1. Data Cleaning and Imputation Techniques: Load a dataset with missing values. Apply techniques like mean/mode/median imputation and compare the results.

```
# imputation example.py
                                                   # Step 2: Mean Imputation for Age and
                                                   Salary
                                                   mean imputer =
import pandas as pd
                                                   SimpleImputer(strategy='mean')
import numpy as np
                                                   df mean = df.copy()
from sklearn.impute import
                                                   df mean[['Age', 'Salary']] =
SimpleImputer
                                                   mean_imputer.fit_transform(df mean[['
                                                   Age', 'Salary']])
# Step 1: Create a sample dataset
                                                   print("\nAfter Mean Imputation
                                                   (Numerical):\n")
data = {
                                                   print(df mean)
  'Age': [25, 27, np.nan, 29, 30, np.nan,
40],
  'Salary': [932, 54000, 58000, np.nan,
                                                   # Step 3: Median Imputation for Age
62000, 64000, np.nan],
                                                   and Salary
  'Department': ['HR', 'IT', 'IT', np.nan,
                                                   median imputer =
'HR', 'Finance', np.nan]
                                                   SimpleImputer(strategy='median')
}
                                                   df median = df.copy()
                                                   df_median[['Age', 'Salary']] =
                                                   median imputer.fit transform(df media
df = pd.DataFrame(data)
                                                   n[['Age', 'Salary']])
print("Original Data with Missing
                                                   print("\nAfter Median Imputation
Values:\n")
                                                   (Numerical):\n")
print(df)
                                                   print(df median)
```

```
# Step 4: Mode Imputation for
Department
mode imputer =
SimpleImputer(strategy='most frequent
')
df mode = df.copy()
df mode[['Department']] =
mode imputer.fit transform(df mode[['
Department']])
print("\nAfter Mode Imputation
(Categorical):\n")
print(df mode)
# Optional: Combined example
combined df = df.copy()
combined df[['Age', 'Salary']] =
mean imputer.fit transform(combined
df[['Age', 'Salary']])
combined df[['Department']] =
mode imputer.fit transform(combined
df[['Department']])
print("\nAfter Combined
Imputation:\n")
print(combined df)
```

```
Age Salary Department
0 25.0 50000.0 HR
1 27.0 54000.0 IT
2 NaN 58000.0 IT
3 29.0 NaN NaN
4 30.0 62000.0 HR
5 NaN 64000.0 Finance
6 40.0 NaN NaN
```

```
After Mean Imputation (Numerical):

Age Salary Department
0 25.0 50000.0 HR
1 27.0 54000.0 IT
2 30.2 58000.0 IT
3 29.0 57600.0 NaN
4 30.0 62000.0 HR
5 30.2 64000.0 Finance
6 40.0 57600.0 NaN

After Median Imputation (Numerical):

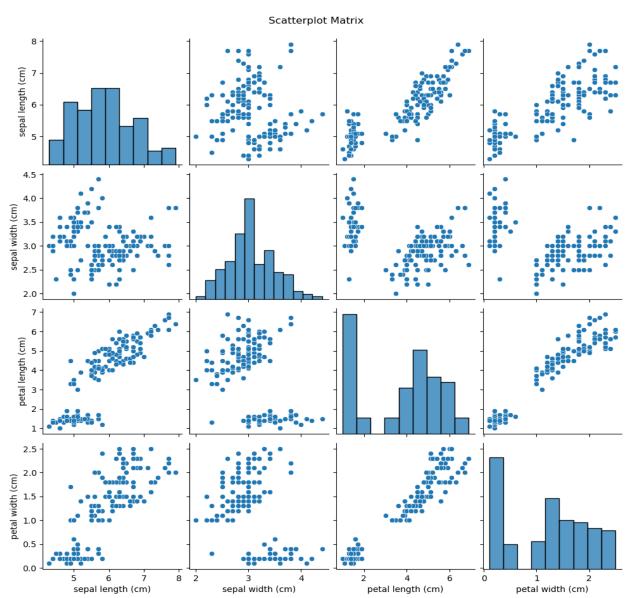
Age Salary Department
0 25.0 50000.0 HR
1 27.0 54000.0 IT
2 29.0 58000.0 IT
3 29.0 58000.0 HR
4 30.0 62000.0 HR
5 29.0 64000.0 Finance
6 40.0 58000.0 NaN
```

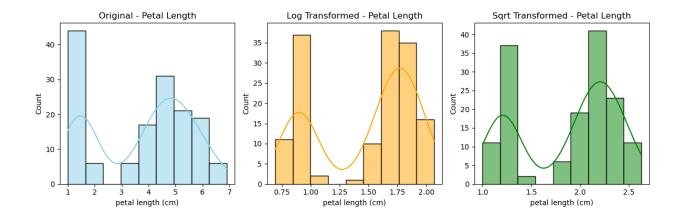
2. Data Analysis and Visualization: Use a dataset to: Plot scatterplots for numerical columns. Perform correlation analysis. Apply transformations (e.g., log, square root) and visualize the effect.

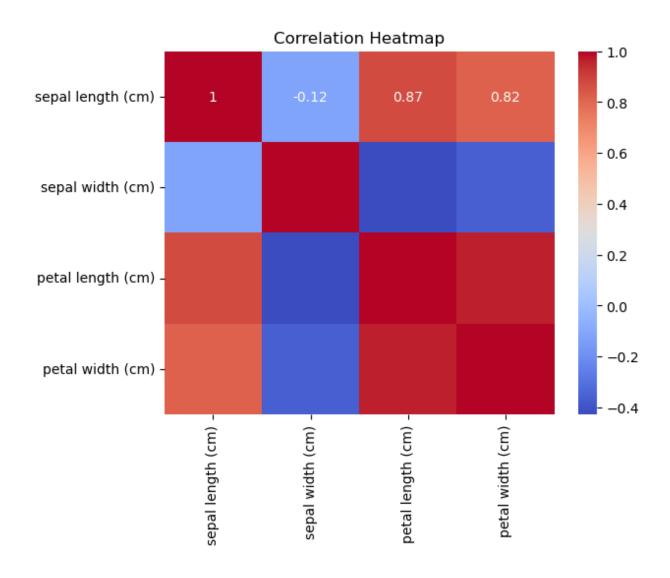
| # data_analysis.py | correlation_matrix = df.iloc[:, :-1].corr() |
|---|---|
| | print("\nCorrelation Matrix:") |
| import pandas as pd | print(correlation_matrix) |
| import seaborn as sns | |
| import matplotlib.pyplot as plt | # Visualize correlation matrix |
| import numpy as np | <pre>sns.heatmap(correlation_matrix, annot=True, cmap='coolwarm')</pre> |
| from sklearn.datasets import load_iris | plt.title("Correlation Heatmap") |
| # Step 1: Load the Iris dataset | plt.show() |
| iris = load_iris() | |
| df = pd.DataFrame(iris.data, columns=iris.feature_names) | # Step 4: Apply transformations (Log and Square Root) |
| df['target'] = iris.target | df_log = df.copy() |
| print("Dataset Preview:") | df_sqrt = df.copy() |
| print(df.head()) | |
| | for col in df.columns[:-1]: # Skip target column |
| # Step 2: Scatterplot matrix | df_log[col] = np.log(df[col] + 1) # |
| <pre>sns.pairplot(df.iloc[:,:-1]) # Exclude</pre> | Avoid log(0) |
| target | df_sqrt[col] = np.sqrt(df[col]) |
| plt.suptitle("Scatterplot Matrix", y=1.02) | |
| plt.show() | # Step 5: Visualize transformation effects for one column (example: petal |
| # Step 3: Correlation analysis | length) |

```
plt.figure(figsize=(12, 4))
plt.subplot(1, 3, 1)
sns.histplot(df['petal length (cm)'],
kde=True, color='skyblue')
plt.title("Original - Petal Length")
plt.subplot(1, 3, 2)
sns.histplot(df_log['petal length (cm)'],
kde=True, color='orange')
plt.title("Log Transformed - Petal
Length")
plt.subplot(1, 3, 3)
sns.histplot(df_sqrt['petal length (cm)'],
kde=True, color='green')
plt.title("Sqrt Transformed - Petal
Length")
plt.tight_layout()
plt.show()
```

