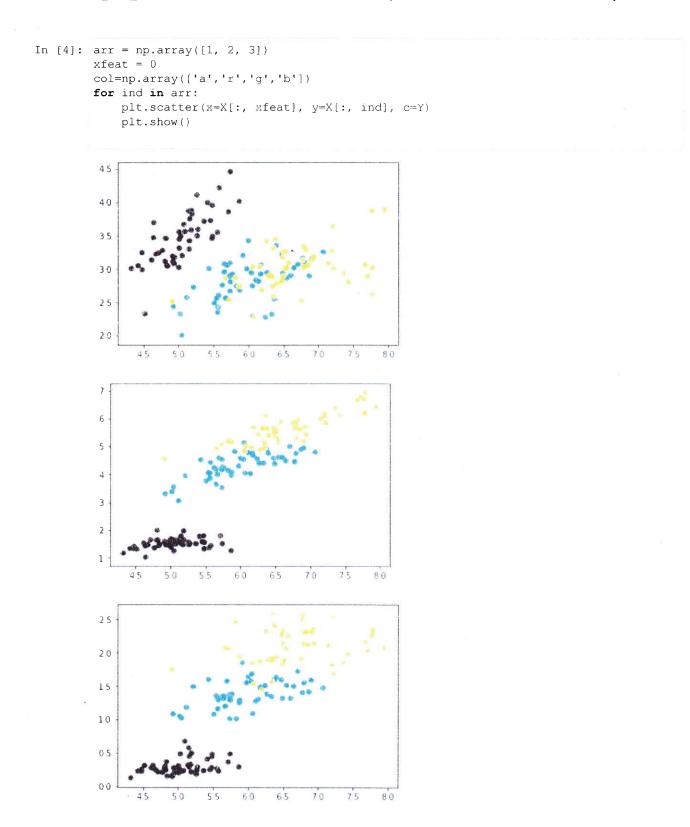
histogram for feature 4



Third question: Compute the mean & standard deviation of the data points for each feature (np.mean,np.std).(5 points)

```
In [4]: for ind in range(4):
            print('mean for feature ' + str(ind+1))
            print(str(np.mean(X[:, ind])))
            print('standard deviation for feature ' + str(ind+1))
            print(str(np.std(X[:, ind])))
            print()
        mean for feature 1
        5.90010376419
        standard deviation for feature 1
        0.833402066775
        mean for feature 2
        3.09893091689
        standard deviation for feature 2
        0.436291838001
        mean for feature 3
        3.81955484054
        standard deviation for feature 3
        1.75405710934
        mean for feature 4
        1.25255548459
        standard deviation for feature 4
        0.758772457026
```

Fourth question: For each pair of features (1,2), (1,3), and (1,4), plot a scatterplot (see plt.plot or plt.scatter) of the feature values, colored according to their target value (class). (For example, plot all data points with y=0 as blue,y=1 as green, andy=2 as red.)(5 points)

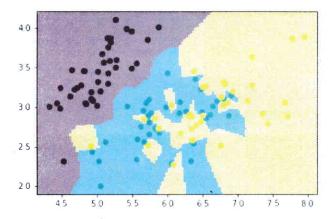


Problem 2: k-nearest-neighbor predictions (25 points)

