Random Forest - Classification

The Data

We will be using the same dataset through our discussions on classification with tree-methods (Decision Tree,Random Forests, and Gradient Boosted Trees) in order to compare performance metrics across these related models.

We will work with the "Palmer Penguins" dataset, as it is simple enough to help us fully understand how changing hyperparameters can change classification results.

Data were collected and made available by Dr. Kristen Gorman and the Palmer Station, Antarctica LTER, a member of the Long Term Ecological Research Network.

Gorman KB, Williams TD, Fraser WR (2014) Ecological Sexual Dimorphism and Environmental Variability within a Community of Antarctic Penguins (Genus Pygoscelis). PLoS ONE 9(3): e90081. doi:10.1371/journal.pone.0090081

Summary: The data folder contains two CSV files. For intro courses/examples, you probably want to use the first one (penguins_size.csv).

- penguins size.csv: Simplified data from original penguin data sets. Contains variables:
 - species: penguin species (Chinstrap, Adélie, or Gentoo)
 - culmen_length_mm: culmen length (mm)
 - culmen_depth_mm: culmen depth (mm)
 - flipper_length_mm: flipper length (mm)
 - body_mass_g: body mass (g)
 - island: island name (Dream, Torgersen, or Biscoe) in the Palmer Archipelago (Antarctica)
 - sex: penguin sex
- (Not used) penguins_lter.csv: Original combined data for 3 penguin species

Note: The culmen is "the upper ridge of a bird's beak"

Our goal is to create a model that can help predict a species of a penguin based on physical attributes, then we can use that model to help researchers classify penguins in the field, instead of needing an experienced biologist

Imports

```
In [1]:
```

```
import numpy as np
import pandas as pd
import matplotlib.pyplot as plt
import seaborn as sns
```

```
In [2]:
```

df=pd.read_csv("D:\\Study\\Programming\\python\\Python course from udemy\\Udemy - 2022 P
ython for Machine Learning & Data Science Masterclass\\01 - Introduction to Course\\1UNZI
P-FOR-NOTEBOOKS-FINAL\\DATA\\penguins_size.csv")
df.head()

Out[2]:

	species	island	culmen_length_mm	culmen_depth_mm	flipper_length_mm	body_mass_g	sex
0	Adelie	Torgersen	39.1	18.7	181.0	3750.0	MALE
1	Adelie	Torgersen	39.5	17.4	186.0	3800.0	FEMALE

```
Torgersen 40.3 18.0 195.0 3250.0 island culmen_length_mm culmen_depth_mm flipper_length_mm body_mass_g
                                                                                           3250.0 FEMALE
 Adelie
species
Adelie
         Torgersen
                                     NaN
                                                         NaN
                                                                             NaN
                                                                                             NaN
                                                                                                        NaN
                                     36.7
                                                          19.3
                                                                             193.0
                                                                                           3450.0 FEMALE
 Adelie Torgersen
```

```
In [3]:
```

```
# Here we are droping all null values
df = df.dropna()
df.head()
```

Out[3]:

	species	island	culmen_length_mm	culmen_depth_mm	flipper_length_mm	body_mass_g	sex
0	Adelie	Torgersen	39.1	18.7	181.0	3750.0	MALE
1	Adelie	Torgersen	39.5	17.4	186.0	3800.0	FEMALE
2	Adelie	Torgersen	40.3	18.0	195.0	3250.0	FEMALE
4	Adelie	Torgersen	36.7	19.3	193.0	3450.0	FEMALE
5	Adelie	Torgersen	39.3	20.6	190.0	3650.0	MALE

Train | Test Split

```
In [4]:
```

```
X=pd.get_dummies(df.drop('species',axis=1),drop_first=True)
```

```
In [5]:
```

```
y=df['species']
```

In [6]:

```
from sklearn.model_selection import train_test_split
```

In [7]:

```
X train, X test, y train, y test = train test split(X, y, test size=0.30, random state=101)
```

Random Forest Classification

```
In [8]:
```

```
from sklearn.ensemble import RandomForestClassifier
```

In [9]:

```
#help(RandomForestClassifier)
```

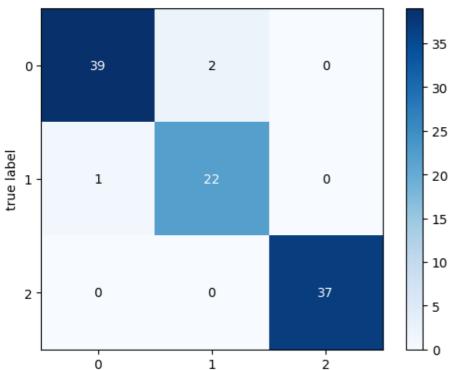
```
In [10]:
```

In [11]:

```
rfc.fit(X train,y train)
```

C:\Users\Chromsy\AppData\Roaming\Python\Python39\site-packages\sklearn\ensemble_forest.p y:424: FutureWarning: `max_features='auto'` has been deprecated in 1.1 and will be remove d in 1.3. To keep the past behaviour, explicitly set `max_features='sqrt'` or remove this parameter as it is also the default value for RandomForestClassifiers and ExtraTreesClass ifiers.

```
warn(
Out[11]:
                             RandomForestClassifier
RandomForestClassifier(max features='auto', n estimators=10, random state=101)
In [12]:
preds = rfc.predict(X test)
In [13]:
from sklearn.metrics import confusion_matrix,classification_report
In [14]:
from mlxtend.plotting import plot_confusion_matrix
In [15]:
d=confusion matrix(y test,preds)
Out[15]:
array([[39, 2, 0],
       [ 1, 22, 0],
       [0, 0, 37], dtype=int64)
In [16]:
plot confusion matrix(d,colorbar=True);
                                                       35
   0 -
            39
                           2
                                         0
                                                       - 30
```



In [17]:

print(classification_report(y_test,preds))

	precision	recall	f1-score	support
Adelie	0.97	0.95	0.96	41
Chinstrap	0.92	0.96	0.94	23
Gentoo	1.00	1.00	1.00	37

1 predicted label

```
accuracy 0.97 101 macro avg 0.96 0.97 0.97 101 weighted avg 0.97 0.97 0.97 101
```

Feature Importance

Very useful attribute of the trained model!

