# CW Part A

April 24, 2021

#### 1 Coursework A

This notebook describes the first part of the ML coursework (code). This part builds on work that you have already done during the labs.

- Get familiar with **common python modules / functions** used for ML in python Get practical experience **implementing** ML methods in python
- Get practical experience regarding **parameter selection** for ML methods Get practical experience on **evaluating** ML methods and applying cross-validation

Notes: - Don't use libraries that implement kNN or cross-validation - your code should be as low-level as possible.

- For more details, read the coursework PDF

```
[1]: import numpy as np
import random
#PRESS <Ctrl>+<Enter> to execute this cell

#%matplotlib inline
mySeed=1234567
#initialize random seed generator
np.random.seed(mySeed)

#In this cell, we load the iris/flower dataset we talked about in class
from sklearn import datasets
import matplotlib.pyplot as plt

iris = datasets.load_iris()

X=iris.data
y=iris.target
```

### 1.1 1. Exploratory Data Analysis

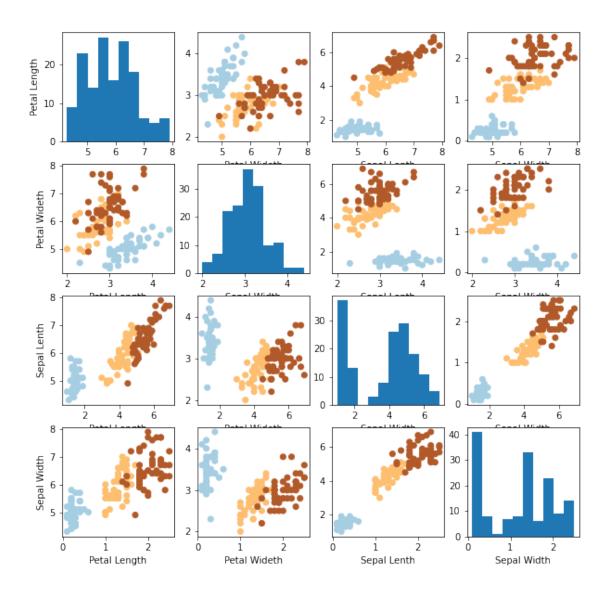
Create a function that given data X and labels y plots the 4x4 grid. The function should be invoked as

```
myplotGrid(X,y)
```

where X is your training data and y are the labels. The final plot should look similarly to the plot from previous lab notebooks, that is:

```
[2]: def myplotGrid(X,y):
         fig = plt.figure(figsize=(10,10))
         labels=['petal length','Petal Wideth','Sepal Lenth','Sepal Width']
         for i in range (X.shape[1]):
             for j in range (X.shape[1]):
                 ax = plt.subplot(X.shape[1],X.shape[1],t)
                 if (i==j \text{ and } i!=0):
                     ax.hist(X[:,i])
                     plt.xlabel('Sepal Width')
                 elif(i==j and i==0):
                     ax.hist(X[:,i])
                     plt.ylabel('Petal Length')
                 elif(i!=j and j==0):
                     ax.scatter(X[:, i], X[:, j], c=y, cmap=plt.cm.Paired)
                     plt.xlabel('Petal Length')
                     plt.ylabel(labels[i])
                 else:
                     ax.scatter(X[:, i], X[:, j], c=y, cmap=plt.cm.Paired)
                     plt.xlabel(labels[j])
                 t=t+1
         return plt.show
```

- [3]: myplotGrid(X,y)
- [3]: <function matplotlib.pyplot.show(close=None, block=None)>



### 1.1.1 1.2. Exploratory Data Analysis under noise

When data are collected under real-world settings (e.g., from webcams or other errors) they usually contain some amount of noise that makes classification more challenging. In the cell below, invoke your exploratory data analysis function above on a noisy version of your data X.

Try to perturb your data with some Gaussian noise,

np.random.seed(mySeed) # initialize random seed to replicate results over different runs XN=X+np.random.normal(0,0.5,X.shape)

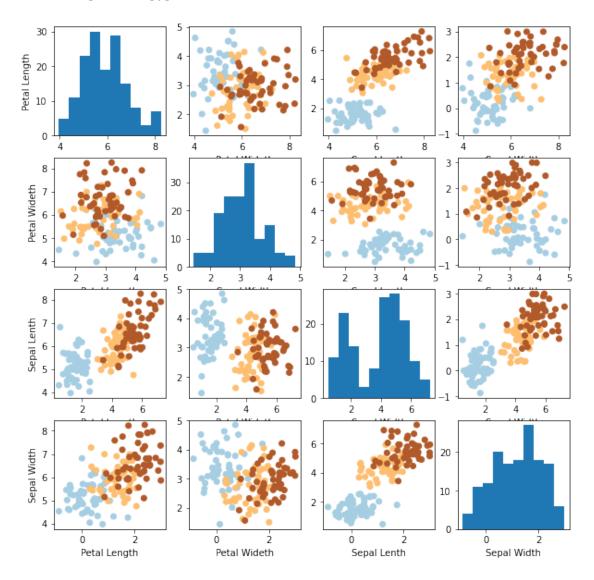
and then invoke

myplotGrid(XN,y)

```
[4]: import numpy as np
mySeed=1234567
np.random.seed(mySeed) # initialize random seed to replicate results over

different runs
XN=X+np.random.normal(0,0.5,X.shape)
myplotGrid(XN,y)
```

# [4]: <function matplotlib.pyplot.show(close=None, block=None)>



## 1.2 2. Implementing kNN

In the cell below, develop your own code for performing k-Nearest Neighbour classification. You are guided to do this by following the notebook accompanying the coursework. Define a function

that performs k-NN given a set of data. Your function should be invoked similary to:

```
y_ = mykNN(X,y,X_,options)
```

where X is your training data, y is your training outputs, X\_ are your testing data and y\_ are your predicted outputs for X\_. The options argument (can be a list or a set of separate arguments depending on how you choose to implement the function) should at least contain the number of neighbours to consider as well as the distance function employed.