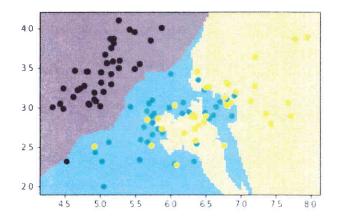
First problem: Modify the code in the next cell to use only the first two features of X(e.g., let X be only the first two columns of iris, instead of the first four), and visualize (plot) the classification boundary for varying values of K=[1, 5, 10, 50] using plotClassify2D.(10 points)

```
In [6]: | iris = np.genfromtxt("data/iris.txt", delimiter=None) # load the data
        Y = iris[:,-1]
       X = iris[:, 0:-1]
        # Note: indexing with ":" indicates all values (in this case, all rows);
        # indexing with a value ("0", "1", "-1", etc.) extracts only that value (here, col
        # indexing rows/columns with a range ("l:-1") extracts any row/column in that rang
        e.
        import mltools as ml
        # We'll use some data manipulation routines in the provided class code
        # Make sure the "mltools" directory is in a directory on your Python path, e.g.,
          export PYTHONPATH=$\$${PYTHONPATH}:/path/to/parent/dir
        # or add it to your path inside Python:
          import sys
          sys.path.append('/path/to/parent/dir/');
        np.random.seed(0)
                                    # set the random number seed
        X,Y = ml.shuffleData(X[:, 0:2],Y); # shuffle data randomly TAKING ONLY FIRST TWO C
        OLUMNS OF X
        # (This is a good idea in case your data are ordered in some systematic way.)
        Xtr,Xva,Ytr,Yva = ml.splitData(X,Y, 0.75); # split data into 75/25 train/validatio
        #-----
        #knn = ml.knn.knnClassify() # create the object and train it
        #knn.train(Xtr, Ytr, 1)  # where K is an integer, e.g. 1 for nearest neighbor p
        rediction
        #YvaHat = knn.predict(Xva)  # get estimates of y for each data point in Xva
        # Alternatively, the constructor provides a shortcut to "train":
        #knn = ml.knn.knnClassify( Xtr, Ytr, K );
        #YvaHat = predict( knn, Xva );
        #-----
        print("k = 1")
        knn = ml.knn.knnClassify()
        knn.train(Xtr, Ytr, 1)
        ml.plotClassify2D( knn, Xtr, Ytr ); # make 2D classification plot with data (Xtr,Y
        tr)
        plt.show()
        print("k = 5")
        knn = ml.knn.knnClassify()
        knn.train(Xtr, Ytr, 5)
        ml.plotClassify2D( knn, Xtr, Ytr ); # make 2D classification plot with data (Xtr,Y
        t.r)
        plt.show()
        print("k = 10")
        knn = ml.knn.knnClassify()
        knn.train(Xtr, Ytr, 10)
        ml.plotClassify2D( knn, Xtr, Ytr ); # make 2D classification plot with data (Xtr,Y
        tr)
        plt.show()
        print("k = 50")
        knn = ml.knn.knnClassify()
        knn.train(Xtr, Ytr, 50)
        plt.show()
        ml.plotClassify2D( knn, Xtr, Ytr ); # make 2D classification plot with data (Xtr,Y
        plt.show()
```