Out[19]:

	Feature Importance
culmen_length_mm	0.318677
culmen_depth_mm	0.101849
flipper_length_mm	0.173434
body_mass_g	0.213170
island_Dream	0.145121
island_Torgersen	0.037201
sex_FEMALE	0.006323
sex_MALE	0.004226

Choosing correct number of trees

Let's explore if continually adding more trees improves performance...

```
In [20]:
```

```
test_error = []
from sklearn.metrics import accuracy_score

for n in range(1,40):
    # Use n random trees
    model = RandomForestClassifier(n_estimators=n,max_features='auto')
    model.fit(X_train,y_train)
    test_preds = model.predict(X_test)
    test_error.append(1-accuracy_score(test_preds,y_test))
```

C:\Users\Chromsy\AppData\Roaming\Python\Python39\site-packages\sklearn\ensemble_forest.p
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```

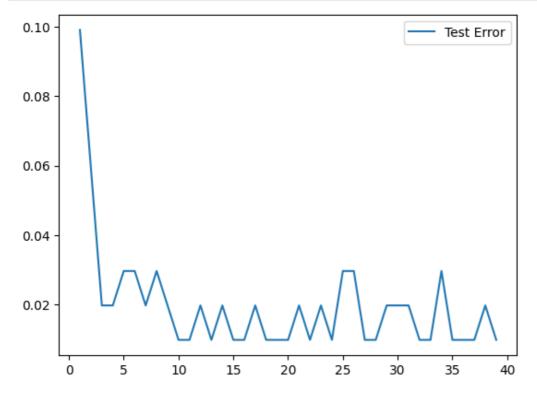
C:\Users\Chromsy\AppData\Roaming\Python\Python39\site-packages\sklearn\ensemble\ forest.p

warn(

```
warn(
C:\Users\Chromsy\AppData\Roaming\Python\Python39\site-packages\sklearn\ensemble\_forest.p
y:424: FutureWarning: `max_features='auto'` has been deprecated in 1.1 and will be remove
d in 1.3. To keep the past behaviour, explicitly set `max_features='sqrt'` or remove this
parameter as it is also the default value for RandomForestClassifiers and ExtraTreesClass
ifiers.
  warn(
```

In [21]:

```
plt.plot(range(1,40),test_error,label='Test Error')
plt.legend();
```



Clearly there are diminishing returns, on such a small dataset, we've pretty much extracted all the information we can after about 5 trees

Random Forest - HyperParameter Exploration

https://archive.ics.uci.edu/ml/datasets/banknote+authentication

Data

In [22]:

df=pd.read_csv("D:\\Study\\Programming\\python\\Python course from udemy\\Udemy - 2022 P
ython for Machine Learning & Data Science Masterclass\\01 - Introduction to Course\\1UNZI
P-FOR-NOTEBOOKS-FINAL\\DATA\\data_banknote_authentication.csv")
df.head()

Out[22]:

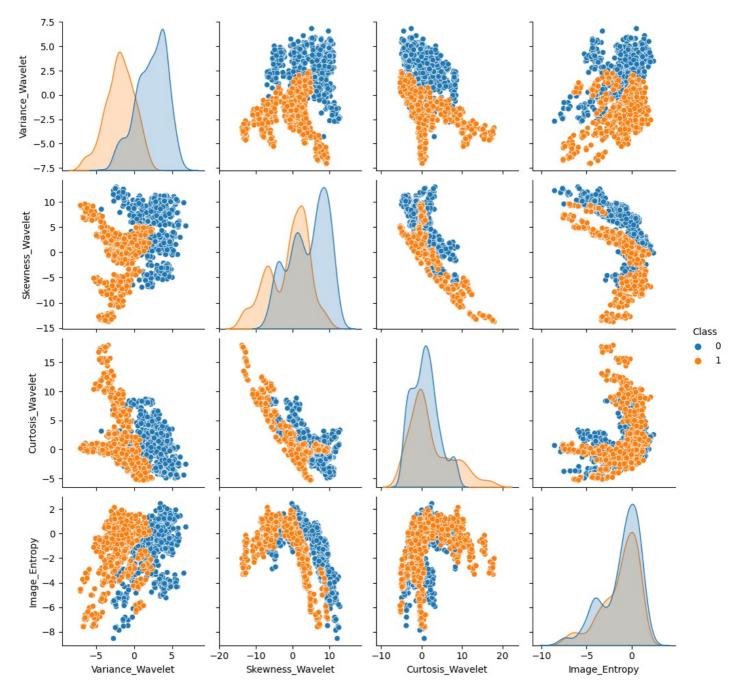
	Variance_Wavelet	Skewness_Wavelet	Curtosis_Wavelet	Image_Entropy	Class
0	3.62160	8.6661	-2.8073	-0.44699	0
1	4.54590	8.1674	-2.4586	-1.46210	0
2	3.86600	-2.6383	1.9242	0.10645	0
3	3.45660	9.5228	-4.0112	-3.59440	0
4	0.32924	-4.4552	4.5718	-0.98880	0

```
In [23]:
```

```
sns.pairplot(df,hue='Class')
```

