	region tenure age marital address income ed employ retire gender reside custcat 0 2 13 44 1 9 64.0 4 5 0.0 0 2 1 1 3 11 33 1 7 136.0 5 5 0.0 0 6 4 2 3 68 52 1 24 116.0 1 29 0.0 1 2 3 3 2 33 33 0 12 33.0 2 0 0.0 1 1 1
[6]:	4
[7]:	<pre>df['custcat'].value_counts() 3 281 1 266 4 236 2 217 Name: custcat, dtype: int64 df.hist(column='income', bins=50) array([[<axessubplot:title={'center':'income'}>]], dtype=object) income</axessubplot:title={'center':'income'}></pre>
	400 300 200
[8]: [[8]: [[9]: [<pre>df.columns Index(['region', 'tenure', 'age', 'marital', 'address', 'income', 'ed',</pre>
L0]:	[3., 11., 33., 1., 7., 136., 5., 5., 0., 0., 6.], [3., 68., 52., 1., 24., 116., 1., 29., 0., 1., 2.], [2., 33., 33., 0., 12., 33., 2., 0., 0., 1., 1.], [2., 23., 30., 1., 9., 30., 1., 2., 0., 0., 4.]]) y = df['custcat'].values y[0:5] array([1, 4, 3, 1, 3])
11]:	<pre>X = preprocessing.StandardScaler().fit(X).transform(X.astype(float)) X[0:5] array([[-0.02696767, -1.055125 , 0.18450456, 1.0100505 , -0.25303431,</pre>
	[-0.02696767, -0.58672182, -0.93080797, 1.0100505, -0.25303431, -0.44429125, -1.36767088, -0.89182893, -0.22207644, -1.03459817, 1.16300577]]) from sklearn.model_selection import train_test_split X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.2, random_state=4) print ('Train set:', X_train.shape, y_train.shape) print ('Test set:', X_test.shape, y_test.shape) Train set: (800, 11) (800,) Test set: (200, 11) (200,)
.4].	<pre>#K nearest neighbor (KNN) from sklearn.neighbors import KNeighborsClassifier k = 4 #Train Model and Predict neigh = KNeighborsClassifier(n_neighbors = k).fit(X_train, y_train) neigh KNeighborsClassifier(n_neighbors=4)</pre>
L6]:	<pre>#predicting yhat = neigh.predict(X_test) yhat[0:5] array([1, 1, 3, 2, 4]) #Accuracy evaluation from sklearn import metrics print("Train set Accuracy: ", metrics.accuracy_score(y_train, neigh.predict(X_train))) print("Test set Accuracy: ", metrics.accuracy_score(y_test, yhat))</pre> Train set Accuracy: 0.5475
17]:	<pre>Test set Accuracy: 0.32 #Building the algorithm with k = 6 k = 6 neigh = KNeighborsClassifier(n_neighbors = k).fit(X_train, y_train) neigh KNeighborsClassifier(n_neighbors=6) yhat = neigh.predict(X_test) yhat[0:5]</pre>
19]:	<pre>array([3, 3, 3, 4, 4]) print("Train set Accuracy: ", metrics.accuracy_score(y_train, neigh.predict(X_train))) print("Test set Accuracy: ", metrics.accuracy_score(y_test, yhat)) Train set Accuracy: 0.51625 Test set Accuracy: 0.31 #For other values of k Ks = 10 mean_acc = np.zeros((Ks-1)) std acc = np.zeros((Ks-1))</pre>
201.	<pre>for n in range(1,Ks): #Train Model and Predict neigh = KNeighborsClassifier(n_neighbors = n).fit(X_train,y_train) yhat=neigh.predict(X_test) mean_acc[n-1] = metrics.accuracy_score(y_test, yhat) std_acc[n-1]=np.std(yhat==y_test)/np.sqrt(yhat.shape[0]) mean_acc array([0.3 , 0.29 , 0.315, 0.32 , 0.315, 0.31 , 0.335, 0.325, 0.34])</pre>
20]:	<pre>plt.plot(range(1,Ks), mean_acc, 'g') plt.fill_between(range(1,Ks), mean_acc - 1 * std_acc, mean_acc + 1 * std_acc, alpha=0.10) plt.fill_between(range(1,Ks), mean_acc - 3 * std_acc, mean_acc + 3 * std_acc, alpha=0.10, color="green") plt.legend(('Accuracy ', '+/- 1xstd','+/- 3xstd')) plt.ylabel('Accuracy ') plt.xlabel('Number of Neighbors (K)') plt.tight_layout() plt.show()</pre>
	0.40
	Number of Neighbors (K) print("The best accuracy was with", mean_acc.max(), "with k=", mean_acc.argmax()+1) The best accuracy was with 0.34 with k= 9 Decision Tree import numpy as np import pandas as pd from sklearn.tree import DecisionTreeClassifier
[3]: [[4]: [[4]: _	# The data: A set of patients, all of whom suffered from the same illness. #During their course of treatment, each patient responded to one of 5 medications: #Drug A, Drug B, Drug c, Drug x and y. df = pd.read_csv('https://cf-courses-data.s3.us.cloud-object-storage.appdomain.cloud/IBMDeveloperSkillsNetw df.head(10) Age Sex BP Cholesterol Na_to_K Drug
	0 23 F HIGH HIGH 25.355 drugY 1 47 M LOW HIGH 13.093 drugC 2 47 M LOW HIGH 10.114 drugX 3 28 F NORMAL HIGH 7.798 drugX 4 61 F LOW HIGH 18.043 drugY 5 22 F NORMAL HIGH 8.607 drugX 6 49 F NORMAL HIGH 11.037 drugY 7 41 M LOW HIGH 11.037 drugY 8 60 M NORMAL HIGH 15.171 drugY
[5]: [5]:	8 60 M NORMAL HIGH 15.171 drugY 9 43 M LOW NORMAL 19.368 drugY df.shape (200, 6) """"Using df as the Drug.csv data read by pandas, declare the following variables: X as the Feature Matrix (data of df) y as the response vector (target) Remove the column containing the target name since it doesn't contain numeric values."""
[7]: [8]:	
[8]: [9]: [9]:	<pre>array([[23, 0, 0, 0, 25.355],</pre>
L0]: [L1]: [<pre>3 drugX 4 drugY Name: Drug, dtype: object #Setting up the decision tree from sklearn.model_selection import train_test_split """train_test_split will return 4 different parameters. Name them: X_trainset, X_testset, y_trainset, y_testset The train_test_split will need the parameters: X, y, test_size=0.3, and random_state=3.</pre>
12]: [The X and y are the arrays required before the split, the test_size represents the ratio of the testing dat X_trainset, X_testset, y_trainset, y_testset = train_test_split(X, y, test_size=0.3, random_state=3) print('Shape of X training set {}'.format(X_trainset.shape),'&',' Size of Y training set {}'.format(y_train Shape of X training set (140, 5) & Size of Y training set (140,) print('Shape of X training set {}'.format(X_testset.shape),'&',' Size of Y training set {}'.format(y_testset) Shape of X training set (60, 5) & Size of Y training set (60,)
L5]: L5]:	<pre>#Modeling #First create an instance of the DecisionTreeClassifier called drugTree. #Inside of the classifier, specify criterion="entropy" so that the information gain of each node can be see drugTree = DecisionTreeClassifier(criterion="entropy", max_depth = 4) drugTree # it shows the default parameters DecisionTreeClassifier(criterion='entropy', max_depth=4) #fit the data with the training feature matrix X_trainset and training response vector y_trainset drugTree.fit(X_trainset,y_trainset)</pre>
L6]: L8]: [L9]: [<pre>DecisionTreeClassifier(criterion='entropy', max_depth=4) #Prediction predTree = drugTree.predict(X_testset) print (predTree [0:5]) print (y_testset [0:5]) ['drugY' 'drugX' 'drugX' 'drugX' 'drugX'] 40</pre>
	<pre>51 drugX 139 drugX 197 drugX 170 drugX Name: Drug, dtype: object #Evaluation from sklearn import metrics import matplotlib.pyplot as plt print("DecisionTrees's Accuracy: ", metrics.accuracy_score(y_testset, predTree)) DecisionTrees's Accuracy: 0.98333333333333333333333333333333333333</pre>
28]:	<pre>#Visualization !conda install -c conda-forge python-graphviz -y !conda install -c conda-forge pydotplus -y from io import StringIO import pydotplus import matplotlib.image as mpimg from sklearn import tree %matplotlib inline</pre>
30]:	<pre>dot_data = StringIO() filename = "drugtree.png" featureNames = df.columns[0:5] out=tree.export_graphviz(drugTree, feature_names=featureNames, out_file=dot_data, class_names= np.unique(y_t graph = pydotplus.graph_from_dot_data(dot_data.getvalue()) graph.write_png(filename) img = mpimg.imread(filename) plt.figure(figsize=(100, 200)) plt.imshow(img,interpolation='nearest')</pre>
	<pre>InvocationException</pre>
	/var/folders/cn/0zbpg_4n7556t3y1p2lb9wlw0000gn/T/ipykernel_7694/3096224755.py in <module></module>
	<pre>/var/folders/cn/0zbpg_4n7556t3y1p21b9wlw0000gn/T/ipykernel_7694/3096224755.py in <module> 4 out=tree.export_graphviz(drugTree, feature_names=featureNames, out_file=dot_data, class_names= np.uni (y_trainset), filled=True, special_characters=True, rotate=False) 5 graph = pydotplus.graph_from_dot_data(dot_data.getvalue())> 6 graph.write_png(filename) 7 img = mpimg.imread(filename) 8 plt.figure(figsize=(100, 200)) ~/opt/anaconda3/lib/python3.9/site-packages/pydotplus/graphviz.py in <lambda>(path, f, prog) 1808</lambda></module></pre>
	<pre>/var/folders/cn/Ozbpg_4n7556t3ylp2lb9wlw0000qn/r/lpykernel_ 1684/3096224755.py in cmodule> 4 out=tree export graphytz (drugTree, feature names=featureNames, out_file=dot_data, class_names= np.uni (y_trainset), filled=True, special_characters=True,rotate=False)</pre>