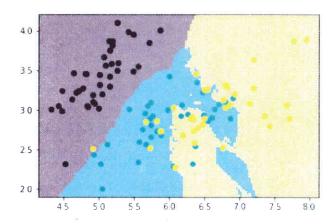




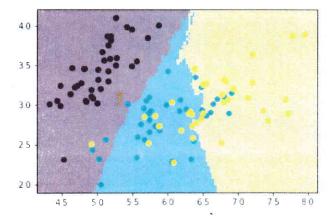
k = 5



k = 10

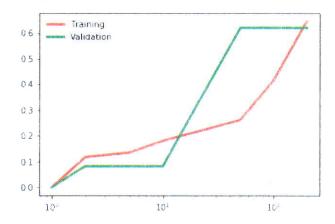


k = 50



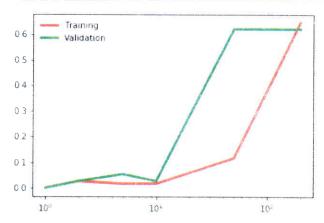
Second question: Again using only the first two features, compute the error rate (number of misclassifications) on both the training and validation data as a function of K=[1,2,5,10,50,100,200]. You can do this most easily with a for-loop:

```
In [9]: iris = np.genfromtxt("data/iris.txt", delimiter=None) # load the data
        Y = iris[:,-1]
        X = iris[:, 0:-1]
        import mltools as ml
        np.random.seed(0)
                                     # set the random number seed
        X,Y = ml.shuffleData(X[:, 0:2],Y); # shuffle data randomly TAKING ONLY FIRST TWO C
        OLUMNS OF X
        # (This is a good idea in case your data are ordered in some systematic way.)
        Xtr, Xva, Ytr, Yva = ml.splitData(X, Y, 0.75); # split data into 75/25 train/validatio
        K=[1,2,5,10,50,100,200];
        errTrain = [None] *len(K)
                                    . # (preallocate storage for training error)
        errVal = [None] *len(K)
        for i,k in enumerate(K):
            knn = ml.knn.knnClassify(Xtr, Ytr, k) # TODO: complete code to train model
            Ytrhat = knn.predict(Xtr) # TODO: predict results on training data
            errTrain[i] = funky(Ytr, Ytrhat)
                                                 # TODO: count what fraction of prediction
        s are wrong
            #TODO: repeat prediction / error evaluation for validation data
            knn2 = ml.knn.knnClassify(Xva, Yva, k)
            Yvahat = knn2.predict(Xva)
            errVal[i] = funky(Yva, Yvahat)
        plt.semilogx(K, errTrain,'r-', lw=3, label='Training')
        plt.semilogx(K, errVal, 'g-', lw=3, label='Validation')
                                                                    #TODO: average and plot
        results on semi-log scale
        plt.legend()
        plt.show()
```



Third question: Create the same error rate plots as the previous part, but with all the features in the dataset. Are the plots very different? Is your recommendation for the best K different? (5 points)

```
In [10]: iris = np.genfromtxt("data/iris.txt",delimiter=None) # load the data
         Y = iris[:,-1]
         X = iris[:, 0:-1]
         import mltools as ml
         np.random.seed(0)
                                     # set the random number seed
         X,Y = ml.shuffleData(X[:,:],Y); # shuffle data randomly TAKING ONLY FIRST TWO COLU
         MNS OF X
         # (This is a good idea in case your data are ordered in some systematic way.)
         Xtr, Xva, Ytr, Yva = ml.splitData(X, Y, 0.75); # split data into 75/25 train/validatio
         K=[1,2,5,10,50,100,200];
         errTrain = [None] *len(K)
                                    * # (preallocate storage for training error)
         errVal = [None] *len(K)
         for i,k in enumerate(K):
             knn = ml.knn.knnClassify(Xtr, Ytr, k)  # TODO: complete code to train model
             Ytrhat = knn.predict(Xtr) # TODO: predict results on training data
             errTrain[i] = funky(Ytr, Ytrhat)
                                                   # TODO: count what fraction of prediction
         s are wrong
             #TODO: repeat prediction / error evaluation for validation data
             knn2 = ml.knn.knnClassify(Xva, Yva, k)
             Yvahat = knn2.predict(Xva)
             errVal[i] = funky(Yva, Yvahat)
         plt.semilogx(K, errTrain,'r-', lw=3, label='Training')
         plt.semilogx(K, errVal,'g-', lw=3, label='Validation')
                                                                     #TODO: average and plot
         results on semi-log scale
         plt.legend()
         plt.show()
```



Across both graphs, the shape (and datapoints) of the validation data's plot remains mostly the same. The only difference is that the value K=5 has a higher number of errors in the second graph than the first. The training data, on the other hand, has far fewer errors in the second graph than in the first leading up to the value of K=100, at which point the two graphs share the same data. With regard to which K value causes the fewest errors, it can be seen that K=1 has the lowest return value, both in the validation data as well as the training data in both plots.

