```
In [1]: import pandas as pd
        import numpy as np
        from sklearn.tree import DecisionTreeClassifier
        import sklearn.tree as tree
In [2]: import subprocess
        def runcmd(cmd, verbose = False, *args, **kwargs):
            process = subprocess.Popen(
                cmd,
                stdout = subprocess.PIPE,
                stderr = subprocess.PIPE,
                text = True,
                shell = True
            std_out, std_err = process.communicate()
            if verbose:
                print(std_out.strip(), std_err)
            pass
```

"Hello, World!"

## Downloading the data

runcmd('echo "Hello, World!"', verbose = True)

```
In [3]: runcmd("wget https://cf-courses-data.s3.us.cloud-object-storage.appdomain.clou
         --2022-10-20 23:47:45-- https://cf-courses-data.s3.us.cloud-object-storage.
        appdomain.cloud/IBMDeveloperSkillsNetwork-ML0101EN-SkillsNetwork/labs/Module%
        203/data/drug200.csv (https://cf-courses-data.s3.us.cloud-object-storage.appd
        omain.cloud/IBMDeveloperSkillsNetwork-ML0101EN-SkillsNetwork/labs/Module%203/
        data/drug200.csv)
        Resolving cf-courses-data.s3.us.cloud-object-storage.appdomain.cloud (cf-cour
        ses-data.s3.us.cloud-object-storage.appdomain.cloud)... 169.45.118.108
        Connecting to cf-courses-data.s3.us.cloud-object-storage.appdomain.cloud (cf-
        courses-data.s3.us.cloud-object-storage.appdomain.cloud) | 169.45.118.108 | :44
        connected.
        HTTP request sent, awaiting response... 200 OK
        Length: 5827 (5,7K) [text/csv]
        Saving to: 'drug200.csv'
             0K .....
                                                                       100% 1,70G=0s
        2022-10-20 23:47:48 (1,70 GB/s) - 'drug200.csv' saved [5827/5827]
```

Exploring the data

```
In [4]: df=pd.read_csv('drug200.csv')
In [5]: df.head()
```

#### Out[5]:

|   | Age | Sex | ВР     | 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 |

The data consists of several information regarding the age, sex, blood pressure(BP), cholesterol, drugs, and how the patients respond to the drugs.

```
In [6]: df.shape
Out[6]: (200, 6)
```

The data consists of 200 patients with 6 informations.

Sklearn decision trees does not read categorical data such as Sex, BP, and Cholesterol. Thus we must change the data into numerical values by using pd.get\_dummies() or preprocessing.LabelEncoder to convert the category into dummy numerics.

```
In [10]: df["Cholesterol"].value_counts()
Out[10]: HIGH
                    103
         NORMAL
                     97
          Name: Cholesterol, dtype: int64
          So F=0 and M=1, continue for BP and Cholesterol.
In [11]: le_BP = preprocessing.LabelEncoder()
         le_BP.fit([ 'LOW', 'NORMAL', 'HIGH'])
         X[:,2] = le_BP.transform(X[:,2])
In [12]: le Chol = preprocessing.LabelEncoder()
         le_Chol.fit([ 'NORMAL', 'HIGH'])
         X[:,3] = le\_Chol.transform(X[:,3])
         X[0:5]
Out[12]: array([[23, 0, 0, 0, 25.355],
                 [47, 1, 1, 0, 13.093],
                 [47, 1, 1, 0, 10.114],
                 [28, 0, 2, 0, 7.798],
                 [61, 0, 1, 0, 18.043]], dtype=object)
         Target Variable
In [13]: |y=df["Drug"]
         y[0:5]
Out[13]: 0
                 drugY
          1
                 drugC
          2
                 drugC
          3
                 drugX
          4
                 drugY
                 . . .
          195
                 drugC
          196
                 drugC
          197
                 drugX
          198
                 drugX
          199
                 drugX
          Name: Drug, Length: 200, dtype: object
```

## Setting the decision tree

```
In [14]:
        from sklearn.model_selection import train_test_split
In [15]: ??train_test_split
In [16]: |X_trainset, X_testset, y_trainset, y_testset=train_test_split(X, y, test_size=
```