Out[23]:

<seaborn.axisgrid.PairGrid at 0x1ed3da63f40>



In [24]:

```
X= df.drop('Class',axis=1)
y=df['Class']
```

In [25]:

```
from sklearn.model_selection import train_test_split
```

In [26]:

```
X_train, X_test, y_train, y_test= train_test_split(X, y, test_size=0.15, random_state=101)
```

In [27]:

```
from sklearn.model selection import GridSearchCV
```

In [28]:

```
param_grid={ 'n_estimators': [64,100,128,200],
           'max_features': [2,3,4],
           'bootstrap':[True, False],
           'oob score':[True,False]}
# Note, oob score only makes sense when bootstrap=True!
In [29]:
rfc = RandomForestClassifier()
In [30]:
grid = GridSearchCV(rfc,param grid)
In [31]:
grid.fit(X train, y train)
C:\Users\Chromsy\AppData\Roaming\Python\Python39\site-packages\sklearn\model selection\ v
alidation.py:378: FitFailedWarning:
60 fits failed out of a total of 240.
The score on these train-test partitions for these parameters will be set to nan.
If these failures are not expected, you can try to debug them by setting error score='rai
Below are more details about the failures:
60 fits failed with the following error:
Traceback (most recent call last):
 File "C:\Users\Chromsy\AppData\Roaming\Python\Python39\site-packages\sklearn\model sele
ction\ validation.py", line 686, in fit and score
   estimator.fit(X_train, y_train, **fit params)
  File "C:\Users\Chromsy\AppData\Roaming\Python\Python39\site-packages\sklearn\ensemble\_
forest.py", line 434, in fit
    raise ValueError("Out of bag estimation only available if bootstrap=True")
ValueError: Out of bag estimation only available if bootstrap=True
  warnings.warn(some fits failed message, FitFailedWarning)
C:\Users\Chromsy\AppData\Roaming\Python\Python39\site-packages\sklearn\model selection\ s
earch.py:952: UserWarning: One or more of the test scores are non-finite: [0.99314038 0.9
9314038 0.99399875 0.99314038 0.99314038 0.99314038
 0.99399875 \ 0.99399875 \ 0.99399875 \ 0.99228201 \ 0.99314038 \ 0.99142365
 0.99314038 0.99314038 0.99314038 0.99399875 0.98799017 0.98799017
 0.98884854 \ 0.98799017 \ 0.9871318 \ 0.98799017 \ 0.98799017 \ 0.98970691
        nan 0.99313305
                            nan 0.99313672
                                                    nan 0.99399142
        nan 0.99313672
                             nan 0.99228201
                                                   nan 0.98970691
        nan 0.99227835
                             nan 0.99056528
                                                   nan 0.97684604
       nan 0.98112688
                             nan 0.97770441
                                                   nan 0.97770441]
 warnings.warn(
Out[31]:
             GridSearchCV
 ▶ estimator: RandomForestClassifier
       RandomForestClassifier
In [32]:
grid.best params
Out[32]:
{'bootstrap': True, 'max features': 2, 'n_estimators': 100, 'oob_score': True}
In [33]:
predis = grid.predict(X test)
```

In [34]:

print(classification_report(y_test,predis)) precision recall f1-score support 0 1.00 0.98 0.99 124 1 0.98 1.00 0.99 82 0.99 206 accuracy 0.99 0.99 macro avg 0.99 206 weighted avg 0.99 0.99 0.99 206

In [35]:

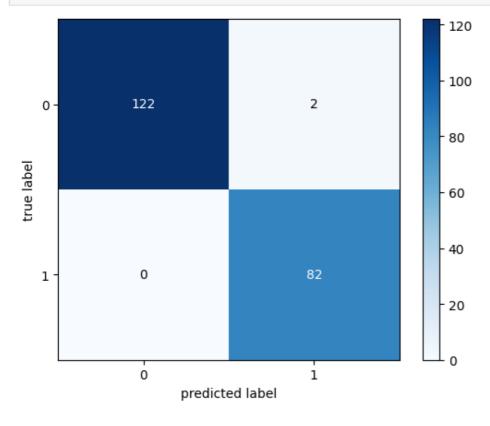
```
p=confusion_matrix(y_test,predis)
p
```

Out[35]:

```
array([[122, 2], [ 0, 82]], dtype=int64)
```

In [36]:

```
plot_confusion_matrix(p,colorbar=True);
```



In [37]:

```
# No underscore, reports back original oob_score parameter
grid.best_estimator_.oob_score
# If it true then it would generate value
```

Out[37]:

True

In [44]:

```
# With underscore, reports back fitted attribute of oob_score
grid.best_estimator_.oob_score_
# That value is only generate when .oob_score is True , show error if its False
```

Out[44]:

Understanding Number of Estimators (Trees)

Let's plot out error vs. Number of Estimators

```
In [39]:
```

```
from sklearn.metrics import accuracy_score
```

