

Decision Trees

Estimated time needed: 15 minutes

Objectives

After completing this lab you will be able to:

• Develop a classification model using Decision Tree Algorithm

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

Table of contents

- 1. About the dataset
- 2. Downloading the Data
- 3. Pre-processing
- 4. Setting up the Decision Tree
- 5. Modeling
- 6. Prediction
- 7. Evaluation
- 8. Visualization

Import the Following Libraries:

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

In [20]:

```
import numpy as np
import pandas as pd
from sklearn.tree import DecisionTreeClassifier
```

About the dataset

Imagine that you are a medical researcher compiling data for a study. You have collected data about 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.

Part of your job is to build a model to find out which drug might be appropriate for a future patient with the same illness. The features of this dataset are Age, Sex, Blood Pressure, and the Cholesterol of the patients, and the target is the drug that each patient responded to.

It is a sample of multiclass classifier, and you can use the training part of the dataset to build a decision tree, and then use it to predict the class of a unknown patient, or to prescribe a drug to a new patient.

Downloading the Data

To download the data, we will use !wget to download it from IBM Object Storage.

!wget -O drug200.csv https://cf-courses-data.s3.us.cloud-object-storage.appdomain.cloud/IBMDeveloperSkillsNetwork-ML0101EN-SkillsNetwork/labs/Module%203/data/drug200.csv

Did you know? When it comes to Machine Learning, you will likely be working with large datasets. As a business, where can you host your data? IBM is offering a unique opportunity for businesses, with 10 Tb of IBM Cloud Object Storage: Sign up now for free

Now, read the data using pandas dataframe:

```
In [22]:     my_data = pd.read_csv(myfile, delimiter=",")
     my_data[0:5]
```

Out[22]:		Age	Sex	BP	Cholesterol	Na_to_K	Drug	
	0	23	F	HIGH	HIGH	25.355	drugY	
	1	47	М	LOW	HIGH	13.093	drugC	
	2	47	М	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 [23]:  # write your code here

my_data.shape

Out[23]:  (200, 6)
```

▶ Click here for the solution

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.

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

```
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_Chol = preprocessing.LabelEncoder()
le_Chol.fit([ 'NORMAL', 'HIGH'])
X[:,3] = le_Chol.transform(X[:,3])
X[0:5]
```

Now we can fill the target variable.

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 [27]: 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 [28]:
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 [29]: # your code
    print('Shape of X training set {}'.format(X_trainset.shape),'&',' Size of Y training set {}'.format(y_trainset.shape))
    Shape of X training set (140, 5) & Size of Y training set (140,)
```

▶ Click here for the solution

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

```
# your code
print('Shape of X training set {}'.format(X_testset.shape),'&',' Size of Y training set {}'.format(y_testset.shape))
```

Shape of X training set (60, 5) & Size of Y training set (60,)

▶ Click here for the solution

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.

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

Out[31]: DecisionTreeClassifier(criterion='entropy', max_depth=4)

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

```
In [32]: drugTree.fit(X_trainset,y_trainset)
PacisionTreeClassifien(spitenien='entropy' max depth=4)
```

Out[32]: DecisionTreeClassifier(criterion='entropy', max_depth=4)

Prediction

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

```
In [33]: predTree = drugTree.predict(X_testset)
```

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

```
In [34]:
    print (predTree [0:5])
    print (y_testset [0:5])

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

Evaluation

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

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

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.

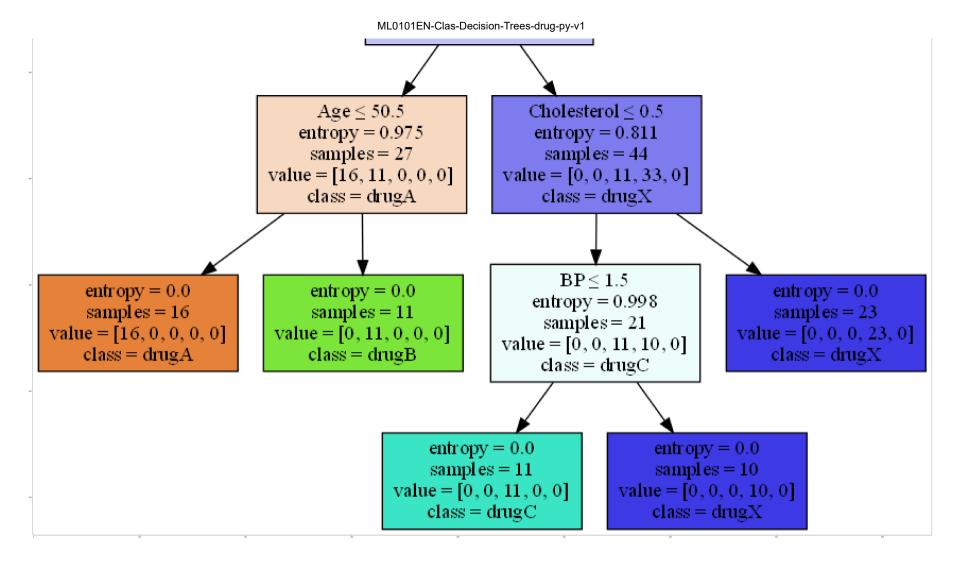
In multilabel classification, the function returns the subset accuracy. If the entire set of predicted labels for a sample strictly match with the true set of labels, then the subset accuracy is 1.0; otherwise it is 0.0.

Visualization

Let's visualize the tree

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 [38]:
          from io import StringIO
          import pydotplus
          import matplotlib.image as mpimg
          from sklearn import tree
          %matplotlib inline
In [39]:
          dot data = StringIO()
          filename = "drugtree.png"
          featureNames = my data.columns[0:5]
          out=tree.export graphviz(drugTree,feature names=featureNames, out file=dot data, class names= np.unique(y trainset), filled=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')
         <matplotlib.image.AxesImage at 0x1945b45d8b0>
Out[39]:
                                                                            Na to K \le 14.615
                                                                              entropy = 1.929
                                                                               samples = 140
                                                                       value = [16, 11, 11, 33, 69]
                                                                               class = drugY
                                                                        True
                                                                                                False
                                                                 BP \le 0.5
                                                                                                entropy = 0.0
                                                              entropy = 1.832
                                                                                                samples = 69
                                                               samples = 71
                                                                                           value = [0, 0, 0, 0, 69]
                                                        value = [16, 11, 11, 33, 0]
                                                                                               class = drugY
                                                               class = drugX
```



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

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