Mustofa Yesilyurf CS 178 HWI

Problem 3: Naive Bayes Classifier

1)
$$\hat{p}(y=-1) = \frac{6}{10} = \frac{3}{5} = 1 - \hat{p}(y=1) = - \hat{p}(y=1) = \frac{4}{10} = \frac{2}{5}$$

$$\hat{p}(x_1|y=1) = \frac{3}{4}, \ \hat{p}(x_4|y=1) = \frac{2}{4} \qquad \hat{p}(x_1|y=-1) = \frac{3}{6}, \ \hat{p}(x_4|y=-1) = \frac{5}{6}, \ \hat{p}(x_2|y=-1) = \frac{5}{6}, \ \hat{p}(x_3|y=-1) = \frac{3}{6}, \ \hat{p}(x_3|y=-1) = \frac{3}{6}, \ \hat{p}(x_3|y=-1) = \frac{4}{6}, \$$

2) Given
$$\vec{x} = (0,0,0,0,0)$$

$$\vec{p}(y = -1) | \vec{x} = 00000) = \frac{(1 - \frac{2}{6})(1 - \frac{5}{6})(1 - \frac{7}{6})(1 -$$

$$\hat{\rho}(y=1|\hat{x}=00000) = \frac{(1-\frac{2}{4})(1-\frac{2$$

$$\Rightarrow \hat{X} = (o, o, o, o, o) \text{ would predict } y = 1$$

Given
$$\vec{x} = (1, 1, 0, 1, 0)$$

$$\vec{p}(y = -1 | \vec{x} = 11010) = \frac{(\frac{2}{5})(\frac{5}{5})(1 - \frac{2}{5})(\frac{5}{5})(1 - \frac{2}{5})(\frac{3}{5})}{(\frac{3}{5})(\frac{7}{5})(1 - \frac{2}{5})(\frac{5}{5})(1 - \frac{2}{5})(\frac{3}{5})} + (\frac{3}{4})(0)(1 - \frac{3}{4})(\frac{3}{4})(1 - \frac{1}{4})(\frac{3}{5})$$

$$= \frac{(5/108)}{\frac{5}{108}+0}$$

$$\hat{\rho}(y=1|\vec{x}=11010) = (\frac{2}{4})(0)(1-\frac{1}{4})(\frac{2}{3})(1-\frac{1}{4})(\frac{2}{3})$$

$$= \frac{(\frac{2}{4})(0)(1-\frac{1}{4})(\frac{2}{3})(1-\frac{1}{4})(\frac{2}{3})}{\frac{2}{108}+0} = \frac{(\frac{2}{3})(0)(1-\frac{1}{4})(\frac{2}{3})(1-\frac{1}{4})(\frac{2}{3})}{\frac{2}{108}+0}$$

=>
$$\vec{x}$$
 = (1,1,0,1,0) would predict $y = -1$

3)
$$\hat{\rho}(y=-1|\hat{x}=00000) = \frac{(1/540)}{\frac{1}{540} + \frac{3}{320}} = [0.16495]$$

$$\hat{\rho}(y=1|\hat{x}=00000) = \frac{(3/320)}{\frac{1}{540} + \frac{3}{320}} = [0.83505]$$

$$\hat{\rho}(y=-1|\hat{x}=11010) = \frac{(5/108)}{\frac{5}{108} + 0} = [1]$$

$$\hat{\rho}(y=1|\hat{x}=11010) = \frac{0}{\frac{5}{108} + 0} = [0]$$

- 4) A joint Bayes classifier would not be suited for this data set because of the feature X2. There is no instance when the feature is true and the class y is 1, meaning there would be no instances of the class being 1 if the features' probabilities were not mutually exclusive. Hence, we use naive Bayes as that method assumes independence for each feature.
- 5) Because the naive Bayes model assumes the features to be independent of each other. The training Later sothered for features Xz through Xs do not need to change or be regathered. Just removing the data and the probabilities associated with X, is enough to allow for new predictions without X, to be made.

4) Statement of Collaboration:

- Nabhan Khan
 - Effect of number of bins in histograms
 - Error counting in 2.2 and 2.3: What makes an error
- Bisher Kudaimi
 - Effect of number of kins in histograms
 - Effect of shuffleData function call

- Sarah andeer
 - Effect of number of like in histograms
 - Effect of shaffle Data function call
- Hibah Wasim
 - Effect of number of bihs in histograms
 - Bayes classifiers: naive versus joint