In [23]:

```
#LIBRARIES
import mglearn as mglearn
import os
import graphviz
import matplotlib.pyplot as plt
import sklearn.datasets as datasets
import pandas as pd
import numpy as np
import tensorflow as tf
from sklearn.metrics import accuracy score
from sklearn.tree import DecisionTreeClassifier
from IPython.display import Image
from sklearn.tree import export graphviz
from sklearn.model selection import train test split
from sklearn.neural_network import MLPRegressor
from sklearn.tree import DecisionTreeRegressor
from sklearn.linear model import LinearRegression
from sklearn.ensemble import RandomForestClassifier
from sklearn.datasets import make moons
from sklearn.ensemble import GradientBoostingClassifier
from sklearn.datasets import make blobs
from sklearn.svm import LinearSVC
import sklearn.svm as linear svm
from mpl toolkits.mplot3d import Axes3D, axes3d
from sklearn.svm import SVC
from sklearn.neural network import MLPClassifier
from tensorflow.examples.tutorials.mnist import input data
from sklearn.datasets import make circles
from sklearn.datasets import load wine
from sklearn.datasets import fetch california housing
from sklearn.neural network import MLPRegressor
```

In [2]:

Wine Data Database

Notes

Data Set Characteristics:

- :Number of Instances: 178 (50 in each of three classes)
- :Number of Attributes: 13 numeric, predictive attributes and the class

:Attribute Information:

- 1) Alcohol
- 2) Malic acid
- 3) Ash
- 4) Alcalinity of ash
- 5) Magnesium
- 6) Total phenols
- 7) Flavanoids
- 8) Nonflavanoid phenols
- 9) Proanthocyanins
- 10)Color intensity
- 11)Hue
- 12)OD280/OD315 of diluted wines
- 13)Proline
- class:
- class 0
- class 1
- class_2

:Summary Statistics:

| ======================================= | ==== | ===== | ====== | ===== |
|---|------|-------|--------|-------|
| | Min | Max | Mean | SD |
| ======================================= | ==== | ===== | ====== | ===== |
| Alcohol: | 11.0 | 14.8 | 13.0 | 0.8 |
| Malic Acid: | 0.74 | 5.80 | 2.34 | 1.12 |
| Ash: | 1.36 | 3.23 | 2.36 | 0.27 |
| Alcalinity of Ash: | 10.6 | 30.0 | 19.5 | 3.3 |
| Magnesium: | 70.0 | 162.0 | 99.7 | 14.3 |
| Total Phenols: | 0.98 | 3.88 | 2.29 | 0.63 |
| Flavanoids: | 0.34 | 5.08 | 2.03 | 1.00 |
| Nonflavanoid Phenols: | 0.13 | 0.66 | 0.36 | 0.12 |
| Proanthocyanins: | 0.41 | 3.58 | 1.59 | 0.57 |
| Colour Intensity: | 1.3 | 13.0 | 5.1 | 2.3 |
| Hue: | 0.48 | 1.71 | 0.96 | 0.23 |
| OD280/OD315 of diluted wines: | 1.27 | 4.00 | 2.61 | 0.71 |
| Proline: | 278 | 1680 | 746 | 315 |
| ======================================= | ==== | ===== | ====== | ===== |

:Missing Attribute Values: None

:Class Distribution: class 0 (59), class 1 (71), class 2 (48)

:Creator: R.A. Fisher

:Donor: Michael Marshall (MARSHALL%PLU@io.arc.nasa.gov)

:Date: July, 1988

This is a copy of UCI ML Wine recognition datasets. https://archive.ics.uci.edu/ml/machine-learning-databases/wine/wine.data

The data is the results of a chemical analysis of wines grown in the

region in Italy by three different cultivators. There are thirteen d ifferent

measurements taken for different constituents found in the three types of

wine.

Original Owners:

Forina, M. et al, PARVUS -

An Extendible Package for Data Exploration, Classification and Corre lation.

Institute of Pharmaceutical and Food Analysis and Technologies, Via Brigata Salerno, 16147 Genoa, Italy.

Citation:

Lichman, M. (2013). UCI Machine Learning Repository
[http://archive.ics.uci.edu/ml]. Irvine, CA: University of California,

School of Information and Computer Science.