In [45]:

```
error = []
misclassifications = []

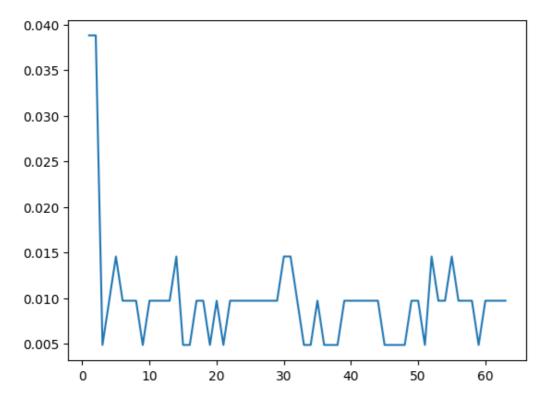
for n in range(1,64):
    rfc = RandomForestClassifier(n_estimators=n,bootstrap=True,max_features=2)
    rfc.fit(X_train,y_train)
    preds = rfc.predict(X_test)
    err = 1- accuracy_score(y_test,preds)
    n_missed=np.sum(preds != y_test) # watch the video to understand this line!!
    error.append(err)
    misclassifications.append(n_missed)
```

In [46]:

```
plt.plot(range(1,64),error)
```

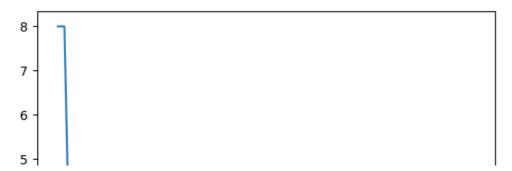
Out[46]:

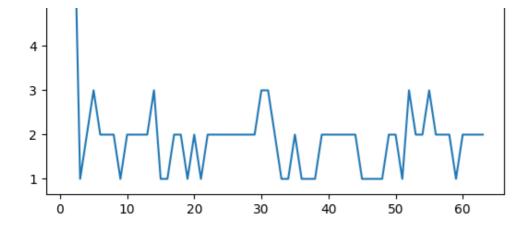
[<matplotlib.lines.Line2D at 0x1ed4ef6bdc0>]



In [47]:

```
plt.plot(range(1,64), misclassifications);
```





Random Forest - Regression

Plus: An Additional Analysis of Various Regression Methods!

The Data

We just got hired by a tunnel boring company which uses X-rays in an attempt to know rock density, ideally this will allow them to switch out boring heads on their equipment before having to mine through the rock!

They have given us some lab test results of signal strength returned in nHz to their sensors for various rock density types tested. You will notice it has almost a sine wave like relationship, where signal strength oscillates based off the density, the researchers are unsure why this is, but

```
In [48]:
```

```
import numpy as np
import pandas as pd
import matplotlib.pyplot as plt
import seaborn as sns
```

Data

In [49]:

df = pd.read_csv("D:\\Study\\Programming\\python\\Python course from udemy\\Udemy - 2022
Python for Machine Learning & Data Science Masterclass\\01 - Introduction to Course\\1UNZ
IP-FOR-NOTEBOOKS-FINAL\\DATA\\rock_density_xray.csv")
df.head()

Out[49]:

Rebound Signal Strength nHz Rock Density kg/m3

0	72.945124	2.456548
1	14.229877	2.601719
2	36.597334	1.967004
3	9.578899	2.300439
4	21.765897	2.452374

```
In [50]:
```

```
df.columns=['Signal','Density']
```

In [51]:

```
df.head()
```

```
        Signal
        Density

        0
        72.945124
        2.456548

        1
        14.229877
        2.601719

        2
        36.597334
        1.967004

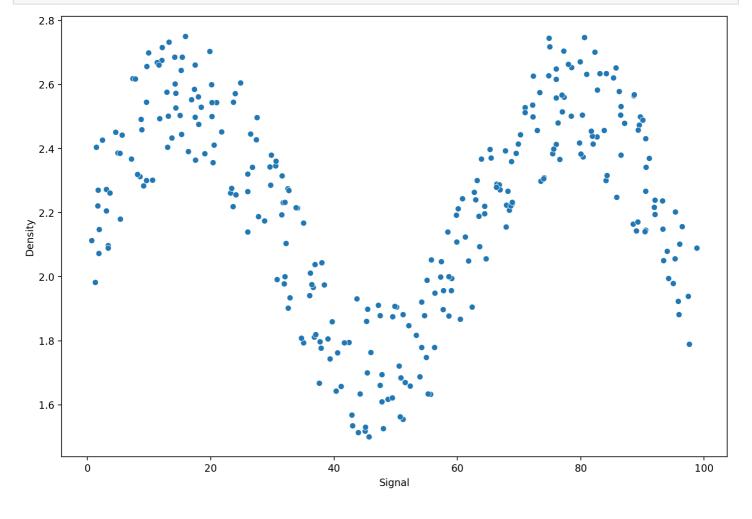
        3
        9.578899
        2.300439
```

21.765897 2.452374

In [56]:

Out[51]:

```
plt.figure(figsize=(12,8),dpi=200)
sns.scatterplot(x='Signal',y='Density',data=df);
```



Splitting the Data

Let's split the data in order to be able to have a Test set for performance metric evaluation.

```
In [98]:

X = df['Signal'].values.reshape(-1,1)
y = df['Density']

In [68]:
```

```
from sklearn.model_selection import train_test_split
```

