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Decision Trees

In []:

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In this lab exercise, you will learn a popular machine learning algorithm, Decision Tree. You will use this classification algorithm to build a model from historical data of patients, and their response to different medications. Then you use the trained decision tree to predict the class of a unknown patient, or to find a proper drug for a new patient.

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Import the Following Libraries:

- **numpy (as np)**
- **pandas**
- **DecisionTreeClassifier** from **sklearn.tree**

In [3]:

```
my_data = pd.read_csv("drug200.csv", delimiter=",")
my_data[0:5]
```

Out[3]:

	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	drugC
3	28	F	NORMAL	HIGH	7.798	drugX
4	61	F	LOW	HIGH	18.043	drugY

Practice

What is the size of data?

In []:

```
# write your code here
```

Pre-processing

Using **my_data** as the Drug.csv data read by pandas, declare the following variables:

- **X** as the **Feature Matrix** (data of my_data)
- **y** as the **response vector (target)**

Remove the column containing the target name since it doesn't contain numeric values.

In [4]:

```
X = my_data[['Age', 'Sex', 'BP', 'Cholesterol', 'Na_to_K']].values
X[0:5]
```

Out[4]:

```
array([[23, 'F', 'HIGH', 'HIGH', 25.355],
       [47, 'M', 'LOW', 'HIGH', 13.093],
       [47, 'M', 'LOW', 'HIGH', 10.113999999999999],
       [28, 'F', 'NORMAL', 'HIGH', 7.797999999999999],
       [61, 'F', 'LOW', 'HIGH', 18.043]], dtype=object)
```

As you may figure out, some features in this dataset are categorical such as **Sex** or **BP**. Unfortunately, Sklearn Decision Trees do not handle categorical variables. But still we can convert these features to numerical values. **pandas.get_dummies()** Convert categorical variable into dummy/indicator variables.

In [5]:

```
from sklearn import preprocessing
le_sex = preprocessing.LabelEncoder()
le_sex.fit(['F', 'M'])
X[:,1] = le_sex.transform(X[:,1])

le_BP = preprocessing.LabelEncoder()
le_BP.fit(['LOW', 'NORMAL', 'HIGH'])
X[:,2] = le_BP.transform(X[:,2])

le_Cholesterol = preprocessing.LabelEncoder()
le_Cholesterol.fit(['NORMAL', 'HIGH'])
X[:,3] = le_Cholesterol.transform(X[:,3])

X[0:5]
```

Out[5]:

```
array([[23, 0, 0, 0, 25.355],
       [47, 1, 1, 0, 13.093],
       [47, 1, 1, 0, 10.113999999999999],
       [28, 0, 2, 0, 7.797999999999999],
       [61, 0, 1, 0, 18.043]], dtype=object)
```

Now we can fill the target variable.

In [6]:

```
y = my_data["Drug"]  
y[0:5]
```

Out[6]:

```
0    drugY
1    drugC
2    drugC
3    drugX
4    drugY
Name: Drug, dtype: object
```

Setting up the Decision Tree

We will be using **train/test split** on our **decision tree**. Let's import **train_test_split** from **sklearn.cross_validation**.

In [7]:

```
from sklearn.model_selection import train_test_split
```

Now **train_test_split** will return 4 different parameters. We will 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.

The **X** and **y** are the arrays required before the split, the **test_size** represents the ratio of the testing dataset, and the **random_state** ensures that we obtain the same splits.

In [8]:

```
X_trainset, X_testset, y_trainset, y_testset = train_test_split(X, y, test_size=0.3, random_state=3)
```

Practice

Print the shape of `X_trainset` and `y_trainset`. Ensure that the dimensions match.

In []:

```
# your code
```

Print the shape of X_testset and y_testset. Ensure that the dimensions match

In []:

```
# your code
```

Modeling

We will first create an instance of the **DecisionTreeClassifier** called **drugTree**.

Inside of the classifier, specify `criterion="entropy"` so we can see the information gain of each node.

In [9]:

```
drugTree = DecisionTreeClassifier(criterion="entropy", max_depth = 4)
drugTree # it shows the default parameters
```

Out[9]:

```
DecisionTreeClassifier(class_weight=None, criterion='entropy', max_dept
h=4,
                        max_features=None, max_leaf_nodes=None,
                        min_impurity_decrease=0.0, min_impurity_split=None,
                        min_samples_leaf=1, min_samples_split=2,
                        min_weight_fraction_leaf=0.0, presort=False, random_state=N
one,
                        splitter='best')
```

Next, we will fit the data with the training feature matrix **X_trainset** and training response vector **y_trainset**

In [10]:

```
drugTree.fit(X_trainset,y_trainset)
```

Out[10]:

```
DecisionTreeClassifier(class_weight=None, criterion='entropy', max_dept
h=4,
                        max_features=None, max_leaf_nodes=None,
                        min_impurity_decrease=0.0, min_impurity_split=None,
                        min_samples_leaf=1, min_samples_split=2,
                        min_weight_fraction_leaf=0.0, presort=False, random_state=N
one,
                        splitter='best')
```

Prediction

Let's make some **predictions** on the testing dataset and store it into a variable called **predTree**.

In [11]:

```
predTree = drugTree.predict(X_testset)
```

You can print out **predTree** and **y_testset** if you want to visually compare the prediction to the actual values.

In [12]:

```
print (predTree [0:5])
print (y_testset [0:5])

['drugY' 'drugX' 'drugX' 'drugX' 'drugX']
40      drugY
51      drugX
139     drugX
197     drugX
170     drugX
Name: Drug, dtype: object
```

Evaluation

Next, let's import **metrics** from sklearn and check the accuracy of our model.

