```
# Import libraries
import pandas as pd
import numpy as np
import matplotlib.pyplot as plt
import seaborn as sns
from sklearn.preprocessing import LabelEncoder
from sklearn.model selection import train test split
from sklearn.decomposition import PCA
from xgboost import XGBRegressor
from xgboost import XGBRFRegressor
from sklearn.metrics import mean squared error
from math import sqrt
import warnings
warnings.filterwarnings('ignore')
# import train dataset
mdz train df = pd.read csv('train.csv')
mdz train df.head()
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[5 rows x 378 columns]
mdz train df.shape
(4209, 378)
# import test dataset
mdz test df = pd.read csv('test.csv')
mdz test df.head()
```

```
X0 X1 X2 X3 X4 X5 X6 X8 X10
                                                   X375
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3
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4
       0
              0
                     0
                            0
[5 rows x 377 columns]
mdz test df.shape
(4209, 377)
In both datasets, we dont need the ID column. Lets remove it.
mdz_train_df.drop(['ID'], axis=1, inplace=True)
mdz_test_df.drop(['ID'], axis=1, inplace=True)
Task 1. If for any column(s), the variance is equal to zero, then you need to remove those
variable(s)
Now lets remove the columns having variance equal to zero in train dataset.
var_zero_cols = mdz_train_df.var()[mdz_train_df.var() ==
0].index.values.tolist()
var zero cols
['X11',
 'X93'
 'X107',
 'X233',
 'X235'
 'X268'
 'X289',
 'X290',
```

'X293', 'X297',

```
'X330',
 'X347'1
# Remove above columns from train and test
mdz train df.drop(var zero cols, axis=1, inplace=True)
mdz test df.drop(var zero cols, axis=1, inplace=True)
print(mdz train df.shape)
print(mdz test df.shape)
(4209, 365)
(4209, 364)
Task 2. Check for null and unique values for test and train sets
mdz train df.isnull().sum().sum()
0
mdz test df.isnull().sum().sum()
0
There is no NaN values in both train and test dataset.
mdz train df.nunique().sum()
3452
mdz_test_df.nunique().sum()
908
There are 3452 unique values in train and 908 unique values in test datasets.
Task 3. Apply label encoder
object cols = mdz train df.describe(include=[object]).columns.values
object cols
array(['X0', 'X1', 'X2', 'X3', 'X4', 'X5', 'X6', 'X8'], dtype=object)
Above columns are of object datatype.
mdz train df[object cols].head()
   X0 X1 X2 X3 X4 X5 X6 X8
0
    k v
          at a d u
                        j
                           0
1
    k t av e d y l
                           0
  az w
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3
  az
       t
           n
              f
                  d x l
                           е
           n f
                  d h d n
  az
le = LabelEncoder()
```

```
for col in object cols:
    le.fit(mdz train df[col].append(mdz test df[col]).values)
    mdz_train_df[col] = le.transform(mdz_train_df[col])
    mdz test df[col] = le.transform(mdz test df[col])
Now we can proceed with the dimensionality reduction and ML model.
x = mdz train df.drop(['y'], axis=1)
y = mdz train df.y
# train test split
x_train, x_test, y_train, y_test = train_test_split(x, y,
test size=0.3, random state=0)
print(x train.shape)
print(x test.shape)
print(y_train.shape)
print(y_test.shape)
(2946, 364)
(1263, 364)
(2946,)
(1263,)
Task 4. Perform dimensionality reduction
pca = PCA(0.98)
pca.fit(x)
pca.n components
12
So we need 12 components generated from all 364 columns to get 98% variance and
project all the data to lower dimension.
pca.explained variance ratio
array([0.40868988, 0.21758508, 0.13120081, 0.10783522, 0.08165248,
       0.0140934 , 0.00660951, 0.00384659, 0.00260289, 0.00214378,
       0.00209857, 0.00180388])
Above are the explained varance ratio/percentage (high to low).
# Convert the data into lower dimension (ld)
x train ld = pca.transform(x train)
x test ld = pca.transform(x test)
x_test_set_ld = pca.transform(mdz_test_df)
print(x train ld.shape)
```

```
print(x test ld.shape)
print(x test set ld.shape)
(2946.12)
(1263, 12)
(4209, 12)
```

