

1. Use your own dataset (Heart.csv is acceptable), create a train and a test set, and build 2 models: Logistic Regression and Decision Tree (shallow). Compare the test results using classification_report and confusion_matrix. Explain which algorithm is optimal

```
In [1]: import numpy as np
import pandas as pd
import seaborn as sns
import matplotlib.pyplot as plt

from sklearn.linear_model import LogisticRegression
from sklearn import preprocessing
from sklearn.tree import DecisionTreeClassifier
from sklearn.metrics import classification_report, confusion_matrix
from sklearn.model_selection import train_test_split
```

I used a dataset from UCI: Wine Quality (Red wine)

(<https://archive.ics.uci.edu/dataset/186/wine+quality> (<https://archive.ics.uci.edu/dataset/186/wine+quality>))

- This data shows scores of wine quality(from 1 to 10) with 11 related features
- Good wine : score ≥ 6 , normal wine : score < 6

1-A. Preprocessing

```
In [2]: # read the dataset as 'df'
df = pd.read_csv("winequality-red.csv")
df.head()
```

Out[2]:

	fixed acidity	volatile acidity	citric acid	residual sugar	chlorides	free sulfur dioxide	total sulfur dioxide	density	pH	sulphates	alcohol	quality
0	7.4	0.70	0.00	1.9	0.076	11.0	34.0	0.9978	3.51	0.56	9.4	5
1	7.8	0.88	0.00	2.6	0.098	25.0	67.0	0.9968	3.20	0.68	9.8	5
2	7.8	0.76	0.04	2.3	0.092	15.0	54.0	0.9970	3.26	0.65	9.8	5
3	11.2	0.28	0.56	1.9	0.075	17.0	60.0	0.9980	3.16	0.58	9.8	6
4	7.4	0.70	0.00	1.9	0.076	11.0	34.0	0.9978	3.51	0.56	9.4	5

- As variable 'quality' has values from 1 to 10, transforming the values into binary values is necessary
- values from 6 to 10 are transformed into value 1(good wine), and values from 1 to 5 are transformed into value 0(normal wine)

```
In [3]: # transform values from 6 to 10 into value 1
df['quality'] = df.quality.between(6,10).astype(int)
df.head()
```

Out[3]:

	fixed acidity	volatile acidity	citric acid	residual sugar	chlorides	free sulfur dioxide	total sulfur dioxide	density	pH	sulphates	alcohol	quality
0	7.4	0.70	0.00	1.9	0.076	11.0	34.0	0.9978	3.51	0.56	9.4	0
1	7.8	0.88	0.00	2.6	0.098	25.0	67.0	0.9968	3.20	0.68	9.8	0
2	7.8	0.76	0.04	2.3	0.092	15.0	54.0	0.9970	3.26	0.65	9.8	0
3	11.2	0.28	0.56	1.9	0.075	17.0	60.0	0.9980	3.16	0.58	9.8	1
4	7.4	0.70	0.00	1.9	0.076	11.0	34.0	0.9978	3.51	0.56	9.4	0

In [4]: #do train/test split

```
x_train, x_test, y_train, y_test = train_test_split(df.drop(['quality'], axis=1), df.quality, test_size=.2)

x_train.shape, x_test.shape, y_train.shape, y_test.shape
```

Out[4]: ((1279, 11), (320, 11), (1279,), (320,))

In [5]: y_test.value_counts()

Out[5]:

```
quality
1    179
0    141
Name: count, dtype: int64
```

1-B. Logistic Regression

In [6]: #generate logistic regression model as 'model1'

```
model1 = LogisticRegression()
model1.fit(x_train, y_train)
```

C:\Users\Wkcosm\Anaconda3\Lib\site-packages\sklearn\linear_model\logistic.py:469: ConvergenceWarning: lbfgs failed to converge (status=1):
STOP: TOTAL NO. of ITERATIONS REACHED LIMIT.

