RAJALAKSHMI ENGINEERING COLLEGE RAJALAKSHMI NAGAR, THANDALAM – 602 105



CS23333 BUILD AND DEPLOYMENT OF MACHINE LEARNING APPLICATIONS

Laboratory Observation Note Book

Name : Guṇayazḥagaṇ B
Year / Branch / Section : 3 nd Year/ AIML
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Exp No: 1	Setting Up the Environment And Preprocessing the Data
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To set up a fully functional machine learning development environment and to perform data preprocessing operations like handling missing values, encoding categorical variables, feature scaling, and splitting datasets.

Algorithm:

- 1. Install Required Libraries:
 - Install numpy, pandas, matplotlib, seaborn, and scikit-learn using pip.
- 2. Import Libraries.
- 3. Load Dataset:
 - Load any dataset (e.g., Titanic or Iris) using pandas.
- 4. Data Exploration:
 - Use df.info(), df.describe(), df.isnull().sum() to understand the data.
- 5. Handle Missing Values:
 - Use .fillna() or .dropna() depending on the strategy.
- 6. Encode Categorical Data:
 - Use pd.get dummies() or LabelEncoder.
- 7. Feature Scaling:
 - Normalize or standardize the numerical features using StandardScaler or MinMaxScaler.
- 8. Split Dataset:
 - Use train test split() from sklearn to create training and testing sets.
- 9. Display the Preprocessed Data.

```
Code:
```

```
# 1. Install necessary libraries (if not already installed)
#!pip install numpy pandas matplotlib seaborn scikit-learn
# 2. Import libraries
import pandas as pd
import numpy as np
from sklearn.model selection import train test split
from sklearn.preprocessing import StandardScaler, LabelEncoder
import seaborn as sns
import matplotlib.pyplot as plt
# 3. Load dataset
df = sns.load dataset('titanic') # Titanic dataset
df.head()
# 4. Explore the dataset
print(df.info())
print(df.describe())
print(df.isnull().sum())
# 5. Handle missing values
# Fill age with median, embark town with mode
df['age'].fillna(df['age'].median(), inplace=True)
df['embark town'].fillna(df['embark town'].mode()[0], inplace=True)
df.drop(columns=['deck'], inplace=True) # too many missing values
# 6. Encode categorical variables
# Convert 'sex' and 'embark town' using LabelEncoder
le = LabelEncoder()
```

```
df['sex'] = le.fit transform(df['sex'])
df['embark town'] = le.fit transform(df['embark town'])
# Drop non-informative or redundant columns
df.drop(columns=['embarked', 'class', 'who', 'alive', 'adult male', 'alone'], inplace=True)
#7. Feature Scaling
scaler = StandardScaler()
numerical cols = ['age', 'fare']
df[numerical cols] = scaler.fit transform(df[numerical cols])
# 8. Split dataset
# Define features (X) and label (y)
X = df.drop(`survived', axis=1)
y = df['survived']
X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.2, random_state=42)
# 9. Show final preprocessed data
print("Training Data Shape:", X train.shape)
print("Test Data Shape:", X_test.shape)
X train.head()
```

Output:

<class 'pandas.core.frame.DataFrame'> RangeIndex: 891 entries, 0 to 890 Data columns (total 15 columns): Column Non-Null Count Dtype 0 survived 891 non-null int64 891 non-null 1 pclass int64 2 891 non-null object 3 age 714 non-null float64 int64 4 sibsp 891 non-null 891 non-null int64 5 parch float64 6 fare 891 non-null 7 embarked 889 non-null object 8 class 891 non-null category 9 891 non-null who object adult_male 10 891 non-null bool 11 deck 203 non-null category 12 embark_town 889 non-null object 13 alive 891 non-null object alone 891 non-null bool dtypes: bool(2), category(2), float64(2), int64(4), object(5) memory usage: 80.7+ KB None survived pclass sibsp parch fare age count 891.000000 891.000000 714.000000 891.000000 891.000000 891.000000 mean 0.383838 2.308642 29.699118 0.523008 0.381594 32.204208 std 0.486592 0.836071 14.526497 1.102743 0.806057 49.693429 min 0.000000 1.000000 0.420000 0.000000 0.000000 0.000000 25% 0.000000 2.000000 20.125000 0.000000 0.000000 7.910400 50% 0.000000 3.000000 28.000000 0.000000 0.000000 14.454200 75% 1.000000 3.000000 38.000000 1.000000 0.000000 31.000000 1.000000 3.000000 80.000000 8.000000 6.000000 512.329200 max survived 0 pclass 0 sex 0 age 177 sibsp 0 parch 0 fare 0 embarked 2 class 0 who 0 adult male 0 deck 688 embark town 2 alive 0 alone 0 dtype: int64

Training Data Shape: (712, 7) Test Data Shape: (179, 7)

/tmp/ipython-input-4068659829.py:3: FutureWarning: A value is trying to be set on a copy of a DataFrame or Series through chained assignment using an inplace method.

The behavior will change in pandas 3.0. This inplace method will never work because the intermediate object on which we are setting values always behaves as a copy.

For example, when doing 'df[col].method(value, inplace=True)', try using 'df.method({col: value}, inplace=True)' or df[col] = df[col].method(value) instead, to perform the operation inplace on the original object.

df['age'].fillna(df['age'].median(), inplace=True)
/tmp/ipython-input-4068659829.py:4: FutureWarning: A value is trying to be set on a copy of a DataFrame or Series through chained assignment using an inplace method.
The behavior will change in pandas 3.0. This inplace method will never work because the intermediate object on which we are setting values always behaves as a copy.