References

(1)

S. Aeberhard, D. Coomans and O. de Vel,

Comparison of Classifiers in High Dimensional Settings,

Tech. Rep. no. 92-02, (1992), Dept. of Computer Science and Dept. of Mathematics and Statistics, James Cook University of North Queensland.

(Also submitted to Technometrics).

The data was used with many others for comparing various classifiers. The classes are separable, though only RDA has achieved 100% correct classification.

(RDA: 100%, QDA 99.4%, LDA 98.9%, 1NN 96.1% (z-transformed data))

(All results using the leave-one-out technique)

(2)

S. Aeberhard, D. Coomans and O. de Vel,

"THE CLASSIFICATION PERFORMANCE OF RDA"

Tech. Rep. no. 92-01, (1992), Dept. of Computer Science and Dept. of Mathematics and Statistics, James Cook University of North Queenslan d.

(Also submitted to Journal of Chemometrics).

Out[2]:

| | Class label | Alcohol | Malic acid | Ash | Alcalinity of ash | Magnesium | Total phenols | Flavanoids | Nonflavanoi pheno |
|---|----------------|---------|---------------|------|----------------------|-----------|------------------|------------|----------------------|
| 0 | 14.23 | 1.71 | 2.43 | 15.6 | 127.0 | 2.80 | 3.06 | 0.28 | 2.29 |
| 1 | 13.20 | 1.78 | 2.14 | 11.2 | 100.0 | 2.65 | 2.76 | 0.26 | 1.28 |
| 2 | 13.16 | 2.36 | 2.67 | 18.6 | 101.0 | 2.80 | 3.24 | 0.30 | 2.81 |
| 3 | 14.37 | 1.95 | 2.50 | 16.8 | 113.0 | 3.85 | 3.49 | 0.24 | 2.18 |
| 4 | 13.24 | 2.59 | 2.87 | 21.0 | 118.0 | 2.80 | 2.69 | 0.39 | 1.82 |

In [3]:

```
X train, X test, y train, y test = train test split(wine.data,
                                                     wine.target,
                                                     stratify = wine.target,
                                                     random state = 42)
tree = DecisionTreeClassifier(criterion = 'entropy',
                             min samples split = 20,
                             min samples leaf = 15,
                             max features = 'sqrt',
                             max leaf nodes = 12,
                             random state = 0)
tree.fit(X_train, y_train)
y pred = tree.predict(X test)
print('#Training data points: %d' % X train.shape[0])
print('#Testing data points: %d' % X test.shape[0])
print("")
print("Accuracy on training set: {:,.3f}".format(tree.score(X_train, y_train)))
print("Accuracy on test set: {:,.3f}".format(tree.score(X_test, y_test)))
print("")
print('Class labels:', np.unique(wine.target))
print('Misclassified samples: %d' % (y_test != y_pred).sum())
#Training data points: 133
```

#Testing data points: 45

Accuracy on training set: 0.895

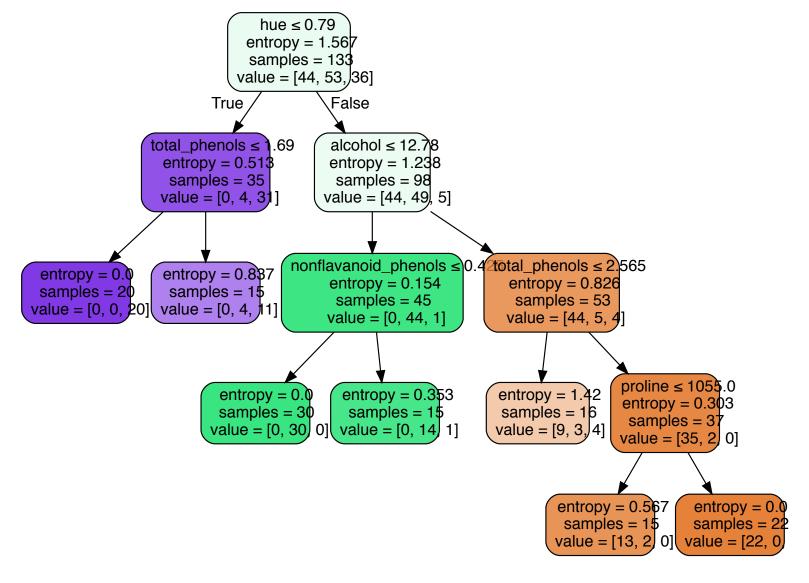
Accuracy on test set: 0.844

Class labels: [0 1 2]

