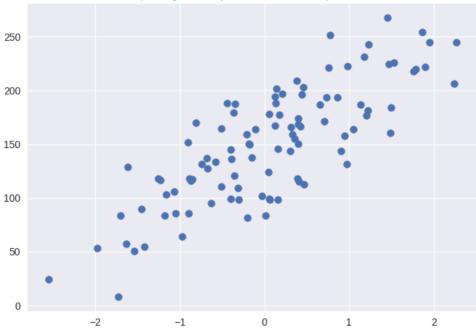
You are currently looking at **version 1.0** of this notebook. To download notebooks and datafiles, as well as get help on Jupyter notebooks in the Coursera platform, visit the <u>Jupyter Notebook FAQ</u> (https://www.coursera.org/learn/python-machine-learning/resources/bANLa) course resource.

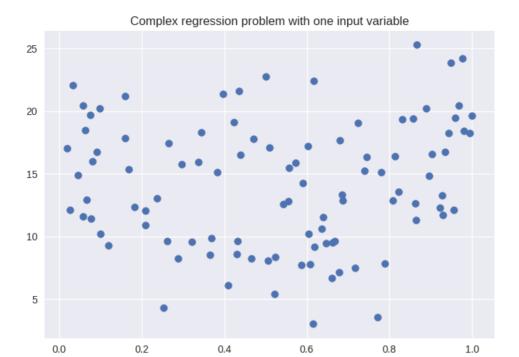
Applied Machine Learning: Module 4 (Supervised Learning, Part II) ¶

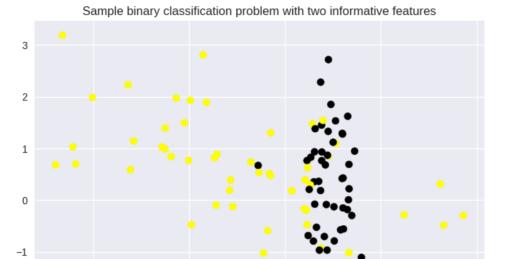
Preamble and Datasets

```
%matplotlib notebook
import numpy as np
import pandas as pd
import seaborn as sn
import matplotlib.pyplot as plt
from sklearn.model selection import train test split
from sklearn.datasets import make classification, make blobs
from matplotlib.colors import ListedColormap
from sklearn.datasets import load breast cancer
from adspy shared utilities import load crime dataset
cmap bold = ListedColormap(['#FFFF00', '#00FF00', '#0000FF', '#000000'])
# fruits dataset
fruits = pd.read table('readonly/fruit data with colors.txt')
feature_names_fruits = ['height', 'width', 'mass', 'color score']
X fruits = fruits[feature names fruits]
y fruits = fruits['fruit label']
target_names_fruits = ['apple', 'mandarin', 'orange', 'lemon']
X_fruits_2d = fruits[['height', 'width']]
y fruits 2d = fruits['fruit label']
# synthetic dataset for simple regression
from sklearn.datasets import make regression
plt.figure()
plt.title('Sample regression problem with one input variable')
X R1, y R1 = make regression(n samples = 100, n features=1,
                            n informative=1, bias = 150.0,
                            noise = 30, random state=0)
plt.scatter(X R1, y R1, marker= 'o', s=50)
plt.show()
# synthetic dataset for more complex regression
from sklearn.datasets import make friedman1
plt.figure()
plt.title('Complex regression problem with one input variable')
X_F1, y_F1 = make_friedman1(n_samples = 100, n_features = 7,
                           random state=0)
plt.scatter(X_F1[:, 2], y_F1, marker= 'o', s=50)
plt.show()
# synthetic dataset for classification (binary)
plt.figure()
plt.title('Sample binary classification problem with two informative features')
X C2, y C2 = make classification(n samples = 100, n features=2,
                                n_redundant=0, n_informative=2,
                                n_clusters_per_class=1, flip_y = 0.1,
                                class_sep = 0.5, random state=0)
plt.scatter(X C2[:, 0], X C2[:, 1], marker= 'o',
           c=y_C2, s=50, cmap=cmap_bold)
plt.show()
# more difficult synthetic dataset for classification (binary)
# with classes that are not linearly separable
```









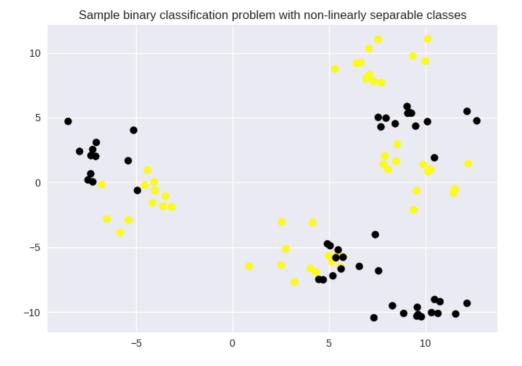
0

1

-2

-2

-1



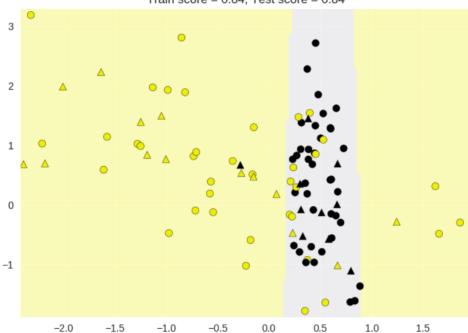
04-01: Naive Bayes classifiers

from sklearn.naive_bayes import GaussianNB # No additional models to control mod
el complexity.

