Overfitting and Underfitting

https://scikit-learn.org/stable/auto_examples/model_selection/plot underfitting overfitting.html

Program for understanding Overfitting and Underfitting

```
import numpy as np
import matplotlib.pyplot as plt
from sklearn.pipeline import Pipeline
from sklearn.preprocessing import PolynomialFeatures
from sklearn.linear model import LinearRegression
from sklearn.model selection import cross val score
def true_fun(X):
    return np.cos(1.5 * np.pi * X)
np.random.seed(0)
n \text{ samples} = 30
degrees = [1, 4, 15]
X = np.sort(np.random.rand(n samples))
y = true fun(X) + np.random.randn(n samples) * 0.1
plt.figure(figsize=(14, 5))
for i in range(len(degrees)):
    ax = plt.subplot(1, len(degrees), i + 1)
    plt.setp(ax, xticks=(), yticks=())
    polynomial features = PolynomialFeatures(degree=degrees[i],
include bias=False)
    linear regression = LinearRegression()
    pipeline = Pipeline(
            ("polynomial_features", polynomial_features),
            ("linear regression", linear regression),
        1
    pipeline.fit(X[:, np.newaxis], y)
    # Evaluate the models using crossvalidation
    scores = cross val score(
        pipeline, X[:, np.newaxis], y,
scoring="neg_mean_squared error", cv=10
```

```
X \text{ test} = \text{np.linspace}(0, 1, 100)
    plt.plot(X test, pipeline.predict(X test[:, np.newaxis]),
label="Model")
    plt.plot(X test, true fun(X test), label="True function")
    plt.scatter(X, y, edgecolor="b", s=20, label="Samples")
    plt.xlabel("x")
    plt.ylabel("y")
    plt.xlim((0, 1))
    plt.ylim((-2, 2))
    plt.legend(loc="best")
    plt.title(
          "Degree {}\nMSE = {:.2e}(+/- {:.2e})".format(
               degrees[i], -scores.mean(), scores.std()
plt.show()
            Degree 1
                                       Degree 4
                                                                  Degree 15
      MSE = 4.08e-01(+/-4.25e-01)
                                 MSE = 4.32e-02(+/-7.08e-02)
                                                            MSE = 1.81e + 08(+/-5.42e + 08)
                                              Model
                                                                         Model
                   True function
                                              True function
                                                                         True function
                   Samples
                                              Samples
                                                                         Samples
```

Overfitting (Printing accuracy at different steps)

https://machinelearningmastery.com/overfitting-machine-learning-models/ # evaluate decision tree performance on train and test sets with different tree depths from sklearn.datasets import make classification

```
from sklearn.model_selection import train_test_split
from sklearn.metrics import accuracy_score
from sklearn.tree import DecisionTreeClassifier
from matplotlib import pyplot

# define dataset
X, y = make_classification(n_samples=10000, n_features=20,
n_informative=5, n_redundant=15, random_state=1)
# summarize the dataset
print(X.shape, y.shape)

(10000, 20) (10000,)
```

```
# split into train test sets
X train, X test, y train, y test = train test split(X, y,
test size=0.3)
# summarize the shape of the train and test sets
print(X train.shape, X test.shape, y train.shape, y test.shape)
(7000, 20) (3000, 20) (7000,) (3000,)
train_scores, test_scores = list(), list()
# define the tree depths to evaluate
values = [i for i in range(1, 31)]
# evaluate a decision tree for each depth
for i in values:
    # configure the model
    model = DecisionTreeClassifier(max depth=i)
    # fit model on the training dataset
    model.fit(X train, y_train)
    # evaluate on the train dataset
    train yhat = model.predict(X train)
    train_acc = accuracy_score(y_train, train_yhat)
    train scores.append(train_acc)
    # evaluate on the test dataset
    test_yhat = model.predict(X_test)
    test acc = accuracy score(y test, test yhat)
    test scores.append(test acc)
    # summarize progress
    print('>%d, train: %.3f, test: %.3f' % (i, train acc, test acc))
>1, train: 0.763, test: 0.767
>2, train: 0.804, test: 0.805
>3, train: 0.871, test: 0.868
>4, train: 0.906, test: 0.890
>5, train: 0.924, test: 0.901
>6, train: 0.937, test: 0.912
>7, train: 0.947, test: 0.917
>8, train: 0.956, test: 0.914
>9, train: 0.966, test: 0.917
>10, train: 0.975, test: 0.911
>11, train: 0.981, test: 0.913
>12, train: 0.985, test: 0.909
>13, train: 0.990, test: 0.909
>14, train: 0.993, test: 0.907
>15, train: 0.995, test: 0.905
>16, train: 0.996, test: 0.910
>17, train: 0.997, test: 0.908
>18, train: 0.998, test: 0.904
>19, train: 0.999, test: 0.905
>20, train: 0.999, test: 0.903
>21, train: 1.000, test: 0.902
```

```
>22, train: 1.000, test: 0.905
>23, train: 1.000, test: 0.903
>24, train: 1.000, test: 0.901
>25, train: 1.000, test: 0.901
>26, train: 1.000, test: 0.906
>27, train: 1.000, test: 0.905
>28, train: 1.000, test: 0.900
>29, train: 1.000, test: 0.900
>30, train: 1.000, test: 0.908
```