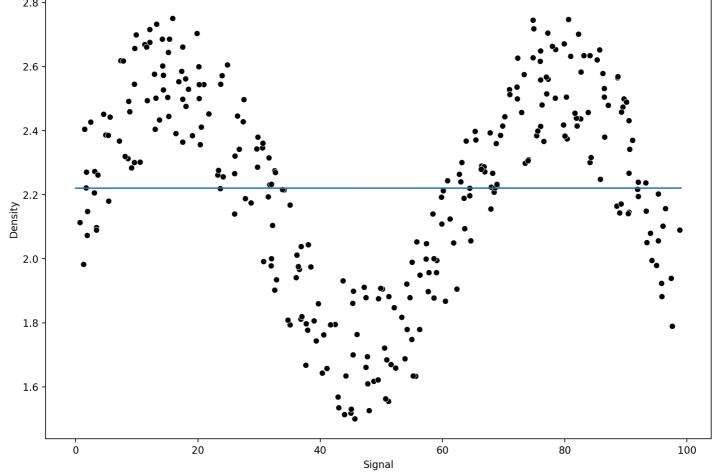
```
In [69]:
```

```
X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.1, random_state=101)
```

Linear Regression

```
In [70]:
from sklearn.linear_model import LinearRegression
In [71]:
lr model = LinearRegression()
In [72]:
lr_model.fit(X_train,y_train)
Out[72]:
▼ LinearRegression
LinearRegression()
In [73]:
lr pred = lr model.predict(X test)
In [89]:
# All number are close to 2.22
lr_pred
Out[89]:
array([2.22029657, 2.22047771, 2.22035637, 2.22034337, 2.22039737,
       2.22050555, 2.22042659, 2.22028877, 2.22034673, 2.22029714,
       2.22041506, 2.22050153, 2.22043891, 2.22042003, 2.22047022,
       2.22032403, 2.22033377, 2.22030628, 2.22035154, 2.22035373,
       2.22029266, 2.22036798, 2.22033018, 2.22030611, 2.22042754,
       2.22044019, 2.2204142 , 2.22040303, 2.22048946, 2.22047495])
In [90]:
from sklearn.metrics import mean squared error, mean absolute error
In [91]:
mean absolute error(y test, lr pred)
Out[91]:
0.211198973318633
In [75]:
np.sqrt(mean_squared_error(y_test,lr_pred))
Out[75]:
0.2570051996584629
What does the fit look like?
In [80]:
signal range = np.arange(0,100)
```

```
In [82]:
lr output = lr model.predict(signal range.reshape(-1,1))
In [92]:
lr_output
Out[92]:
array([2.22028446, 2.22028673, 2.22028899, 2.22029126, 2.22029353,
       2.22029579, 2.22029806, 2.22030032, 2.22030259, 2.22030485,
       2.22030712, 2.22030938, 2.22031165, 2.22031391, 2.22031618,
       2.22031844, 2.22032071, 2.22032297, 2.22032524, 2.2203275 ,
       2.22032977, 2.22033204, 2.2203343 , 2.22033657, 2.22033883,
       2.2203411 , 2.22034336, 2.22034563, 2.22034789, 2.22035016,
       2.22035242, 2.22035469, 2.22035695, 2.22035922, 2.22036148,
       2.22036375, 2.22036602, 2.22036828, 2.22037055, 2.22037281,
       2.22037508, 2.22037734, 2.22037961, 2.22038187, 2.22038414,
       2.2203864 , 2.22038867, 2.22039093, 2.2203932 , 2.22039546,
       2.22039773, 2.22039999, 2.22040226, 2.22040453, 2.22040679,
       2.22040906, 2.22041132, 2.22041359, 2.22041585, 2.22041812,
       2.22042038, 2.22042265, 2.22042491, 2.22042718, 2.22042944,
       2.22043171, 2.22043397, 2.22043624, 2.2204385 , 2.22044077,
       2.22044304, 2.2204453 , 2.22044757, 2.22044983, 2.2204521 ,
       2.22045436, 2.22045663, 2.22045889, 2.22046116, 2.22046342,
       2.22046569, 2.22046795, 2.22047022, 2.22047248, 2.22047475,
       2.22047701, 2.22047928, 2.22048155, 2.22048381, 2.22048608,
       2.22048834, 2.22049061, 2.22049287, 2.22049514, 2.2204974,
       2.22049967, 2.22050193, 2.2205042 , 2.22050646, 2.22050873])
In [84]:
plt.figure(figsize=(12,8),dpi=200)
sns.scatterplot(x='Signal',y='Density',data=df,color='black')
plt.plot(signal_range, lr_output);
  2.8
```



Polynomial Regression

Attempting with a Polynomial Regression Model

Let's explore why our standard regression approach of a polynomial could be difficult to fit here, keep in mind, we're in a fortunate situation where we can easily visualize results of y vs x.

Function to Help Run Models

```
In [85]:
```

```
from sklearn.linear_model import LinearRegression
model = LinearRegression()
```

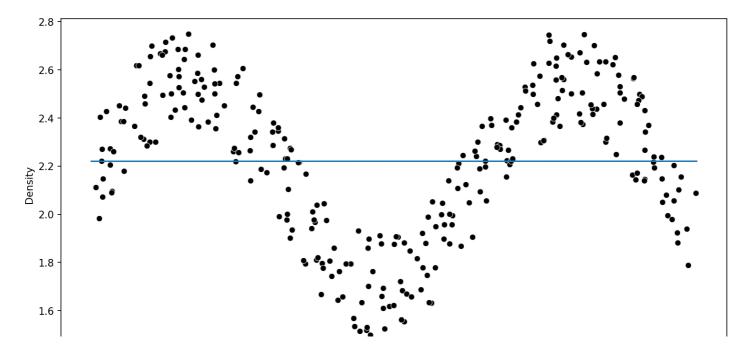
```
In [101]:
```

```
def run_model(model, X_train, y_train, X_test, y_test):
    # Fit Model
   model.fit(X_train,y_train)
    # Get Metrics
    preds = model.predict(X test)
    # Evaluation
    rmse = np.sqrt(mean squared error(y test,preds))
   mae = mean absolute_error(y_test,preds)
    print(f"MAE : {mae}")
   print(f"RMSE: {rmse}")
    #Plot results
    signal range = np.arange(0,100)
    output = model.predict(signal range.reshape(-1,1))
    plt.figure(figsize=(12,6),dpi=150)
    sns.scatterplot(x='Signal',y='Density',data=df,color='black')
   plt.plot(signal range,output)
```