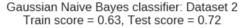
from adspy shared utilities import plot class regions for classifier

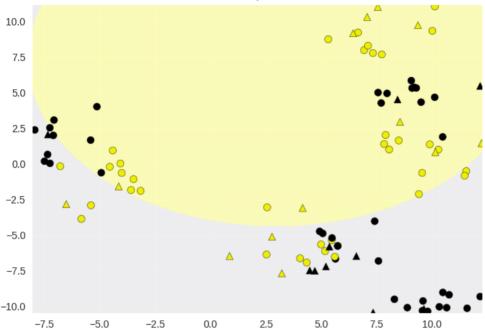
X_train, X_test, y_train, y_test = train_test_split(X_C2, y_C2, random_state=0)
Create the classifier and fit the parameters - do your usual routine
nbclf = GaussianNB().fit(X_train, y_train)
plot class regions for classifier(nbclf, X train, y train, X test, y test.

Gaussian Naive Bayes classifier: Dataset 1 Train score = 0.84, Test score = 0.84



In [4]:





Application to a real-world dataset

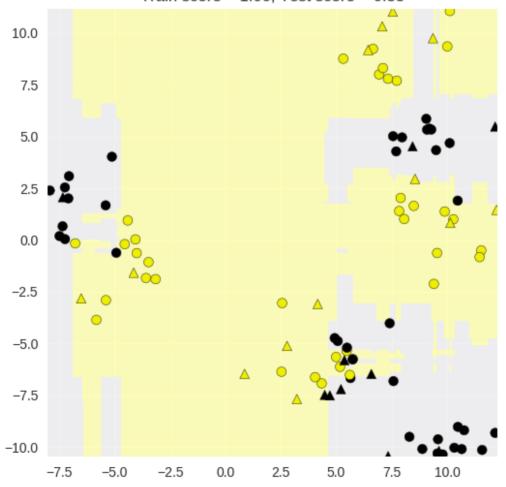
In [5]:

Breast cancer dataset
Accuracy of GaussianNB classifier on training set: 0.95
Accuracy of GaussianNB classifier on test set: 0.94

04-02: Ensembles of Decision Trees - Random Forests

In [6]:

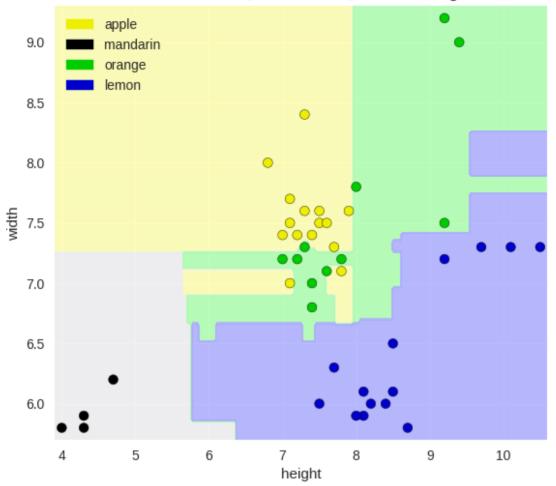
Random Forest Classifier, complex binary dataset, default settings Train score = 1.00, Test score = 0.88



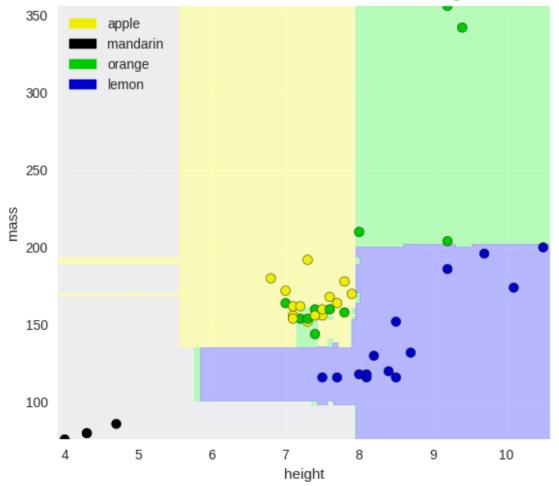
Random forest: Fruit dataset

```
from sklearn.ensemble import RandomForestClassifier
from sklearn.model selection import train test split
from adspy shared utilities import plot class regions for classifier subplot
X train, X test, y train, y test = train test split(X fruits.as matrix(),
                                                   y fruits.as matrix(),
                                                   random state = 0)
fig, subaxes = plt.subplots(6, 1, figsize=(6, 32))
title = 'Random Forest, fruits dataset, default settings'
# After creating the train-test split on the data, we iterate through each pair
of feature columns
# on the training set.
pair list = [[0,1], [0,2], [0,3], [1,2], [1,3], [2,3]]
##### ----- To plot the Random Forest on the fruit dataset ------
for pair, axis in zip(pair list, subaxes):
   X = X train[:, pair]
   y = y train
   # Create the random forest classifier and fit the model
    # For each pair of features we call the fit method on that subset of the tra
ining data X
   # using the labels y
   clf = RandomForestClassifier().fit(X, y)
   # Use the utility function to plot this graph
   plot class regions for classifier subplot(clf, X, y, None,
                                             None, title, axis,
                                             target names fruits)
   axis.set xlabel(feature names fruits[pair[0]])
   axis.set_ylabel(feature_names_fruits[pair[1]])
plt.tight layout()
plt.show()
### ------ To see the score of the classifier on default parameters --
clf = RandomForestClassifier(n estimators = 10,
                            random state=0).fit(X train, y train)
print('Random Forest, Fruit dataset, default settings')
print('Accuracy of RF classifier on training set: {:.2f}'
     .format(clf.score(X_train, y_train)))
print('Accuracy of RF classifier on test set: {:.2f}'
     .format(clf.score(X test, y test)))
```