Task 5. Predict your test_df values using xgboost

```
We can build xgboost (XGBRegeessor, XGBRFRegressor) models with different
configuration of learning rates.
xgbRegressor 1 = XGBRegressor(objective='reg:squarederror',
learning rate = 0.1)
xgbRegressor 1.fit(x_train_ld, y_train)
XGBRegressor(base score=0.5, booster=None, colsample bylevel=1,
             colsample bynode=1, colsample bytree=1, gamma=0, gpu id=-
1,
             importance type='gain', interaction constraints=None,
             learning rate=0.1, max delta step=0, max depth=6,
             min child weight=1, missing=nan,
monotone_constraints=None,
             n_estimators=100, n jobs=0, num parallel tree=1,
random state=0,
             reg alpha=0, reg lambda=1, scale pos weight=1,
subsample=1,
             tree method=None, validate parameters=False,
verbosity=None)
xgbRegressor 1 y pred = xgbRegressor 1.predict(x test ld)
print('RMSE', sqrt(mean squared error(y test, xgbRegressor 1 y pred)))
RMSE 9.996668252109297
xgbRegressor 2 = XGBRegressor(objective='reg:squarederror',
learning rate = 1)
xgbRegressor 2.fit(x train ld, y train)
XGBRegressor(base score=0.5, booster=None, colsample bylevel=1,
             colsample bynode=1, colsample bytree=1, gamma=0, gpu id=-
1,
             importance type='gain', interaction constraints=None,
             learning rate=1, max delta step=0, max depth=6,
min child weight=1,
             missing=nan, monotone_constraints=None, n estimators=100,
n jobs=0,
             num parallel tree=1, random state=0, reg alpha=0,
reg lambda=1,
             scale pos weight=1, subsample=1, tree method=None,
             validate parameters=False, verbosity=None)
```

```
xgbRegressor_2_y_pred = xgbRegressor 2.predict(x test ld)
print('RMSE', sqrt(mean squared error(y test, xgbRegressor 2 y pred)))
RMSE 12.141280477327937
Here we can choose xgbRegressor_1 model. It gave less RMSE.
xgbRFRegressor 1 = XGBRFRegressor(objective='reg:squarederror',
learning rate = 0.1)
xgbRFRegressor 1.fit(x train ld, y train)
XGBRFRegressor(base score=0.5, booster=None, colsample bylevel=1,
               colsample bytree=1, gamma=0, gpu id=-1,
importance type='gain',
               interaction constraints=None, learning rate=0.1,
               max delta step=0, max depth=6, min child weight=1,
missing=nan,
               monotone constraints=None, n estimators=100, n jobs=0,
               num parallel tree=100, objective='reg:squarederror',
               random state=0, reg alpha=0, scale pos weight=1,
               tree method=None, validate parameters=False,
verbosity=None)
xgbRFRegressor 1 y pred = xgbRFRegressor 1.predict(x test ld)
print('RMSE', sqrt(mean squared error(y test,
xgbRFRegressor 1 y pred)))
RMSE 91.53699087619204
xqbRFReqressor 2 = XGBRFReqressor(objective='req:squarederror',
learning rate = 1)
xqbRFRegressor 2.fit(x_train_ld, y_train)
XGBRFRegressor(base score=0.5, booster=None, colsample bylevel=1,
               colsample bytree=1, gamma=0, gpu id=-1,
importance type='gain',
               interaction constraints=None, max delta step=0,
max depth=6,
               min child weight=1, missing=nan,
monotone_constraints=None,
               n estimators=100, n jobs=0, num parallel tree=100,
               objective='reg:squarederror', random state=0,
reg alpha=0,
               scale pos weight=1, tree method=None,
validate_parameters=False,
               verbosity=None)
xgbRFRegressor 2 y pred = xgbRFRegressor 2.predict(x test ld)
print('RMSE', sqrt(mean squared error(y test,
xgbRFRegressor 2 y pred)))
RMSE 10.313216428415938
```

XGBRegressor model with learning rate 0.1 gave the least RMSE. So we can choose this model for prediction.