For example, when doing 'df[col].method(value, inplace=True)', try using 'df.method({col: value}, inplace=True)' or df[col] = df[col].method(value) instead, to perform the operation inplace on the original object.

df['embark_town'].fillna(df['embark_town'].mode()[0], inplace=True)

	pclass	sex	age	sibsp	parch	fare	embark_town
331	1	1	1.240235	0	0	-0.074583	2
733	2	1	-0.488887	0	0	-0.386671	2
382	3	1	0.202762	0	0	-0.488854	2
704	3	1	-0.258337	1	0	-0.490280	2
813	3	0	-1.795334	4	2	-0.018709	2

Exp No: 2

Support Vector Machine (SVM) and Random Forest for Binary & Multiclass Classification

Aim

To build classification models using **Support Vector Machines (SVM)** and **Random Forest**, apply them to a dataset, and evaluate the models using performance metrics like accuracy and confusion matrix.

Algorithm

Part A: SVM Model

- 1. Import necessary libraries
- 2. Load and explore the dataset
- 3. Handle missing values if any
- 4. Encode categorical variables
- 5. Split dataset into training and testing sets
- 6. Build SVM classifier using SVC()
- 7. Train and predict
- 8. Evaluate the model using accuracy and confusion matrix

Part B: Random Forest Model

- 1. Initialize Random Forest using RandomForestClassifier()
- 2. Train and predict
- 3. Evaluate and compare with SVM

Code:

1. Import libraries

import pandas as pd

from sklearn.datasets import load iris

from sklearn.model selection import train test split

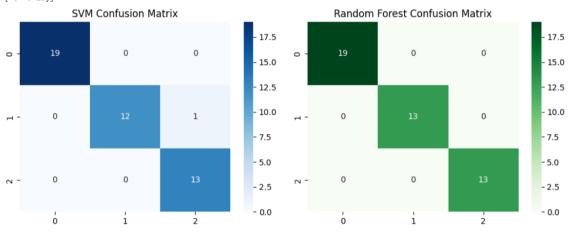
from sklearn.preprocessing import StandardScaler

from sklearn.svm import SVC

from sklearn.ensemble import RandomForestClassifier

```
from sklearn.metrics import accuracy score, confusion matrix
import seaborn as sns
import matplotlib.pyplot as plt
# 2. Load dataset
iris = load iris()
X = iris.data
y = iris.target
#3. Feature scaling
scaler = StandardScaler()
X scaled = scaler.fit transform(X)
# 4. Train-test split
X train, X test, y train, y test = train test split(X scaled, y, test size=0.3, random state=42)
# -----
# Part A: SUPPORT VECTOR MACHINE
# -----
# 5. Initialize and train SVM
svm model = SVC(kernel='linear') # You can also try 'rbf', 'poly'
svm_model.fit(X_train, y_train)
# 6. Predict and evaluate SVM
y pred svm = svm model.predict(X test)
print("SVM Accuracy:", accuracy score(y test, y pred svm))
print("SVM Confusion Matrix:\n", confusion matrix(y test, y pred svm))
```

```
# -----
# Part B: RANDOM FOREST
# -----
#7. Initialize and train Random Forest
rf model = RandomForestClassifier(n estimators=100, random state=42)
rf model.fit(X train, y train)
#8. Predict and evaluate Random Forest
y pred rf = rf model.predict(X test)
print("Random Forest Accuracy:", accuracy score(y test, y pred rf))
print("Random Forest Confusion Matrix:\n", confusion matrix(y test, y pred rf))
# -----
# 9. Visual comparison using seaborn heatmap
# -----
plt.figure(figsize=(10, 4))
plt.subplot(1, 2, 1)
sns.heatmap(confusion matrix(y test, y pred svm), annot=True, cmap='Blues', fmt='d')
plt.title("SVM Confusion Matrix")
plt.subplot(1, 2, 2)
sns.heatmap(confusion matrix(y test, y pred rf), annot=True, cmap='Greens', fmt='d')
plt.title("Random Forest Confusion Matrix")
plt.tight layout()
plt.show()
```



ExpNo:3

Classification with Decision Trees

Aim

To implement a Decision Tree classifier and evaluate its performance using **accuracy score** and **confusion matrix** on a real-world dataset.

Algorithm

- 1. Import necessary libraries
- 2. Load a classification dataset (e.g., Iris or Titanic)
- 3. Split the dataset into training and test sets
- 4. Preprocess data if needed
- 5. Train a DecisionTreeClassifier from sklearn.tree
- 6. Predict on test data
- 7. Evaluate using:
 - Confusion Matrix
 - Accuracy Score
- 8. Visualize the Decision Tree (optional)

Code:

Step 1: Import Libraries

from sklearn.datasets import load iris

from sklearn.tree import DecisionTreeClassifier, plot_tree

from sklearn.model selection import train test split

from sklearn.metrics import confusion matrix, accuracy score

import matplotlib.pyplot as plt

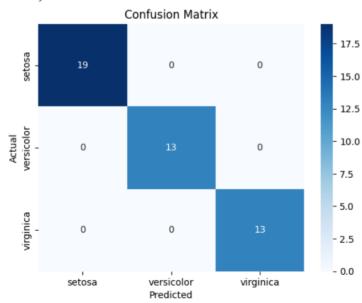
import seaborn as sns

```
# Step 2: Load Dataset
iris = load iris()
X = iris.data
y = iris.target
# Step 3: Split the dataset
X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.3, random_state=42)
# Step 4: Train the Decision Tree Classifier
dt model = DecisionTreeClassifier(criterion='gini', random state=0)
dt_model.fit(X_train, y_train)
# Step 5: Predict
y pred = dt model.predict(X test)
# Step 6: Evaluate the Model
cm = confusion matrix(y test, y pred)
acc = accuracy score(y test, y pred)
print("Confusion Matrix:\n", cm)
print("Accuracy Score:", acc)
# Step 7: Visualize Confusion Matrix
sns.heatmap(cm,
                          annot=True,
                                               cmap="Blues",
                                                                        xticklabels=iris.target names,
yticklabels=iris.target names)
plt.xlabel("Predicted")
plt.ylabel("Actual")
plt.title("Confusion Matrix")
plt.show()
# Step 8: Visualize the Decision Tree
plt.figure(figsize=(12,8))
plot tree(dt model, filled=True, feature names=iris.feature names, class names=iris.target names)
```

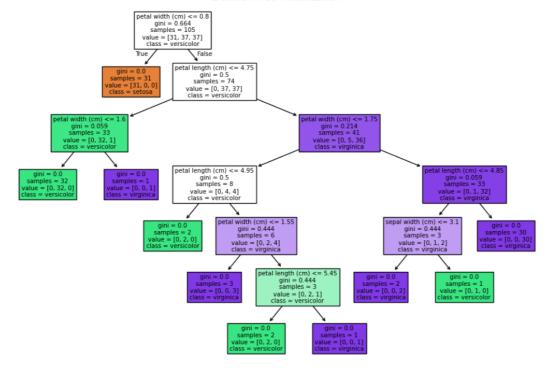
plt.title("Decision Tree Visualization") plt.show()

OUTPUT:

Confusion Matrix: [[19 0 0] [0 13 0] [0 0 13]] Accuracy Score: 1.0



Decision Tree Visualization



Exp No: 4A Support Vector Machines (SVM)

Aim:

To build an SVM model for a binary classification task, tune its hyperparameters, and evaluate it using accuracy, precision, recall, F1-score, confusion matrix, and ROC-AUC.

Algorithm:

- 1. Import libraries: numpy, pandas, matplotlib, sklearn.
- 2. Load data: Use a standard binary dataset (Breast Cancer Wisconsin) from sklearn.datasets.
- 3. Train/Test split: 80/20 split with a fixed random_state.
- 4. Preprocess: Standardize features (StandardScaler).
- 5. SVMs are sensitive to feature scale.
- 6. Model selection: Use SVC (RBF kernel).
- 7. Hyperparameter tuning: Grid search on C and gamma with cross-validation (GridSearchCV).
- 8. Train final model: Fit on training data using best parameters.
- 9. Evaluate: Predict on test set; compute metrics and plot ROC curve.
- 10. Report: Best params, metrics, and brief observations.

from sklearn.datasets import load_breast_cancer

```
from sklearn.model selection import train test split, GridSearchCV
from sklearn.preprocessing import StandardScaler
from sklearn.svm import SVC
from sklearn.metrics import (
  accuracy score, precision score, recall score, f1 score,
  confusion matrix, classification report, roc auc score, roc curve
)
#2) Load dataset (binary classification)
data = load breast cancer()
X = pd.DataFrame(data.data, columns=data.feature names)
y = pd.Series(data.target, name="target") # 0 = malignant, 1 = benign
#3) Train/test split
X_train, X_test, y_train, y_test = train test split(
  X, y, test size=0.20, random state=42, stratify=y
# 4) Standardize features (important for SVMs)
scaler = StandardScaler()
X_train_sc = scaler.fit_transform(X_train)
X \text{ test } sc = scaler.transform(X \text{ test})
#5) Define model
svm = SVC(kernel='rbf', probability=True, random state=42)
# 6) Hyperparameter grid & tuning
param grid = {
  "C": [0.1, 1, 10, 100],
  "gamma": ["scale", 0.01, 0.001, 0.0001]
```

```
grid = GridSearchCV(
  estimator=svm,
  param grid=param grid,
  scoring='f1', # You can change to 'accuracy' or 'roc auc'
  cv=5,
  n jobs=-1,
  verbose=0
)
grid.fit(X train sc, y train)
print("Best Parameters from Grid Search:", grid.best_params_)
best svm = grid.best estimator
#7) Train final model & predict
best svm.fit(X train sc, y train)
y pred = best svm.predict(X test sc)
y prob = best svm.predict proba(X test sc)[:, 1]
#8) Evaluation
acc = accuracy score(y test, y pred)
prec = precision score(y test, y pred, zero division=0)
rec = recall score(y test, y pred)
f1 = f1 score(y test, y pred)
auc = roc_auc_score(y_test, y_prob)
cm = confusion matrix(y test, y pred)
print("\n=== SVM (RBF) — Test Metrics ====")
print(f"Accuracy : {acc:.4f}")
print(f"Precision: {prec:.4f}")
```

```
print(f"Recall : {rec:.4f}")
print(f"F1-Score : {f1:.4f}")
print(f"ROC-AUC : {auc:.4f}")
print("\nConfusion Matrix:\n", cm)
print("\nClassification Report:\n", classification report(y test, y pred, zero division=0))
#9) Plot ROC Curve
fpr, tpr, thresholds = roc curve(y test, y prob)
plt.figure()
plt.plot(fpr, tpr, label=f"SVM (AUC = {auc:.3f})")
plt.plot([0, 1], [0, 1], linestyle="--", color='gray')
plt.xlabel("False Positive Rate")
plt.ylabel("True Positive Rate")
plt.title("ROC Curve — SVM (RBF)")
plt.legend()
plt.grid(True)
plt.show()
```

