Capstone Project 1

(BY MULA SHASHANK)

ANOMA DATA -Automated Anomaly Detection for Predictive Maintenance

Problem Statement:

Many different industries need predictive maintenance solutions to reduce risks and gain actionable insights through processing data from their equipment. Although system failure is a very general issue that can occur in any machine, predicting the failure and taking steps to prevent such failure is most important for any machine or software application. Predictive maintenance evaluates the condition of equipment by performing online monitoring. The goal is to perform maintenance before the equipment degrades or breaks down. This Capstone project is aimed at predicting the machine breakdown by identifying the anomalies in the data. The data we have contains about 18000+ rows collected over few days. The column 'y' contains the binary labels, with 1 denoting there is an anomaly. The rest of the columns are predictors.

Loading the required libraries

```
In [1]:
import pandas as pd
import numpy as np
import matplotlib.pyplot as plt
%matplotlib inline
import seaborn as sns
from sklearn.model selection import train test split, cross val score
from sklearn.preprocessing import StandardScaler, LabelEncoder
from sklearn.ensemble import RandomForestClassifier
from sklearn.metrics import classification report, confusion matrix,
accuracy score
                                                                          In [2]:
# Loading the dataset
data = pd.read excel('AnomaData.xlsx')
# Displaying the first few rows of the dataset
data.head()
      Out[2]
5 rows * 62 colums
                                                                       In[3]
data.columns
                                                                        Out[3]:
Index(['time', 'y', 'x1', 'x2', 'x3', 'x4', 'x5', 'x6', 'x7', 'x8', 'x9',
```

```
'x10', 'x11', 'x12', 'x13', 'x14', 'x15', 'x16', 'x17', 'x18', 'x19'

'x20', 'x21', 'x22', 'x23', 'x24', 'x25', 'x26', 'x27', 'x28', 'x29'

'x30', 'x31', 'x32', 'x33', 'x34', 'x35', 'x36', 'x37', 'x38', 'x39'

'x40', 'x41', 'x42', 'x43', 'x44', 'x45', 'x46', 'x47', 'x48', 'x49'

'x50', 'x51', 'x52', 'x54', 'x55', 'x56', 'x57', 'x58', 'x59', 'x60'

'y.1'],
dtype='object')

In [4]:
```

data.drop(['y.1','time'],axis=1,inplace=True)

Exploratory Data Analysis

```
In [ ]:
from ydata profiling import ProfileReport
rep=ProfileReport(data)
rep
Summarize dataset: 0%| | 0/5 [00:00<?, ?it/s]
                                                                      In [7]:
# Checking null values
data.isna().sum()
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No null values present
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data.describe()

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```

8 rows × 60 columns

```
# Separate features and target variables
X = data.drop('y',axis=1)
y = data['y'] # Binary target variable for anomaly detection
In [10]:
```

Checking the skewness of the columns

from scipy.stats import skew