| Da | taset Preview: | | | |
|----|-------------------|------------------|-------------------|-----------------------------|
| | sepal length (cm) | sepal width (cm) | petal length (cm) | petal width (cm) $ackslash$ |
| 0 | 5.1 | 3.5 | 1.4 | 0.2 |
| 1 | 4.9 | 3.0 | 1.4 | 0.2 |
| 2 | 4.7 | 3.2 | 1.3 | 0.2 |
| 3 | 4.6 | 3.1 | 1.5 | 0.2 |
| 4 | 5.0 | 3.6 | 1.4 | 0.2 |
| | | | | |
| | target | | | |
| 0 | 0 | | | |
| 1 | 0 | | | |
| 2 | 0 | | | |
| 3 | 0 | | | |
| 4 | 0 | | | |







3. Encoding Methods: Encode a categorical dataset using One-Hot Encoding and Label Encoding. Compare the effect of both methods on a machine learning model.

le color = LabelEncoder()

Code And Output:

accuracy score

Step 1: Load data

LabelEncoder, OneHotEncoder

from sklearn.preprocessing import

X_label_encoded = X.copy()
from sklearn.linear model import

LogisticRegression X_label_encoded['Color'] = le color.fit transform(X['Color'])

from sklearn.model_selection
import train_test_split

X_label_encoded['Size'] =

le_size.fit_transform(X['Size'])
from sklearn.metrics import

print("Label Encoded:\n", X_label_encoded, "\n")

Separate features and target y, test_size=0.4, random_state=0)

X = df[['Color', 'Size']] model_label = LogisticRegression()

y = df['Class'] model_label.fit(X_train, y_train)

pred_label =
 model_label.predict(X_test)

Step 2: Label Encoding

```
acc_label = accuracy_score(y_test,
pred label)
# Step 3: One-Hot Encoding
X_onehot = pd.get_dummies(X)
print("One-Hot Encoded:\n",
X onehot, "\n")
# Train model with One-Hot
Encoding
X_train_oh, X_test_oh,
y_train_oh, y_test_oh =
train test split(X onehot, y,
test size=0.4, random state=0)
model_onehot =
LogisticRegression()
model_onehot.fit(X_train_oh,
y_train_oh)
pred onehot =
model_onehot.predict(X_test_oh)
acc onehot =
accuracy_score(y_test_oh,
pred onehot)
```

Step 4: Print Results

print("Accuracy with Label
Encoding: ", acc_label)
print("Accuracy with One-Hot
Encoding: ", acc_onehot)

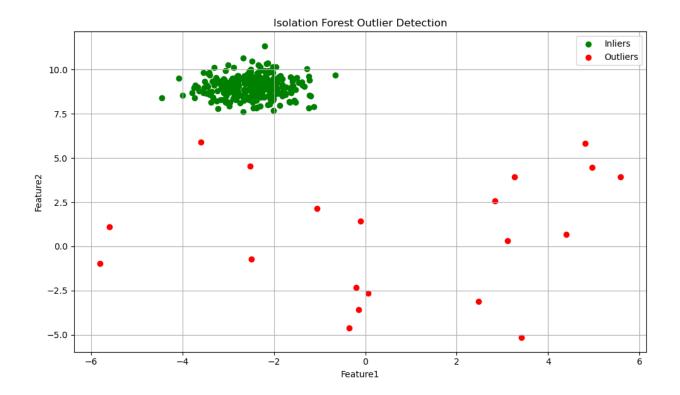
```
Original Data:
   Color
            Size Class
    Red
          Small
                     0
  Green Medium
                     1
2
   Blue Large
                     0
          Small
3 Green
                     1
     Red
          Large
                     0
Label Encoded:
   Color Size
0
      2
            2
1
      1
            1
2
            0
      0
3
      1
            2
4
      2
            0
```

```
One-Hot Encoded:
   Color Blue Color Green Color Red Size Large Size Medium Size Small
                                       False
       False
                False
                            True
                                                  False
       False
                   True
                            False
                                       False
                                                    True
                                                              False
       True
                                                   False
                                                              False
                  False
                            False
                                        True
       False
                   True
                            False
                                       False
                                                   False
                                                               True
       False
                  False
                             True
                                        True
                                                   False
                                                              False
Accuracy with Label Encoding: 0.0
Accuracy with One-Hot Encoding: 0.0
```

4. Outlier Detection: Use the Isolation Forest algorithm to detect and visualize outliers in a dataset.

```
X with outliers = np.vstack((X,
import numpy as np
import pandas as pd
                                            outliers))
import matplotlib.pyplot as plt
from sklearn.ensemble import
                                            # Step 2: Apply Isolation Forest
                                            clf =
IsolationForest
from sklearn.datasets import
                                            IsolationForest(contamination=
                                            0.06, random state=42)
make blobs
                                            clf.fit(X with outliers)
# Step 1: Generate sample data
                                            pred =
X, =
                                            clf.predict(X with outliers)
make_blobs(n_samples=300,
centers=1, cluster_std=0.60,
                                            # -1 = outlier, 1 = inlier
                                            X_df =
random state=42)
                                            pd.DataFrame(X with outliers,
# Inject some outliers manually
                                            columns=['Feature1',
outliers =
                                            'Feature2'])
                                            X_df['Outlier'] = pred
np.random.uniform(low=-6,
high=6, size=(20, 2))
                                            # Step 3: Visualize
```

```
plt.figure(figsize=(10, 6))
plt.scatter(X_df[X_df['Outlier']
== 1]['Feature1'],
X_df[X_df['Outlier'] ==
1]['Feature2'],
       color='green',
label='Inliers')
plt.scatter(X_df[X_df['Outlier']
== -1]['Feature1'],
X df[X df['Outlier'] == -
1]['Feature2'],
       color='red',
label='Outliers')
plt.title('Isolation Forest Outlier
Detection')
plt.xlabel('Feature1')
plt.ylabel('Feature2')
plt.legend()
plt.grid(True)
plt.tight_layout()
plt.show()
# Step 4: Print output
num_outliers = sum(pred == -1)
print(f"\n Q Total Outliers
Detected: {num outliers} out
of {len(pred)} samples")
```