Cross-validation

```
https://scikit-learn.org/stable/modules/cross_validation.html
import numpy as np
from sklearn.model selection import train test split
from sklearn import datasets
from sklearn import svm
X, y = datasets.load iris(return X y=True)
X.shape, y.shape
((150, 4), (150,))
Basic method to compute score
X train, X test, y train, y test = train test split(
    X, y, test size=0.4, random state=0)
X train.shape, y train.shape
X_test.shape, y_test.shape
clf = svm.SVC(kernel='linear', C=1).fit(X train, y train)
clf.score(X_test, y_test)
0.96666666666666
Estimate the accuracy by splitting the data, computing the score 5 consecutive times (with
different splits each time)
from sklearn.model selection import cross val score
clf = svm.SVC(kernel='linear', C=1, random state=42)
scores = cross val score(clf, X, y, cv=5)
scores
```

```
array([0.96666667, 1. , 0.96666667, 0.96666667, 1.
                                                                   1)
print("%0.2f accuracy with a standard deviation of %0.2f" %
(scores.mean(), scores.std()))
0.98 accuracy with a standard deviation of 0.02
Using the different scoring parameter
from sklearn import metrics
scores = cross val score(
    clf, X, y, cv=5, scoring='f1 macro')
scores
array([0.96658312, 1.
                             , 0.96658312, 0.96658312, 1.
                                                              1)
Specified multiple metrics of predefined scorer names
from sklearn.model selection import cross validate
from sklearn.metrics import recall score
scoring = ['precision_macro', 'recall_macro']
clf = svm.SVC(kernel='linear', C=1, random_state=0)
scores = cross validate(clf, X, y, scoring=scoring)
sorted(scores.keys())
scores['test recall macro']
array([0.96666667, 1. , 0.96666667, 0.96666667, 1.
                                                                   ])
Calculate cross validation score by passing a cross validation iterator
from sklearn.model selection import ShuffleSplit
n \text{ samples} = X.shape[0]
cv = ShuffleSplit(n splits=5, test size=0.3, random state=0)
cross val score(clf, X, y, cv=cv)
array([0.97777778, 0.97777778, 1. , 0.95555556, 1.
                                                                   1)
Use an iterable yielding (train, test) splits as arrays of indices
def custom cv 2folds(X):
    n = X.shape[0]
    i = 1
    while i <= 2:
        idx = np.arange(n * (i - 1) / 2, n * i / 2, dtype=int)
        vield idx, idx
        i += 1
custom cv = custom cv 2 folds(X)
cross_val_score(clf, X, y, cv=custom_cv)
array([1. , 0.97333333])
```