```
for col in X.columns:
    print(col+" --> Skew=", skew(data[col]))
x1 \longrightarrow Skew = -1.247625732759167
x2 --> Skew= 0.24618128774105086
x3 --> Skew= 0.4055454880638477
x4 \longrightarrow Skew = 0.12514637378313767
x5 \longrightarrow Skew = 0.6837495582949324
x6 \longrightarrow Skew = -0.7783583177473792
x7 \longrightarrow Skew = 2.4357311339796475
x8 --> Skew= -1.6026471752443343
x9 \longrightarrow Skew = 9.012488007639806
x10 --> Skew= 8.943562386353957
x11 \longrightarrow Skew = -7.185593183352757
x12 \longrightarrow Skew = -8.361297553646873
x13 --> Skew= -0.12379559435046963
x14 --> Skew= 10.143169032655202
x15 --> Skew= 7.49124549847556
x16 --> Skew= -7.889659645806755
x17 --> Skew= 0.8767548887172883
x18 --> Skew= 0.6193348339124728
x19 --> Skew= -14.642912582263218
x20 \longrightarrow Skew = -1.2036063560137629
x21 --> Skew= 0.25536084410168486
x22 --> Skew= 0.5803436878881983
x23 --> Skew= -3.1418220516846533
x24 \longrightarrow Skew = 0.554933106032238
x25 \longrightarrow Skew = -9.498804786346405
x26 --> Skew= 0.7062128721051979
x27 \longrightarrow Skew = -1.09724465412704
x28 --> Skew= 0.8690546825826609
x29 \longrightarrow Skew = -0.3144094618837163
x30 \longrightarrow Skew = -0.38429639300540624
x31 --> Skew= -0.06350061244827468
x32 \longrightarrow Skew = -9.61533865007797
x33 --> Skew= 0.27416672685782983
x34 \longrightarrow Skew = -0.7427083382729173
x35 \longrightarrow Skew = -0.05333477661865808
x36 --> Skew= 4.796638845977901
x37 \longrightarrow Skew = -2.01988817570014
x38 --> Skew= 9.851534132930421
x39 --> Skew= 3.7708045847451492
x40 \longrightarrow Skew = -4.413442472540411
x41 --> Skew= 0.29427380773473477
x42 --> Skew= 1.51225912751579
x43 \longrightarrow Skew = 7.522535603813891
x44 \longrightarrow Skew = -0.751908745142232
x45 --> Skew= 9.14525391875508
x46 --> Skew= 2.085440032593785
x47 \longrightarrow Skew = -0.9776387034518712
x48 --> Skew= -1.8096800837955427
x49 \longrightarrow Skew = -0.41343459474009875
x50 --> Skew= 0.23970568808172968
x51 \longrightarrow Skew = -10.384495189618546
x52 \longrightarrow Skew = -1.3077363575331327
x54 \longrightarrow Skew = -0.9314628850541723
x55 --> Skew= -0.46067152341753304
```

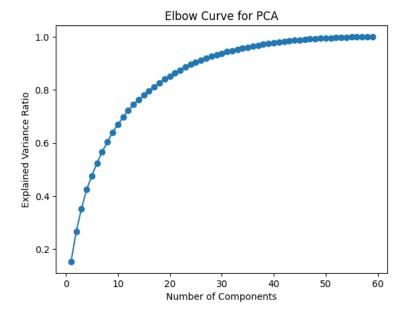
```
x56 --> Skew= 0.5252953172596814
x57 \longrightarrow Skew = 0.20301850665091772
x58 \longrightarrow Skew = -2.460119223931329
x59 \longrightarrow Skew = -9.49916162199559
x60 --> Skew= 0.4793085411594726
\# Applying yeo-johnson transformation as there are both negative and
positive skew
from sklearn.preprocessing import PowerTransformer
pt=PowerTransformer(method='yeo-johnson')
                                                                          In [12]:
# Performing train test split first to prevent data leakage later
X_train, X_test, y_train, y_test = train_test_split(X, y,
test_size=0.2,random_state=42,stratify=y)
                                                                          In [13]:
print("Shape of X train: ",X train.shape)
print("Shape of y train: ",y train.shape)
print("Shape of X test: ",X test.shape)
print("Shape of y test: ",y test.shape)
                                                                         Out[13]
Shape of X train: (14718, 59)
Shape of y train: (14718,)
Shape of X test: (3680, 59)
Shape of y test: (3680,)
# Transforming the X train and X test
X_train=pt.fit_transform(X_train)
X test=pt.transform(X test)
                                                                         In [15]:
# Checking the value counts to see if there is imbalance in target column
y train.value counts()
                                                                         Out[15]:
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    14619
        99
Name: count, dtype: int64
```

Dataset is imbalanced. So we will apply oversampling method SMOTE