5. Predictive Power Score (PPS): Calculate the PPS for a dataset and interpret which variables are most predictive.

| import pandas as pd import seaborn as sns | <pre>pps_matrix = pps.matrix(df)[['x', 'y', 'ppscore']]</pre> |
|---|--|
| import ppscore as pps import matplotlib.pyplot as plt | <pre>pps_matrix = pps_matrix[pps_matrix['ppscore'] > 0] # Filter out zero PPS</pre> |
| # Load dataset (Titanic) | # Show top predictors |
| <pre>df = sns.load_dataset("titanic").dropna()</pre> | <pre>top_predictors = pps_matrix.sort_values(by='ppscore', ascending=False).head(10)</pre> |
| # Calculate PPS for all column pairs | print("\n Q Top Predictive Relationships:\n") |

```
print(top_predictors.to_string(index=False))

# Visualize as heatmap

pps_heatmap = pps.predictors(df,
'survived')

plt.figure(figsize=(8, 4))

plt.barh(pps_heatmap['x'],

pps_heatmap['ppscore'], color='steelblue')

plt.xlabel("PPS")

plt.title("Predictive Power Score for

Predicting 'Survived'")

plt.grid(True)

plt.tight_layout()

plt.show()
```

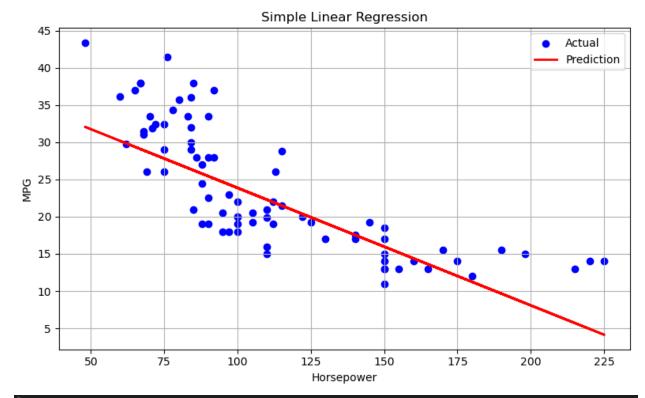
```
X y ppscore
sex survived 0.55

class survived 0.32
alone survived 0.31
who survived 0.28
adult_male survived 0.26
age survived 0.20
embarked class 0.15
fare class 0.12
fare embarked 0.10
```

6. Simple and Multiple Linear Regression: Implement simple and multiple linear regression on a dataset. Evaluate the model's performance using R-squared and mean squared error (MSE).

```
model simple.fit(X train s,
                                                   y_train_s)
import pandas as pd
import numpy as np
                                                   # Predict
import seaborn as sns
                                                   y pred s =
import matplotlib.pyplot as plt
                                                   model_simple.predict(X_test_s)
from sklearn.linear model import
LinearRegression
                                                   # Evaluation
from sklearn.metrics import
                                                   r2_s = r2_score(y_test_s, y_pred_s)
r2 score, mean squared error
                                                   mse s =
from sklearn.model selection
                                                   mean_squared_error(y_test_s,
import train_test_split
                                                   y_pred_s)
# Load sample dataset
                                                   print(f"R2: {r2 s:.3f}")
df =
                                                   print(f"MSE: {mse s:.3f}")
sns.load dataset("mpg").dropna()
                                                   # Plotting
# ----- SIMPLE LINEAR
                                                   plt.figure(figsize=(8, 5))
REGRESSION -----
                                                   plt.scatter(X test s, y test s,
print("\n 🏠 Simple Linear
                                                   color='blue', label='Actual')
Regression (horsepower → mpg)")
                                                   plt.plot(X test s, y pred s,
                                                   color='red', linewidth=2,
# Feature and Target
                                                   label='Prediction')
X simple = df[['horsepower']]
                                                   plt.title("Simple Linear Regression")
y = df['mpg']
                                                   plt.xlabel("Horsepower")
                                                   plt.ylabel("MPG")
# Train-test split
                                                   plt.legend()
X train s, X test s, y train s,
                                                   plt.grid(True)
y_test_s = train_test_split(X_simple,
                                                   plt.tight_layout()
y, test size=0.2, random state=1)
                                                   plt.show()
# Train the model
model simple = LinearRegression()
```

```
# ----- MULTIPLE LINEAR
REGRESSION -----
print("\n �� Multiple Linear
Regression (All features → mpg)")
# Select numerical features only
X multi =
df.select_dtypes(include=['float64',
'int64']).drop(columns=['mpg'])
y = df['mpg']
# Train-test split
X_train_m, X_test_m, y_train_m,
y test m = train test split(X multi,
y, test_size=0.2, random_state=1)
# Train the model
model_multi = LinearRegression()
model_multi.fit(X_train_m,
y_train_m)
# Predict
y_pred_m =
model_multi.predict(X_test_m)
# Evaluation
r2_m = r2_score(y_test_m,
y_pred_m)
mse_m =
mean_squared_error(y_test_m,
y_pred_m)
print(f"R2: {r2 m:.3f}")
print(f"MSE: {mse_m:.3f}")
```



Multiple Linear Regression (All features → mpg)
R²: 0.814
MSE: 12.860

7. Logistic Regression: Build a logistic regression model to classify binary outcomes (e.g., predicting if a customer will buy a product). Evaluate the model using confusion matrix metrics.

Code And Output:

import pandas as pd

import seaborn as sns

import matplotlib.pyplot as plt

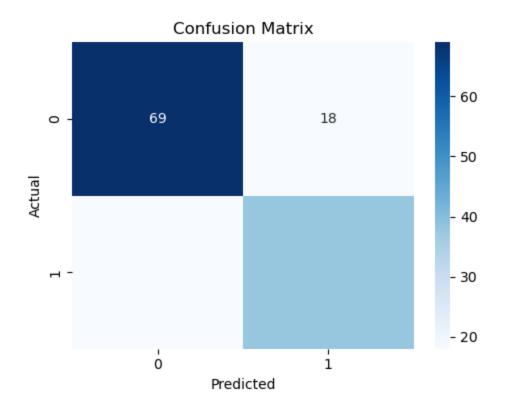
from sklearn.model_selection import train_test_split