Increase the number of iterations (max_iter) or scale the data as shown in:

<https://scikit-learn.org/stable/modules/preprocessing.html> (<https://scikit-learn.org/stable/modules/preprocessing.html>)

Please also refer to the documentation for alternative solver options:

https://scikit-learn.org/stable/modules/linear_model.html#logistic-regression (https://scikit-learn.org/stable/modules/linear_model.html#logistic-regression)

```
n_iter_i = _check_optimize_result(
```

Out[6]:

▼ LogisticRegression

```
LogisticRegression()
```

In [7]: #As the warning sign indicates total number of iterations reached limit,
#I increased maximum number of iterations from 100(defalut) to 1000

```
model1 = LogisticRegression(max_iter=1000)
model1.fit(x_train, y_train)
```

Out[7]:

▼ LogisticRegression

```
LogisticRegression(max_iter=1000)
```

In [8]: #Print the test results

```
prediction1 = model1.predict(x_test)
print(classification_report(y_test, prediction1))
print(confusion_matrix(y_test, prediction1))
```

```
              precision    recall  f1-score   support

     0       0.70      0.74      0.72         141
     1       0.79      0.74      0.76         179

 accuracy          0.74
 macro avg       0.74      0.74      0.74
weighted avg       0.75      0.74      0.74

[[105  36]
 [ 46 133]]
```

1-C. Decision Tree

In [9]: #build a shallow decision tree model(max_depth=3) as 'model2'

```
model2 = DecisionTreeClassifier(criterion='entropy', max_depth=3)
model2.fit(x_train, y_train)
```

Out[9]:

DecisionTreeClassifier
DecisionTreeClassifier(criterion='entropy', max_depth=3)

In [10]: #Print the test results

```
prediction2 = model2.predict(x_test)
print(classification_report(y_test, prediction2))
print(confusion_matrix(y_test, prediction2))
```

	precision	recall	f1-score	support
0	0.58	0.92	0.71	141
1	0.89	0.47	0.62	179
accuracy			0.67	320
macro avg	0.73	0.70	0.67	320
weighted avg	0.75	0.67	0.66	320


```
[[130 11]
 [ 94 85]]
```

Result : Comparing cell 8 and 10, Logistic Regression model is more optimal.

- Logistic Regression model shows higher f1-score in predicting good wine (value 1, 0.76 > 0.62) and normal wine (value 0, 0.72 > 0.71)

2. Repeat 1. but let the Decision Tree be much deeper to allow over-fitting. Compare the two models' test results again, and explain which is optimal

2-A. Deeper Decision Tree(max_depth=10)

In [11]: #build a deeper decision tree model(max_depth=10) as 'model3'

```
model3 = DecisionTreeClassifier(criterion='entropy', max_depth=10)
model3.fit(x_train, y_train)
```

Out[11]:

DecisionTreeClassifier
DecisionTreeClassifier(criterion='entropy', max_depth=10)

```
In [12]: #Print the test results
prediction3 = model3.predict(x_test)
print(classification_report(y_test, prediction3))
print(confusion_matrix(y_test, prediction3))
```

	precision	recall	f1-score	support
0	0.71	0.75	0.73	141
1	0.79	0.75	0.77	179
accuracy			0.75	320
macro avg	0.75	0.75	0.75	320
weighted avg	0.76	0.75	0.75	320

```
[[106 35]
 [ 44 135]]
```

2-B. Much Deeper Decision Tree(max_depth=50)

```
In [13]: #build a much deeper decision tree model(max_depth=50) as 'model4'

model4 = DecisionTreeClassifier(criterion='entropy', max_depth=50)
model4.fit(x_train, y_train)
```

Out[13]:

DecisionTreeClassifier
DecisionTreeClassifier(criterion='entropy', max_depth=50)

(https://scikit-learn.org/1.4/modules/generated/sklearn.tree.DecisionTreeClassifier.html)

```
In [14]: #Print the test results
prediction4 = model4.predict(x_test)
print(classification_report(y_test, prediction4))
print(confusion_matrix(y_test, prediction4))
```

	precision	recall	f1-score	support
0	0.70	0.79	0.74	141
1	0.82	0.73	0.77	179
accuracy			0.76	320
macro avg	0.76	0.76	0.76	320
weighted avg	0.77	0.76	0.76	320

```
[[112 29]
 [ 48 131]]
```

Result : Comparing the cell 8, 12, and 14, deeper Decision Tree model outperforms Logistic Regression model

- both Decision Tree model shows higher f1-score than Logistic Regression model
- f1-score improved as max depth of Decision Tree increased