In [13]:

```
from sklearn import metrics
import matplotlib.pyplot as plt
print("DecisionTrees's Accuracy: ", metrics.accuracy_score(y_testset, predTree))
```

DecisionTrees's Accuracy: 0.9833333333333333

Accuracy classification score computes subset accuracy: the set of labels predicted for a sample must exactly match the corresponding set of labels in `y_true`.

In multilabel classification, the function returns the subset accuracy. If the entire set of predicted labels for a sample strictly match with the true set of labels, then the subset accuracy is 1.0; otherwise it is 0.0.

Practice

Can you calculate the accuracy score without sklearn ?

In []:

```
# your code here
```

Visualization

Lets visualize the tree

In []:

```
# Notice: You might need to uncomment and install the pydotplus and graphviz libraries if you have not installed these before  
# !conda install -c conda-forge pydotplus -y  
# !conda install -c conda-forge python-graphviz -y
```

In [14]:

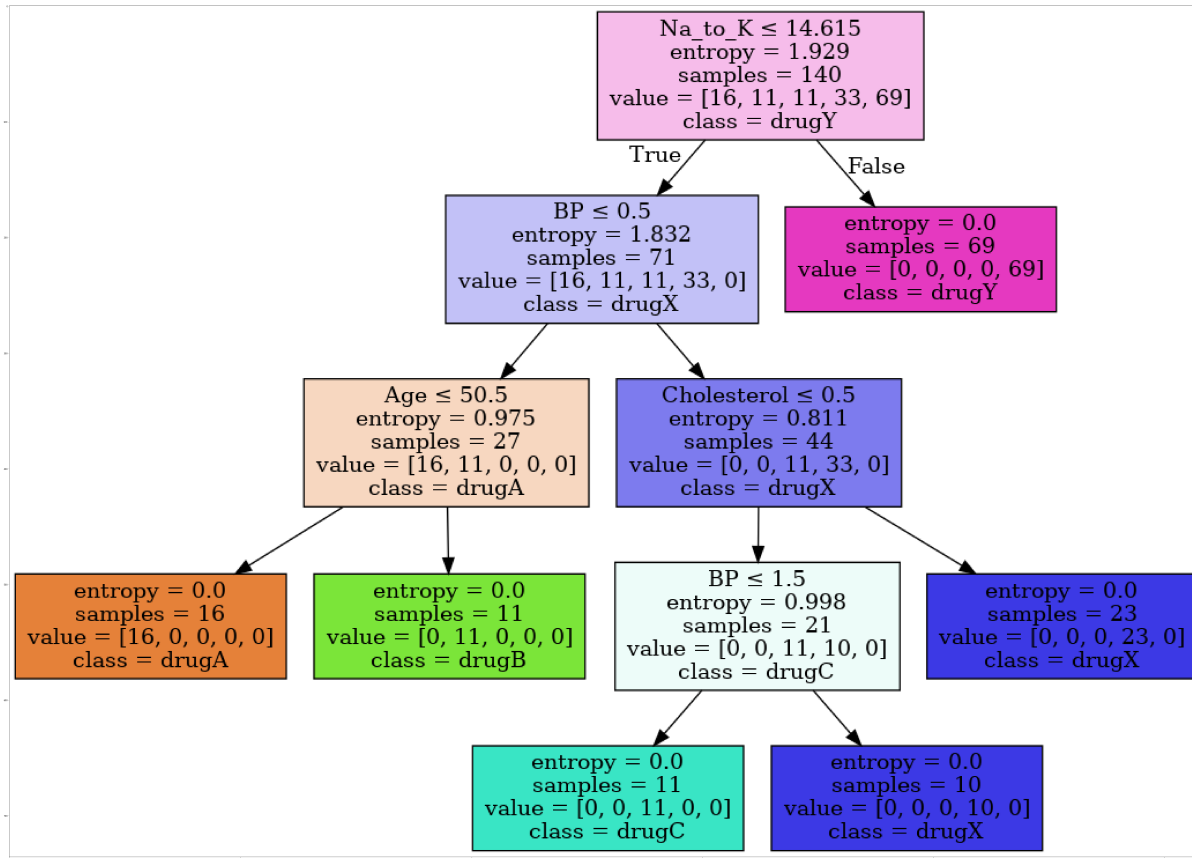
```
from sklearn.externals.six import StringIO  
import pydotplus  
import matplotlib.image as mpimg  
from sklearn import tree  
%matplotlib inline
```


In [15]:

```
dot_data = StringIO()
filename = "drugtree.png"
featureNames = my_data.columns[0:5]
targetNames = my_data["Drug"].unique().tolist()
out=tree.export_graphviz(drugTree,feature_names=featureNames, out_file=dot_data, class_names= np.unique(y_trainset), filled=True, special_characters=True,rotate=False)
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')
```

Out[15]:

<matplotlib.image.AxesImage at 0x7f8600d36d30>



Want to learn more?

IBM SPSS Modeler is a comprehensive analytics platform that has many machine learning algorithms. It has been designed to bring predictive intelligence to decisions made by individuals, by groups, by systems – by your enterprise as a whole. A free trial is available through this course, available here: [SPSS Modeler \(http://cocl.us/ML0101EN-SPSSModeler\)](http://cocl.us/ML0101EN-SPSSModeler)

Also, you can use Watson Studio to run these notebooks faster with bigger datasets. Watson Studio is IBM's leading cloud solution for data scientists, built by data scientists. With Jupyter notebooks, RStudio, Apache Spark and popular libraries pre-packaged in the cloud, Watson Studio enables data scientists to collaborate on their projects without having to install anything. Join the fast-growing community of Watson Studio users today with a free account at [Watson Studio \(https://cocl.us/ML0101EN_DSX\)](https://cocl.us/ML0101EN_DSX)

Thanks for completing this lesson!

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