Best Parameters from Grid Search: {'C': 10, 'gamma': 0.01}

=== SVM (RBF) - Test Metrics ===

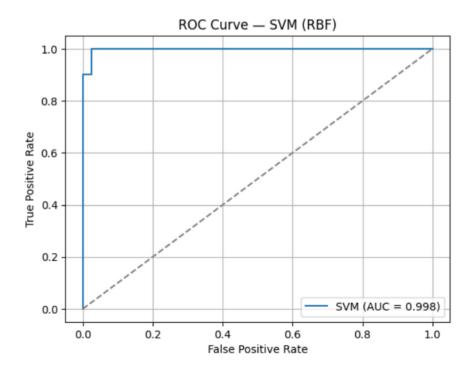
Accuracy : 0.9825 Precision: 0.9861 Recall : 0.9861 F1-Score : 0.9861 ROC-AUC : 0.9977

Confusion Matrix:

[[41 1] [1 71]]

Classification Report:

	precision	recall	f1-score	support
0	0.98	0.98	0.98	42
1	0.99	0.99	0.99	72
accuracy			0.98	114
macro avg	0.98	0.98	0.98	114
weighted avg	0.98	0.98	0.98	114



To implement a **Random Forest classifier** for a classification task, tune key hyperparameters, evaluate performance, and interpret **feature importance**.

Algorithm:

- 1. Import libraries.
- 2. Load data (use same dataset to compare with SVM).
- 3. Train/Test split with stratification.
- 4. (Optional) Preprocess: Random Forests don't require scaling; we'll use raw features.
- 5. Model: RandomForestClassifier.
- 6. Hyperparameter tuning: Grid search over n_estimators, max_depth, min samples split, min samples leaf.
- 7. Train the best model on training data.
- 8. Evaluate with accuracy, precision, recall, F1, confusion matrix, ROC-AUC.
- 9. Interpretation: Plot top feature importances.

from sklearn.model selection import train test split, GridSearchCV

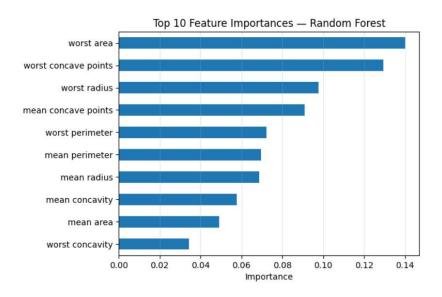
from sklearn.ensemble import RandomForestClassifier

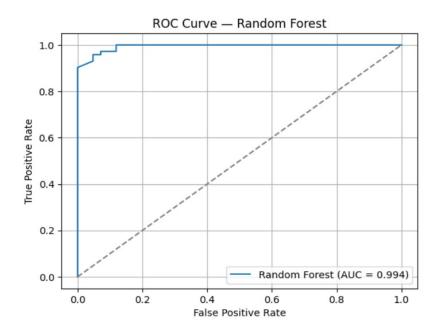
```
from sklearn.metrics import (
  accuracy score, precision score, recall score, f1 score,
  confusion matrix, classification report, roc auc score, roc curve
)
#2) Load dataset (same as 4A for comparison)
data = load breast cancer()
X = pd.DataFrame(data.data, columns=data.feature names)
y = pd.Series(data.target, name="target")
# 3) Train/test split (no scaling needed for RF)
X train, X test, y train, y test = train test split(
  X, y, test size=0.20, random state=42, stratify=y
)
#4) Define model
rf = RandomForestClassifier(random_state=42, n_jobs=-1)
# 5) Hyperparameter grid & tuning
param grid = {
  "n estimators": [100],
  "max depth": [None, 10],
  "min samples split": [2],
  "min samples leaf": [1]
}
grid = GridSearchCV(
  estimator=rf,
  param grid=param grid,
  scoring="f1",
  cv=3,
  n jobs=-1,
  verbose=0)
```

```
grid.fit(X train, y train)
print("Best Parameters (CV):", grid.best params )
best rf = grid.best estimator
#6) Train final model & predict
best rf.fit(X train, y train)
y pred = best rf.predict(X test)
y prob = best_rf.predict_proba(X_test)[:, 1]
#7) Evaluate
acc = accuracy score(y test, y pred)
prec = precision score(y test, y pred, zero division=0)
rec = recall score(y test, y pred)
f1 = f1 score(y test, y pred)
auc = roc auc_score(y_test, y_prob)
cm = confusion_matrix(y_test, y_pred)
print("\n=== Random Forest — Test Metrics ===")
print(f"Accuracy: {acc:.4f}")
print(f"Precision: {prec:.4f}")
print(f"Recall : {rec:.4f}")
print(f"F1-Score : {f1:.4f}")
print(f"ROC-AUC : {auc:.4f}")
print("\nConfusion Matrix:\n", cm)
print("\nClassification Report:\n", classification report(y test, y pred, zero division=0))
#8) Feature Importance (Top 10)
importances = pd.Series(best rf.feature importances , index=X.columns)
top10 = importances.sort values(ascending=False).head(10)
```

```
plt.figure()
top10[::-1].plot(kind="barh")
plt.xlabel("Importance")
plt.title("Top 10 Feature Importances — Random Forest")
plt.grid(axis="x", alpha=0.3)
plt.show()
#9) ROC Curve
fpr, tpr, thresholds = roc curve(y test, y prob)
plt.figure()
plt.plot(fpr, tpr, label=f"Random Forest (AUC = {auc:.3f})")
plt.plot([0, 1], [0, 1], linestyle="--", color='gray')
plt.xlabel("False Positive Rate")
plt.ylabel("True Positive Rate")
plt.title("ROC Curve — Random Forest")
plt.legend()
plt.grid(True)
plt.show()
```

```
Best Parameters (CV): {'max_depth': None, 'min_samples_leaf': 1, 'min_samples_split': 2, 'n_estimators': 100}
=== Random Forest - Test Metrics ===
Accuracy: 0.9561
Precision: 0.9589
Recall : 0.9722
F1-Score : 0.9655
ROC-AUC : 0.9937
Confusion Matrix:
[ 2 70]]
Classification Report:
                         recall f1-score support
             precision
                 0.95
                          0.93
                                   0.94
                 0.96
                          0.97
                                   0.97
                                   0.96
                                             114
   accuracy
  macro avg
                 0.96
                          0.95
                                   0.95
                                             114
weighted avg
                 0.96
                          0.96
                                   0.96
                                             114
```