Misclassified samples: 7

```
In [4]:
```

```
X train, X test, y train, y test = train test split(wine.data,
                                                     wine.target,
                                                     stratify = wine.target,
                                                     random state = 42)
tree = DecisionTreeClassifier(criterion = 'entropy',
                             min samples split = 20,
                             min samples leaf = 15,
                             max features = 'sqrt',
                             max leaf_nodes = 12,
                             random state = 0)
tree.fit(X_train, y_train)
export graphviz(tree, out file = "tree.dot",
                feature names = wine.feature names,
                filled = True, rounded = True,
               special characters = True) #[writes data into .dot file]
with open("tree.dot") as f:
    dot graph = f.read()
display(graphviz.Source(dot graph))
print("Feature importance: \n{}".format(tree.feature importances ))
def plot feature importances wine(model):
    n features = wine.data.shape[1]
    plt.barh(range(n_features), model.feature_importances_, align = 'center')
    plt.yticks(np.arange(n features), wine.feature names)
    plt.xlabel("Feature Importance")
    plt.ylabel("Feature")
    plt.ylim(-1, n features)
plot feature importances wine(tree)
tree = mglearn.plots.plot tree not monotone()
display(tree)
```



Feature importance:

[0.44311648 0.

0.

0.

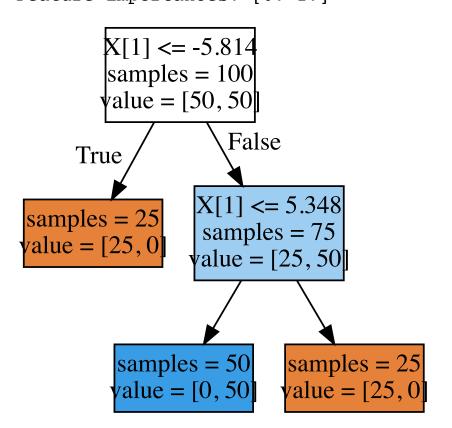
0.

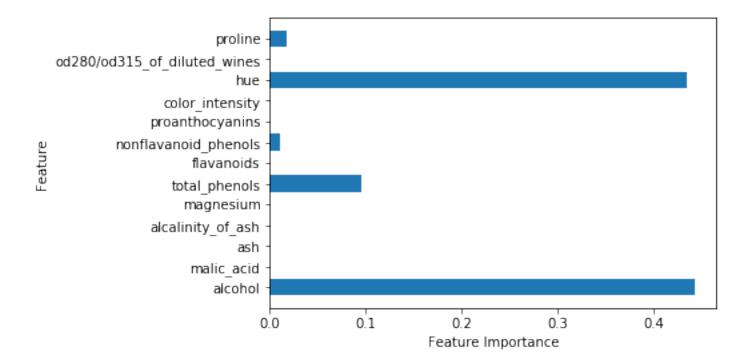
0.09543221

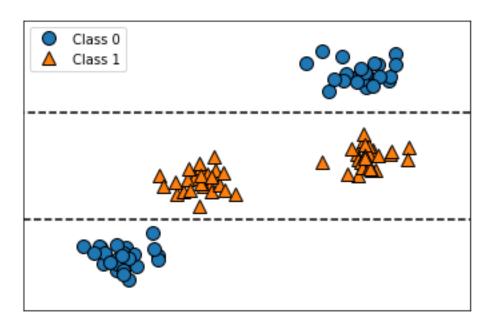
0.01015208 0.

0.43418745 0.

0.01711179]
Feature importances: [0. 1.]