Random Forest, fruits dataset, default settings

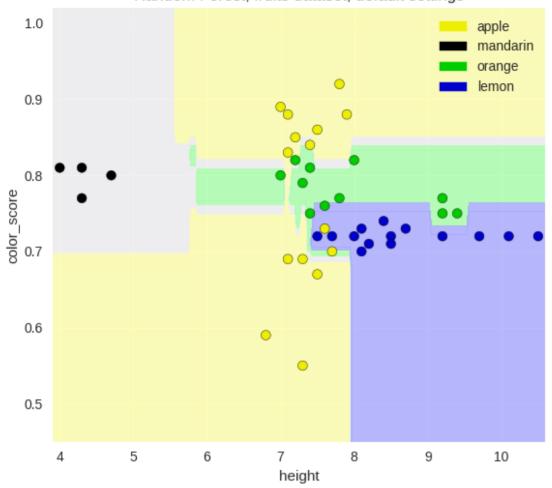


Random Forest, fruits dataset, default settings

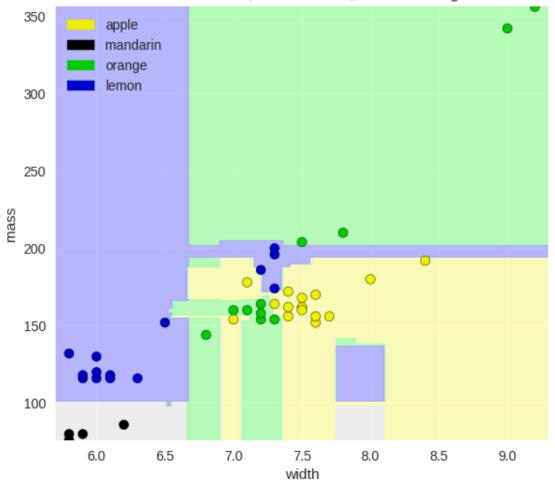


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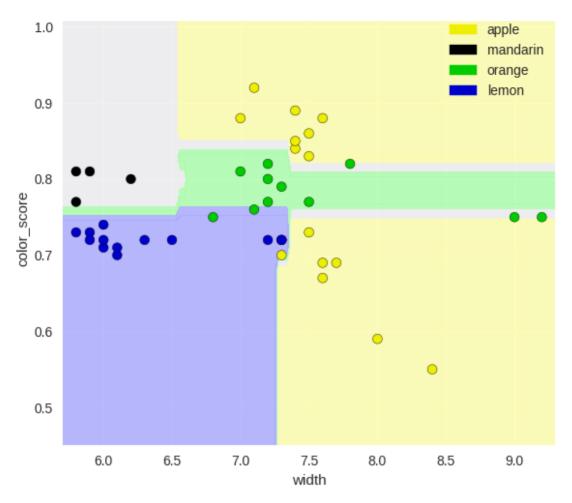
ranuom rorest, muits uataset, detauit settings

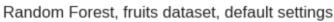


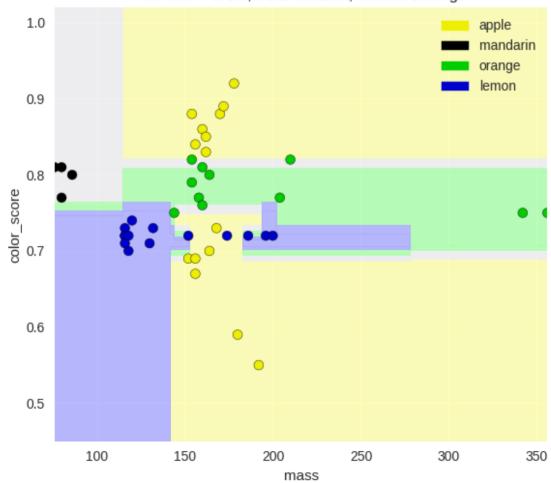
Random Forest, fruits dataset, default settings



Random Forest, fruits dataset, default settings







```
Random Forest, Fruit dataset, default settings
Accuracy of RF classifier on training set: 1.00
Accuracy of RF classifier on test set: 0.80
```

Random Forests on a real-world dataset

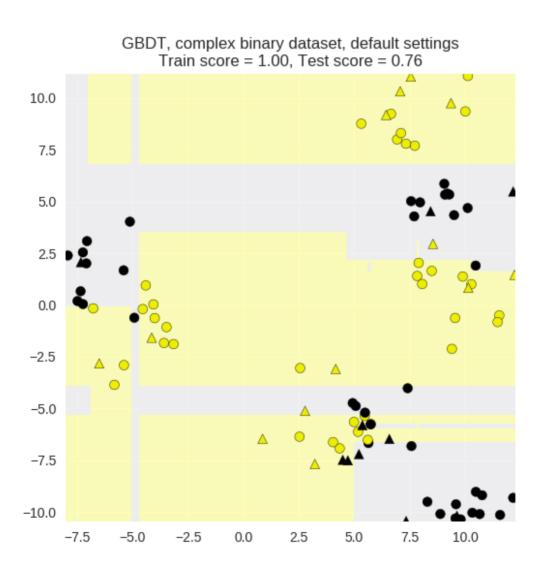
In [8]:

Breast cancer dataset
Accuracy of RF classifier on training set: 1.00
Accuracy of RF classifier on test set: 0.99

We see that this random forest - with no feature scaling or parameter tuning - achieves very good test set performance on this dataset. In fact, it is as good or better than all the other supervised methods we've seen so far including kernelised SVMs and neural networks.