Exp No: 5	Clustering with K-Means and Dimensionality Reduction with PCA

To demonstrate the application of Unsupervised Learning models, specifically K-Means clustering for grouping data points and Principal Component Analysis (PCA) for dimensionality reduction and visualization, using a suitable dataset.

Algorithm:

1. K-Means Clustering

K-Means is an iterative clustering algorithm that aims to partition \$n\$ observations into \$k\$ clusters, where each observation belongs to the cluster with the nearest mean (centroid).

Steps:

- 1. **Initialization:** Choose \$k\$ initial centroids randomly from the dataset.
- 2. **Assignment:** Assign each data point to the cluster whose centroid is closest (e.g., using Euclidean distance).
- 3. Update: Recalculate the centroids as the mean of all data points assigned to that cluster.
- 4. **Iteration:** Repeat steps 2 and 3 until the centroids no longer move significantly or a maximum number of iterations is reached.

2. Principal Component Analysis (PCA)

PCA is a statistical procedure that uses an orthogonal transformation to convert a set of observations of possibly correlated variables into a set of values of linearly uncorrelated variables called principal components.

Steps:

- 1. **Standardization:** Standardize the dataset (mean = 0, variance = 1).
- 2. Covariance Matrix Calculation: Compute the covariance matrix of the standardized data.
- 3. **Eigenvalue Decomposition:** Calculate the eigenvalues and eigenvectors of the covariance matrix.
- 4. **Feature Vector Creation:** Sort the eigenvectors by decreasing eigenvalues and select the top \$k\$ eigenvectors to form a feature vector (projection matrix).
- 5. **Projection:** Project the original data onto the new feature space using the feature vector.