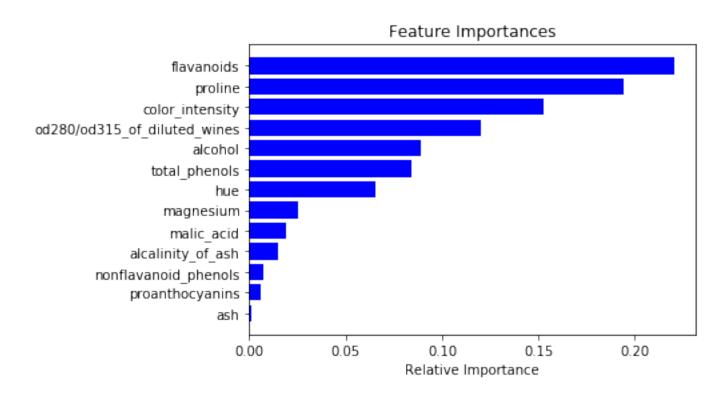






```
#RANDOM FOREST
forest = RandomForestClassifier(criterion='entropy',
                                n = 50,
                                 min samples split = 20,
                                min samples leaf = 15,
                                max features = 'sqrt',
                                max leaf nodes = 12,
                                random state = 0)
forest.fit(X train, y train)
y pred = forest.predict(X test)
print("Accuracy on training set: {:.3f}".format(forest.score(X_train, y_train)))
print("Accuracy on test set: {:.3f}".format(forest.score(X_test, y_test)))
importances = forest.feature importances
features = wine['feature names']
indices = np.argsort(importances)
plt.title('Feature Importances')
plt.barh(range(len(indices)), importances[indices], color='b', align='center')
plt.yticks(range(len(indices)), [features[i] for i in indices])
plt.xlabel('Relative Importance')
plt.show()
```

Accuracy on training set: 0.985 Accuracy on test set: 1.000



```
In [6]:
estimator = forest.estimators [29]
export_graphviz(estimator, out_file = "rftree.dot",
                 feature names = wine.feature names,
                class names = wine.target names,
                rounded = True, proportion = False,
                precision = 2, filled = True)
with open("rftree.dot") as f:
    dot graph = f.read()
display(graphviz.Source(dot graph))
                    hue <= 0.83
                   entropy = 1.56
                   samples = 84
                 value = [46, 53, 34]
                   class = class_1
                               False
                True
                            alcohol <= 13.1
          entropy = 0.43
                            entropy = 1.16
          samples = 21
                             samples = 63
         value = [0, 3, 31]
                           √alue = [46, 50, 3]
          class = class
                            class = class_1
                 proline <= 541.0 color_intensity <= 5.22
                  entropy = 0.39
                                       entropy = 0.6
```

entropy = 0.0 samples = 15 value = [0, 29, 0] class = class 1

entropy = 0.77 samples = 15 value = [1, 16, 2] class = class 1

samples = 30

/alue = [1, 45, 2]

class = class_

entropy = 0.85 samples = 15 value = [13, 5, 0] class = class_0

samples = 33

value = [45, 5, 1]

class = class_0

entropy = 0.2 samples = 18 value = [32, 0, 1] class = class 0

In [16]:

| 1) | Flavanoids | 0.220670 |
|-----|------------------------------|----------|
| 2) | Proline | 0.194247 |
| 3) | Color intensity | 0.152337 |
| 4) | OD280/OD315 of diluted wines | 0.120117 |
| 5) | Alcohol | 0.088855 |
| 6) | Total phenols | 0.084423 |
| 7) | Hue | 0.065712 |
| 8) | Magnesium | 0.025427 |
| 9) | Malic acid | 0.018752 |
| 10) | Alcalinity of ash | 0.015239 |
| 11) | Nonflavanoid phenols | 0.007226 |
| 12) | Proanthocyanins | 0.005836 |
| 13) | Ash | 0.001158 |
| | | |

In [21]:

```
#Gradient Boosted Trees MaxDepth = 0.01
X train, X test, y train, y test = train test split(wine.data, wine.target, rand
om state=0)
gbrt=GradientBoostingClassifier(random state=0, learning rate=0.01)
gbrt.fit(X_train, y_train)
print("Accuracy on training set: {:.3f}".format(gbrt.score(X_train, y_train)))
print("Accuracy on test set: {:.3f}".format(gbrt.score(X test, y test)))
importances = gbrt.feature importances
features = wine['feature names']
indices = np.argsort(importances)
def plot feature importances wine(model):
       n features = wine.data.shape[1]
       plt.barh(range(n features), model.feature importances , align='center')
       plt.yticks(np.arange(n features), wine.feature names)
       plt.xlabel("Feature importance")
       plt.ylabel("Fetaure")
       plt.ylim(-1, n features)
       plot_feature_importances_wine(gbrt)
```