```
У
0
     14619
1
     14619
Name: count, dtype: int64
Now the dataset is balanced
print("Shape of X train: ",X train.shape)
print("Shape of y train: ",y train.shape)
Shape of X train: (29238, 59)
Shape of y train:
                   (29238,)
Featuring Scaling
                                                                        In [19]:
# Scaling the data using Standard Scaler
sc=StandardScaler()
X train sc=sc.fit transform(X train)
X test sc=sc.transform(X test)
                                                                        In [20]:
train data processed=pd.concat([pd.DataFrame(X train sc,columns=X.columns),
pd.DataFrame(y train,columns=['y'])],axis=1)
train data processed.to csv('/content/train data processed.csv')
                                                                        In [21]:
test data processed=pd.concat([pd.DataFrame(X test sc,columns=X.columns),
                                 pd.DataFrame(y test, columns=['y'])], axis=1)
                                                                        In [22]:
test data processed.to csv('/content/test data processed.csv')
                                                                        In [23]:
# Applying PCA on n number of components to see how many components needed
# to achieve more than 90% explained variance ratio
from sklearn.decomposition import PCA
n components range=np.arange(1, X train sc.shape[1]+1)
explained variances=[]
for n component in n components range:
    pca=PCA(n components=n component)
    X_train_pca=pca.fit_transform(X_train_sc)
    explained variance=pca.explained variance ratio
    explained variances.append(np.sum(pca.explained variance ratio ))
plt.plot(n_components_range,explained_variances,marker='o')
plt.xlabel('Number of Components')
plt.ylabel('Explained Variance Ratio')
```

plt.title('Elbow Curve for PCA')

plt.show()



We see that for n_component=30, we can reach 90% of the explained variance ratio. So we choose n_components=30

```
In [24]:
pca=PCA(n components=30)
X train pca=pca.fit transform(X train sc)
                                                                   In [25]:
print("Shape of X train before PCA: ",X train sc.shape)
print("Shape of X train after PCA: ",X train pca.shape)
Shape of X train before PCA: (29238, 59)
Shape of X train after PCA: (29238, 30)
                                                                   In [26]:
X test pca=pca.transform(X test sc)
                                                                   In [27]:
print("Shape of X test before PCA: ",X_test_sc.shape)
print("Shape of X test after PCA: ", X test pca.shape)
Shape of X test before PCA: (3680, 59)
Shape of X test after PCA: (3680, 30)
                                                                   In[28]
# Predict and evaluate the binary classifier
def evaluate model(clf, X train, X test):
   clf.fit(X train, y train)
   y pred = clf.predict(X test)
   print ("============"")
   print("Binary Classification - Predictictive Maintenance")
   print("============="")
   print("Accuracy:", accuracy_score(y_test, y_pred))
   print("Classification Report:\n", classification report(y test, y pred,
target names=['No Anomaly','Anomaly']))
   sns.heatmap(confusion matrix(y test, y pred),annot=True,fmt='g',
          yticklabels=['No Anomaly','Anomaly'],xticklabels=['No
Anomaly','Anomaly'])
```