CODE:

```
# EXPERIMENT — K-Means & PCA
# Import necessary libraries
import numpy as np
import pandas as pd
import matplotlib.pyplot as plt
import seaborn as sns
from sklearn.datasets import make blobs
from sklearn.preprocessing import StandardScaler
from sklearn.cluster import KMeans
from sklearn.decomposition import PCA
from sklearn.metrics import silhouette score
# --- Part 1: K-Means Clustering ---
print("--- Part 1: K-Means Clustering ---")
# 1. Generate dataset
X, y = make blobs(n samples=300, centers=3, cluster std=0.60, random state=42)
df kmeans = pd.DataFrame(X, columns=['Feature 1', 'Feature 2'])
print("\nOriginal K-Means Dataset Head:")
print(df kmeans.head())
# 2. Elbow Method
wcss = []
for i in range(1, 11):
                 KMeans(n clusters=i, init='k-means++', max iter=300,
  kmeans
                                                                               n init=10,
random state=42)
  kmeans.fit(X)
  wcss.append(kmeans.inertia)
plt.figure(figsize=(10, 6))
plt.plot(range(1, 11), wcss, marker='o', linestyle='--')
plt.title('Elbow Method for Optimal K (K-Means)')
plt.xlabel('Number of Clusters (K)')
plt.ylabel('WCSS')
plt.grid(True)
plt.show()
# 3. Apply K-Means with chosen K
optimal k = 3
```

```
kmeans = KMeans(n clusters=optimal k, init='k-means+++', max iter=300, n init=10,
random state=42)
clusters = kmeans.fit predict(X)
df kmeans['Cluster'] = clusters
# 4. Visualize K-Means clusters
plt.figure(figsize=(10, 8))
sns.scatterplot(x='Feature 1', y='Feature 2', hue='Cluster', data=df kmeans, palette='viridis',
s=100, alpha=0.8)
plt.scatter(kmeans.cluster centers [:, 0], kmeans.cluster centers [:, 1], s=300, c='red',
marker='X', label='Centroids')
plt.title(f'K-Means Clustering with K=\{\text{optimal } k\}'\}
plt.xlabel('Feature 1')
plt.ylabel('Feature 2')
plt.legend()
plt.grid(True)
plt.show()
# 5. Silhouette Score
silhouette avg = silhouette score(X, clusters)
print(f"\nSilhouette Score for K-Means (K={optimal k}): {silhouette avg:.3f}")
# --- Part 2: Dimensionality Reduction with PCA ---
print("\n--- Part 2: Dimensionality Reduction with PCA ---")
# 1. Generate 4D dataset
X pca, y pca = make blobs(n samples=500, n features=4, centers=4, cluster std=1.0,
random state=25)
df pca original
                 = pd.DataFrame(X pca,
                                               columns=[fFeature {i+1}'
range(X pca.shape[1])])
df pca original['True Cluster'] = y pca
print("\nOriginal PCA Dataset Head:")
print(df pca original.head())
print(f"Original PCA Dataset Shape: {df pca original.shape}")
# 2. Standardize
scaler = StandardScaler()
X pca scaled = scaler.fit transform(X pca)
# 3. PCA (4D \rightarrow 2D)
pca = PCA(n components=2)
principal components = pca.fit transform(X pca scaled)
df principal components
                                                     pd.DataFrame(principal components,
columns=['Principal Component 1', 'Principal Component 2'])
df principal components['True Cluster'] = y pca
```

```
explained variance = pca.explained variance ratio
print("\nPrincipal Components Head:")
print(df principal components.head())
print(f"\nExplained Variance Ratio: {explained variance}")
print(f"Total Explained Variance by 2 PCs: {explained variance.sum():.3f}")
# 4. Visualize PCA result
plt.figure(figsize=(10, 8))
sns.scatterplot(x='Principal Component 1', y='Principal Component 2', hue='True Cluster',
         data=df principal components, palette='Paired', s=100, alpha=0.8)
plt.title('PCA - Dimensionality Reduction to 2 Components')
plt.xlabel(fPC1 ({explained variance[0]*100:.2f}%)')
plt.ylabel(fPC2 ({explained variance[1]*100:.2f}%)')
plt.grid(True)
plt.show()
# 5. K-Means on PCA-reduced data
                  KMeans(n clusters=4, init='k-means++',
kmeans pca
              =
                                                              max iter=300,
                                                                               n init=10,
random state=42)
clusters pca = kmeans pca.fit predict(principal components)
df principal components['KMeans Cluster on PCA'] = clusters pca
plt.figure(figsize=(10, 8))
sns.scatterplot(x='Principal Component 1',
                                                              y='Principal Component 2',
hue='KMeans Cluster on PCA',
         data=df principal components, palette='viridis', s=100, alpha=0.8)
plt.scatter(kmeans pca.cluster centers [:, 0], kmeans pca.cluster centers [:, 1], s=300,
c='red', marker='X', label='Centroids')
plt.title('K-Means Clustering on PCA-Reduced Data')
plt.xlabel('Principal Component 1')
plt.ylabel('Principal Component 2')
plt.legend()
plt.grid(True)
plt.show()
# 6. Silhouette Score for PCA-reduced KMeans
silhouette avg pca = silhouette score(principal components, clusters pca)
print(f"\nSilhouette
                      Score
                               for
                                      K-Means
                                                         PCA-Reduced
                                                                          Data
                                                  on
                                                                                   (K=4):
{silhouette avg pca:.3f}")
```

--- Part 1: K-Means Clustering ---

Original K-Means Dataset Head:

Feature_1 Feature_2

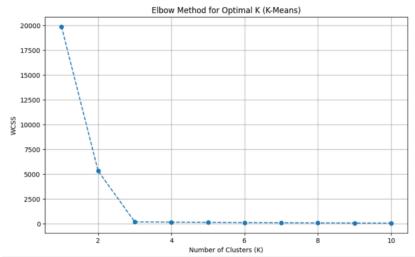
0 -7.155244 -7.390016

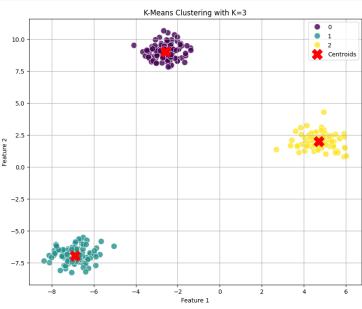
1 -7.395875 -7.110843

2 -2.015671 8.281780

3 4.509270 2.632436

4 -8.102502 -7.484961





Silhouette Score for K-Means (K=3): 0.908

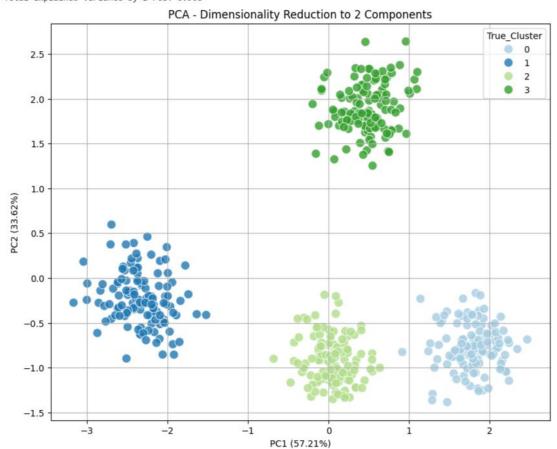
--- Part 2: Dimensionality Reduction with PCA ---