Note that we **did not** have to do scaling or any preprocessing of the data. This is an advantage of using random forests. Also note that we passed in a random_state = 0 parameter in order to make the results reproducible. If we didn't set the random_state parameter, the model will likely be different each time due to the randomised nature of the random forest algorithm.

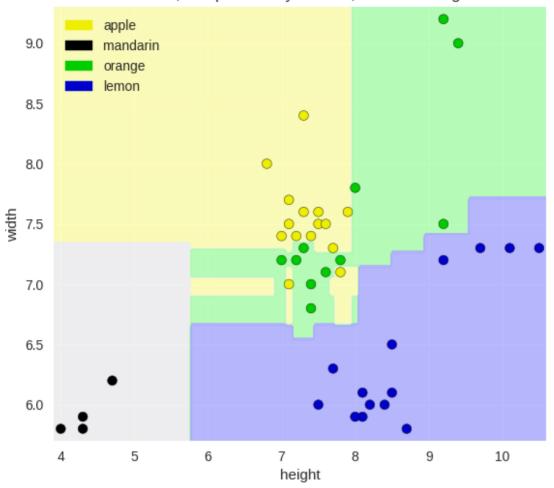
04-03: Ensemble Trees: Gradient-boosted decision trees



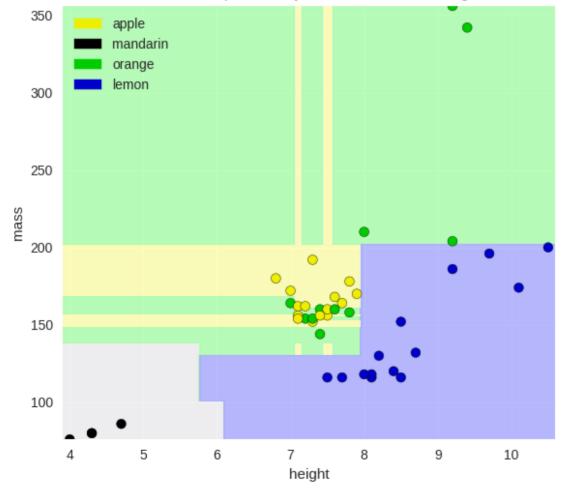
Gradient boosted decision trees on the fruit dataset

```
X_train, X_test, y_train, y_test = train_test_split(X_fruits.as_matrix(),
                                                   y_fruits.as_matrix(),
                                                    random state = 0)
fig, subaxes = plt.subplots(6, 1, figsize=(6, 32))
pair_list = [[0,1], [0,2], [0,3], [1,2], [1,3], [2,3]]
for pair, axis in zip(pair_list, subaxes):
    X = X_train[:, pair]
    y = y train
    clf = GradientBoostingClassifier().fit(X, y)
    plot_class_regions_for_classifier_subplot(clf, X, y, None,
                                             None, title, axis,
                                             target names fruits)
    axis.set xlabel(feature names fruits[pair[0]])
    axis.set ylabel(feature names fruits[pair[1]])
plt.tight layout()
plt.show()
clf = GradientBoostingClassifier().fit(X train, y train)
print('GBDT, Fruit dataset, default settings')
print('Accuracy of GBDT classifier on training set: {:.2f}'
     .format(clf.score(X_train, y_train)))
print('Accuracy of GBDT classifier on test set: {:.2f}'
     .format(clf.score(X_test, y_test)))
```

GBDT, complex binary dataset, default settings

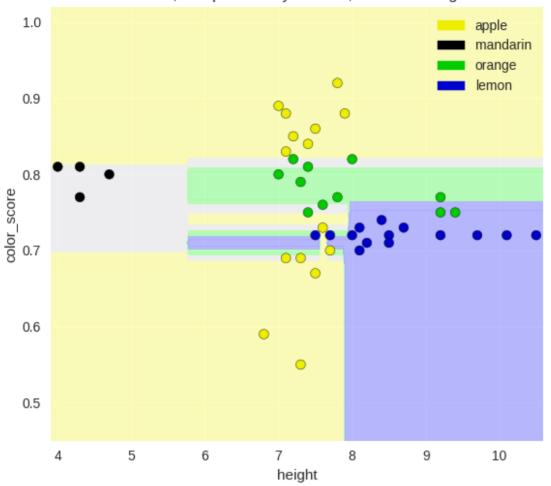


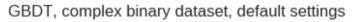
GBDT, complex binary dataset, default settings

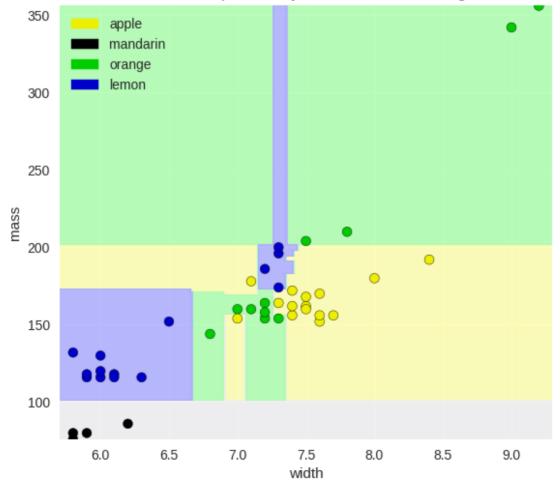


CRDT complex hinary datacet default cettings

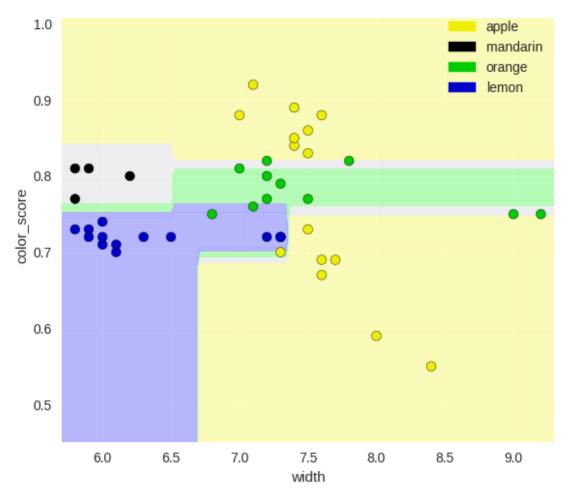
ז עסט ד, complex binary ualasel, delault sellings



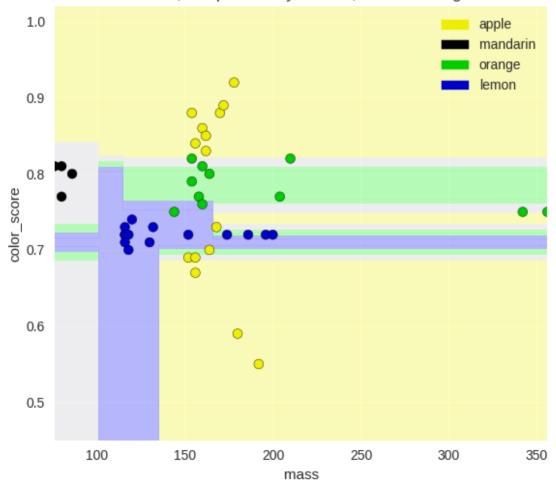




GBDT, complex binary dataset, default settings



GBDT, complex binary dataset, default settings



```
GBDT, Fruit dataset, default settings
Accuracy of GBDT classifier on training set: 1.00
Accuracy of GBDT classifier on test set: 0.80
```

Gradient-boosted decision trees on a real-world dataset

In [11]:

```
from sklearn.ensemble import GradientBoostingClassifier
X train, X test, y train, y test = train test split(X cancer, y cancer, random s
tate = 0)
clf = GradientBoostingClassifier(random state = 0)
clf.fit(X train, y train)
print('Breast cancer dataset (learning rate=0.1, max depth=3)')
print('Accuracy of GBDT classifier on training set: {:.2f}'
     .format(clf.score(X train, y train)))
print('Accuracy of GBDT classifier on test set: {:.2f}\n'
     .format(clf.score(X_test, y_test)))
clf = GradientBoostingClassifier(learning rate = 0.01, max depth = 2, random sta
te = 0)
clf.fit(X_train, y_train)
print('Breast cancer dataset (learning_rate=0.01, max_depth=2)')
print('Accuracy of GBDT classifier on training set: {:.2f}'
     .format(clf.score(X train, y train)))
print('Accuracy of GBDT classifier on test set: {:.2f}'
     .format(clf.score(X_test, y_test)))
Breast cancer dataset (learning_rate=0.1, max_depth=3)
Accuracy of GBDT classifier on training set: 1.00
Accuracy of GBDT classifier on test set: 0.96
Breast cancer dataset (learning rate=0.01, max depth=2)
Accuracy of GBDT classifier on training set: 0.97
Accuracy of GBDT classifier on test set: 0.97
```

Notice that the first run on the Breast cancer dataset has an accuracy of 1 on the training set - which implies that the model is overfitting.

- So we go about this by **reducing the learning_rate**, such that a **less** complex model is produced, and hence less likely to fix the mistakes of its predecessor.
- In addition, we can reduce the max_depth parameter for individual trees in the ensemble.

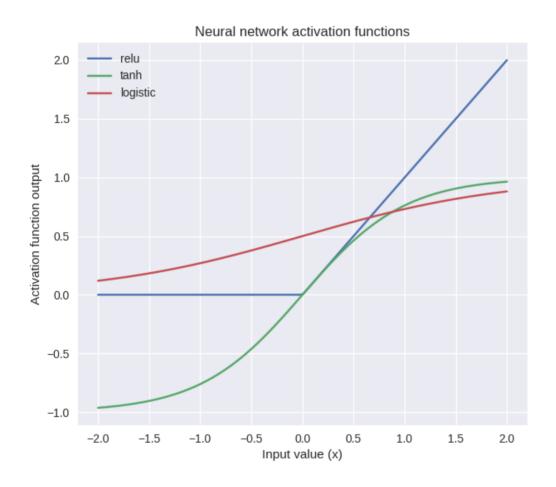
The 2nd classifier example makes these changes in the parameters, and you can see that the training set accuracy decreases slightly, with an increase in the test set accuracy.