] [] [31]:	/war/folders/on/deppg_dn356fdyp2plobelw0000gn/r/pykernel 7684/308622475.pp in concules 4 out-tree.export_graphividerure_newer_enewer=featureNames_out_file=dot_data, class_names= np.uni (y_trainset), filed=True, special_characters=True.rotate=Raise) 5 graph_write_pog(file=newe) 7 img = ppimg_interes_uraph_from_dot_data(dot_data.getvalue()) 7 img = ppimg_intered 8 plt.figure[figsize-(10, 200)) -/opt/anacconds]/ip/python3.8/site-packages/pydotplus/praphviz.py in <lambda*(peth. -="" 1809="" 1810="" 1811="" 1811<="" 1812="" 1813="" 1814="" 1815="" 1816="" 1817="" 1818="" 1819="" anacconds]="" f,="" format)="" graphviz.py="" iip="" in="" opt="" path,="" prog)="" prog,="" pydotplus="" python3.9="" site-packages="" th="" xrite(self,=""></lambda*(peth.>
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In [59]:	<pre>#Split dataset into train and test set X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.2, random_state=4) print ('Train set:', X_train.shape, y_train.shape) print ('Test set:', X_test.shape, y_test.shape) Train set: (546, 9) (546,) Test set: (137, 9) (137,)</pre>
In [60]:	"""Modeling (SVM with Scikit-learn) The SVM algorithm offers a choice of kernel functions for performing its processing. Basically, mapping data in The mathematical function used for the transformation is known as the kernel function, and can be of different 1.Linear 2.Polynomial 3.Radial basis function (RBF) 4.Sigmoid Each of these functions has its characteristics, its pros and cons, and its equation, but as there's no easy was We usually choose different functions in turn and compare the results. Using the default, RBF (Radial Basis Function)."""
Out[60]: In [61]: Out[61]:	<pre>from sklearn import svm clf = svm.SVC(kernel='rbf') clf.fit(X_train, y_train)</pre> SVC()
Out[61]: In [62]: In [63]:	<pre>#Evaluation from sklearn.metrics import classification_report, confusion_matrix import itertools def plot_confusion_matrix(cm, classes,</pre>
	<pre>cmap=plt.cm.Blues): """ This function prints and plots the confusion matrix. Normalization can be applied by setting `normalize=True`. """ if normalize: cm = cm.astype('float') / cm.sum(axis=1)[:, np.newaxis] print("Normalized confusion matrix") else: print('Confusion matrix, without normalization')</pre>
	<pre>plt.imshow(cm, interpolation='nearest', cmap=cmap) plt.title(title) plt.colorbar() tick_marks = np.arange(len(classes)) plt.xticks(tick_marks, classes, rotation=45) plt.yticks(tick_marks, classes) fmt = '.2f' if normalize else 'd' thresh = cm.max() / 2. for i, j in itertools.product(range(cm.shape[0]), range(cm.shape[1])): plt.text(j, i, format(cm[i, j], fmt),</pre>
In [64]:	<pre>plt.text(j, i, format(cm[i, j], fmt),</pre>
	<pre>np.set_printoptions(precision=2) print (classification_report(y_test, yhat)) # Plot non-normalized confusion matrix plt.figure() plot_confusion_matrix(cnf_matrix, classes=['Benign(2)','Malignant(4)'],normalize= False, title='Confusion matrix</pre>
	accuracy 0.96 137 macro avg 0.95 0.97 0.96 137 weighted avg 0.97 0.96 0.96 137 Confusion matrix, without normalization [[85 5] [0 47]] Confusion matrix Confusion matrix Sequence 85 5 5 5 5 5 5 5 5 5 5 5 5 5 5 5 5 5 5
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<pre>In [65]: Out[65]:</pre>	# Using f1_score from sklearn from sklearn.metrics import f1_score f1_score(y_test, yhat, average='weighted') 0.9639038982104676
In [66]: Out[66]: In [67]:	<pre>from sklearn.metrics import jaccard_score jaccard_score(y_test, yhat,pos_label=2) 0.944444444444444 #Rebuilding the model, with a _linear_ kernel. # usingkernel='linear' option, when defining the svm. How does the accuracy change with the new kernel fur</pre>
In []:	<pre>clf2 = svm.SVC(kernel='linear') clf2.fit(X_train, y_train) yhat2 = clf2.predict(X_test) print("Avg F1-score: %.4f" % f1_score(y_test, yhat2, average='weighted')) print("Jaccard score: %.4f" % jaccard_score(y_test, yhat2,pos_label=2)) Avg F1-score: 0.9639 Jaccard score: 0.9444</pre>