```
plt.xlabel('Predicted')
    plt.ylabel('True')
    plt.tight_layout()
    plt.show()
                                                                         In[29]
from sklearn.linear model import LogisticRegression
from sklearn.svm import SVC
from sklearn.tree import DecisionTreeClassifier
from sklearn.metrics import accuracy score, recall score, f1 score
                                                                         In [30]:
from sklearn.ensemble import AdaBoostClassifier,GradientBoostingClassifier
from xgboost import XGBClassifier
                                                                         In [31]:
model dict={'Support Vector Classifier':SVC(),
      'Random Forest': RandomForestClassifier(),
      'Gradient Boosting Classifier' : GradientBoostingClassifier(),
      'XGBoost':XGBClassifier()
          }
                                                                         In [32]:
def model results(model dict):
    results=[]
    for modelname, model in model dict.items():
        model.fit(X train pca,y train)
        acc=accuracy_score(y_test, model.predict(X_test_pca))
        rec=recall score(y test, model.predict(X test pca))
        f1=f1 score(y test, model.predict(X test pca))
crossvalscore=np.mean(cross val score(model, X train pca, y train, cv=3))
        results.append({'Model Name':modelname,
                         'Accuracy':acc,
                         'Recall':rec,
                         'F1 Score':f1,
                         'Cross Val Score': crossvalscore})
    results df=pd.DataFrame(results)
    return results df
model results(model dict)
                                                                        Out[33]:
                Model Name Accuracy Recall F1 Score Cross Val Score
 0
      Support Vector Classifier 0.982609
                                          0.36 0.219512
                                                                0.990184
 1
               Random Forest 0.994837
                                          0.32 0.457143
                                                                0.998940
```

0.52 0.111588

0.943750

0.965045

2 Gradient Boosting Classifier

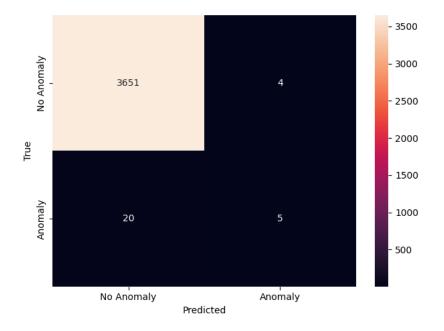
3 XGBoost 0.992935 0.32 0.380952 0.997093

For all the models, recall is extremely low, which means model isn't performing well on test data

In[33]

rf clf=RandomForestClassifier()

II_CII—Kanaomi	OTCDCCTGDDTT	TCT ()			
evaluate_model	(rf_clf,X_tr	ain_pca,X	_test_pca)		Out[33]
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Binary Classif	ication - Pr	edicticti	ve Maintena	ance	
=======================================				====	
Accuracy: 0.99	347826086956	53			
Classification	Report:				
	precision	recall	f1-score	support	
	precipion	rcoarr	11 00010	cappore	
No Anomaly	0.99	1.00	1.00	3655	
2					
Anomaly	0.56	0.20	0.29	25	
accuracy			0.99	3680	
macro avg	0.78	0.60	0.65	3680	
weighted avg	0.99	0.99	0.99	3680	



In[34]

gb_clf=GradientBoostingClassifier()
evaluate_model(gb_clf,X_train_pca,X_test_pca)

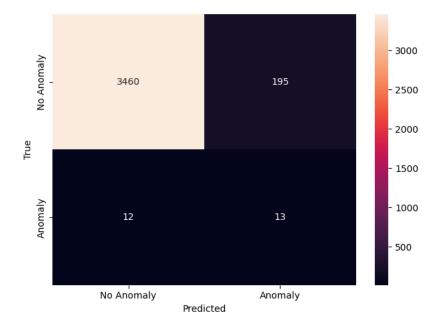
<pre>evaluate_model(gb_clf,X_train_pca,X_test_pca)</pre>	evaluate	_model(gb_	_clf,X_	_train_	pca,X	_test_	pca)
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Binary Classification - Predictictive Maintenance

Accuracy: 0.94375

Classification Report:

	precision	recall	f1-score	support
No Anomaly	1.00	0.95	0.97	3655
Anomaly	0.06	0.52	0.11	25
accuracy			0.94	3680
macro avg	0.53	0.73	0.54	3680
weighted avg	0.99	0.94	0.97	3680



For both the models, we see from the confusion matrix and classification report that the recall is extremely low. So we use a different method instead of PCA.

Using Feature Selection

```
In [35]:

from sklearn.feature_selection import SelectFromModel

In [36]:

# Function to select the best features and show the evaluation metrics for that best model

def select_from_model(clf):
    sfm=SelectFromModel(clf)
    sfm.fit(X_train_sc, y_train)
    sfm.transform(X_train_sc)
    X_train_sc_df=pd.DataFrame(X_train_sc,columns=X.columns)
    X_train_sc_df=X_train_sc_df[X_train_sc_df.columns[sfm.get_support()]]
    X_test_sc_df=pd.DataFrame(X_test_sc,columns=X.columns)
```

evaluate_model(clf,X_train_sc_df,X_test_sc_df)

gb_clf2=GradientBoostingClassifier()
gb_clf2.fit(X_train_sc,y_train)
select_from_model(gb_clf2)

In [37]:

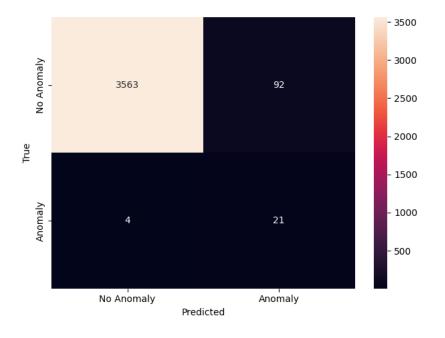
OUT[37]

Binary Classification - Predictictive Maintenance

Accuracy: 0.9739130434782609

Classification Report:

	precision	recall	f1-score	support
No Anomaly	1.00	0.97	0.99	3655
Anomaly	0.19	0.84	0.30	25
accuracy			0.97	3680
macro avg	0.59	0.91	0.65	3680
weighted avg	0.99	0.97	0.98	3680



In[38]

Out[38]

rf_clf2=RandomForestClassifier()
rf_clf2.fit(X_train_sc,y_train)
select from model(rf clf2).

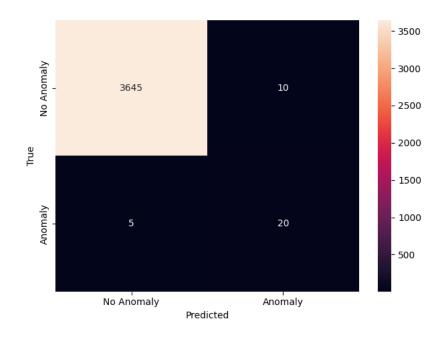
Binary Classification - Predictictive Maintenance

Accuracy: 0.9959239130434783

Classification Report:

	precision	recall	f1-score	support
No Anomaly Anomaly	1.00 0.67	1.00	1.00 0.73	3655 25
accuracv			1.00	3680

macro	avg	0.83	0.90	0.86	3680
weighted	avα	1.00	1.00	1.00	3680



In[39]

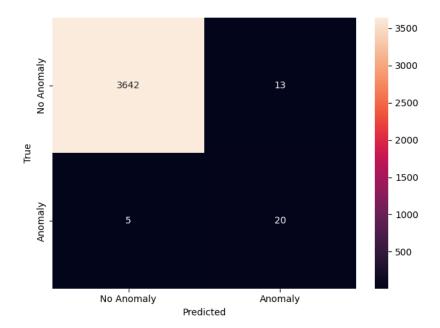
xgb_clf=XGBClassifier()
xgb_clf.fit(X_train_sc,y_train)
select_from_model(xgb_clf).

Out[39]

Binary Classification - Predictictive Maintenance

Classification Report:

	precision	recall	f1-score	support
No Anomaly	1.00	1.00	1.00	3655
Anomaly	0.61	0.80	0.69	25
accuracy			1.00	3680
macro avg	0.80	0.90	0.84	3680
weighted avg	1.00	1.00	1.00	3680



As we can see, the Gradient Boosting Classifier has excellent recall on the 'Anomaly'(1) class, but overall accuracy and F1 score is higher for Random Forest Model. The Random Forest Model is also slightly performing better than the XGBoost model

We see that the test data is highly imbalanced but still model performs very well with feature selection.

Hyperparameter tuning for the Random Forest Model

```
In[40]
# Selecting the best features and building the Random forest model
  sfm=SelectFromModel(rf clf2)
  sfm.fit(X train sc, y train)
  sfm.transform(X train sc)
  X train sc df=pd.DataFrame(X train sc,columns=X.columns)
 X_train_sc_df=X_train_sc_df[X_train_sc_df.columns[sfm.get_support()]]
 X test sc df=pd.DataFrame(X test sc,columns=X.columns)
 X test sc df=X test sc df[X test sc df.columns[sfm.qet support()]]
  rf clf2=sfm.estimator
                                                                        In [41]:
from sklearn.model_selection import GridSearchCV
# Defining the parameter grid
param grid = {
    'n estimators': [300,400],
    'max_depth': [None, 10, 20],
    'min_samples_split': [2, 5,10],
    'min samples leaf': [1, 2, 4],
}
# Initialize GridSearchCV
```