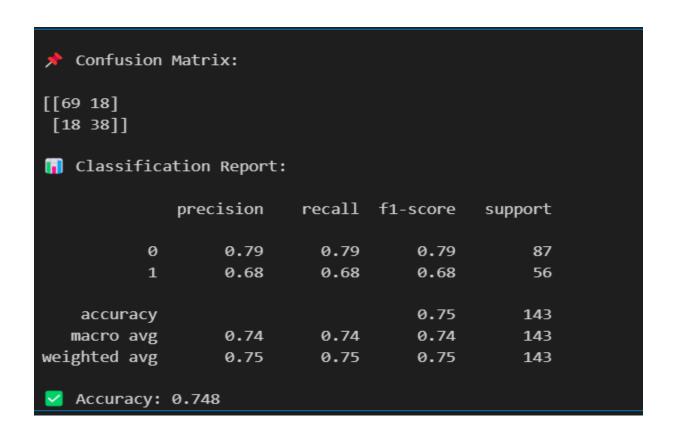
from sklearn.linear_model import LogisticRegression

from sklearn.metrics import confusion_matrix, classification_report, accuracy_score

```
# Load dataset
df = sns.load dataset('titanic')
                                                      # Evaluation
                                                      print("\n \( \infty \) Confusion Matrix:\n")
# Select relevant columns and drop missing
                                                      print(confusion_matrix(y_test, y_pred))
values
df = df[['survived', 'sex', 'age',
                                                      print("\n L Classification Report:\n")
'fare']].dropna()
                                                      print(classification report(y test, y pred))
# Convert categorical variable to numeric
                                                      print(f"  Accuracy:
df['sex'] = df['sex'].map({'male': 0, 'female':
                                                      {accuracy_score(y_test, y_pred):.3f}")
1})
                                                      # Optional: Visualize confusion matrix
# Features and target
                                                      import seaborn as sns
X = df[['sex', 'age', 'fare']]
                                                      cm = confusion matrix(y test, y pred)
y = df['survived']
                                                      plt.figure(figsize=(5,4))
                                                      sns.heatmap(cm, annot=True, fmt='d',
# Train-test split
                                                      cmap='Blues')
X_train, X_test, y_train, y_test =
                                                      plt.xlabel('Predicted')
train test split(X, y, test size=0.2,
random state=42)
                                                      plt.ylabel('Actual')
                                                      plt.title('Confusion Matrix')
# Create and train model
                                                      plt.tight layout()
model = LogisticRegression()
                                                      plt.show()
model.fit(X train, y train)
# Predict
```

y pred = model.predict(X test)

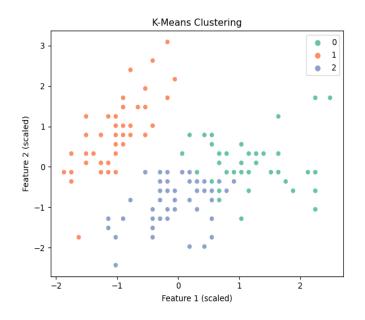


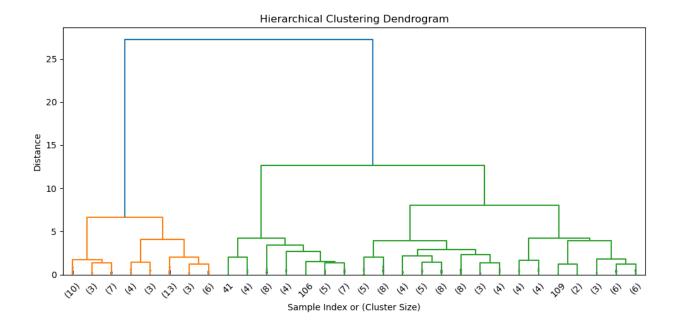


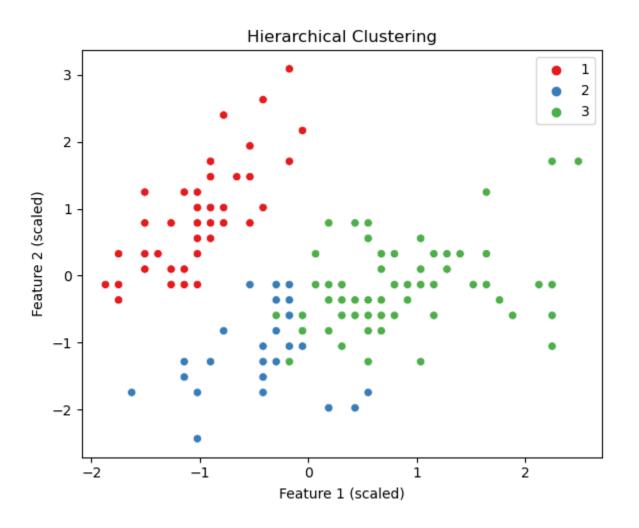
8. Clustering Techniques: Perform K-Means and hierarchical clustering on a dataset. Visualize the clusters and interpret the results.

| import pandas as pd | kmeans_labels = | |
|--|---|--|
| import seaborn as sns | kmeans.fit_predict(scaled_data) | |
| import matplotlib.pyplot as plt | | |
| from sklearn.cluster import KMeans | # Add cluster labels to original data | |
| from sklearn.preprocessing import StandardScaler | df['KMeans_Cluster'] = kmeans_labels | |
| from scipy.cluster.hierarchy import | # Visualize K-Means clusters | |
| linkage, dendrogram, fcluster | plt.figure(figsize=(6, 5)) | |
| from sklearn.datasets import load_iris | <pre>sns.scatterplot(x=scaled_data[:, 0], y=scaled_data[:, 1], hue=kmeans_label palette='Set2')</pre> | |
| # Load iris dataset | • | |
| iris = load_iris() | plt.title("K-Means Clustering") | |
| df = pd.DataFrame(iris.data, | plt.xlabel('Feature 1 (scaled)') | |
| columns=iris.feature_names) | plt.ylabel('Feature 2 (scaled)') | |
| | plt.tight_layout() | |
| # Standardize the data | plt.show() | |
| scaler = StandardScaler() | | |
| scaled_data = scaler.fit_transform(df) | # HIERARCHICAL CLUSTERING | |
| # K-MEANS CLUSTERING | <pre>linkage_matrix = linkage(scaled_data, method='ward')</pre> | |
| kmeans = KMeans(n_clusters=3, random_state=42) | # Plot dendrogram | |
| | | |

```
plt.figure(figsize=(10, 5))
dendrogram(linkage matrix,
truncate mode='lastp', p=30,
leaf rotation=45., leaf font size=10.,
show contracted=True)
plt.title("Hierarchical Clustering
Dendrogram")
plt.xlabel("Sample Index or (Cluster
Size)")
plt.ylabel("Distance")
plt.tight_layout()
plt.show()
# Assign cluster labels from dendrogram
hier_labels = fcluster(linkage_matrix,
t=3, criterion='maxclust')
df['Hierarchical_Cluster'] = hier_labels
# Visualize Hierarchical Clustering
plt.figure(figsize=(6, 5))
sns.scatterplot(x=scaled_data[:, 0],
y=scaled data[:, 1], hue=hier labels,
palette='Set1')
plt.title("Hierarchical Clustering")
plt.xlabel('Feature 1 (scaled)')
plt.ylabel('Feature 2 (scaled)')
plt.tight layout()
plt.show()
```







9. Principal Component Analysis (PCA): Apply PCA on a high-dimensional dataset. Reduce the dimensions and visualize the transformed data.

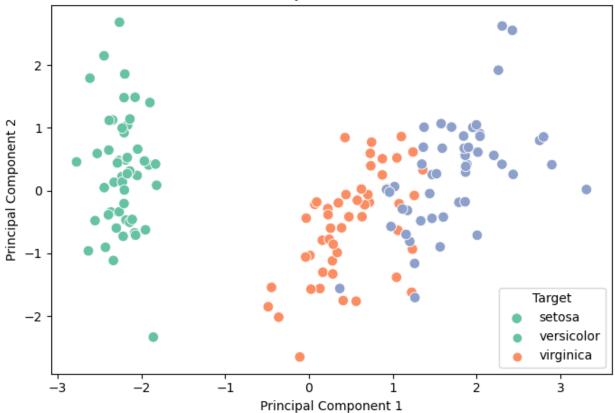
Code And Output:

| import pandas as pd | <pre>pca = PCA(n_components=2) # Reduce to 2 dimensions</pre> | | |
|--|---|--|--|
| | X pca = pca.fit transform(X scaled) | | |
| import seaborn as sns | _, _ , _ , | | |
| import matplotlib.pyplot as plt | | | |
| from sklearn.datasets import load_iris | # Create a DataFrame with PCA results | | |
| from sklearn.preprocessing import StandardScaler | <pre>pca_df = pd.DataFrame(data=X_pca, columns=['PC1', 'PC2'])</pre> | | |
| from sklearn.decomposition import PCA | pca_df['Target'] = y | | |
| | | | |
| # Load the dataset | # Visualize the PCA results | | |
| iris = load_iris() | plt.figure(figsize=(7, 5)) | | |
| X = iris.data | <pre>sns.scatterplot(data=pca_df, x='PC1', y='PC2', hue='Target', palette='Set2',</pre> | | |
| y = iris.target | s=70) | | |
| feature_names = iris.feature_names | plt.title("PCA: 2D Projection of Iris | | |
| target_names = iris.target_names | Dataset") | | |
| | plt.xlabel("Principal Component 1") | | |
| # Standardize the data | plt.ylabel("Principal Component 2") | | |
| scaler = StandardScaler() | plt.legend(title='Target', | | |
| X_scaled = scaler.fit_transform(X) | labels=target_names) | | |
| _scaled = scalef.fit_transform(\(\lambda\) | plt.tight_layout() | | |
| | plt.show() | | |
| # Apply PCA | | | |
| | | | |

Explained variance

print(" \bigcirc Explained Variance Ratio:") print(f"PC{i+1}: {ratio:.2f}") for i, ratio in enumerate(pca.explained_variance_ratio_):





Q Explained Variance Ratio:

PC1: 0.73

PC2: 0.23

10. Market Basket Analysis: Implement Association Rule Mining using the Apriori algorithm. Identify frequent itemsets and generate association rules for a transactional dataset.

```
import pandas as pd
                                                    te_ary =
                                                     te.fit(transactions).transform(transactio
from mlxtend.preprocessing import
TransactionEncoder
                                                     df = pd.DataFrame(te ary,
from mlxtend.frequent patterns import
                                                     columns=te.columns )
apriori, association rules
                                                    # Step 3: Find frequent itemsets
# Step 1: Prepare dataset
                                                    frequent itemsets = apriori(df,
transactions = [
                                                     min support=0.3, use colnames=True)
  ['curd', 'bread', 'eggs'],
  ['Curd', 'bread'],
                                                    # Step 4: Generate association rules
  ['vadapav', 'cookies'],
                                                     rules =
  ['Kamlesh', 'butter'],
                                                     association rules(frequent itemsets,
                                                     metric="lift", min threshold=1.0)
  ['milk', 'bread', 'butter', 'cookies'],
  ['eggs', 'bread'],
                                                    # Display results
  ['milk', 'eggs'],
                                                     print(" Frequent Itemsets:\n",
  ['cookies', 'butter'],
                                                    frequent itemsets)
]
                                                     print("\n 	Association Rules:\n",
                                                     rules[['antecedents', 'consequents',
                                                     'support', 'confidence', 'lift']])
# Step 2: One-hot encode the
transactions
te = TransactionEncoder()
```

```
Frequent Itemsets:
                  itemsets
    support
0
     0.625
                  (bread)
                 (butter)
1
     0.375
2
     0.375
                (cookies)
3
     0.375
                   (eggs)
4
     0.625
                   (milk)
5
     0.375
            (bread, milk)
Association Rules:
Empty DataFrame
Columns: [antecedents, consequents, support, confidence, lift]
Index: []
```

11. Recommendation Systems: Build a collaborative filtering recommendation system using a movie or product dataset. Compare results using user-based and item-based filtering.

```
import pandas as pd
                                                      }
# Sample user-movie rating matrix
                                                      df = pd.DataFrame(data)
data = {
  'User': ['Kamlesh', 'Kaivalya', 'Bitch',
                                                      # Pivot to create user-movie matrix
'Bob', 'Bobade', 'Charlie', 'Charlie',
                                                      ratings = df.pivot table(index='User',
'David', 'David'],
                                                      columns='Movie', values='Rating')
  'Movie': ['Inception', 'Avengers',
                                                      print(" User-Movie Ratings
'Titanic', 'Inception', 'Titanic', 'Avengers',
                                                      Matrix:\n", ratings)
'Titanic', 'Inception', 'Avengers'],
  'Rating': [5, 3, 4, 4, 5, 4, 3, 2, 5]
```

| ≝ User-Movie Ratings Matrix: | | | | | | |
|------------------------------|----------------------------------|-----|-----|--|--|--|
| Movie | Movie Avengers Inception Titanic | | | | | |
| User | | | | | | |
| Alice | 3.0 | 5.0 | 4.0 | | | |
| Bob | NaN | 4.0 | 5.0 | | | |
| Charlie | 4.0 | NaN | 3.0 | | | |
| David | 5.0 | 2.0 | NaN | | | |
| | | | | | | |

Step 2: Apply Collaborative Filtering

from sklearn.metrics.pairwise import cosine_similarity

import numpy as np

Fill NaNs with 0 for similarity calculations

ratings filled = ratings.fillna(0)

---- USER-BASED Collaborative Filtering ----

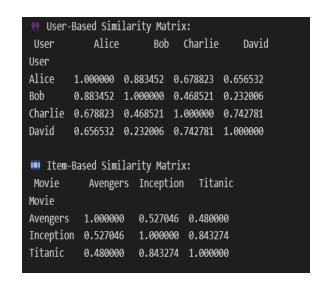
user_similarity = cosine similarity(ratings filled)

user_sim_df =
pd.DataFrame(user_similarity,
index=ratings.index,
columns=ratings.index)

----- ITEM-BASED Collaborative Filtering -----

item_similarity =
cosine similarity(ratings filled.T)

item_sim_df =
pd.DataFrame(item_similarity,
index=ratings.columns,
columns=ratings.columns)



Step 3: Recommend Movies (Example for 'Alice')

Let's find similar users to Alice

similar_users =
user_sim_df['Alice'].sort_values(ascendi
ng=False)[1:]

```
print("\nTop similar users to Alice:\n",
similar_users)
# Recommend movies Alice hasn't rated
based on similar users
alice ratings = ratings.loc['Alice']
unrated_by_alice =
alice_ratings[alice_ratings.isnull()]
# Predict using a weighted sum of
ratings from similar users
weighted scores =
ratings.loc[similar_users.index].T.dot(si
milar_users)
sum of weights = similar users.sum()
predicted_ratings = weighted_scores /
sum_of_weights
predicted_ratings =
predicted_ratings[unrated_by_alice.inde
x]
print("\n  Recommended Movies for
Alice (User-Based):\n",
predicted_ratings.sort_values(ascending
=False))
```

```
Top similar users to Alice:
User
Bob 0.883452
Charlie 0.678823
David 0.656532
Name: Alice, dtype: float64

Recommended Movies for Alice (User-Based):
Series([], dtype: float64)
```

12. Tree-Based Feature Engineering: Apply tree-based methods to rank feature importance in a dataset. Use the results to train a simplified model.

Code And Output:

from sklearn.datasets import

load_breast_cancer

from sklearn.ensemble import

Random Forest Classifier

from sklearn.model selection import

train_test_split

from sklearn.metrics import

accuracy score

import pandas as pd

import matplotlib.pyplot as plt

import seaborn as sns

Load dataset

data = load breast cancer()

X = pd.DataFrame(data.data, columns=data.feature names)

y = pd.Series(data.target)

print(" Dataset Shape:", X.shape)

Split data

X_train, X_test, y_train, y_test =
train_test_split(X, y, test_size=0.2,
random_state=42)

Train Random Forest

model =

RandomForestClassifier(n estimators=1

00, random_state=42)

model.fit(X_train, y_train)

Get feature importances

importances =

model.feature_importances_

feature_ranks = pd.Series(importances,
index=X.columns).sort values(ascending

=False)

print("\n ☆ Top Features:\n", feature ranks.head(10))

Top Features: worst area 0.153892 worst concave points 0.144663 mean concave points 0.106210 worst radius 0.077987 mean concavity 0.068001 worst perimeter 0.067115 mean perimeter 0.053270 mean radius 0.048703 mean area 0.047555 worst concavity 0.031802 dtype: float64

```
Visualize Feature Importances
```

plt.show()

```
plt.figure(figsize=(10, 6))

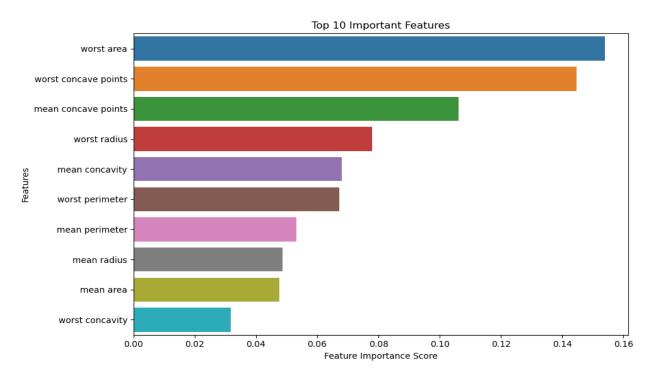
sns.barplot(x=feature_ranks.head(10),
y=feature_ranks.head(10).index)

plt.title("Top 10 Important Features")

plt.xlabel("Feature Importance Score")

plt.ylabel("Features")

plt.tight_layout()
```



13. Recursive Feature Elimination (RFE): Perform feature selection using RFE. Evaluate the performance of a machine learning model before and after feature selection.

Code And Output:

Initialize Logistic Regression

from sklearn.datasets import model = load breast cancer LogisticRegression(max iter=10000, solver='liblinear') from sklearn.linear model import LogisticRegression from sklearn.feature selection import # ----- Before RFE -----RFE model.fit(X train, y train) from sklearn.model_selection import y pred full = model.predict(X test) train_test_split full acc = accuracy score(y test, from sklearn.metrics import y pred full) accuracy score, classification report print(" Full Model Accuracy (All import pandas as pd features):", full acc) # Load dataset # ----- Apply RFE ----data = load breast cancer() rfe = RFE(estimator=model, X = pd.DataFrame(data.data, n_features_to_select=10) columns=data.feature names) rfe.fit(X_train, y_train) y = pd.Series(data.target) # Transform datasets # Split dataset X train rfe = rfe.transform(X train) X train, X test, y train, y test = X test rfe = rfe.transform(X test) train test split(X, y, test size=0.2, random state=42) # ----- After RFE ------

```
model.fit(X_train_rfe, y_train) # Show selected features

y_pred_rfe = model.predict(X_test_rfe) selected_features =

rfe_acc = accuracy_score(y_test,

y_pred_rfe) print("\n ✓ Selected Features by RFE:")

print(" ☆ RFE Model Accuracy (Top 10 print(selected_features)

features):", rfe_acc)
```

| ■ Before RFE: Accuracy: 0.956140350877193 Classification Report: | | | | | | |
|---|-----------|--------|----------|---------|--|--|
| | precision | recall | f1-score | support | | |
| Ø | 0.97 | 0.91 | 0.94 | 43 | | |
| 1 | 0.95 | 0.99 | 0.97 | 71 | | |
| accuracy | | | 0.96 | 114 | | |
| macro avg | 0.96 | 0.95 | 0.95 | 114 | | |
| weighted avg | 0.96 | 0.96 | 0.96 | 114 | | |
| | | | | | | |
| Selected Features by RFE: | | | | | | |
| ['mean radius' 'mean compactness' 'mean concavity' 'texture error' | | | | | | |
| 'worst radius' 'worst smoothness' 'worst compactness' 'worst concavity' | | | | | | |
| 'worst concave points' 'worst symmetry'] | | | | | | |

| ✓ After RFE: Accuracy: 0.9736842105263158 Classification Report: | | | | | | |
|--|-----|-----------|--------|----------|---------|--|
| | | precision | recall | f1-score | support | |
| | ^ | 0.00 | 0.05 | 0.00 | 43 | |
| | 0 | 0.98 | 0.95 | 0.96 | 43 | |
| | 1 | 0.97 | 0.99 | 0.98 | 71 | |
| | | | | | | |
| accura | асу | | | 0.97 | 114 | |
| macro a | avg | 0.97 | 0.97 | 0.97 | 114 | |
| weighted a | avg | 0.97 | 0.97 | 0.97 | 114 | |

14. Train-Test Split and Cross-Validation: Split a dataset into train-test sets. Use Shuffle Cross-Validation to evaluate a model and compare the results.

Code And Output:

Test Split

model =

LogisticRegression(max iter=200)

model.fit(X train, y train)

from sklearn.datasets import load iris y pred = model.predict(X test) from sklearn.linear_model import print(" Accuracy (Train-Test Split):", LogisticRegression accuracy_score(y_test, y_pred)) from sklearn.model selection import train_test_split, ShuffleSplit, # Step 4: Shuffle Cross-Validation cross_val_score shuffle split = ShuffleSplit(n splits=5, from sklearn.metrics import test size=0.3, random state=42) accuracy score cv scores = cross val score(LogisticRegression(max # Step 1: Load dataset _iter=200), X, y, cv=shuffle_split) data = load_iris() X = data.data print("\n♥ Shuffle Cross-Validation Scores:") y = data.target print(cv scores) print(" Mean Accuracy (Shuffle CV):", # Step 2: Train-Test Split cv_scores.mean()) X train, X test, y train, y test = train_test_split(X, y, test_size=0.3, random state=42) # Step 3: Train and Evaluate using Train-Shuffle Cross-Validation Scores:

0.91111111 0.9555556 0.93333333]

Mean Accuracy (Shuffle CV): 0.96

15. Bagging and Random Forest: Build and evaluate a Random Forest model. Visualize the decision trees and feature importance.

| # Import libraries import pandas as pd import numpy as np | <pre>X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.3, random_state=42)</pre> |
|---|---|
| import matplotlib.pyplot as plt from sklearn.datasets import load_iris from sklearn.ensemble import RandomForestClassifier, BaggingClassifier from sklearn.tree import plot_tree from sklearn.model_selection import train_test_split from sklearn.metrics import accuracy_score | # # A Random Forest Classifier # rf = RandomForestClassifier(n_estimators=5, random_state=42) rf.fit(X_train, y_train) y_pred_rf = rf.predict(X_test) print(" Random Forest Accuracy:", accuracy_score(y_test, y_pred_rf)) |
| <pre># Load the dataset iris = load_iris() X = iris.data y = iris.target feature_names = iris.feature_names class_names = iris.target_names</pre> | # # Bagging Classifier (with Decision Trees) # from sklearn.tree import DecisionTreeClassifier |
| # Split the data | bagging = BaggingClassifier(base_estimator=Decisi |

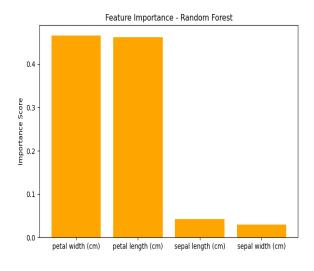
```
onTreeClassifier(), n_estimators=5,
random state=42)
bagging.fit(X_train, y_train)
y_pred_bag = bagging.predict(X_test)
print(" ( Bagging Accuracy:",
accuracy_score(y_test, y_pred_bag))
# [1] Feature Importance from Random
Forest
# -----
importances = rf.feature_importances_
indices = np.argsort(importances)[::-1]
plt.figure(figsize=(8, 5))
plt.title("Feature Importance - Random
Forest")
plt.bar([feature_names[i] for i in
indices], importances[indices],
color='orange')
plt.ylabel("Importance Score")
plt.show()
# -----
# 
 Visualize one Decision Tree from
Random Forest
# -----
```

plt.figure(figsize=(15, 8))

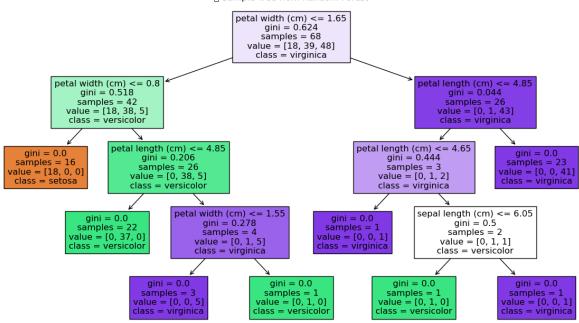
plot_tree(rf.estimators_[0], filled=True, feature_names=feature_names, class_names=class_names)

plt.title(" Sample Tree from Random Forest")

plt.show()



☐ Sample Tree from Random Forest



16 Boosting Methods: Implement AdaBoost and XGBoost on classification task. Compare their accuracy and runtime performance.

```
xgb_model.fit(X_train, y_train)
xgb pred = xgb model.predict(X test)
end xgb = time.time()
xgb_accuracy = accuracy_score(y_test,
xgb_pred)
xgb_time = end_xgb - start_xgb
print("\n X XGBoost Accuracy:",
xgb accuracy)
print(" XGBoost Runtime:", xgb time,
"seconds")
# -----
# [1] Performance Summary
# -----
print("\n | Performance Comparison:")
print(f"AdaBoost -> Accuracy:
{ada accuracy:.4f}, Runtime:
{ada_time:.4f} sec")
print(f"XGBoost -> Accuracy:
{xgb_accuracy:.4f}, Runtime:
{xgb_time:.4f} sec")
```

- AdaBoost Accuracy: 1.0
 AdaBoost Runtime: 0.0532 seconds

 XGBoost Accuracy: 1.0
 XGBoost Runtime: 0.1687 seconds

 Performance Comparison:
- AdaBoost -> Accuracy: 1.0000, Runtime: 0.0532 sec XGBoost -> Accuracy: 1.0000, Runtime: 0.1687 sec

17. K-Nearest Neighbors (KNN): Implement a KNN classifier for a classification task. Experiment with different values of K and analyze their impact on accuracy.

Code And OutPut:

Try different values of K

| import numpy as np | k_values = range(1, 21) |
|---|--|
| import matplotlib.pyplot as plt | accuracies = [] |
| from sklearn.datasets import load_iris | |
| from sklearn.model_selection import train_test_split | for k in k_values: knn = |
| from sklearn.neighbors import | KNeighborsClassifier(n_neighbors=k) |
| KNeighborsClassifier | knn.fit(X_train, y_train) |
| from sklearn.metrics import accuracy_score | predictions = knn.predict(X_test) |
| 4004.407_500.C | <pre>acc = accuracy_score(y_test, predictions)</pre> |
| # Load Iris dataset | accuracies.append(acc) |
| iris = load_iris() | |
| X = iris.data | # Print accuracy for each K |
| y = iris.target | for k, acc in zip(k_values, accuracies): |
| | print(f"K={k}: Accuracy={acc:.4f}") |
| # Split into train and test sets | |
| <pre>X_train, X_test, y_train, y_test = train test split(X, y, test size=0.3,</pre> | # Plotting accuracy vs K |
| random_state=42) | plt.figure(figsize=(10, 5)) |
| # Try different values of K | <pre>plt.plot(k_values, accuracies, marker='o', linestyle='-')</pre> |

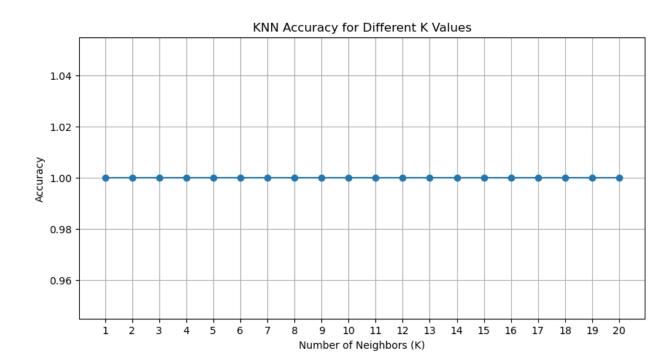
plt.title('KNN Accuracy for Different K plt.show()
Values')

plt.xlabel('Number of Neighbors (K)')

plt.ylabel('Accuracy')

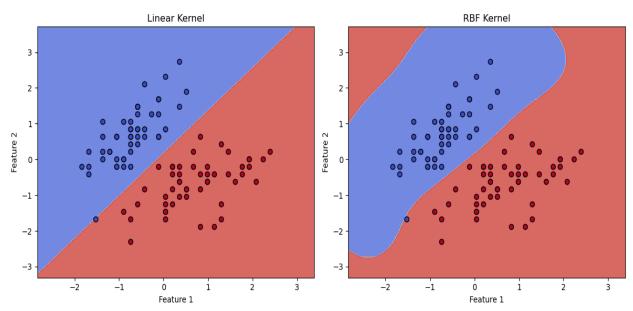
plt.xticks(k_values)

plt.grid(True)