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```
In [17]: X_trainset, X_testset, y_trainset, y_testset
Out[17]: (array([[26, 0, 0, 1, 19.161],
                  [41, 0, 2, 1, 22.905],
                  [28, 0, 2, 0, 19.675],
                  [19, 0, 0, 0, 13.313],
                  [50, 1, 2, 1, 15.79],
                  [24, 1, 2, 0, 25.786],
                  [72, 1, 1, 0, 16.31],
                  [74, 0, 1, 0, 20.942],
                  [37, 0, 1, 1, 12.006],
                  [31, 1, 0, 1, 17.069],
                  [22, 0, 2, 0, 8.607],
                  [20, 0, 2, 1, 9.281],
                  [28, 0, 1, 0, 13.127],
                  [59, 0, 2, 0, 13.884],
                  [15, 1, 0, 1, 17.206],
                  [51, 0, 1, 1, 23.003],
                  [45, 1, 1, 1, 10.017],
                  [33, 0, 1, 0, 33.486],
                  [39, 1, 0, 0, 9.664],
In [18]: | print('Shape of X test {}'.format(X_testset.shape),'&','Shape of y test {}'.fo
         Shape of X test (60, 5) & Shape of y test (60,)
         Modelling
In [19]:
         drugtree=DecisionTreeClassifier(criterion='entropy', max_depth=4)
         drugtree
```

### **Prediction**

Let's make prediction using the model above on the test data set and store it into a variable predTree.

```
In [21]: predTree=drugtree.predict(X_testset)
```

61

175

drugA

drugY

```
In [22]:
         print(predTree)
         print(y_testset)
         ['drugY' 'drugX'
                          'drugX' 'drugX' 'drugC' 'drugY' 'drugA' 'drugB'
           'drugA' 'drugY'
                          'drugA' 'drugY' 'drugY' 'drugX'
                                                           'drugY' 'drugX'
                                                                           'drugX'
           'drugB' 'drugX' 'drugY' 'drugY' 'drugY' 'drugX' 'drugB' 'drugY'
           'drugY' 'drugA'
                           'drugX' 'drugB' 'drugC' 'drugX' 'drugX' 'drugC'
          'drugY' 'drugX'
                          'drugX' 'drugA' 'drugY' 'drugC' 'drugY' 'drugA'
           'drugY' 'drugY' 'drugY' 'drugY' 'drugB' 'drugX' 'drugY' 'drugX'
           'drugY' 'drugY' 'drugA' 'drugX' 'drugY' 'drugX']
         40
                drugY
         51
                drugX
         139
                drugX
         197
                drugX
         170
                drugX
         82
                drugC
         183
                drugY
         46
                drugA
         70
                drugB
         100
                drugA
         179
                drugY
         83
                drugA
         25
                drugY
         190
                drugY
         159
                drugX
         173
                drugY
         95
                drugX
         3
                drugX
         41
                drugB
         58
                drugX
         14
                drugX
         143
                drugY
         12
                drugY
         6
                drugY
         182
                drugX
         161
                drugB
         128
                drugY
         122
                drugY
         101
                drugA
         86
                drugX
         64
                drugB
         47
                drugC
         158
                drugC
         34
                drugX
         38
                drugX
         196
                drugC
         4
                drugY
         72
                drugX
         67
                drugX
         145
                drugX
         156
                drugA
                drugY
         115
         155
                drugC
         15
                drugY
```

```
120
       drugY
130
       drugY
23
       drugY
153
       drugX
31
       drugB
103
       drugX
89
       drugY
132
       drugX
109
       drugY
126
       drugY
17
       drugA
30
       drugX
178
       drugY
162
       drugX
Name: Drug, dtype: object
```

The first 5 are okay as we could see, but let's evaluate the overall performance of the model.

#### **Evaluate The Model**

For model evaluation, we will use metrics from sklearn to check the accuracy of the model.

The closer the value to 1, the closer it matches with the testset, implying better accuracy.

# **Visualizing The Tree**

First, let's install the packages and import them

```
In [25]: #!conda install -c conda-forge pydotplus -y
#!conda install -c conda-forge python-graphviz -y
In [26]: ?tree.plot_tree
```