```
grid_search = GridSearchCV(estimator=rf_clf2, param_grid=param_grid, cv=3,
scoring='accuracy', verbose=2, n_jobs=-1)

# Fit GridSearchCV
grid_search.fit(X_train_sc_df, y_train)

# Get the best parameters
best_params = grid_search.best_params_
print("Best Parameters:", best_params)

Fitting 3 folds for each of 54 candidates, totalling 162 fits
Best Parameters: {'max_depth': None, 'min_samples_leaf': 1, 'min_samples_sp
lit': 2, 'n_estimators': 300}

In [42]:
# The best Random Forest Model
best rf clf=grid search.best estimator
```

evaluate_model(best_rf_clf, X_train_sc_df, X_test_sc_df)

Out[43]

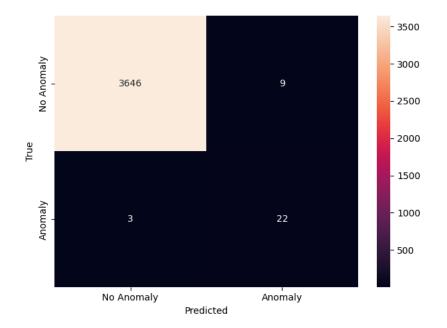
In [43]:

Binary Classification - Predictictive Maintenance

Accuracy: 0.9967391304347826

Classification Report:

	precision	recall	f1-score	support
No Anomaly	1.00	1.00	1.00	3655
Anomaly	0.71	0.88	0.79	25
accuracy			1.00	3680
macro avg	0.85	0.94	0.89	3680
weighted avg	1.00	1.00	1.00	3680



As we can see that for our tuned Random Forest model, the accuracy, F1 score has increased compared to the untuned model.

Saving the model as a pickle file

```
import pickle

In [44]:

import pickle

In [45]:

with open('/content/random_forest_model.pkl', 'wb') as file:
    pickle.dump(best_rf_clf, file)

In [46]:

from google.colab import files

# Download the file
files.download('/content/random_forest_model.pkl')
```

Using Deep Learning

```
In [47]:
from tensorflow.keras.models import Sequential
from tensorflow.keras.layers import Dense, Dropout

In [48]:
model=Sequential()
model.add(Dense(128,activation='relu',input_shape=(59,)))
model.add(Dropout(0.5))
model.add(Dense(64,activation='relu'))
model.add(Dropout(0.5))
model.add(Dense(1,activation='relu'))
model.add(Dense(1,activation='sigmoid'))

model.compile(optimizer='adam',loss='binary_crossentropy',metrics=['accurac y'])
```