Different type of Cross validation iterators

```
K-fold
import numpy as np
from sklearn.model selection import KFold
X = ["a", "b", "c", "d"]
kf = KFold(n splits=2)
for train, test in kf.split(X):
    print("%s %s" % (train, test))
[2 3] [0 1]
[0 1] [2 3]
Repeated K-Fold
import numpy as np
from sklearn.model_selection import RepeatedKFold
X = np.array([[1, \overline{2}], [3, 4], [1, 2], [3, 4]])
random state = 12883823
rkf = RepeatedKFold(n splits=2, n repeats=2,
random state=random state)
for train, test in rkf.split(X):
    print("%s %s" % (train, test))
[2 3] [0 1]
[0 1] [2 3]
[0 2] [1 3]
[1 3] [0 2]
Leave One Out (LOO)
from sklearn.model selection import LeaveOneOut
X = [1, 2, 3, 4]
loo = LeaveOneOut()
for train, test in loo.split(X):
    print("%s %s" % (train, test))
[1 2 3] [0]
[0 2 3] [1]
[0 1 3] [2]
[0 1 2] [3]
Leave P Out (LPO)
from sklearn.model selection import LeavePOut
X = np.ones(4)
lpo = LeavePOut(p=2)
for train, test in lpo.split(X):
     print("%s %s" % (train, test))
```

```
[2 3] [0 1]
[1 3] [0 2]
[1 2] [0 3]
[0 3] [1 2]
[0 2] [1 3]
[0 1] [2 3]
Stratified k-fold
from sklearn.model selection import StratifiedKFold, KFold
import numpy as np
X, y = np.ones((50, 1)), np.hstack(([0] * 45, [1] * 5))
skf = StratifiedKFold(n splits=3)
for train, test in skf.split(X, y):
    print('train - {} | test - {}'.format(
         np.bincount(y[train]), np.bincount(y[test])))
train -
         [30
              31
                       test - [15
                                     21
                       test - [15 2]
train - [30
              31
train - [30 4]
                       test - [15 1]
kf = KFold(n splits=3)
for train, test in kf.split(X, y):
     print('train - {} | test - {}'.format(
                  np.bincount(y[train]), np.bincount(y[test])))
train -
         [28 5]
                       test - [17]
         [28 5]
train -
                       test - [17]
         [34] | test - [11 5]
train -
GroupShuffleSplit
from sklearn.model selection import GroupShuffleSplit
X = [0.1, 0.2, 2.2, 2.4, 2.3, 4.55, 5.8, 0.001]

y = ["a", "b", "b", "c", "c", "c", "a"]
groups = [1, 1, 2, 2, 3, 3, 4, 4]
gss = GroupShuffleSplit(n splits=4, test size=0.5, random state=0)
for train, test in gss.split(X, y, groups=groups):
    print("%s %s" % (train, test))
[0 1 2 3] [4 5 6 7]
[2 3 6 7] [0 1 4 5]
[2 3 4 5] [0 1 6 7]
[4 5 6 7] [0 1 2 3]
```