```
In [27]: tree.plot_tree(drugtree)
         plt.show()
In [28]: with open("drug200.csv", 'w') as f:
             f = tree.export_graphviz(clf, out_file=f)
In [29]: import os
         os.unlink('drug200.csv')
In [30]: import pydotplus
         dot_data=tree.export_graphviz(clf, out_file=None)
         graph2=pydotplus.graph_from_dot_data(dot_data)
         graph2.write("drug200.csv.pdf")
Out[30]: True
```

In [31]: feature\_names=['Age', 'Sex', 'BP', 'Cholesterol', 'Na\_to\_K']

In [32]: target\_names=['drugA','drugB', 'drugC', 'drugX', 'drugY']

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```
In [33]:
            from IPython.display import Image
            dot_data=tree.export_graphviz(clf, out_file=None,
                                       feature_names=feature_names, class_names=target_names, fi
                                       special_characters=True)
            graph2 = pydotplus.graph_from_dot_data(dot_data)
            nodes = graph2.get_node_list()
            Image(graph2.create_png() )
Out[33]:
                                                                   Na_to_K ≤ 14.615
                                                                    entropy = 1.929
                                                                    samples = 140
                                                               value = [16, 11, 11, 33, 69]
                                                                     class = drugY
                                                                True
                                                                                   False
                                                          BP ≤ 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
                                           Age ≤ 50.5
                                                                   Cholesterol ≤ 0.5
                                         entropy = 0.975
                                                                    entropy = 0.811
                                          samples = 27
                                                                     samples = 44
                                      value = [16, 11, 0, 0, 0]
                                                                value = [0, 0, 11, 33, 0]
                                                                    class = drugX
                                          class = drugA
                                                                       BP ≤ 1.5
                                                                                              entropy = 0.0
                  entropy = 0.0
                                           entropy = 0.0
                                                                   entropy = 0.998
                                           samples = 11
                                                                                              samples = 23
                  samples = 16
                                                                    samples = 21
              value = [16, 0, 0, 0, 0]
                                       value = [0, 11, 0, 0, 0]
                                                                                          ∨alue = [0, 0, 0, 23, 0]
                                                                value = [0, 0, 11, 10, 0]
                 class = drugA
                                          class = drugB
                                                                                              class = drugX
                                                                    class = drugC
                                                                                 entropy = 0.0
                                                        entropy = 0.0
                                                        samples = 11
                                                                                 samples = 10
                                                                             value = [0, 0, 0, 10, 0]
                                                    value = [0, 0, 11, 0, 0]
                                                        class = drugC
                                                                                 class = drugX
```

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#### Notes:

- 1) Everytime a split happens, we try to increase the purity of the prediction by using a certain criteria in each nodes.
- 2) With the accuracy of 0.98 is quite accurate. Below I also make sure that the classification is correct by filtering out data via conditions.
- 3) Drug A is for those who have Na\_to\_K lower than or equals to 14.6, High BP, and younger or at age 50, while Drug Y is for those who have Na\_to\_K higher than 14.6.

Numeric value interpretation:

```
BP: High=0, Low=1, Normal=2
Cholesterol: High=1, Normal=0
```

# Making sure that our labelling of drug classification is correct

```
In [34]: newdf = df[(df.Drug == "drugA")] #reading the graph, check whether BP of patien
drugXBP=newdf['BP'].value_counts()
newdf2 = df[(df.Drug == "drugB")]
drugBBP=newdf2['BP'].value_counts()
print("BP of Drug X:",drugXBP,"& BP of Drug B", drugBBP)
```

```
BP of Drug X: HIGH 23
Name: BP, dtype: int64 & BP of Drug B HIGH 16
Name: BP, dtype: int64
```

```
newdf = df[(df.Drug == "drugA")] #reading the graph, check whether Age of patie
In [36]:
          drugXAge=newdf['Age'].value_counts()
          newdf2 = df[(df.Drug == "drugB")]
          drugBAge=newdf2['Age'].value_counts()
          print("Age of Drug X:",
                drugXAge,
                "& Age of Drug B", drugBAge)
          Age of Drug X: 31
          49
                2
          32
                2
          43
                1
          35
                1
          47
                1
          48
                1
          42
                1
          20
                1
          26
                1
          39
                1
          50
                1
          45
                1
          38
                1
          19
                1
          36
                1
          29
                1
          24
                1
          37
                1
          23
                1
          Name: Age, dtype: int64 & Age of Drug B 60
          68
                2
          70
                2
          74
                1
          58
                1
          65
                1
          59
                1
          72
                1
          53
                1
          55
                1
          51
                1
          57
          Name: Age, dtype: int64
          It is as classified by the model.
```

In [ ]:

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