Accuracy on training set: 1.000 Accuracy on test set: 0.933

```
#Gradient Boosted Trees MaxDepth = 0.05
X train, X test, y train, y test = train test split(wine.data, wine.target, rand
om state=0)
gbrt=GradientBoostingClassifier(random state=42, learning rate=0.05)
gbrt.fit(X_train, y_train)
print("Accuracy on training set: {:.3f}".format(gbrt.score(X_train, y_train)))
print("Accuracy on test set: {:.3f}".format(gbrt.score(X test, y test)))
importances = gbrt.feature importances
features = wine['feature names']
indices = np.argsort(importances)
def plot feature importances wine(model):
       n features = wine.data.shape[1]
       plt.barh(range(n features), model.feature importances , align='center')
       plt.yticks(np.arange(n features), wine.feature names)
       plt.xlabel("Feature importance")
       plt.ylabel("Fetaure")
       plt.ylim(-1, n features)
       plot_feature_importances_wine(gbrt)
```

Accuracy on training set: 1.000 Accuracy on test set: 0.956

In [10]:

```
#Gradient Boosted Trees MaxDepth = 0.1
X train, X test, y train, y test = train test split(wine.data, wine.target, rand
om state=0)
gbrt=GradientBoostingClassifier(random state=42, learning rate=0.1)
gbrt.fit(X_train, y_train)
print("Accuracy on training set: {:.3f}".format(gbrt.score(X_train, y_train)))
print("Accuracy on test set: {:.3f}".format(gbrt.score(X test, y test)))
importances = gbrt.feature importances
features = wine['feature names']
indices = np.argsort(importances)
def plot feature importances wine(model):
       n features = wine.data.shape[1]
       plt.barh(range(n features), model.feature importances , align='center')
       plt.yticks(np.arange(n features), wine.feature names)
       plt.xlabel("Feature importance")
       plt.ylabel("Fetaure")
       plt.ylim(-1, n features)
       plot_feature_importances_wine(gbrt)
```

Accuracy on training set: 1.000 Accuracy on test set: 0.956

```
In [11]:
```

```
#Gradient Boosted Trees MaxDepth = 1
X train, X test, y train, y test = train test split(wine.data,
                                                     wine.target,
                                                     stratify = wine.target,
                                                     random state = 0)
gbrt=GradientBoostingClassifier(random state=42, max depth=1)
gbrt.fit(X train, y train)
print("Accuracy on training set: {:.3f}".format(gbrt.score(X train, y train)))
print("Accuracy on test set: {:.3f}".format(gbrt.score(X test, y test)))
importances = gbrt.feature importances
features = wine['feature names']
indices = np.argsort(importances)
def plot feature importances wine(model):
       n features = wine.data.shape[1]
       plt.barh(range(n features), model.feature importances , align='center')
       plt.yticks(np.arange(n features), wine.feature names)
       plt.xlabel("Feature importance")
       plt.ylabel("Fetaure")
       plt.ylim(-1, n_features)
       plot_feature_importances wine(gbrt)
Accuracy on training set: 1.000
Accuracy on test set: 1.000
In [24]:
```

```
ch = fetch_california_housing()
df = pd.DataFrame(ch.data,columns=ch.feature_names)
df.head()
```

Out[24]:

| | MedInc | HouseAge | AveRooms | AveBedrms | Population | AveOccup | Latitude | Longi |
|---|--------|----------|----------|-----------|------------|----------|----------|--------|
| 0 | 8.3252 | 41.0 | 6.984127 | 1.023810 | 322.0 | 2.555556 | 37.88 | -122.2 |
| 1 | 8.3014 | 21.0 | 6.238137 | 0.971880 | 2401.0 | 2.109842 | 37.86 | -122.2 |
| 2 | 7.2574 | 52.0 | 8.288136 | 1.073446 | 496.0 | 2.802260 | 37.85 | -122.2 |
| 3 | 5.6431 | 52.0 | 5.817352 | 1.073059 | 558.0 | 2.547945 | 37.85 | -122.2 |
| 4 | 3.8462 | 52.0 | 6.281853 | 1.081081 | 565.0 | 2.181467 | 37.85 | -122.2 |

In [27]:

```
print("Adjusted R-square on training set: {:.3f}".format(mlp.score(X_train, y_tr
ain)))
print("Adjusted R-square on test set: {:.3f}".format(mlp.score(X_test, y_test)))
```

Adjusted R-square on training set: 0.466 Adjusted R-square on test set: 0.451