18. Support Vector Machines (SVM): Train an SVM model with both linear and RBF kernels on a dataset. Visualize the decision boundaries.

```
# Create a meshgrid for plotting decision
import numpy as np
                                                      boundaries
import matplotlib.pyplot as plt
                                                      h = .02
from sklearn import svm, datasets
                                                      x_{min}, x_{max} = X[:, 0].min() - 1, X[:,
from sklearn.preprocessing import
                                                      0].max() + 1
StandardScaler
                                                      y min, y max = X[:, 1].min() - 1, X[:, 1]
                                                      1].max() + 1
# Load the dataset (2 classes and 2
                                                      xx, yy = np.meshgrid(np.arange(x min,
features for simplicity)
                                                      x max, h),
iris = datasets.load iris()
                                                                  np.arange(y min, y max, h))
X = iris.data[:, :2] # Using only 2
features for visualization
                                                      # Models with Linear and RBF Kernels
y = iris.target
                                                      models = [
                                                        ('Linear Kernel',
# Use only 2 classes (binary
                                                      svm.SVC(kernel='linear', C=1.0)),
classification)
                                                        ('RBF Kernel', svm.SVC(kernel='rbf',
X = X[y != 2]
                                                      gamma=0.7, C=1.0))
y = y[y != 2]
                                                      ]
# Scale features
                                                      plt.figure(figsize=(12, 5))
scaler = StandardScaler()
X = scaler.fit transform(X)
                                                      for i, (title, clf) in enumerate(models):
                                                        clf.fit(X, y)
```



19. Regularization Techniques: Use Lasso and Ridge regression on a dataset. Analyze how they handle multicollinearity and reduce model complexity.

```
import numpy as np
                                                   X train, X test, y train, y test =
                                                   train test split(X, y, test size=0.2,
import pandas as pd
                                                   random_state=42)
import matplotlib.pyplot as plt
from sklearn.linear model import
                                                   # Step 3: Train models
LinearRegression, Ridge, Lasso
                                                   models = {
from sklearn.model selection import
                                                     "Linear Regression":
train test split
                                                   LinearRegression(),
from sklearn.metrics import
                                                     "Ridge Regression": Ridge(alpha=1.0),
mean_squared_error
                                                     "Lasso Regression": Lasso(alpha=0.1)
# Step 1: Generate synthetic data with
                                                   }
multicollinearity
                                                   coefficients = {}
np.random.seed(0)
                                                   for name, model in models.items():
n_samples = 100
                                                     model.fit(X_train, y_train)
X1 = np.random.rand(n samples)
                                                     y_pred = model.predict(X_test)
X2 = X1 + np.random.normal(0, 0.1,
                                                     mse = mean squared error(y test,
n samples) # Highly correlated with X1
                                                   y_pred)
X3 = np.random.rand(n samples)
                                                     coefficients[name] = model.coef
                                                     print(f"{name} MSE: {mse:.4f}")
X = np.vstack([X1, X2, X3]).T
v = 4 * X1 + 2 * X2 + 3 * X3 +
                                                   # Step 4: Compare Coefficients
np.random.normal(0, 0.1, n samples)
                                                   coef df = pd.DataFrame(coefficients,
                                                   index=["X1", "X2", "X3"])
# Step 2: Split dataset
```

print("\nQ Coefficients
Comparison:\n", coef_df)

Ridge vs Lasso")

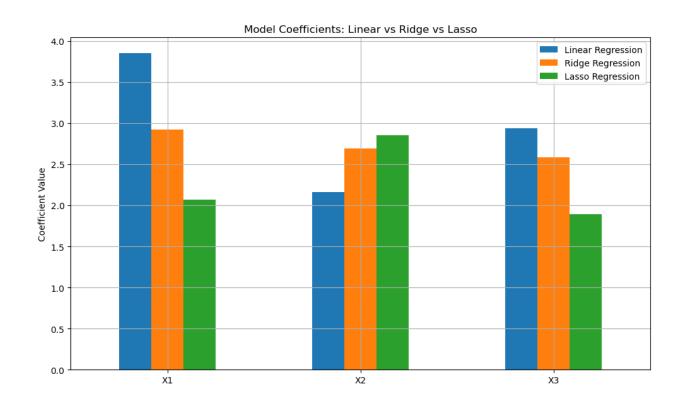
Step 5: Visualize Coefficients

coef_df.plot(kind='bar', figsize=(10, 6))

plt.title("Model Coefficients: Linear vs

plt.ylabel("Coefficient Value")
plt.xticks(rotation=0)
plt.grid(True)
plt.tight_layout()
plt.show()

Linear Regression MSE: 0.015/ Ridge Regression MSE: 0.0483 Lasso Regression MSE: 0.2287 Coefficients Comparison: Linear Regression Ridge Regression Lasso Regression X1 3.848380 2.924626 2.072207 X2 2.162766 2.854834 2.689867 2.581383 **X**3 2.939791 1.891608



20. Introduction to Neural Networks: Build a simple Artificial Neural Network (ANN) to classify data. Use optimization algorithms (Gradient Descent, SGD) and visualize the loss during training.

Code And Output:

X_train, X_test, y_train, y_test = train test split(X, y, test size=0.2, # Imports random state=42) import numpy as np import matplotlib.pyplot as plt # Feature scaling from sklearn.datasets import scaler = StandardScaler() load_breast_cancer X train = scaler.fit transform(X train) from sklearn.model selection import train test split X test = scaler.transform(X test) from sklearn.preprocessing import StandardScaler # Build ANN model from tensorflow.keras.models import model = Sequential([Sequential Dense(16, input shape=(X.shape[1],), from tensorflow.keras.layers import activation='relu'), Dense Dense(8, activation='relu'), from tensorflow.keras.optimizers import Dense(1, activation='sigmoid') # SGD Output layer for binary classification]) # Load dataset data = load breast cancer() # Compile model X = data.data optimizer = SGD(learning_rate=0.01) y = data.target model.compile(optimizer=optimizer, loss='binary_crossentropy', # Train-test split metrics=['accuracy'])

Train model

history = model.fit(X_train, y_train, validation_data=(X_test, y_test), epochs=100, batch size=16, verbose=0)

Evaluate model

loss, accuracy = model.evaluate(X_test,
y_test, verbose=0)

print(f" Test Accuracy:
{accuracy:.4f}")

plt.plot(history.history['loss'],
label='Train Loss')

plt.plot(history.history['val_loss'],
label='Validation Loss')

plt.title('Loss over Epochs')

plt.xlabel('Epoch')

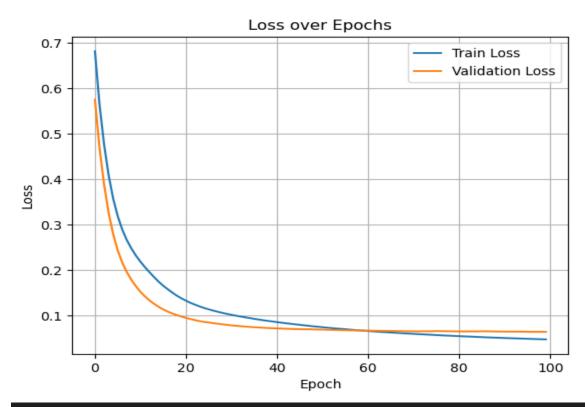
plt.ylabel('Loss')

plt.legend()

plt.grid(True)

plt.show()

Visualize loss during training



super().__init__(activity_regularizer=activity_regularizer, **kwargs)

🔽 Test Accuracy: 0.9737