04-04: Neural networks

Activation functions

```
xrange = np.linspace(-2, 2, 200)
plt.figure(figsize=(7,6))

plt.plot(xrange, np.maximum(xrange, 0), label = 'relu')
plt.plot(xrange, np.tanh(xrange), label = 'tanh')
plt.plot(xrange, 1 / (1 + np.exp(-xrange)), label = 'logistic')
plt.legend()
plt.title('Neural network activation functions')
plt.xlabel('Input value (x)')
plt.ylabel('Activation function output')
```

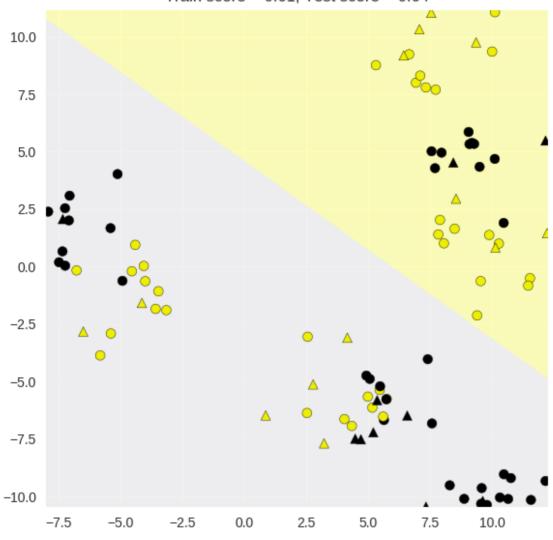


Neural networks: Classification

Synthetic dataset 1: single hidden layer

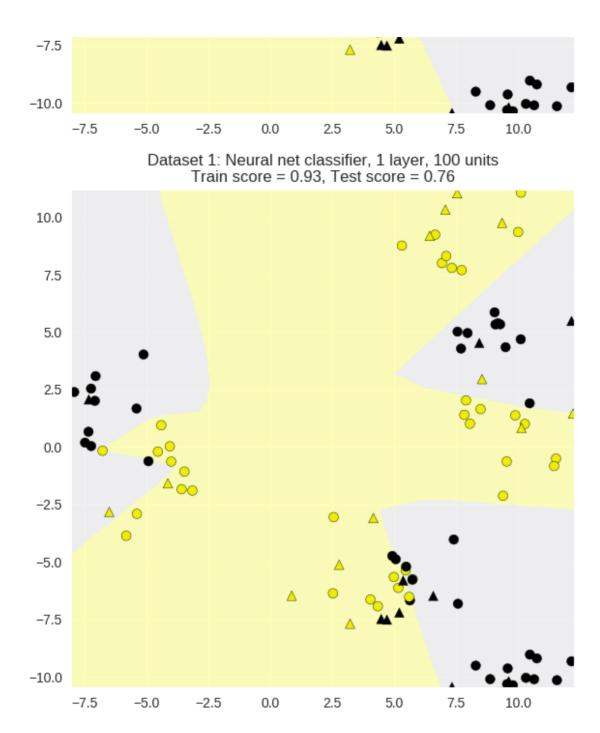
```
from sklearn.neural network import MLPClassifier
from adspy shared utilities import plot class regions for classifier subplot
X train, X test, y train, y test = train test split(X D2, y D2, random state=0)
fig, subaxes = plt.subplots(3, 1, figsize=(6,18))
for units, axis in zip([1, 10, 100], subaxes):
    # Create the classifier
    # note the parameter hidden layer sizes -> this parameter is a list, with 1
 element for each hidden layer
    # that gives the number of hidden units to use in this layer.
    # In this example, we would like 1 hidden layer, using the number in the var
iable called "units"
    # By default, if you don't specify the hidden layer sizes parameter, scikit 1
earn will create a single hidden layer
    # with 100 hidden units.
    # While a setting of 10 may work well like the one we use as examples here,
 for really complex data sets,
    # the number of hidden units could be in the thousands.
    # It's also possible to create an MLP ( in a later example) with more than 1
hidden layer, by passing in the
    # hidden_layer_sizes parameter with multiple entries.
    # Note the use of the extra parameter here, called "Solver". This specifies
 the algorithm to use for learning
    # the weights of the network. Here we use the LBFGS algorithm, we'll discuss
the solver parameter setting further,
    # at the end of the lecture.
    # Also note that we are passing in a random state parameter when creating th
e MLP Classifier object, like we did
    # for the train-test split function, and we set its value to 0.
    # This is because for neural networks, their weights are initialised randoml
y, which can affect the model that
    # is learned. Because of this, even without changing the key parameters on t
he dataset, the same neural network
    # algorithm might learn two different models, depending on the value of the
 internal random seed chosen. By
    # choosing the same random seed, we can then guarantee reproducible results.
    # OKAY, so create the classifier with appropriate params, and fit the model
 with the training data.
    nnclf = MLPClassifier(hidden layer sizes = [units], solver='lbfgs',
                         random state = 0).fit(X train, y train)
    title = 'Dataset 1: Neural net classifier, 1 layer, {} units'.format(units)
    # MPL Support
    plot_class_regions_for_classifier_subplot(nnclf, X_train, y_train,
                                             X test, y test, title, axis)
    plt.tight layout()
```

Dataset 1: Neural net classifier, 1 layer, 1 units Train score = 0.61, Test score = 0.64

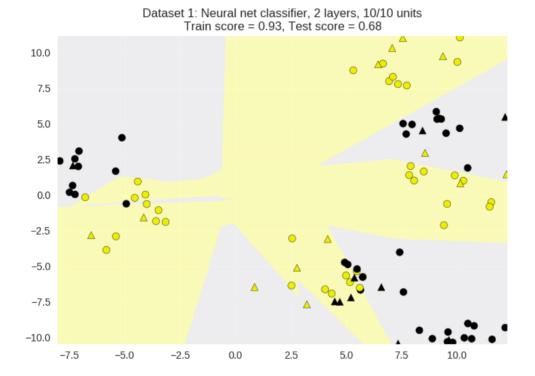


Dataset 1: Neural net classifier, 1 layer, 10 units Train score = 0.77, Test score = 0.64