```
Epoch 5/20
914/914 [============== ] - 2s 2ms/step - loss: 0.0318 - acc
uracy: 0.9903 - val loss: 0.8399 - val accuracy: 0.9913
Epoch 6/20
914/914 [============= ] - 2s 3ms/step - loss: 0.0267 - acc
uracy: 0.9921 - val loss: 0.9352 - val_accuracy: 0.9908
Epoch 7/20
914/914 [============= ] - 3s 3ms/step - loss: 0.0246 - acc
uracy: 0.9926 - val loss: 1.0304 - val accuracy: 0.9918
Epoch 8/20
914/914 [============= ] - 4s 4ms/step - loss: 0.0185 - acc
uracy: 0.9945 - val loss: 1.1935 - val accuracy: 0.9867
Epoch 9/20
914/914 [============ ] - 3s 3ms/step - loss: 0.0204 - acc
uracy: 0.9934 - val loss: 1.1871 - val accuracy: 0.9927
Epoch 10/20
914/914 [============== ] - 3s 3ms/step - loss: 0.0160 - acc
uracy: 0.9952 - val loss: 1.3342 - val accuracy: 0.9932
Epoch 11/20
uracy: 0.9952 - val_loss: 1.5859 - val_accuracy: 0.9935
Epoch 12/20
914/914 [============ ] - 3s 3ms/step - loss: 0.0166 - acc
uracy: 0.9956 - val loss: 1.5284 - val_accuracy: 0.9927
Epoch 13/20
914/914 [============= ] - 4s 4ms/step - loss: 0.0151 - acc
uracy: 0.9961 - val loss: 1.6516 - val accuracy: 0.9932
Epoch 14/20
914/914 [============= ] - 3s 3ms/step - loss: 0.0140 - acc
uracy: 0.9961 - val loss: 1.6380 - val accuracy: 0.9940
Epoch 15/20
914/914 [============= ] - 2s 3ms/step - loss: 0.0112 - acc
uracy: 0.9969 - val loss: 1.8362 - val accuracy: 0.9937
Epoch 16/20
uracy: 0.9953 - val_loss: 1.8580 - val_accuracy: 0.9924
Epoch 17/20
uracy: 0.9963 - val loss: 1.9435 - val accuracy: 0.9929
Epoch 18/20
uracy: 0.9965 - val_loss: 2.0000 - val_accuracy: 0.9940
Epoch 19/20
uracy: 0.9962 - val loss: 1.9488 - val accuracy: 0.9929
Epoch 20/20
914/914 [============== ] - 2s 3ms/step - loss: 0.0108 - acc
uracy: 0.9970 - val loss: 1.8913 - val accuracy: 0.9935
                                                     Out[49]:
<keras.src.callbacks.History at 0x7c809023a6b0>
```

```
In [50]:
loss,accuracy=model.evaluate(X test sc,y test)
                                                           Out[50]
uracy: 0.9937
                                                             In [51]:
threshold = 0.5
y pred prob=model.predict(X test sc)
y_pred_classes = (y_pred_prob >= threshold).astype(int)
115/115 [=========== ] - 0s lms/step
                                                             In [52]:
print("Predicted probabilities:", y_pred_prob[:10].flatten())
print("Predicted classes:", y pred classes[:10].flatten())
                                                          Out[52]
Predicted probabilities: [1.2075962e-25 6.9271626e-25 1.4649272e-07 8.66433
89e-08 6.5011552e-09
1.6045803e-30 0.0000000e+00 2.7664161e-20 2.7629006e-15 2.2056058e-19]
Predicted classes: [0 0 0 0 0 0 0 0 0]
                                                             In [53]:
print("========="")
print("Binary Classification - Predictictive Maintenance")
print("======="")
print("Accuracy:", accuracy score(y test, y pred classes))
print("Classification Report:\n", classification report(y test,
y pred classes,
                                                 target names=['No
Anomaly','Anomaly']))
sns.heatmap(confusion matrix(y test, y pred classes),annot=True,fmt='g',
     yticklabels=['No Anomaly','Anomaly'],xticklabels=['No
Anomaly','Anomaly'])
plt.xlabel('Predicted')
plt.ylabel('True')
plt.tight layout()
plt.show()
                                                           Out[53]
Binary Classification - Predictictive Maintenance
_____
Accuracy: 0.9923913043478261
Classification Report:
             precision recall f1-score
                                         support
                                  1.00
                                           3655
 No Anomaly
                1.00
                         1.00
                0.43
                         0.40
                                  0.42
    Anomaly
                                            25
   accuracy
                                  0.99
                                           3680
                0.72
                        0.70
                                  0.71
                                          3680
  macro avg
```

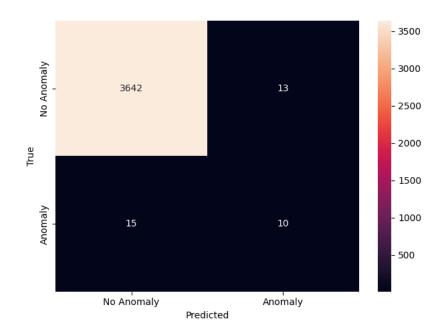
weighted avg

0.99

0.99

0.99

3680



Deep learning model not performing as good as our Random Forest model.