Validation curve

https://keeeto.github.io/blog/bias_variance/

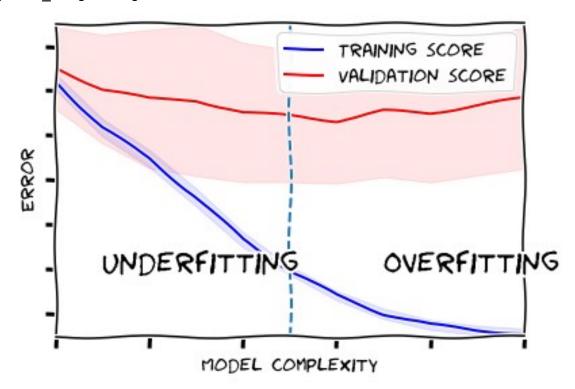
```
https://scikit-learn.org/stable/modules/learning_curve.html
import pandas as pd
import numpy as np
import matplotlib.pyplot as plt
from sklearn.ensemble import RandomForestClassifier
from sklearn.model selection import GridSearchCV
from sklearn.model selection import cross val score, learning curve,
validation curve
df train = pd.read csv('train.csv')
df_test = pd.read_csv('test.csv')
df comb = df train.append(df test)
X = pd.DataFrame()
def encode sex(x):
    return 1 if x == 'female' else 0
def family size(x):
    size = x.SibSp + x.Parch
    return 4 if size > 3 else size
X['Sex'] = df comb.Sex.map(encode sex)
X['Pclass'] = df comb.Pclass
X['FamilySize'] = df comb.apply(family size, axis=1)
fare median = df train.groupby(['Sex', 'Pclass']).Fare.median()
fare median.name = 'FareMedian'
age mean = df train.groupby(['Sex', 'Pclass']).Age.mean()
age mean.name = 'AgeMean'
def join(df, stat):
    return pd.merge(df, stat.to frame(), left on=['Sex', 'Pclass'],
right index=True, how='left')
```

```
X['Fare'] = df comb.Fare.fillna(join(df comb, fare median).FareMedian)
X['Age'] = df comb.Age.fillna(join(df comb, age mean).AgeMean)
def quantiles(series, num):
    return pd.qcut(series, num, retbins=True)[1]
def discretize(series, bins):
    return pd.cut(series, bins, labels=range(len(bins)-1),
include lowest=True)
X['Fare'] = discretize(X.Fare, quantiles(df comb.Fare, 10))
X['Age'] = discretize(X.Age, quantiles(df comb.Age, 10))
X train = X.iloc[:df train.shape[0]]
X test = X.iloc[df train.shape[0]:]
y_train = df_train.Survived
clf 1 = RandomForestClassifier(n estimators=100, bootstrap=True,
random state=0)
clf 1.fit(X train, y train)
# Number of folds for cross validation
num folds = 7
def plot_curve(ticks, train_scores, test_scores):
    train scores mean = -1 * np.mean(train scores, axis=1)
    train scores std = -1 * np.std(train scores, axis=1)
    test scores mean = -1 * np.mean(test scores, axis=1)
    test scores std = -1 * np.std(test scores, axis=1)
    plt.figure()
    plt.fill between(ticks,
                     train scores mean - train scores std,
                     train scores mean + train scores std, alpha=0.1,
color="b")
    plt.fill between(ticks,
                     test scores mean - test scores std,
                     test scores mean + test scores std, alpha=0.1,
color="r")
    plt.plot(ticks, train_scores_mean, 'b-', label='Training score')
    plt.plot(ticks, test_scores_mean, 'r-', label='Validation score')
    plt.legend(fancybox=True, facecolor='w')
    return plt.gca()
def plot validation curve(clf, X, y, param name, param range,
scoring='roc auc'):
    ax = plot curve(param range, *validation curve(clf, X, y,
cv=num folds,
                                                   scoring=scoring,
```

```
param_name=param_name,

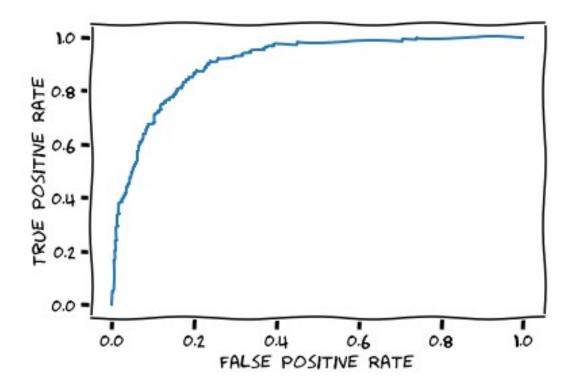
param_range=param_range, n_jobs=-1))
   ax.set_title('')
   ax.set_xticklabels([])
   ax.set_yticklabels([])
   ax.set_xlim(2,12)
   ax.set_ylim(-0.97, -0.83)
   ax.set_ylabel('Error')
   ax.set_xlabel('Model complexity')
   ax.text(9, -0.94, 'Overfitting', fontsize=22)
   ax.text(3, -0.94, 'Underfitting', fontsize=22)
   ax.axvline(7, ls='--')
   plt.tight layout()
```

plot_validation_curve(clf_1, X_train, y_train, param_name='max_depth',
param_range=range(2,13))



ROC

```
https://www.statology.org/plot-roc-curve-python/
import pandas as pd
import numpy as np
from sklearn.model selection import train test split
from sklearn.linear model import LogisticRegression
from sklearn import metrics
import matplotlib.pyplot as plt
#import dataset from CSV file on Github
url =
"https://raw.githubusercontent.com/Statology/Python-Guides/main/defaul
t.csv"
data = pd.read csv(url)
#define the predictor variables and the response variable
X = data[['student', 'balance', 'income']]
y = data['default']
#split the dataset into training (70%) and testing (30%) sets
X train,X test,y train,y test =
train test split(X,y,test size=0.3,random state=0)
#instantiate the model
log regression = LogisticRegression()
#fit the model using the training data
log_regression.fit(X_train,y_train)
LogisticRegression()
#define metrics
y pred proba = log regression.predict proba(X test)[::,1]
fpr, tpr, _ = metrics.roc_curve(y_test, y_pred_proba)
#create ROC curve
plt.plot(fpr,tpr)
plt.ylabel('True Positive Rate')
plt.xlabel('False Positive Rate')
plt.show()
```



Task 1: Perform all of the above codes of Overfitting, Cross Validation, etc. with the help of the given reference link.