In [102]:

```
model = LinearRegression()
run_model(model, X_train, y_train, X_test, y_test)
```

MAE: 0.211198973318633 RMSE: 0.2570051996584629



Pipeline for Poly Orders

In [103]:

from sklearn.pipeline import make pipeline

In [104]:

from sklearn.preprocessing import PolynomialFeatures

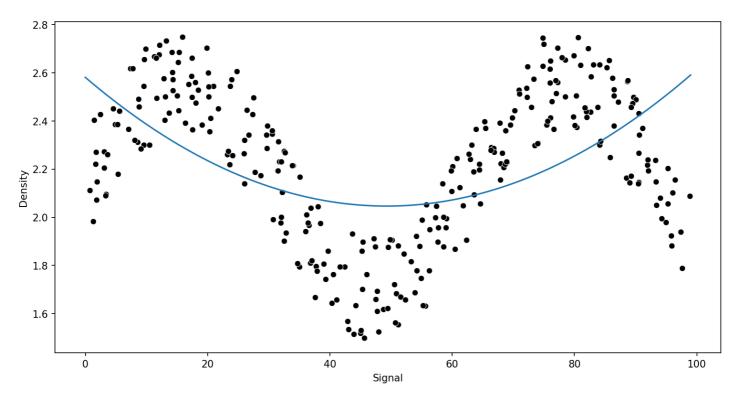
In [105]:

pipe = make_pipeline(PolynomialFeatures(2), LinearRegression())

In [106]:

run_model(pipe, X_train, y_train, X_test, y_test)

MAE : 0.22903105443511335 RMSE: 0.2817309563725596



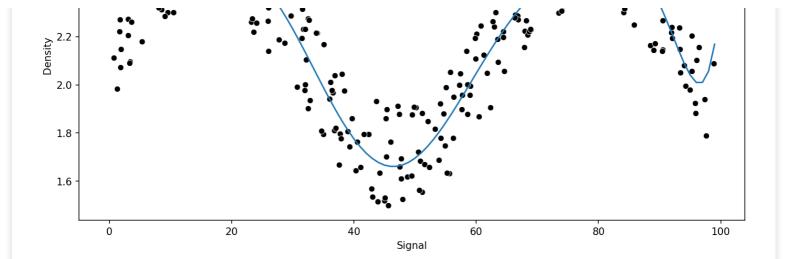
Comparing Various Polynomial Orders

In [107]:

pipe = make_pipeline(PolynomialFeatures(10), LinearRegression())
run_model(pipe, X_train, y_train, X_test, y_test)

MAE : 0.12478026429426053 RMSE: 0.14049911421019984





KNN Regression

In [108]:

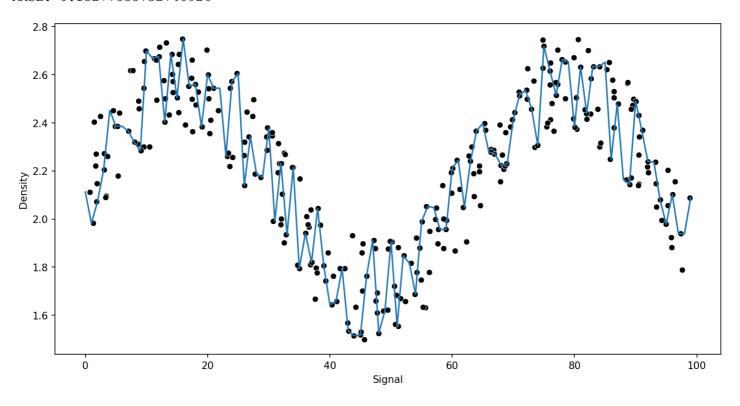
from sklearn.neighbors import KNeighborsRegressor

```
In [109]:
```

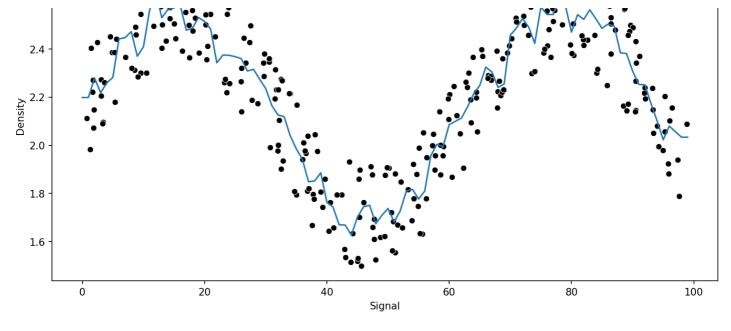
```
preds ={}
k_values = [1,5,10]
for n in k_values:

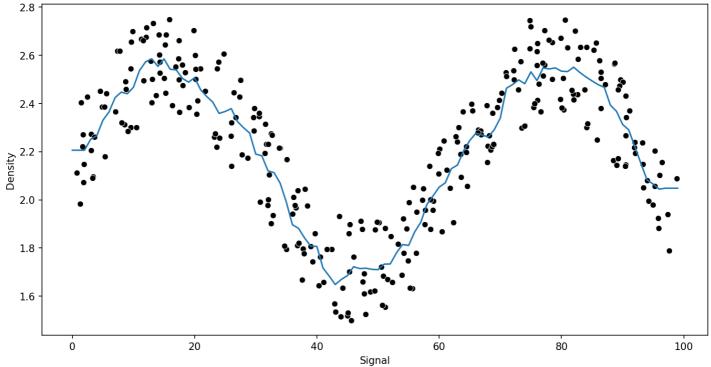
model = KNeighborsRegressor(n_neighbors=n)
   run_model(model,X_train,y_train,X_test,y_test)
```

