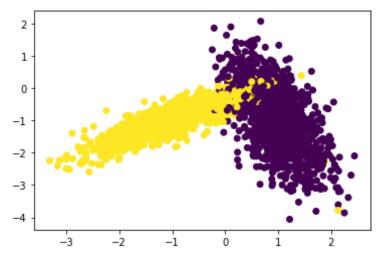
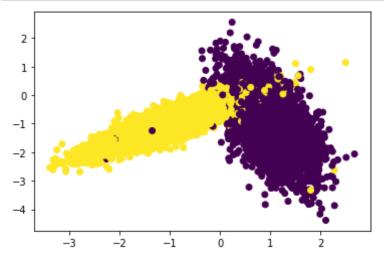
```
In [6]: from sklearn.datasets import make_classification
    from sklearn.model_selection import train_test_split
    from sklearn.model_selection import KFold
    from sklearn.preprocessing import StandardScaler
    import numpy
    from tqdm import tqdm
    import numpy as np
    from sklearn.metrics.pairwise import euclidean_distances

x,y = make_classification(n_samples=10000, n_features=2, n_informative=
2, n_redundant= 0, n_clusters_per_class=1, random_state=60)
X_train, X_test, y_train, y_test = train_test_split(x,y,stratify=y,rand om_state=42)
```





```
In [10]: %matplotlib inline
   import matplotlib.pyplot as plt
   colors = {0:'red', 1:'blue'}
   plt.scatter(X_train[:,0], X_train[:,1],c=y_train)
   plt.show()
```



Implementing Custom RandomSearchCV

```
def RandomSearchCV(x_train,y_train,classifier, param_range, fold
s):
    # x_train: its numpy array of shape, (n,d)
    # y_train: its numpy array of shape, (n,) or (n,1)
    # classifier: its typically KNeighborsClassifier()
    # param_range: its a tuple like (a,b) a < b
    # folds: an integer, represents number of folds we need to d
evide the data and test our model</pre>
```

#1.generate 10 unique values(uniform random distribution) in

the given range "param range" and store them as "params"

ex: if param_range = (1, 50), we need to generate 10 rando m numbers in range 1 to 50

#2.devide numbers ranging from 0 to len(X_train) into group
s= folds

ex: folds=3, and len(x_{train})=100, we can devide numbers f rom 0 to 100 into 3 groups

group 1: 0-33, group 2:34-66, group 3: 67-100

#3.for each hyperparameter that we generated in step 1:

and using the above groups we have created in step 2 y
ou will do cross-validation as follows

first we will keep group 1+group 2 i.e. 0-66 as train
data and group 3: 67-100 as test data, and find train and
 test accuracies

third we will keep group 2+group 3 i.e. 34-100 as trai
n data and group 1: 0-33 as test data, and find train and
 test accuracies

based on the 'folds' value we will do the same procedu
re

find the mean of train accuracies of above 3 steps and store in a list "train scores"

find the mean of test accuracies of above 3 steps and store in a list "test_scores"

#4. return both "train scores" and "test scores"

#5. call function RandomSearchCV(x_train,y_train,classifier, par

am_range, folds) and store the returned values into "train_scor
e", and "cv_scores"
#6. plot hyper-parameter vs accuracy plot as shown in reference
notebook and choose the best hyperparameter
#7. plot the decision boundaries for the model initialized with
the best hyperparameter, as shown in the last cell of reference
notebook

```
In [12]: from sklearn.metrics import accuracy score
         def RandomSearch(x train,y train,classifier, params, folds=3):
             trainscores = []
             cvscores = []
             for k in tqdm(params['n neighbors']):
                 trainscores folds = []
                 cvscores folds = []
                 dist = int(len(x train)/folds) #here it divides len(x train) in
         to 3 equal fields
                 count = 0
                 for i in range(folds): #this for loop will run for 3 folds
                     x_tr = []
                     x cv = []
                     y tr = []
                     y cv = []
                     index = count+dist
                     index cv = list(range(count,index))#this will take 2.5k poi
         nts for cv
                    # print(len(index cv))
                     count = dist+count
                    # print(count)
                     total = list(range(len(x train))) #total of all 7500 points
                     #print(len(total))
                     index tr = set(total)-set(index cv) #for train it will cons
         ider 7500-2500 = 5000
                     index tr = list(index tr)
                     #print(len(index tr))
                     for i cv in index cv:
```

```
x_cv.append(X_train[i_cv])
           y cv.append(y train[i cv])
       for i tr in index tr:
           x tr.append(X train[i tr])
           y tr.append(y train[i tr])
       x cv = np.array(x cv)
       y cv = np.array(y cv)
       x tr = np.array(x tr)
       y tr = np.array(y tr)
       classifier.n neighbors = k
       classifier.fit(x tr,y tr) #fits the module
       Y predicted = classifier.predict(x cv)
       cvscores folds.append(accuracy score(y cv, Y predicted))
       Y predicted = classifier.predict(x tr)
       trainscores folds.append(accuracy score(y tr, Y predicted))
   trainscores.append(np.mean(np.array(trainscores folds)))
   cvscores.append(np.mean(np.array(cvscores folds)))
return trainscores,cvscores
```