Note: if you need help implementing k-NN, you can use Nested Cross-Validation.ipynb

```
[5]: #Euclidean destance
     def EuclideanDistance(X_train,X_test):
         diff=[X test[i]-X train[i] for i in range(len(X test))]
         dist=np.sqrt(sum([diff[i]**2 for i in range(len(diff))]))
         return dist
[6]: # Manhattan Distance
     def ManhattanDistance(X_train, X_test):
         diff=[X_test[i]-X_train[i] for i in range(len(X_test))]
         dist=sum([abs(diff[i]) for i in range(len(diff))])
         return dist
[7]: def getNeighbours1(x_,X,n,T):
         Distance=[EuclideanDistance(X[i],x_) for i in range(T) ]
         indices=[b[0] for b in sorted(enumerate(Distance), key=lambda i:i[1])]
         indic_Neigh=[indices[i] for i in range(n)]
         return indic_Neigh
[8]: #Returning indices of n Neighbours
     def getNeighbours2(x_,X,n,T):
         Distance=[ManhattanDistance(X[i],x_) for i in range(T) ]
         indices=[b[0] for b in sorted(enumerate(Distance), key=lambda i:i[1])]
         indic_Neigh=[indices[i] for i in range(n)]
         return indic_Neigh
[9]: import random
     def assignLabel(nLabels):
         nlabel=np.array(nLabels)
         unique_lab=list(np.unique(nlabel))
         n_unique=[len(np.where(nlabel==unique_lab[i])[0]) for i in_
      →range(len(unique_lab))]
         max unique=[]
         for i in range(len( n_unique)):
             if (n_unique[i] == max(n_unique)):
                 max_unique.append(i)
                 ind=random.choice(max_unique)
```

```
return unique_lab[ind]
```

```
[10]: #Splitting data into training and testing
      #iris_data=[[iris.data[i],iris.target[i]] for i in range(len(iris.data))]
      \#test\_ind=random.sample(range(0,150),30)
      \#ind=random.sample(range(0,150),150)
      #train_ind=list(set(ind)-set(test_ind))
      #Train=[iris data[i] for i in train ind]
      #Test=[iris_data[i] for i in test_ind]
      #X_train=[Train[i][0] for i in range(len(Train))]
      #X_test=[Test[i][0] for i in range(len(Test))]
      #y_train=[Train[i][1] for i in range(len(Train))]
      #y_test=[Test[i][1] for i in range(len(Test))]
[11]: def kNNClassifier(X_train, X_test, y_train, n, dist):
          if (dist=='euclidean'):
              n_indices=[getNeighbours1(X_test[i],X_train,n,len(X_train)) for i in_u
       →range(len(X_test))]
          elif (dist=='manhattan'):
              n indices=[getNeighbours2(X test[i], X train, n, len(X train)) for i in_
       →range(len(X test))]
          else:
              pass
          n_labels=[[y_train[k] for k in n_indices[i]]for i in range(len(n_indices))]
          y_pred=[assignLabel(n_labels[i]) for i in range(len(n_labels)) ]
          return y_pred
[12]: #kNNClassifier(X_train, X_test, y, 10, 'euclidean')
[13]: #y_test
[14]: def kNNClassifier_score(X_train, X_test, y_test, y_train, n, dist):
          if (dist=='euclidean'):
              n_indices=[getNeighbours1(X_test[i],X_train,n,len(X_train)) for i in_
       →range(len(X_test))]
          elif (dist=='manhattan'):
              n_indices=[getNeighbours2(X_test[i],X_train,n,len(X_train)) for i in_
       →range(len(X_test))]
          else:
              pass
          n_labels=[[y_train[k] for k in n_indices[i]]for i in range(len(n_indices))]
          y_pred=[assignLabel(n_labels[i]) for i in range(len(n_labels)) ]
          #Building accuracy_score
          corr_pred=[]
```

```
for i in range(len(y_pred)):
               if (y_pred[i]==y_test[i]):
                   corr_pred.append(i)
               else:
                   pass
          accuracy_score=len(corr_pred)/len(y_pred)
          #Building Confusion matrix
          N_pred_00=[]
          N_pred_01=[]
          N_pred_02=[]
          N_pred_11=[]
          N_pred_10=[]
          N_pred_12=[]
          N_pred_22=[]
          N_pred_20=[]
          N_pred_21=[]
          for i in range(len(y_pred)):
                   (y_test[i] == 0 and y_pred[i] == 0):
                   N_pred_00.append(i)
              elif (y_test[i] == 0 and y_pred[i] == 1):
                   N_pred_01.append(i)
               elif (y_test[i] == 0 and y_pred[i] == 2):
                   N_pred_02.append(i)
               elif (y_test[i]==1 and y_pred[i]==1):
                   N_pred_11.append(i)
               elif (y_test[i]==1 and y_pred[i]==0):
                   N_pred_10.append(i)
              elif (y_test[i] == 1 and y_pred[i] == 2):
                   N_pred_12.append(i)
               elif (y_test[i] == 2 and y_pred[i] == 2):
                   N_pred_22.append(i)
               elif (y_test[i]==2 and y_pred[i]==0):
                   N_pred_20.append(i)
               else:
                   N_pred_21.append(i)
          confusion_matrix=np.
       \rightarrowarray([[len(N_pred_00),len(N_pred_01),len(N_pred_02)],[len(N_pred_10),len(N_pred_11),len(N_pred_10)]
          return confusion_matrix,accuracy_score,
[15]: #kNNClassifier_score(15, 'euclidean')
[16]: #kNNClassifier_score(10, 'manhattan')[1]
```

