Extra Credit Link: https://youtu.be/AYfcrX7Ya2U

```
from ucimlrepo import fetch_ucirepo

cervical_cancer_risk_factors = fetch_ucirepo(id=383)

X = cervical_cancer_risk_factors.data.features
y = cervical_cancer_risk_factors.data.targets

print(cervical_cancer_risk_factors.metadata)

print(cervical_cancer_risk_factors.variables)
```

```
from ucimlrepo import fetch ucirepo
import pandas as pd
from sklearn.model selection import train test split
from sklearn.preprocessing import StandardScaler, LabelEncoder
from imblearn.over sampling import SMOTE
from sklearn.ensemble import RandomForestClassifier
from sklearn.metrics import classification report, accuracy score
cervical cancer risk factors = fetch ucirepo(id=383)
X = cervical cancer risk factors.data.features
y = cervical cancer risk factors.data.targets
data = pd.concat([X, y], axis=1)
print("Missing values in the dataset:")
print(data.isnull().sum())
y = X['Biopsy']
X = X.drop(columns=['Biopsy'])
columns to drop = X.columns[X.isnull().mean() > 0.5]
X.drop(columns=columns to drop, inplace=True)
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= X.fillna(X.median())
categorical columns = X.select dtypes(include=['object']).columns
for col in categorical columns:
    X[col] = X[col].fillna(X[col].mode()[0])
label encoder = LabelEncoder()
for col in ['Smokes', 'Hormonal Contraceptives', 'IUD']:
    if col in X.columns:
        X[col] = label encoder.fit transform(X[col])
scaler = StandardScaler()
X scaled = scaler.fit transform(X)
X train, X test, y train, y test = train_test_split(X_scaled, y,
test size=0.2, random state=42)
smote = SMOTE(random state=42)
X train smote, y train smote = smote.fit resample(X train, y train)
print(f"Training data shape (after SMOTE): {X_train_smote.shape}")
print(f"Testing data shape: {X test.shape}")
model = RandomForestClassifier(random state=42)
model.fit(X train smote, y train smote)
y pred = model.predict(X test)
print("\nModel Evaluation:")
print(f"Accuracy: {accuracy score(y test, y pred)}")
print("Classification Report:")
print(classification report(y test, y pred))
```

```
from sklearn.model selection import GridSearchCV
from sklearn.ensemble import RandomForestClassifier
from sklearn.metrics import classification report, accuracy score
param_grid = {
    'min samples leaf': [1, 2, 4],
    'bootstrap': [True, False],
grid search rf = GridSearchCV(
    RandomForestClassifier(random state=42),
   param grid=param grid,
   scoring='accuracy',
grid search rf.fit(X train smote, y train smote)
print("Best hyperparameters for Random Forest:")
print(grid search rf.best params )
best rf model = grid search rf.best estimator
y pred rf = best rf model.predict(X test)
print("Random Forest with Best Hyperparameters Evaluation:")
print(f"Accuracy: {accuracy score(y test, y pred rf)}")
print("Classification Report:")
print(classification report(y test, y pred rf))
from xgboost import XGBClassifier
from sklearn.metrics import classification report, accuracy score
```

```
class 0 count = sum(y train smote == 0)
class 1 count = sum(y train smote == 1)
xgb_model = XGBClassifier(scale_pos_weight=class_0_count/class_1_count,
random state=42)
xgb model.fit(X train smote, y train smote)
y_pred_xgb = xgb_model.predict(X test)
print("XGBoost Model Evaluation:")
print(f"Accuracy: {accuracy score(y test, y pred xgb)}")
print("Classification Report:")
print(classification report(y test, y pred xgb))
from sklearn.model selection import GridSearchCV
param_grid = {
grid search = GridSearchCV(XGBClassifier(random state=42), param grid,
scoring='accuracy', cv=3)
grid search.fit(X train smote, y train smote)
print("Best parameters found: ", grid search.best params )
xgb model best = XGBClassifier(
    learning rate=0.05,
    \max depth=3,
```

```
n estimators=200,
    subsample=1.0,
    scale pos weight=class 0 count/class 1 count,
    random state=42
xgb model best.fit(X train smote, y train smote)
y pred xgb best = xgb model best.predict(X test)
print("XGBoost Model Evaluation with Best Parameters:")
print(f"Accuracy: {accuracy score(y test, y pred xgb best)}")
print("Classification Report:")
print(classification report(y test, y pred xgb best))
y prob = xgb model best.predict proba(X test)[:, 1]
threshold = 0.3
y pred threshold = (y_prob > threshold).astype(int)
print("XGBoost Model with Adjusted Threshold Evaluation:")
print(f"Accuracy: {accuracy score(y test, y pred threshold)}")
print("Classification Report:")
print(classification report(y test, y pred threshold))
from imblearn.over sampling import ADASYN
from sklearn.metrics import classification report, accuracy score
adasyn = ADASYN(sampling strategy=0.7, random state=42)
X train adasyn, y train adasyn = adasyn.fit resample(X train, y train)
print("Class distribution after ADASYN:")
print(pd.Series(y train adasyn).value counts())
from xgboost import XGBClassifier
```

```
xgb model adasyn = XGBClassifier(
   max depth=3,
    subsample=1.0,
    scale pos weight=class 0 count/class_1_count,
xgb model adasyn.fit(X train adasyn, y train adasyn)
y prob adasyn = xgb model adasyn.predict proba(X test)[:, 1]
threshold = 0.3
y pred threshold adasyn = (y prob adasyn > threshold).astype(int)
print("XGBoost Model with ADASYN and Further Adjusted Threshold
Evaluation:")
print(f"Accuracy: {accuracy score(y test, y pred threshold adasyn)}")
print("Classification Report:")
print(classification_report(y_test, y_pred_threshold_adasyn))
import matplotlib.pyplot as plt
import numpy as np
models = ['XGBoost', 'Random Forest', 'ADASYN + XGBoost']
precision class 1 = [0.53, 0.60, 0.56]
recall class 1 = [0.73, 0.55, 0.82]
f1 class 1 = [0.62, 0.57, 0.67]
fig, ax = plt.subplots(figsize=(10, 6))
ax.plot(models, precision class 1, label='Precision (Class 1)',
color='r', marker='o')
ax.plot(models, recall class 1, label='Recall (Class 1)', color='y',
marker='o')
ax.plot(models, f1 class 1, label='F1-score (Class 1)', color='m',
marker='o')
ax.set xlabel('Models')
ax.set_ylabel('Scores')
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ax.set_title('Performance Comparison of Models for Cancer Prediction
  (Class 1)')
ax.legend()
plt.tight_layout()
plt.show()
```