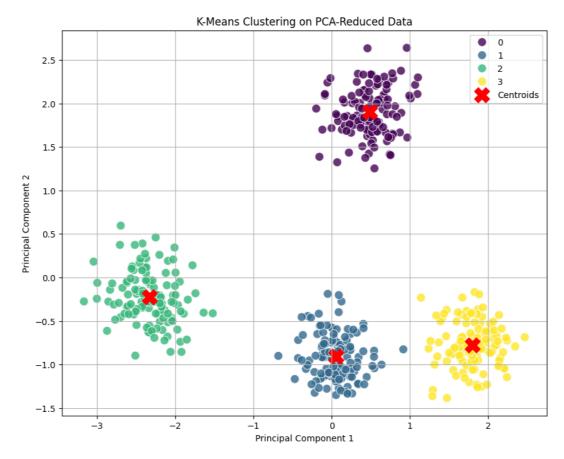
Original PCA Dataset Head: Feature_1 Feature_2 Feature_3 Feature_4 True_Cluster 0 -0.638667 1.110057 -6.400722 -0.204990 3 1 -2.951556 -7.657445 3.844794 0.903589 1 2 -0.253177 2.125103 -7.869801 0.559678 3 3 -2.151209 3.401400 -5.734930 0.965230 3 4 -2.347519 -7.230467 3.478891 -0.443440 1 Original PCA Dataset Shape: (500, 5)

Principal Components Head:

	The That components head		
	Principal_Component_1	Principal_Component_2	True_Cluster
0	0.455305	1.623917	3
1	-2.705622	0.375012	1
2	0.810234	1.966926	3
3	0.427139	2.149626	3
4	-2.407508	0.099250	1

Explained Variance Ratio: [0.57208431 0.33622342] Total Explained Variance by 2 PCs: 0.908





Silhouette Score for K-Means on PCA-Reduced Data (K=4): 0.776

Exp No: 6	Feedforward and Convolutional Neural Networks

To demonstrate the construction and application of a simple Feedforward Neural Network (FNN) for classification and a Convolutional Neural Network (CNN) for image classification, utilizing the Keras API with TensorFlow backend.

Algorithm:

1. Feedforward Neural Network (FNN)

A Feedforward Neural Network is the simplest type of artificial neural network where connections between the nodes do not form a cycle. It consists of an input layer, one or more hidden layers, and an output layer. Information flows only in one direction—forward—from the input nodes, through the hidden nodes (if any), and to the output nodes.

Steps:

- 1. Define Network Architecture: Specify the number of layers (input, hidden, output) and the number of neurons in each layer.
- 2. Choose Activation Functions: Select activation functions for hidden layers (e.g., ReLU) and the output layer (e.g., Sigmoid for binary classification, Softmax for multi-class classification).
- 3. Define Loss Function: Choose a loss function appropriate for the task (e.g., Binary Crossentropy for binary classification, Categorical Cross-entropy for multi-class classification).
- 4. Choose Optimizer: Select an optimization algorithm (e.g., Adam, SGD) to update network weights during training.
- 5. Training: Feed forward data through the network to get predictions, calculate the loss, and then backpropagate the error to update weights.
- 6. Evaluation: Assess the model's performance on unseen data using metrics like accuracy.

2. Convolutional Neural Network (CNN)

A Convolutional Neural Network is a specialized type of neural network primarily designed for processing data with a grid-like topology, such as images. Key components include convolutional layers, pooling layers, and fully connected layers.

Steps:

- 1. Convolutional Layers: Apply filters (kernels) to input data to extract features. Each filter detects a specific pattern (e.g., edges, textures).
- 2. Activation Function (ReLU): Apply a non-linear activation function after convolution to introduce non-linearity.
- 3. Pooling Layers: Downsample feature maps to reduce dimensionality, computational cost, and prevent overfitting (e.g., Max Pooling).
- 4. Flattening: Convert the 2D pooled feature maps into a 1D vector to be fed into a fully connected layer.
- 5. Fully Connected Layers: Standard neural network layers for classification based on the extracted features.
- 6. Output Layer: Final layer with an activation function (e.g., Softmax) to output class probabilities.
- 7. Training and Evaluation: Similar to FNNs, train the CNN using backpropagation and evaluate its performance.