#### 1.3 3. Nested Cross-validation using your implementation of KNN

In the cell below, develop your own code for performing 5-fold nested cross-validation along with your implementation of k-NN above. Again, you are guided to complete this task by following the appropriate notebook accompanying this coursework. Your code for nested cross-validation should invoke your kNN function (see above). You cross validation function should be invoked similary to:

```
accuracy_fold=myNestedCrossVal(X,y,5,list(range(1,11)),['euclidean','manhattan'],mySeed)
```

where X is your data matrix (containing all samples and features for each sample), 5 is the number of folds, y are your known output labels, list(range(1,11) evaluates the neighbour parameter from 1 to 10, and ['euclidean', 'manhattan'] evaluates the two distances on the validation sets. mySeed is simply a random seed to enable us to replicate your results.

**Notes:** - you should perform nested cross-validation on both your original data X, as well as the data pertrubed by noise as shown in the cells above (XN) - you should implement/validate at least three distance functions - you should evaluate number of neighbours from 1 to 10 - your function should return a list of accuracies per fold - for each fold, your function should print: - the accuracy per distinct set of parameters on the validation set - the best set of parameters for the fold after validation - the confusion matrix per fold (on the testing set)

Note: if you need help implementing cv, you can use nested cross-validation.ipynb

```
[17]: import itertools import random
```

```
[18]: def myNestedCrossVal(X,y,foldK,nns,dist,mySeed):
          np.random.seed(mySeed)
           #Reshuffling data
          indices= np.random.permutation(np.array(X).shape[0])
          X=np.array([X[i] for i in indices])
          y=np.array([y[i] for i in indices])
          accuracy_fold=[]
          accuracy_score=[]
          confusion_matrix=[]
          n_list=[]
          dist list=[]
          fold_list=[]
          pred=[]
          Y_test=[]
          fold1_dist=[]
          fold2_dist=[]
          fold3_dist=[]
          fold4_dist=[]
          fold5_dist=[]
          fold1 as=[]
          fold2_as=[]
          fold3_as=[]
          fold4_as=[]
```

```
fold5_as=[]
   fold1_n=[]
   fold2_n=[]
   fold3_n=[]
   fold4_n=[]
   fold5_n=[]
   accuracy_score1=[]
   y_valid=[y[i] for i in range(30)]
  #creating 5 folds and splitting data into training, validation and testing.
   X_{bins}=[[X[i+int(len(X)/foldK)*j]] for i in range(int(len(X)/foldK))] for ju
→in range(foldK) ]
   y_bins=[[y[i+int(len(X)/foldK)*j] for i in range(int(len(X)/foldK))] for ju
→in range(foldK) ]
   lst_test_X_bins=[]
   lst_test_y_bins=[]
   lst_valid_X_bins=[]
   lst_valid_y_bins=[]
   lst_train_X_bins=[]
   lst_train_y_bins=[]
   lst_pred_y_bins_e=[]
   test_X_bins=X_bins[4]
   lst_test_X_bins.append(test_X_bins)
   test_y_bins=y_bins[4]
   lst_test_y_bins.append(test_y_bins)
   valid_X_bins=X_bins[0]
   lst_valid_X_bins.append(valid_X_bins)
   valid y bins=y bins[0]
   lst_valid_y_bins.append(valid_y_bins)
   train_ind=[j for j in range(foldK) if (j!=0 and j!=4)]
   train_X=[X_bins[i] for i in train_ind]
   train_y=[y_bins[i] for i in train_ind]
   train_X_bins=list(itertools.chain(*train_X))
   lst_train_X_bins.append(train_X_bins)
   train_y_bins=list(itertools.chain(*train_y))
   lst_train_y_bins.append(train_y_bins)
   for i in range(foldK-1):
       test_X_bins=X_bins[i]
       lst_test_X_bins.append(test_X_bins)
       test_y_bins=y_bins[i]
       lst_test_y_bins.append(test_y_bins)
       valid_X_bins=X_bins[i+1]
       lst_valid_X_bins.append(valid_X_bins)
       valid_y_bins=y_bins[i+1]
       lst_valid_y_bins.append(valid_y_bins)
       train_ind=[j for j in range(foldK) if (j!=i and j!=i+1)]
       train_X=[X_bins[i] for i in train_ind]
       train_y=[y_bins[i] for i in train_ind]
```

```
train_X_bins=list(itertools.chain(*train_X))
       lst_train_X_bins.append(train_X_bins)
       train_y_bins=list(itertools.chain(*train_y))
       lst_train_y_bins.append(train_y_bins)
   A=[[i,n,d] for i in range(5) for n in nns for d in dist]
                #X_train=lst_train_X_bins[i]
                \#X\_test=lst\_valid\_X\_bins[i]
                #y_train=lst_train_y_bins[i]
                \#y\_test=lst\_valid\_y\_bins[i]
   for i in range(len(A)):
               accuracy_score.append(np.
→round(kNNClassifier_score(lst_train_X_bins[A[i][0]],lst_valid_X_bins[A[i][0]],lst_valid_y_b
               n_list.append(A[i][1])
               dist_list.append(A[i][2])
               fold_list.append(A[i][0])
   for i in range(len(fold_list)):
       if (fold_list[i]==0):
           fold1_dist.append(dist_list[i])
           fold1_n.append(n_list[i])
           fold1_as.append(accuracy_score[i])
   for i in range(len(fold1_as)):
       print (fold1_as[i],(fold1_n[i],fold1_dist[i]))
   for i in range(len(fold1_as)):
       if (fold1_as[i] == max(fold1_as)):
           fold1_bestNN=fold1_n[i]
           fold1_bestDist=fold1_dist[i]
           fold1_bestas=fold1_as[i]
   #print(fold1_dist)
   opt=['\x1b[6;30;46m' +'** End of fold1 validation, best AS,best NN ,best_{\topic}]
\hookrightarrowDist='+ str([fold1_bestas,fold1_bestNN,fold1_bestDist] ) + '\x1b[0m' for i_\_
\rightarrowin range(5)]
       \#A = [print(Happy_New_Year[i]) for i in range(5)]
   print(opt[0])
   for i in range(len(fold_list)):
       if (fold_list[i]==1):
           fold2_dist.append(dist_list[i])
           fold2_n.append(n_list[i])
           fold2_as.append(accuracy_score[i])
   for i in range(len(fold1_as)):
       print (fold2_as[i],(fold2_n[i],fold2_dist[i]))
   for i in range(len(fold2_as)):
       if (fold2_as[i] == max(fold2_as)):
           fold2_bestNN=fold2_n[i]
           fold2_bestDist=fold2_dist[i]
```

```
fold2_bestas=fold2_as[i]
   opt=['\x1b[6;30;46m' +'** End of fold2 validation, best AS,best NN ,best_
→Dist='+ str([fold2_bestas,fold2_bestNN,fold2_bestDist] ) + '\x1b[0m' for i_
\rightarrowin range(5)]
       #A=[print(Happy_New_Year[i]) for i in range(5)]
   print(opt[0])
   for i in range(len(fold_list)):
       if (fold_list[i]==2):
           fold3_dist.append(dist_list[i])
           fold3_n.append(n_list[i])
           fold3_as.append(accuracy_score[i])
   for i in range(len(fold1_as)):
       print (fold3_as[i],(fold3_n[i],fold3_dist[i]))
   for i in range(len(fold3_as)):
       if (fold3_as[i] == max(fold3_as)):
           fold3_bestNN=fold3_n[i]
           fold3_bestDist=fold3_dist[i]
           fold3_bestas=fold3_as[i]
   opt=['\x1b[6;30;46m' +'** End of fold3 validation,best AS, best NN ,best_
→Dist='+ str([fold3_bestas,fold3_bestNN,fold3_bestDist] ) + '\x1b[0m' for i_
\rightarrowin range(5)]
       #A=[print(Happy_New_Year[i]) for i in range(5)]
   print(opt[0])
   for i in range(len(fold list)):
       if (fold_list[i]==3):
           fold4_dist.append(dist_list[i])
           fold4_n.append(n_list[i])
           fold4_as.append(accuracy_score[i])
   for i in range(len(fold1_as)):
       print (fold4_as[i],(fold4_n[i],fold4_dist[i]))
   for i in range(len(fold4_as)):
       if (fold4_as[i] == max(fold4_as)):
           fold4_bestNN=fold4_n[i]
           fold4_bestDist=fold4_dist[i]
           fold4_bestas=fold4_as[i]
   opt=['\x1b[6;30;46m' +'** End of fold4 validation,best AS, best NN ,best_
→Dist='+ str([fold4_bestas,fold4_bestNN,fold4_bestDist] ) + '\x1b[0m' for i_
\rightarrowin range(5)]
       #A=[print(Happy_New_Year[i]) for i in range(5)]
   print(opt[0])
```

```
for i in range(len(fold_list)):
       if (fold list[i]==4):
           fold5_dist.append(dist_list[i])
           fold5_n.append(n_list[i])
           fold5_as.append(accuracy_score[i])
  for i in range(len(fold1_as)):
       print (fold5_as[i],(fold5_n[i],fold5_dist[i]))
  for i in range(len(fold5_as)):
       if (fold5 as[i] == max(fold5 as)):
           fold5_bestNN=fold5_n[i]
           fold5 bestDist=fold5 dist[i]
           fold5_bestas=fold5_as[i]
   opt=['\x1b[6;30;46m' +'** End of fold5 validation,best AS, best NN ,best ∪,
→Dist='+ str([fold5_bestas,fold5_bestNN,fold5_bestDist] ) + '\x1b[0m' for i_
\rightarrowin range(5)]
       #A=[print(Happy_New_Year[i]) for i in range(5)]
  print(opt[0])
   #print(fold_list)
   #print(accuracy score)
  l_bestNN=[fold1_bestNN,fold2_bestNN,fold3_bestNN,fold4_bestNN,fold5_bestNN]
→l_bestDist=[fold1_bestDist,fold2_bestDist,fold3_bestDist,fold4_bestDist,fold5_bestDist]
  l bestas=[fold1_bestas,fold2_bestas,fold3_bestas,fold4_bestas,fold5_bestas]
  for i in range(len(l_bestas)):
       if (l_bestas[i] == max(l_bestas)):
           bestDistance= l_bestDist[i]
           bestNN=l_bestNN[i]
  print('Now out of the optimal parameters of the 5 folds, we chose the ones⊔
⇒that give the best accuracy as shown below and proceed with⊔
opt=['\x1b[6;30;43m' +'** Best NN ,best Dist='+ str([bestNN,bestDistance]_
\rightarrow) + '\x1b[0m' for i in range(5)]
       #A=[print(Happy_New_Year[i]) for i in range(5)]
  print(opt[0])
   #print('** End of val, best NN', bestNN, 'best Dist', bestDistance)
  for i in range(len(lst_test_X_bins)):
      X_train=lst_train_X_bins[i]
```

```
y_train=lst_train_y_bins[i]
              X test=lst_test_X_bins[i]
              y_test=lst_test_y_bins[i]
              Y_test.append(y_test)
       →matrix,score=kNNClassifier_score(X_train,X_test,y_test,y_train,bestNN,bestDistance)
              #matrix=kNNClassifier score(bestNN, bestDistance)
              confusion_matrix.append(matrix)
              accuracy_fold.append(np.round(score,3))
          print('Now using the best parameters obtained in validation, here is the \sqcup
       ⇒score obtained using testing sets in 5 folds')
          accuracy_fold=['\x1b[6;30;42m' +str(accuracy_fold ) + '\x1b[0m' for i in_
       \rightarrowrange(5)]
          accuracy_fold=[print(accuracy_fold[i]) for i in range(1)]
          print('Here is the total Confusion Matrix:')
          print(sum(confusion_matrix))
          return accuracy_fold
[19]: iris = datasets.load_iris()
      X=iris.data
      y=iris.target
      accuracy_fold=myNestedCrossVal(X,y,5,list(range(1,11)),['euclidean','manhattan'],mySeed)
      print(accuracy_fold)
     0.967 (1, 'euclidean')
     0.967 (1, 'manhattan')
     0.967 (2, 'euclidean')
     0.967 (2, 'manhattan')
     0.967 (3, 'euclidean')
     0.967 (3, 'manhattan')
     0.967 (4, 'euclidean')
     0.967 (4, 'manhattan')
     0.967 (5, 'euclidean')
     0.967 (5, 'manhattan')
     0.967 (6, 'euclidean')
     0.967 (6, 'manhattan')
     0.967 (7, 'euclidean')
     0.967 (7, 'manhattan')
     0.967 (8, 'euclidean')
     0.967 (8, 'manhattan')
     0.967 (9, 'euclidean')
     0.967 (9, 'manhattan')
     0.967 (10, 'euclidean')
     0.967 (10, 'manhattan')
```

```
** End of fold1 validation, best AS, best NN , best Dist=[0.967, 10, 'manhattan']
0.933 (1, 'euclidean')
0.933 (1, 'manhattan')
0.933 (2, 'euclidean')
0.9 (2, 'manhattan')
0.967 (3, 'euclidean')
0.967 (3, 'manhattan')
0.933 (4, 'euclidean')
0.9 (4, 'manhattan')
0.933 (5, 'euclidean')
0.933 (5, 'manhattan')
0.933 (6, 'euclidean')
0.933 (6, 'manhattan')
0.933 (7, 'euclidean')
0.933 (7, 'manhattan')
0.9 (8, 'euclidean')
0.933 (8, 'manhattan')
0.933 (9, 'euclidean')
0.933 (9, 'manhattan')
0.933 (10, 'euclidean')
0.933 (10, 'manhattan')
** End of fold2 validation, best AS, best NN , best Dist=[0.967, 3, 'manhattan']
0.933 (1, 'euclidean')
0.9 (1, 'manhattan')
0.9 (2, 'euclidean')
0.867 (2, 'manhattan')
0.967 (3, 'euclidean')
1.0 (3, 'manhattan')
1.0 (4, 'euclidean')
0.933 (4, 'manhattan')
1.0 (5, 'euclidean')
0.933 (5, 'manhattan')
0.933 (6, 'euclidean')
0.933 (6, 'manhattan')
0.933 (7, 'euclidean')
0.933 (7, 'manhattan')
0.967 (8, 'euclidean')
0.967 (8, 'manhattan')
1.0 (9, 'euclidean')
1.0 (9, 'manhattan')
1.0 (10, 'euclidean')
0.967 (10, 'manhattan')
** End of fold3 validation, best AS, best NN , best Dist=[1.0, 10, 'euclidean']
1.0 (1, 'euclidean')
1.0 (1, 'manhattan')
1.0 (2, 'euclidean')
1.0 (2, 'manhattan')
1.0 (3, 'euclidean')
```

```
1.0 (3, 'manhattan')
1.0 (4, 'euclidean')
1.0 (4, 'manhattan')
1.0 (5, 'euclidean')
1.0 (5, 'manhattan')
1.0 (6, 'euclidean')
1.0 (6, 'manhattan')
1.0 (7, 'euclidean')
1.0 (7, 'manhattan')
1.0 (8, 'euclidean')
1.0 (8, 'manhattan')
1.0 (9, 'euclidean')
1.0 (9, 'manhattan')
1.0 (10, 'euclidean')
1.0 (10, 'manhattan')
** End of fold4 validation, best AS, best NN , best Dist=[1.0, 10, 'manhattan']
0.967 (1, 'euclidean')
0.967 (1, 'manhattan')
0.967 (2, 'euclidean')
0.967 (2, 'manhattan')
0.967 (3, 'euclidean')
0.967 (3, 'manhattan')
1.0 (4, 'euclidean')
1.0 (4, 'manhattan')
1.0 (5, 'euclidean')
0.967 (5, 'manhattan')
0.933 (6, 'euclidean')
1.0 (6, 'manhattan')
1.0 (7, 'euclidean')
0.967 (7, 'manhattan')
0.967 (8, 'euclidean')
0.967 (8, 'manhattan')
0.967 (9, 'euclidean')
0.933 (9, 'manhattan')
0.967 (10, 'euclidean')
0.933 (10, 'manhattan')
** End of fold5 validation, best AS, best NN , best Dist=[1.0, 7, 'euclidean']
Now out of the optimal parameters of the 5 folds, we chose the ones that give
the best accuracy as shown below and proceed with cross-validation
** Best NN ,best Dist=[7, 'euclidean']
Now using the best parameters obtained in validation, here is the score obtained
using testing sets in 5 folds
[1.0, 0.967, 0.933, 0.967, 1.0]
Here is the total Confusion Matrix:
[[50 0 0]
[ 0 48 2]
 [ 0 2 48]]
[None]
```

```
[20]: iris = datasets.load_iris()
      X=iris.data
      y=iris.target
      Xn=X+np.random.normal(0,5,np.array(X).shape)
      accuracy_fold=myNestedCrossVal(Xn,y,5,list(range(1,11)),['euclidean','manhattan'],mySeed)
      print(accuracy_fold)
     0.333 (1, 'euclidean')
     0.367 (1, 'manhattan')
     0.333 (2, 'euclidean')
     0.367 (2, 'manhattan')
     0.267 (3, 'euclidean')
     0.3 (3, 'manhattan')
     0.233 (4, 'euclidean')
     0.4 (4, 'manhattan')
     0.4 (5, 'euclidean')
     0.367 (5, 'manhattan')
     0.367 (6, 'euclidean')
     0.367 (6, 'manhattan')
     0.4 (7, 'euclidean')
     0.4 (7, 'manhattan')
     0.4 (8, 'euclidean')
     0.4 (8, 'manhattan')
     0.5 (9, 'euclidean')
     0.4 (9, 'manhattan')
     0.433 (10, 'euclidean')
     0.467 (10, 'manhattan')
     ** End of fold1 validation, best AS, best NN , best Dist=[0.5, 9, 'euclidean']
     0.333 (1, 'euclidean')
     0.333 (1, 'manhattan')
     0.267 (2, 'euclidean')
     0.367 (2, 'manhattan')
     0.233 (3, 'euclidean')
     0.467 (3, 'manhattan')
     0.3 (4, 'euclidean')
     0.367 (4, 'manhattan')
     0.3 (5, 'euclidean')
     0.4 (5, 'manhattan')
     0.233 (6, 'euclidean')
     0.233 (6, 'manhattan')
     0.233 (7, 'euclidean')
     0.267 (7, 'manhattan')
     0.2 (8, 'euclidean')
     0.167 (8, 'manhattan')
     0.167 (9, 'euclidean')
     0.233 (9, 'manhattan')
     0.2 (10, 'euclidean')
```

```
0.233 (10, 'manhattan')
** End of fold2 validation, best AS, best NN , best Dist=[0.467, 3, 'manhattan']
0.467 (1, 'euclidean')
0.433 (1, 'manhattan')
0.3 (2, 'euclidean')
0.333 (2, 'manhattan')
0.4 (3, 'euclidean')
0.467 (3, 'manhattan')
0.4 (4, 'euclidean')
0.4 (4, 'manhattan')
0.467 (5, 'euclidean')
0.4 (5, 'manhattan')
0.367 (6, 'euclidean')
0.367 (6, 'manhattan')
0.333 (7, 'euclidean')
0.367 (7, 'manhattan')
0.333 (8, 'euclidean')
0.4 (8, 'manhattan')
0.333 (9, 'euclidean')
0.367 (9, 'manhattan')
0.3 (10, 'euclidean')
0.367 (10, 'manhattan')
** End of fold3 validation, best AS, best NN , best Dist=[0.467, 5, 'euclidean']
0.3 (1, 'euclidean')
0.267 (1, 'manhattan')
0.367 (2, 'euclidean')
0.2 (2, 'manhattan')
0.333 (3, 'euclidean')
0.333 (3, 'manhattan')
0.4 (4, 'euclidean')
0.367 (4, 'manhattan')
0.3 (5, 'euclidean')
0.3 (5, 'manhattan')
0.333 (6, 'euclidean')
0.433 (6, 'manhattan')
0.333 (7, 'euclidean')
0.333 (7, 'manhattan')
0.367 (8, 'euclidean')
0.433 (8, 'manhattan')
0.4 (9, 'euclidean')
0.467 (9, 'manhattan')
0.433 (10, 'euclidean')
0.533 (10, 'manhattan')
** End of fold4 validation, best AS, best NN , best Dist=[0.533, 10, 'manhattan']
0.267 (1, 'euclidean')
0.333 (1, 'manhattan')
0.233 (2, 'euclidean')
0.3 (2, 'manhattan')
```

```
0.233 (3, 'euclidean')
     0.333 (3, 'manhattan')
     0.333 (4, 'euclidean')
     0.333 (4, 'manhattan')
     0.3 (5, 'euclidean')
     0.3 (5, 'manhattan')
     0.367 (6, 'euclidean')
     0.4 (6, 'manhattan')
     0.4 (7, 'euclidean')
     0.433 (7, 'manhattan')
     0.367 (8, 'euclidean')
     0.433 (8, 'manhattan')
     0.333 (9, 'euclidean')
     0.467 (9, 'manhattan')
     0.433 (10, 'euclidean')
     0.467 (10, 'manhattan')
     ** End of fold5 validation, best AS, best NN , best Dist=[0.467, 10, 'manhattan']
     Now out of the optimal parameters of the 5 folds, we chose the ones that give
     the best accuracy as shown below and proceed with cross-validation
     ** Best NN ,best Dist=[10, 'manhattan']
     Now using the best parameters obtained in validation, here is the score obtained
     using testing sets in 5 folds
     Here is the total Confusion Matrix:
     [[20 20 10]
      [27 14 9]
      [12 14 24]]
     [None]
[21]: #Fuction to calculate mean of data in a list X
      def mean(X):
          mean=sum(X)/len(X)
          return mean
[22]: #Fuction to calculate standard deviation of data in a list X
      def std(X):
          diff=[(X[i]-mean(X))**2 for i in range(len(X)) ]
          std=np.sqrt(sum(
              diff)/(len(X)-1))
          return std
[23]: #Clean data
      accuracy_clean= [96.7, 96.7, 100, 100, 100]
      mean(accuracy_clean)
[23]: 98.6799999999999
```

```
[24]: std(accuracy_clean)
[24]: 1.8074844397670466
[25]: #Noisy Data
      accuracy_noisy= [46.7, 36.7, 53.3, 56.7, 46.7]
      mean(accuracy noisy)
[25]: 48.0199999999999
[26]: std(accuracy_noisy)
[26]: 7.662375610735875
     1.4 Part 3
[27]: def ManhattanDistance(X_train, X_test):
          dist=[sum(list(abs(X_test-X_train[i]))) for i in range(len(X_train))]
          return dist
[28]: X_train=np.array([(2,2),(1,4),(5,5),(5,7),(5,8)])
      X \text{ test=np.array}((1,1))
      y_train=np.array((0,0,1,1,1))
[29]: ManhattanDistance(X_train, X_test)
[29]: [2, 3, 8, 10, 11]
     2 Weighted k-NN
[30]: from sklearn import datasets
      iris = datasets.load_iris()
      X=iris.data
      y=iris.target
[31]: #Euclidean destance
      def EuclideanDistance(X_train,X_test):
          diff=[X_test[i]-X_train[i] for i in range(len(X_test))]
          dist=np.sqrt(sum([diff[i]**2 for i in range(len(diff))]))
          return dist
[32]: # Manhattan Distance
      def ManhattanDistance(X_train,X_test):
```

diff=[X\_test[i]-X\_train[i] for i in range(len(X\_test))]

```
dist=sum([abs(diff[i]) for i in range(len(diff))])
          return dist
[33]: def getNeighbours3(x ,X train,n,T):
          Distance=[EuclideanDistance(X[i],x_) for i in range(T) ]
          indices=[b[0] for b in sorted(enumerate(Distance), key=lambda i:i[1])]
          indic_Neigh=[indices[i] for i in range(n)]
          return indic_Neigh ,Distance
[34]: def getNeighbours4(x_,X,n,T):
          Distance=[ManhattanDistance(X[i],x_) for i in range(T) ]
          indices=[b[0] for b in sorted(enumerate(Distance), key=lambda i:i[1])]
          indic_Neigh=[indices[i] for i in range(n)]
          return indic_Neigh, Distance
[35]: import random
      def kNNClassifier1(X_train, X_test, y_test, n, dist):
          if (dist=='euclidean'):
              n_indices=[getNeighbours3(X_test[i],X_train,n,len(X_train))[0] for i in_
       →range(len(X test))]
              Distance=[getNeighbours3(X_test[i],X_train,n,len(X_train))[1] for i in__
       →range(len(X_test))]
          elif (dist=='manhattan'):
              n_indices=[getNeighbours4(X_test[i],X_train,n,len(X_train))[0] for i in_
       →range(len(X test))]
              Distance=[getNeighbours4(X_test[i],X_train,n,len(X_train))[1] for i in_
       →range(len(X_test))]
          else:
              pass
          return n_indices,Distance
[36]: def kNNClassifier_score1(X_train, X_test, y_test, y_train, n, dist):
          n_indices, Distance=kNNClassifier1(X_train,X_test,y_test,n,dist)
          ind_class_0=[[i for i in n_indices[j] if(y_train[i]==0)] for j in_u
       →range(len(n_indices))]
          ind_class_1=[[i for i in n_indices[j] if(y_train[i]==1)] for j in_
       →range(len(n_indices))]
          ind_class_2=[[i for i in n_indices[j] if(y_train[i]==2)] for j in__
       →range(len(n_indices))]
          wd class 0=[]
          wd_class_1=[]
          wd_class_2=[]
          for k in range(len(ind_class_0)):
```

```
if(len(ind_class_0[k])==0):
            wd class_0.append(0.000001)
        else:
            wd_class_0.append(sum([1/Distance[k][j] for j in ind_class_0[k]]))
   for k in range(len(ind_class_1)):
        if(len(ind_class_1[k])==0):
            wd_class_1.append(0.000001)
        else:
            wd_class_1.append(sum([1/Distance[k][j] for j in ind_class_1[k]]))
   for k in range(len(ind_class_2)):
        if(len(ind class 2[k])==0):
            wd_class_2.append(0.000001)
        else:
            wd_class_2.append(sum([1/Distance[k][j] for j in ind_class_2[k]]))
   y_pred=[]
   for i in range(len( wd_class_0)):
        a=random.choice([0,1])
       b=random.choice([0,2])
       c=random.choice([1,2])
        d=random.choice([0,1,2])
        if ( wd_class_0[i]> wd_class_1[i] and wd_class_0[i]> wd_class_2[i]):
            y_pred.append(0)
        elif ( wd_class_1[i]> wd_class_0[i] and wd_class_1[i]> wd_class_2[i]):
            y pred.append(1)
        elif ( wd_class_2[i]>wd_class_0[i] and wd_class_2[i]> wd_class_1[i]):
            y_pred.append(2)
        elif ( wd_class_0[i] == wd_class_1[i] and wd_class_0[i] > wd_class_2[i]):
            y_pred.append(a)
        elif ( wd_class_0[i] == wd_class_2[i] and wd_class_0[i] > wd_class_1[i]):
            y_pred.append(b)
        elif ( wd_class_1[i] == wd_class_2[i] and wd_class_1[i] > wd_class_0[i]):
            y_pred.append(c)
        else:
            y_pred.append(d)
   corr_pred=[]
   for i in range(len(y_pred)):
        if (y_pred[i]==y_test[i]):
            corr pred.append(i)
        else:
            pass
   accuracy_score=len(corr_pred)/len(y_pred)
#Building Confusion matrix
   N_pred_00=[]
   N_pred_01=[]
   N_pred_02=[]
   N_pred_11=[]
```

```
N_pred_10=[]
  N_pred_12=[]
  N_pred_22=[]
  N_pred_20=[]
  N_pred_21=[]
  for i in range(len(y_pred)):
            (y_test[i] == 0 and y_pred[i] == 0):
           N_pred_00.append(i)
       elif (y_test[i]==0 and y_pred[i]==1):
           N_pred_01.append(i)
       elif (y_test[i] == 0 and y_pred[i] == 2):
           N_pred_02.append(i)
       elif (y_test[i]==1 and y_pred[i]==1):
           N_pred_11.append(i)
       elif (y_test[i]==1 and y_pred[i]==0):
           N_pred_10.append(i)
       elif (y_test[i]==1 and y_pred[i]==2):
           N_pred_12.append(i)
       elif (y_test[i]==2 and y_pred[i]==2):
           N_pred_22.append(i)
       elif (y_test[i]==2 and y_pred[i]==0):
           N_pred_20.append(i)
       else:
           N_pred_21.append(i)
   confusion_matrix=np.
→array([[len(N_pred_00),len(N_pred_01),len(N_pred_02)],[len(N_pred_10),len(N_pred_11),len(N_
  return confusion_matrix, accuracy_score
```

### 2.1 Testing our Weighted k-NN with Clean data

Note that we are using the best parameters found for clean data after running cross validation of k-NN. So our n=10 and distance is Euclidean.

```
[37]: from sklearn import datasets
  import random
  iris = datasets.load_iris()
  X=iris.data
  y=iris.target
  indices= np.random.permutation(np.array(X).shape[0])
  X=np.array([X[i] for i in indices])
  y=np.array([y[i] for i in indices])
  X_train_c=[X[i] for i in range(120)]
  X_test_c=[X[i] for i in range(120,150)]
```

### 2.2 Testing our Weighted k-NN with Noisy data

For noisy data our best parameters as obtained from cross validation is n=10 and distance is manhattan

```
[38]: from sklearn import datasets
      import random
      iris = datasets.load_iris()
      X=iris.data
      y=iris.target
      Xn=X+np.random.normal(10,30,np.array(X).shape)
      y=iris.target
      indices= np.random.permutation(np.array(X).shape[0])
      Xn=np.array([X[i] for i in indices])
      y=np.array([y[i] for i in indices])
      X_train_n=[Xn[i] for i in range(120)]
      X test n=[Xn[i] \text{ for } i \text{ in } range(120,150)]
      y_train_n=[y[i] for i in range(120)]
      y_test_n=[y[i] for i in range(120,150)]
      kNNClassifier_score1(X_train_n,X_test_n,y_test_n,y_train_n,10,'manhattan')
     <ipython-input-36-a9859d259743>:23: RuntimeWarning: divide by zero encountered
     in double scalars
       wd_class_2.append(sum([1/Distance[k][j] for j in ind_class_2[k]]))
[38]: (array([[15, 0, 0],
              [0, 7, 0],
              [0, 0, 8]]),
       1.0)
 []:
 []:
```