MAE : 0.11877297474442378 RMSE: 0.1523487028635337 MAE : 0.12198383614100558 RMSE: 0.13730685016923647 MAE : 0.11635971693292672 RMSE: 0.13277855732740926









Decision Tree Regression

```
In [110]:
```

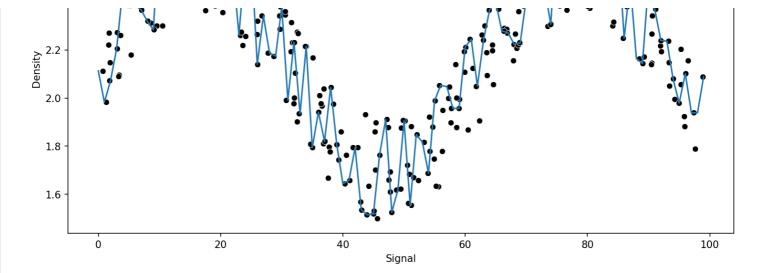
```
from sklearn.tree import DecisionTreeRegressor
```

```
In [111]:
```

```
model = DecisionTreeRegressor()
run_model(model, X_train, y_train, X_test, y_test)
```

MAE : 0.11877297474442378 RMSE: 0.1523487028635337





```
In [112]:
model.get_n_leaves()
Out[112]:
270
```

Support Vector Regression

```
In [113]:
    from sklearn.svm import SVR

In [114]:
    from sklearn.model_selection import GridSearchCV

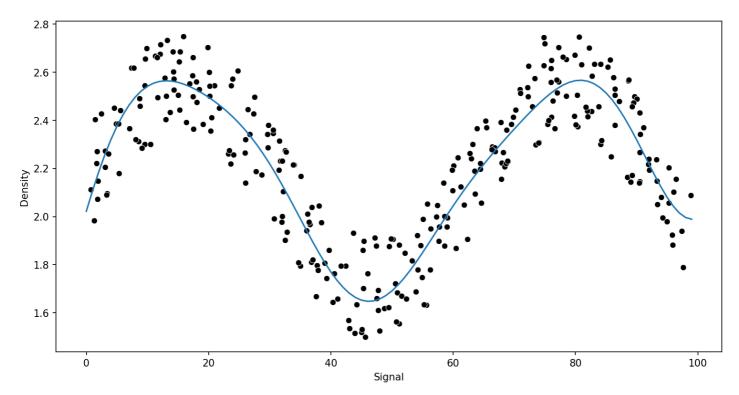
In [115]:
    param_grid = {'C':[0,0.01,.1,1,5,10,100,1000],'gamma':['auto','scale']}
    svr = SVR()

In [116]:
    grid = GridSearchCV(svr,param_grid)

In [117]:
    run_model(grid,X_train,y_train,X_test,y_test)
```

```
C:\Users\Chromsy\AppData\Roaming\Python\Python39\site-packages\sklearn\model selection\ v
alidation.py:378: FitFailedWarning:
10 fits failed out of a total of 80.
The score on these train-test partitions for these parameters will be set to nan.
If these failures are not expected, you can try to debug them by setting error score='rai
Below are more details about the failures:
10 fits failed with the following error:
Traceback (most recent call last):
 File "C:\Users\Chromsy\AppData\Roaming\Python\Python39\site-packages\sklearn\model_sele
ction\ validation.py", line 686, in fit and score
    estimator.fit(X train, y train, **fit params)
  File "C:\Users\Chromsy\AppData\Roaming\Python\Python39\site-packages\sklearn\svm\_base.
py", line 180, in fit
    self. validate params()
  File "C:\Users\Chromsy\AppData\Roaming\Python\Python39\site-packages\sklearn\base.py",
line 581, in validate params
```

MAE : 0.10854210121348368 RMSE: 0.12646999302046696



```
In [118]:
```

grid.best estimator

Out[118]:

▼ SVR SVR (C=1000)

Random Forest Regression

```
In [119]:
```

```
from sklearn.ensemble import RandomForestRegressor
```

In [120]:

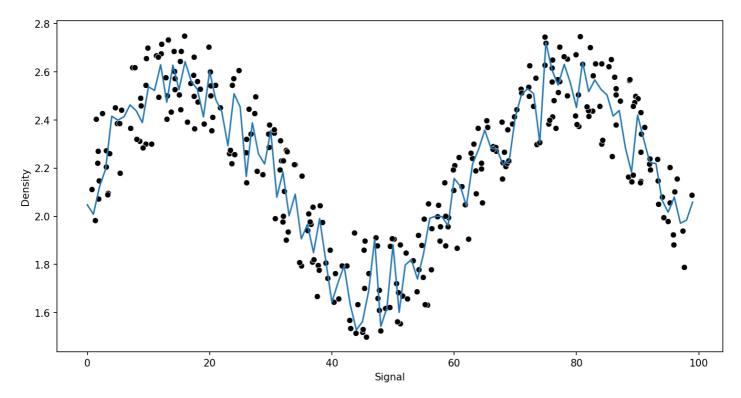
```
# help(RandomForestRegressor)
```

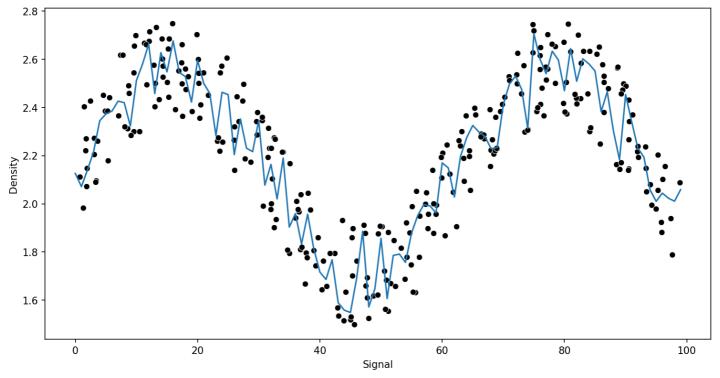
```
In [121]:
```

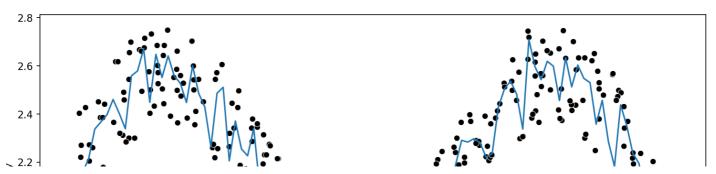
```
trees = [10,50,100]
for n in trees:
```

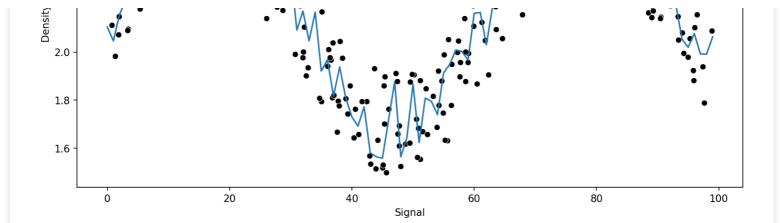
model = RandomForestRegressor(n_estimators=n)
run_model(model, X_train, y_train, X_test, y_test)

MAE : 0.10794614401455976 RMSE: 0.12923933126466158 MAE : 0.11190211837017762 RMSE: 0.13637615617898904 MAE : 0.10990104288555731 RMSE: 0.13096990474747555









Gradient Boosting

We will cover this in more detail in next section.

In [122]:

from sklearn.ensemble import GradientBoostingRegressor

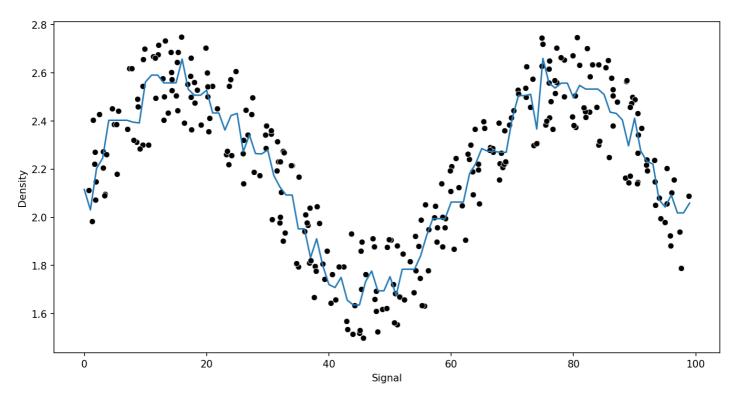
In [123]:

help(GradientBoostingRegressor)

In [124]:

```
model = GradientBoostingRegressor()
run_model(model, X_train, y_train, X_test, y_test)
```

MAE: 0.11318284854800689 RMSE: 0.13294148649584667



Adaboost

In [127]:

from sklearn.ensemble import AdaBoostRegressor

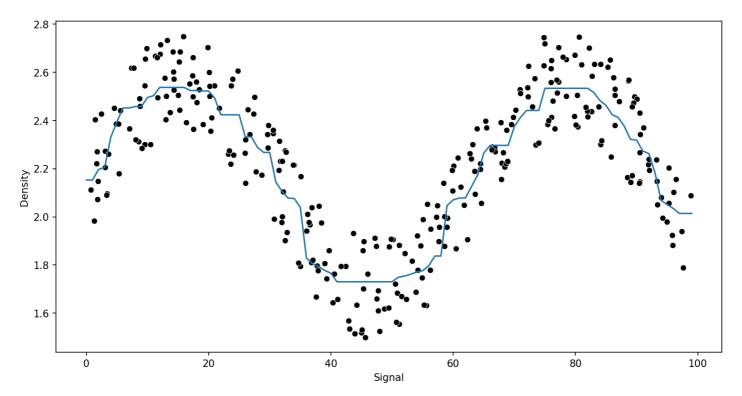
Tn [1281:

______.

model = AdaBoostRegressor()

run_model(model,X_train,y_train,X_test,y_test)

MAE : 0.11711438520863086 RMSE: 0.13457543971524755



In []: