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In [1]: from sklearn.datasets import make_classification
         from sklearn.model_selection import train_test_split
        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,random_state=42)
        # del X_train, X_test
In [2]: %matplotlib inline
        import matplotlib.pyplot as plt
        colors = {0:'red', 1:'blue'}
        plt.scatter(X_test[:,0], X_test[:,1],c=y_test)
        plt.show()
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        Implementing Custom RandomSearchCV
            def RandomSearchCV(x_train, y_train, classifier, param_range, folds):
                # 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</pre>
                # folds: an integer, represents number of folds we need to devide the data and tes
            t our model
                #1.generate 10 unique values(uniform random distribution) in the given range "para
            m_range" and store them as "params"
                # ex: if param_range = (1, 50), we need to generate 10 random numbers in range 1 t
            o 50
                #2.devide numbers ranging from 0 to len(X_train) into groups= folds
                # ex: folds=3, and len(x_train)=100, we can devide numbers from 0 to 100 into 3 gr
            oups
                  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 you will do cross-valid
            ation as follows
                    \# first we will keep group 1+group 2 i.e. 0-66 as train data and group 3: 67-1
            00 as test data, and find train and
                      test accuracies
                    # second we will keep group 1+group 3 i.e. 0-33, 67-100 as train data and grou
            p 2: 34-66 as test data, and find
                      train and test accuracies
                    # third we will keep group 2+group 3 i.e. 34-100 as train 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 procedure
                    # find the mean of train accuracies of above 3 steps and store in a list "trai
            n_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, param_range, folds) and s
            tore the returned values into "train_score", and "cv_scores"
            #6. plot hyper-parameter vs accuracy plot as shown in reference notebook and choose th
            e best hyperparameter
            #7. plot the decision boundaries for the model initialized with the best hyperparamete
            r, as shown in the last cell of reference notebook
In [3]: from sklearn.metrics import accuracy_score
        #Function to generate 10 random values(parameters)
        def randomn_params(params):
            param=random.sample(range(1, params), 10)
            param.sort()
            return param
        def RandomsearchCV(x_train, y_train, classifier, param_range, folds):
            # List to store the train and test scores
            trainscores=[]
            testscores=[]
            #Dividing the points into groups called folds for cross-validation
            dividing_by_folds=int(len(x_train)/folds)
            #Looping through each hyper-parameter k
            for k in tqdm(params_list['n_neighbors']):
                train_folds=[]
                test_folds=[]
                for j in range(0, folds):
                    # Looping to get the required points for each fold CV
                     test_points=list(set(list(range((dividing_by_folds*j),(dividing_by_folds*(j+1
        ))))))
                    # Separating the test and train points
                    train_points=list(set(list(range(1,len(x_train))))-set(test_points))
                    # selecting the data points based on the train_indices and test_indices
                    X_train=x_train[train_points]
                    Y_train=y_train[train_points]
                    X_test=x_train[test_points]
                    Y_test=y_train[test_points]
                    # Fitting the training data points into the model
                    classifier.n_neighbors=k
                    classifier.fit(X_train,Y_train)
                     #Predicting the accuracy of both train and test data points
                     Y_predict=classifier.predict(X_test)
                     test_folds.append(accuracy_score(Y_test,Y_predict))
                    Y_predict1=classifier.predict(X_train)
                    train_folds.append(accuracy_score(Y_train,Y_predict1))
                 #Adding the scores of each fold in training and test datasets for all hyper-paramete
        r and storing them
                 trainscores.append(np.mean(np.array(train_folds)))
                 testscores.append(np.mean(np.array(test_folds)))
            return trainscores, testscores
In [4]: | from sklearn.metrics import accuracy_score
         from sklearn.neighbors import KNeighborsClassifier
        import matplotlib.pyplot as plt
        import random
        import warnings
        warnings.filterwarnings("ignore")
        n=KNeighborsClassifier()
        params=50
        params_list={'n_neighbors':randomn_params(params)}
        folds=3
        print(params_list)
        trainscores, testscores=RandomsearchCV(X_train, y_train, n, params_list, folds)
        plt.plot(params_list['n_neighbors'], trainscores, label='train data')
        plt.plot(params_list['n_neighbors'], testscores, label='test data')
        plt.title('Hyper-parameter vs accuracy')
        plt.legend
        plt.show
        {'n_neighbors': [2, 10, 11, 24, 25, 34, 39, 42, 45, 48]}
                                                     | 10/10 [00:02<00:00, 3.95it/s]
Out[4]: <function matplotlib.pyplot.show(*args, **kw)>
                       Hyper-parameter vs accuracy
         0.965
         0.960
         0.955
         0.950
         0.945
         0.940
                             20
                     10
                                     30
                                             40
In [5]: 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 [6]: from matplotlib.colors import ListedColormap
        neigh = KNeighborsClassifier(n_neighbors = 45)
        neigh.fit(X_train, y_train)
        plot_decision_boundary(X_train[:, 0], X_train[:, 1], y_train, neigh)
                     2-Class classification (k = 45)
         -3
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