Task 2: Explain your analysis of the code. Make a detailed analysis that can also cover the following questions: (Submit the PDF of Report)

- 1) According to you, why do overfitting and underfitting occur, and how resolve them? What is the difference between them?
- 2) What kind of pattern did you analyze in the Train and Test score while running the code of overfitting?
- 3) What is cross-validation, and what did you analyze in a different type of validation that you performed?
- 4) Explain the analysis from generated ROC and validation curve and what they represent?

Task 3: Using the given Cross Validation iterators perform all types of Cross Validations we did in the task:

- 1) K-fold
- 2) Repeated K-Fold
- 3) Leave One Out (LOO)

Apart from this three, try to perform validation using three new iterators.

Task 4: With the help of the given code and references complete all of the following step:

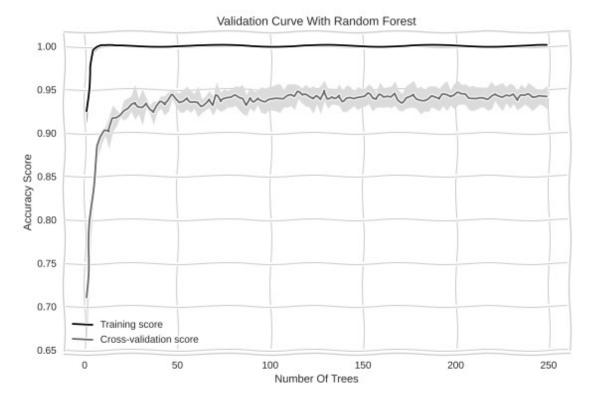
- 1) Choose one new dataset. Train a overfitted model with the help of any machine learning technique, such as KNN, classification, regression.
- 2) Try to resolve the overfitting.
- 3) Calculate the Validation score by any two or three given techniques and Validation iterators.
- 4) Generate the validation curve
- 5) Predict the output of testing data.

```
6) Generate the ROC curve using the predicted data and actual data.
import numpy as np
import matplotlib.pyplot as plt
from sklearn.pipeline import Pipeline
from sklearn.preprocessing import PolynomialFeatures
from sklearn.linear model import LinearRegression
from sklearn.model selection import cross val score
from sklearn.datasets import load wine
X, y = make classification(n samples=5000, n features=20,
n informative=15)
from sklearn.model selection import train test split
X_train, X_test, y_train, y_test = train_test_split(
    X, y, test size=0.2, shuffle=True)
print(len(X),len(X train),len(X test))
5000 4000 1000
from sklearn.neighbors import KNeighborsClassifier
# Create KNN classifier
knn = KNeighborsClassifier(n neighbors = 3)
# Fit the classifier to the data
knn.fit(X train,y train)
KNeighborsClassifier(n neighbors=3)
#show first 5 model predictions on the test data
knn.predict(X test)[0:5]
array([0, 1, 1, 1, 1])
knn.score(X_test, y_test)
0.934
from sklearn.model selection import ShuffleSplit
n \text{ samples} = X.shape[0]
cv = ShuffleSplit(n splits=5, test size=0.3, random state=0)
cross_val_score(clf, X, y, cv=cv)
array([0.74066667, 0.73266667, 0.726 , 0.74866667, 0.72
                                                                   ])
def custom cv 2folds(X):
    n = X.shape[0]
    i = 1
    while i <= 2:
        idx = np.arange(n * (i - 1) / 2, n * i / 2, dtype=int)
        yield idx, idx
        i += 1
```

```
custom_cv = custom cv 2folds(X)
cross_val_score(clf, X, y, cv=custom_cv)
array([0.748 , 0.7364])
import numpy as np
from sklearn.model selection import KFold
X = ["a", "b", "c", "d"]
kf = KFold(n splits=2)
for train, test in kf.split(X):
     print("%s %s" % (train, test))
[2 3] [0 1]
[0 1] [2 3]
import numpy as np
from sklearn.model selection import RepeatedKFold
X = np.array([[1, 2], [3, 4], [1, 2], [3, 4]])
random state = 12883823
rkf = RepeatedKFold(n_splits=2, n_repeats=2,
random state=random state)
for train, test in rkf.split(X):
    print("%s %s" % (train, test))
[2 3] [0 1]
[0 1] [2 3]
[0 2] [1 3]
[1 3] [0 2]
import matplotlib.pyplot as plt
import numpy as np
from sklearn.datasets import load wine