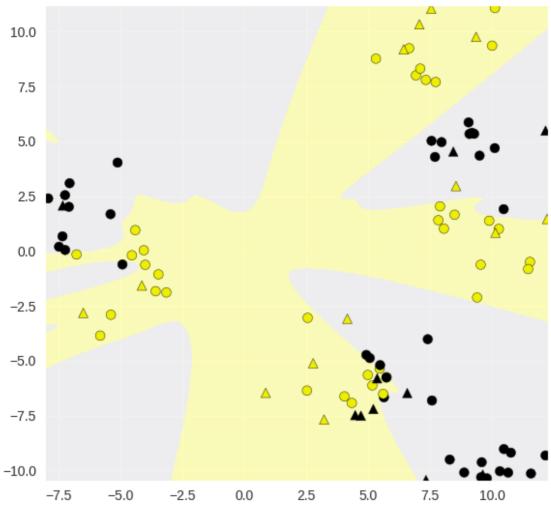
Synthetic dataset 1: two hidden layers



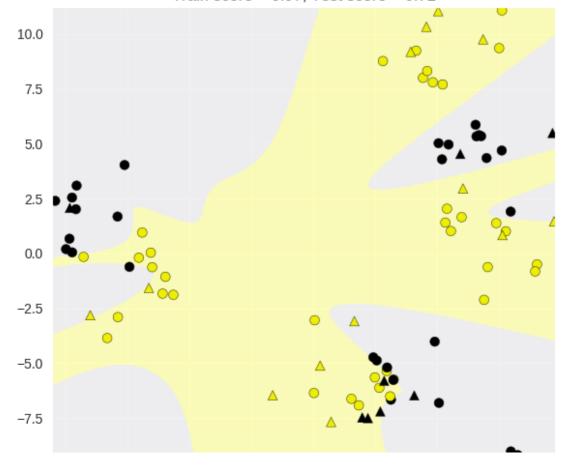
Regularization parameter: alpha

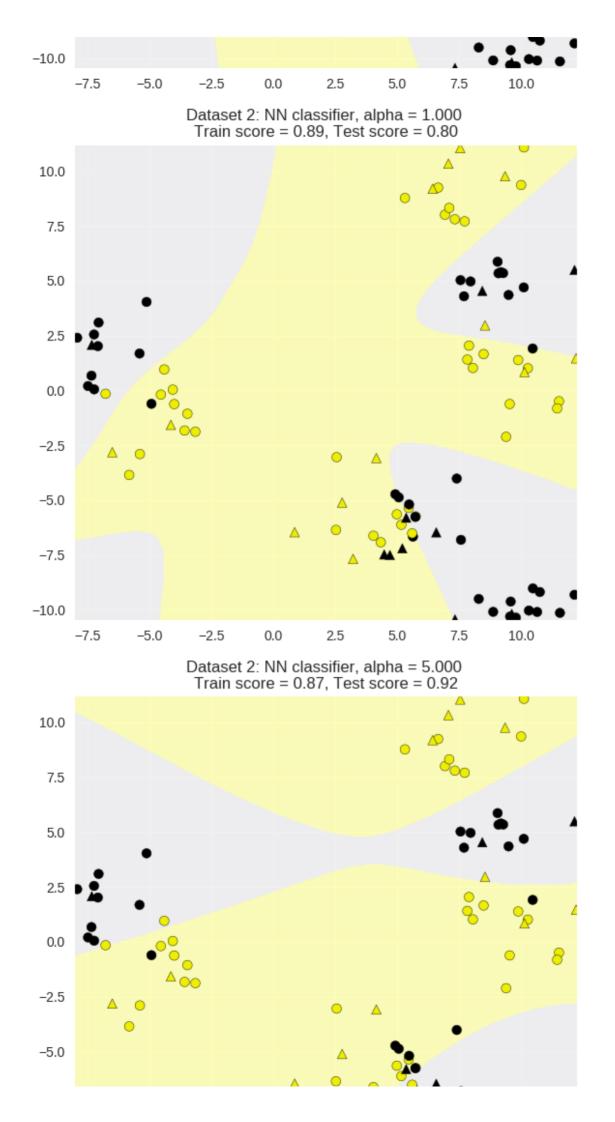
```
X_train, X_test, y_train, y_test = train_test_split(X_D2, y_D2, random_state=0)
fig, subaxes = plt.subplots(4, 1, figsize=(6, 23))
# This code shows an example of the effect of changing alpha on the neural netwo
rk.
for this alpha, axis in zip([0.01, 0.1, 1.0, 5.0], subaxes):
    nnclf = MLPClassifier(solver='lbfgs', activation = 'tanh', # For variety we
 set the activation function to the
                                                                # hyperbolic tan
gent function
                         alpha = this alpha, # by default this is set to 0.0001
                         hidden_layer_sizes = [100, 100],
                         random state = 0).fit(X train, y train)
    title = 'Dataset 2: NN classifier, alpha = {:.3f} '.format(this alpha)
    plot_class_regions_for_classifier_subplot(nnclf, X_train, y_train,
                                             X_test, y_test, title, axis)
    plt.tight layout()
```