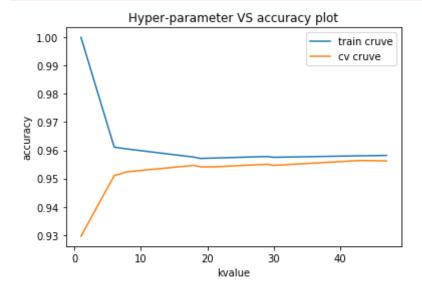
```
In [17]: from sklearn.metrics import accuracy_score
    from sklearn.neighbors import KNeighborsClassifier
    import matplotlib.pyplot as plt
    import random
    import warnings
    import numpy as np
    warnings.filterwarnings("ignore")

neigh = KNeighborsClassifier()
    pram = range(1,50) #it will take the range of values between 1-50

params = {'n_neighbors':np.sort(random.sample(pram,10))} #it will consider 10 values
    folds = 3

trainscores,cvscores = RandomSearch(X_train, y_train, neigh, params, fo
```

```
plt.plot(params['n_neighbors'], trainscores, label='train cruve')
plt.plot(params['n_neighbors'], cvscores, label='cv cruve')
plt.title('Hyper-parameter VS accuracy plot')
plt.xlabel("kvalue")
plt.ylabel("accuracy")
plt.legend()
plt.show()
100%
```



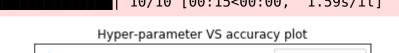
```
In [28]: from sklearn.metrics import accuracy_score
    from sklearn.neighbors import KNeighborsClassifier
    import matplotlib.pyplot as plt
    import random
    import warnings
    import numpy as np
    warnings.filterwarnings("ignore")
```

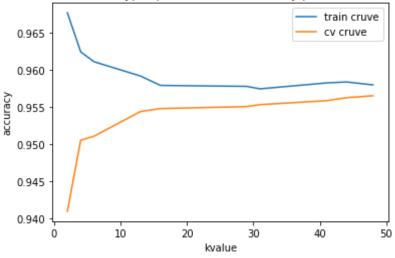
```
neigh = KNeighborsClassifier()
pram = range(1,50)

params = {'n_neighbors':np.sort(random.sample(pram,10))}
folds = 3

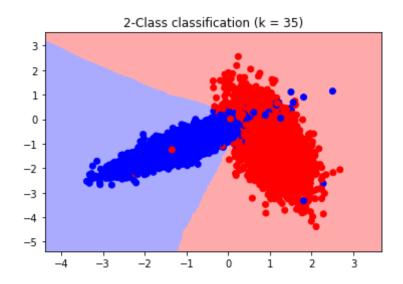
trainscores,cvscores = RandomSearch(X_train, y_train, neigh, params, folds)

plt.plot(params['n_neighbors'],trainscores, label='train cruve')
plt.plot(params['n_neighbors'],cvscores, label='cv cruve')
plt.title('Hyper-parameter VS accuracy plot')
plt.xlabel("kvalue")
plt.ylabel("accuracy")
plt.legend()
plt.show()
100%
```

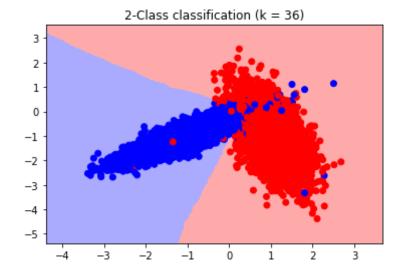




```
In [26]: # understanding this code line by line is not that importent
         def plot decision boundary(X1, X2, y, clf):
                 # Create color maps
             cmap light = ListedColormap(['#FFAAAA', '#AAFFAA', '#AAAAFF'])
             cmap bold = ListedColormap(['#FF0000', '#00FF00', '#0000FF'])
             x \min, x \max = X1.\min() - 1, X1.\max() + 1
             y \min, y \max = X2.\min() - 1, X2.\max() + 1
             xx, yy = np.meshgrid(np.arange(x min, x max, 0.02), np.arange(y min)
          , y \max, 0.02)
             Z = clf.predict(np.c [xx.ravel(), yy.ravel()])
             Z = Z.reshape(xx.shape)
             plt.figure()
             plt.pcolormesh(xx, yy, Z, cmap=cmap light)
             # Plot also the training points
             plt.scatter(X1, X2, c=y, cmap=cmap_bold)
             plt.xlim(xx.min(), xx.max())
             plt.ylim(yy.min(), yy.max())
             plt.title("2-Class classification (k = %i)" % (clf.n neighbors))
             plt.show()
In [75]: from matplotlib.colors import ListedColormap
         neigh = KNeighborsClassifier(n neighbors = 35)
         neigh.fit(X train, y train)
         plot decision boundary(X train[:, 0], X train[:, 1], y train, neigh)
```



In [20]: from matplotlib.colors import ListedColormap
 neigh = KNeighborsClassifier(n_neighbors = 36)
 neigh.fit(X_train, y_train)
 plot_decision_boundary(X_train[:, 0], X_train[:, 1], y_train, neigh)



observation:

since it is random search cv everytime i run seeing the variation while choosing the hyperparameter i had drawn 2 plots which i observed best hyperparameter i.e., k = 36 and k = 35