from sklearn.ensemble import RandomForestClassifier
from sklearn.model selection import validation curve
wine = load wine()
# Create feature matrix and target vector
X, y = wine.data, wine.target
train scores, test scores = list(), list()
# define the tree depths to evaluate
values = [i for i in range(1, 31)]
for i in values:
    # configure the model
    model = RandomForestClassifier(max depth=i)
    # fit model on the training dataset
    model.fit(X train, y train)
```

```
# evaluate on the train dataset
    train yhat = model.predict(X train)
    train_acc = accuracy_score(y_train, train_yhat)
    train scores.append(train acc)
    # evaluate on the test dataset
    test yhat = model.predict(X test)
    test acc = accuracy score(y test, test yhat)
    test scores.append(test acc)
    # summarize progress
    print('>%d, train: %.3f, test: %.3f' % (i, train acc, test acc))
    plt.show()
>1, train: 0.903, test: 0.893
>2, train: 0.922, test: 0.909
>3, train: 0.984, test: 0.960
>4, train: 0.996, test: 0.980
>5, train: 0.998, test: 0.987
>6, train: 0.999, test: 0.993
>7, train: 1.000, test: 0.993
>8, train: 1.000, test: 0.993
>9, train: 1.000, test: 0.996
>10, train: 1.000, test: 0.996
>11, train: 1.000, test: 0.993
>12, train: 1.000, test: 0.993
>13, train: 1.000, test: 0.996
>14, train: 1.000, test: 0.991
>15, train: 1.000, test: 0.991
>16, train: 1.000, test: 0.993
>17, train: 1.000, test: 0.991
>18, train: 1.000, test: 0.991
>19, train: 1.000, test: 0.993
>20, train: 1.000, test: 0.987
>21, train: 1.000, test: 0.993
>22, train: 1.000, test: 0.989
>23, train: 1.000, test: 0.996
>24, train: 1.000, test: 0.993
>25, train: 1.000, test: 0.996
>26, train: 1.000, test: 0.996
>27, train: 1.000, test: 0.993
>28, train: 1.000, test: 0.996
>29, train: 1.000, test: 0.993
>30, train: 1.000, test: 0.991
param range = np.arange(1, 250, 2)
# Calculate accuracy on training and test set using range of parameter
values
train scores, test scores = validation curve(RandomForestClassifier(),
```

```
param name="n estimators",
                                             param range=param range,
                                             cv=3,
                                             scoring="accuracy",
                                             n jobs=-1
# Calculate mean and standard deviation for training set scores
train mean = np.mean(train scores, axis=1)
train std = np.std(train scores, axis=1)
# Calculate mean and standard deviation for test set scores
test mean = np.mean(test scores, axis=1)
test std = np.std(test_scores, axis=1)
# Plot mean accuracy scores for training and test sets
plt.plot(param range, train mean, label="Training score",
color="black")
plt.plot(param range, test mean, label="Cross-validation score",
color="dimgrey")
# Plot accurancy bands for training and test sets
plt.fill between(param range, train mean - train std, train mean +
train std, color="gray")
plt.fill between(param range, test mean - test std, test mean +
test std, color="gainsboro")
# Create plot
plt.title("Validation Curve With Random Forest")
plt.xlabel("Number Of Trees")
plt.ylabel("Accuracy Score")
plt.tight layout()
plt.legend(loc="best")
plt.show()
```



from sklearn.datasets import load_wine
wine = load wine()

```
from sklearn.naive_bayes import GaussianNB
from sklearn.model_selection import train_test_split
X_train, X_test, y_train, y_test = train_test_split(wine.data,
wine.target)
clf = GaussianNB()
clf.fit(X train, y train)
```