Dataset 2: NN classifier, alpha = 0.010 Train score = 0.97, Test score = 0.72



Dataset 2: NN classifier, alpha = 0.100 Train score = 0.97, Test score = 0.72

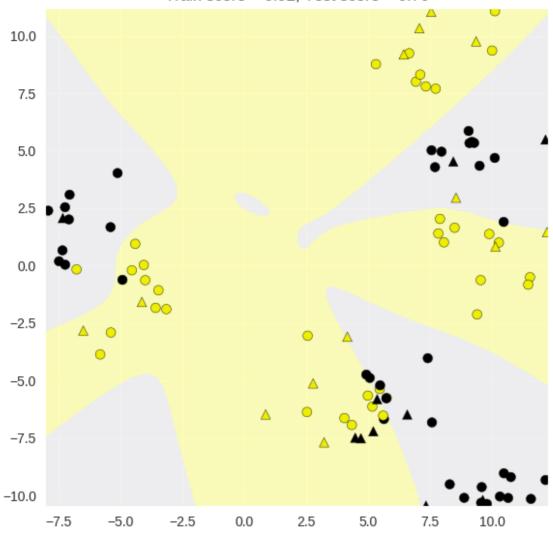




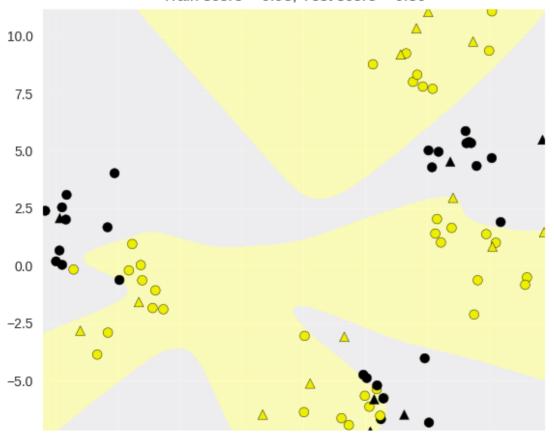


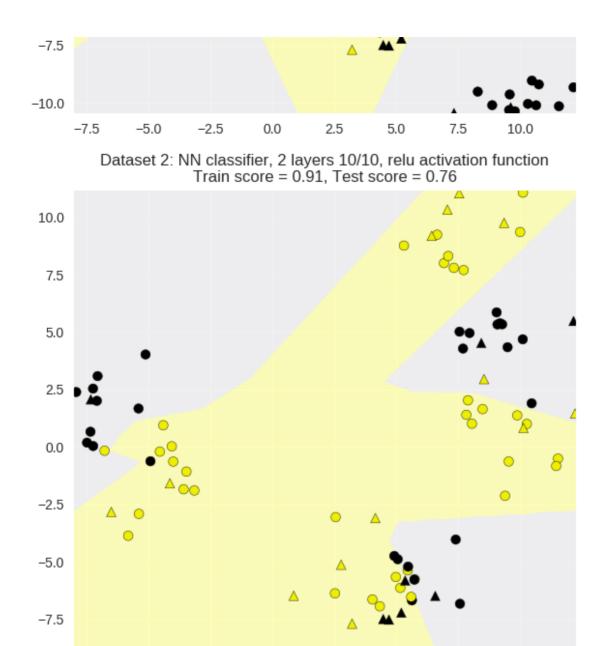
The effect of different choices of activation function

Dataset 2: NN classifier, 2 layers 10/10, logistic activation function Train score = 0.92, Test score = 0.76



Dataset 2: NN classifier, 2 layers 10/10, tanh activation function Train score = 0.93, Test score = 0.80





Neural networks: Regression

-5.0

-2.5

0.0

2.5

5.0

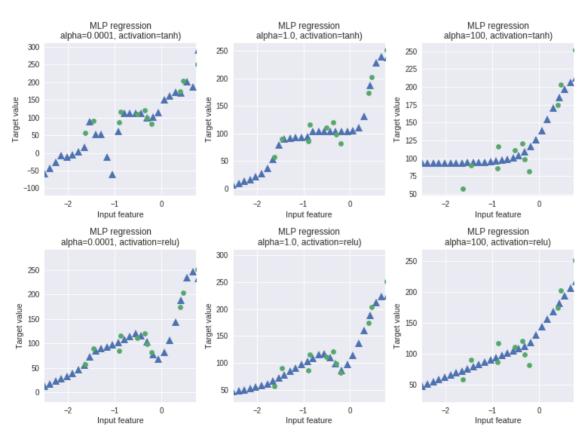
7.5

10.0

-10.0

-7.5

```
from sklearn.neural network import MLPRegressor # of course
fig, subaxes = plt.subplots(2, 3, figsize=(11,8), dpi=70)
X predict input = np.linspace(-3, 3, 50).reshape(-1,1)
X_train, X_test, y_train, y_test = train_test_split(X_R1[0::5], y_R1[0::5], rand
om state = 0)
for thisaxisrow, thisactivation in zip(subaxes, ['tanh', 'relu']):
    for thisalpha, thisaxis in zip([0.0001, 1.0, 100], thisaxisrow):
        # Create the MLP Regression object and fit the training data
        # we create this MLP with 2 hidden layers of 100 nodes
        # a varied activation function
        # and also a varied alpha variable for L2 regularisation.
        mlpreg = MLPRegressor(hidden layer sizes = [100,100],
                             activation = thisactivation,
                             alpha = thisalpha,
                             solver = 'lbfgs').fit(X_train, y_train)
        y predict output = mlpreg.predict(X predict input)
        thisaxis.set_xlim([-2.5, 0.75])
        thisaxis.plot(X_predict_input, y_predict_output,
                        , markersize = 10)
        thisaxis.plot(X_train, y_train, 'o')
        thisaxis.set xlabel('Input feature')
        thisaxis.set ylabel('Target value')
        thisaxis.set_title('MLP regression\nalpha={}, activation={})'
                          .format(thisalpha, thisactivation))
        plt.tight layout()
```



Application to real-world dataset for classification

In [16]:

Breast cancer dataset
Accuracy of NN classifier on training set: 0.98
Accuracy of NN classifier on test set: 0.97

04-05: Data Leakage

No sample code for this section.