GaussianNB()

5 5

predicted = clf.predict(X_test)
expected = y_test
print(predicted)

[5 7 1 5 7 7 7 7 4 7 4 2 0 7 2 7 0 4 3 5 7 7 7 9 9 3 3 3 5 8 5 4 0 5 7 5 9 0 8 7 2 2 1 6 1 4 6 7 9 6 2 7 1 6 8 7 7 7 7 7 3 8 2 8 1 8 4 2 0 5 7 7 3 1 8 1 7 2 7 2 7 2 8 7 7 7 4 6 0 5 8 9 8 7 6 0 6 0 5 6 4 3 5 6 5 0 6 3 7 1 1 3 4 7 6 6 9 8 4 6 6 4 3 2 7 6 6 6 6 3 8 7 8 6 7 4 2 2 1 6 8 1 8 7 5 3 8 2 8 8 4 2 1 2 7 8 0 6 9 9 8 0 5 0 1 3 7 6 7

```
7 7 6 0 6 2 2 1 7 4 9 7 9 1 8 8 7 6 8 2 1 7 8 2 2 6 8 2 3 7 9 4 1 4 3
9 6
8 8 9 0 6 5 9 9 6 6 7 5 6 5 2 8 7 8 7 8 3 7 5 7 8 0 9 4 1 0 3 6 6 0 2
 8 6 0 7 4 9 1 7 7 2 4 8 9 0 8 6 9 2 1 9 2 1 0 2 0 4 5 9 3 7 3 6 3 8 1
 1 0 2 8 5 4 6 6 2 1 9 6 1 4 8 3 9 4 5 5 2 4 7 3 3 6 1 2 9 0 9 5 5 0 1
5 8 4 9 2 8 2 0 5 4 0 7 4 7 3 2 8 0 6 7 6 5 4 8 5 7 6 0 5 2 3 2 1 7 2
5 7
 1 7 6 4 4 7 0 1 0 1 8 6 5 5 5 5 5 2 6 7 6 3 2 7 7 3 9 5 4 1 6 8 3 2 2 0
7 2
8 1 7 4 8 9 2 0 3 1 3 4 8 5 4 8 3 8 1 8 5 7 9 6 3 9 9 9 8 1 3 3 5 0 2
0 4
4 0 8 7 1 3]
from sklearn import datasets
from sklearn.linear model import LogisticRegression
from sklearn.metrics import roc curve, roc auc score
from sklearn.model selection import train test split
import matplotlib.pyplot as plt
dataset = datasets.load wine()
X = dataset.data
y = dataset.target
from sklearn import datasets, ensemble, metrics, model selection,
dummy
import matplotlib.pyplot as plt
wine = datasets.load wine()
# First we work only on sevens:
sevens = (wine.target == 7)
classifier = ensemble.RandomForestClassifier()
most frequent = dummy.DummyClassifier(strategy='most frequent')
X_train, X_test, y_train, y_test = model_selection.train_test_split(
    wine.data, sevens, random_state=0)
y score = classifier.fit(X train, y train).predict proba(X test)
most frequent score = most frequent.fit(X train,
                                        y train).predict proba(X test)
fpr, tpr, = metrics.roc curve(y test, y score[:, 1])
roc auc = metrics.auc(fpr, tpr)
fpr_dummy, tpr_dummy, _ = metrics.roc_curve(y_test,
```

```
most frequent score[:, 1])
roc auc dummy = metrics.auc(fpr dummy, tpr dummy)
plt.figure(figsize=(3, 3))
lw = 2
plt.plot(fpr, tpr, color='darkorange',
         lw=lw, label='RandomForestClassifier\n(area = %0.2f)' %
roc auc)
plt.plot(fpr dummy, tpr dummy, color='.5',
         lw=lw, label='Dummy\n(area = %0.2f)' % roc auc dummy)
plt.xlim([0.0, 1.0])
plt.ylim([0.0, 1.05])
plt.xlabel('False Positive Rate')
plt.ylabel('True Positive Rate')
plt.title('Receiver operating characteristic')
plt.legend(loc="lower left", title='ROC curve')
plt.tight_layout()
plt.show()
```

RECEIVER OPERATING CHARACTERISTIC