CODE:

```
# Import necessary libraries
import numpy as np
import matplotlib.pyplot as plt
import tensorflow as tf
from tensorflow import keras
from tensorflow.keras import layers
from tensorflow.keras.datasets import mnist, fashion mnist
from sklearn.metrics import classification report, confusion matrix
import seaborn as sns
# Suppress TensorFlow warnings for cleaner output
tf.keras.utils.disable interactive logging()
# --- Part 1: Building a Simple Feedforward Neural Network ---
print("--- Part 1: Building a Simple Feedforward Neural Network ---")
# 1. Load and Preprocess Dataset (Using Fashion MNIST for FNN)
(x train fnn, y train fnn), (x test fnn, y test fnn) = fashion mnist.load data()
print(f"\nOriginal FNN training data shape: {x train fnn.shape}")
print(f''Original FNN test data shape: {x test fnn.shape}")
# Flatten images to 1D array
x train fnn flat = x train fnn.reshape(-1, 28 * 28)
```

```
x test fnn flat = x test fnn.reshape(-1, 28 * 28)
# Normalize pixel values
x train fnn norm = x train fnn flat / 255.0
x test fnn norm = x test fnn flat / 255.0
print(f"Flattened & Normalized FNN training data shape: {x train fnn norm.shape}")
print(f"Flattened & Normalized FNN test data shape: {x test fnn norm.shape}")
# 2. Build FNN Model
model fnn = keras.Sequential([
  layers.Dense(128, activation='relu', input shape=(784,)),
  layers.Dropout(0.2),
  layers.Dense(64, activation='relu'),
  layers.Dense(10, activation='softmax')
])
#3. Compile Model
model fnn.compile(optimizer='adam',
          loss='sparse categorical crossentropy',
          metrics=['accuracy'])
print("\n--- FNN Model Summary ---")
model fnn.summary()
#4. Train Model
print("\n--- Training FNN Model ---")
history fnn = model fnn.fit(x train fnn norm, y train fnn, epochs=10,
                validation split=0.1, verbose=1)
# 5. Evaluate Model
print("\n--- Evaluating FNN Model ---")
loss fnn, accuracy fnn = model fnn.evaluate(x test fnn norm, y test fnn, verbose=0)
print(f"FNN Test Loss: {loss fnn:.4f}")
print(f"FNN Test Accuracy: {accuracy fnn:.4f}")
# Classification report & confusion matrix
y pred fnn = np.argmax(model fnn.predict(x test fnn norm), axis=-1)
print("\n--- FNN Classification Report ---")
print(classification report(y test fnn, y pred fnn))
```

```
print("\n--- FNN Confusion Matrix ---")
cm fnn = confusion matrix(y test fnn, y pred fnn)
plt.figure(figsize=(10, 8))
sns.heatmap(cm fnn, annot=True, fmt="d", cmap="Blues", cbar=False)
plt.title("FNN Confusion Matrix")
plt.xlabel("Predicted Label")
plt.ylabel("True Label")
plt.show()
# Plot Accuracy & Loss
plt.figure(figsize=(12, 5))
plt.subplot(1, 2, 1)
plt.plot(history fnn.history['accuracy'], label='Training Accuracy')
plt.plot(history fnn.history['val accuracy'], label='Validation Accuracy')
plt.title('FNN Model Accuracy')
plt.xlabel('Epoch')
plt.ylabel('Accuracy')
plt.legend()
plt.grid(True)
plt.subplot(1, 2, 2)
plt.plot(history fnn.history['loss'], label='Training Loss')
plt.plot(history fnn.history['val loss'], label='Validation Loss')
plt.title('FNN Model Loss')
plt.xlabel('Epoch')
plt.ylabel('Loss')
plt.legend()
plt.grid(True)
plt.tight layout()
plt.show()
# --- Part 2: Convolutional Neural Network (CNN) ---
print("\n--- Part 2: Implementing a CNN ---")
# 1. Load MNIST for CNN
(x train cnn, y train cnn), (x test cnn, y test cnn) = mnist.load data()
print(f"\nOriginal CNN training data shape: {x train cnn.shape}")
print(f"Original CNN test data shape: {x test cnn.shape}")
```

```
# Reshape for channel dimension
x train cnn = x train cnn.reshape(x train cnn.shape[0], 28, 28, 1)
x test cnn = x test cnn.reshape(x test cnn.shape[0], 28, 28, 1)
# Normalize
x train cnn = x train cnn.astype('float32') / 255.0
x test cnn = x test cnn.astype('float32') / 255.0
print(f''Reshaped & Normalized CNN training data shape: {x train cnn.shape}'')
print(f"Reshaped & Normalized CNN test data shape: {x test cnn.shape}")
num classes cnn = 10
# 2. Build CNN Model
model cnn = keras.Sequential([
  layers.Conv2D(32, (3, 3), activation='relu', input shape=(28, 28, 1)),
  layers.MaxPooling2D((2, 2)),
  layers.Conv2D(64, (3, 3), activation='relu'),
  layers.MaxPooling2D((2, 2)),
  layers.Flatten(),
  layers.Dense(128, activation='relu'),
  layers.Dropout(0.5),
  layers.Dense(num classes cnn, activation='softmax')
])
# 3. Compile Model
model cnn.compile(optimizer='adam',
           loss='sparse categorical crossentropy',
          metrics=['accuracy'])
print("\n--- CNN Model Summary ---")
model cnn.summary()
#4. Train Model
print("\n--- Training CNN Model ---")
history cnn = model cnn.fit(x train cnn, y train cnn, epochs=10,
                 validation split=0.1, verbose=1)
# 5. Evaluate Model
print("\n--- Evaluating CNN Model ---")
```

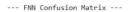
```
loss cnn, accuracy cnn = model cnn.evaluate(x test cnn, y test cnn, verbose=0)
print(f"CNN Test Loss: {loss cnn:.4f}")
print(f"CNN Test Accuracy: {accuracy cnn:.4f}")
# Classification report & confusion matrix
y pred cnn = np.argmax(model cnn.predict(x test cnn), axis=-1)
print("\n--- CNN Classification Report ---")
print(classification report(y test cnn, y pred cnn))
print("\n--- CNN Confusion Matrix ---")
cm cnn = confusion matrix(y test cnn, y pred cnn)
plt.figure(figsize=(10, 8))
sns.heatmap(cm cnn, annot=True, fmt="d", cmap="Blues", cbar=False)
plt.title("CNN Confusion Matrix")
plt.xlabel("Predicted Label")
plt.ylabel("True Label")
plt.show()
# Plot Accuracy & Loss
plt.figure(figsize=(12, 5))
plt.subplot(1, 2, 1)
plt.plot(history cnn.history['accuracy'], label='Training Accuracy')
plt.plot(history cnn.history['val accuracy'], label='Validation Accuracy')
plt.title('CNN Model Accuracy')
plt.xlabel('Epoch')
plt.ylabel('Accuracy')
plt.legend()
plt.grid(True)
plt.subplot(1, 2, 2)
plt.plot(history cnn.history['loss'], label='Training Loss')
plt.plot(history cnn.history['val loss'], label='Validation Loss')
plt.title('CNN Model Loss')
plt.xlabel('Epoch')
plt.ylabel('Loss')
plt.legend()
plt.grid(True)
plt.tight layout()
plt.show()
```

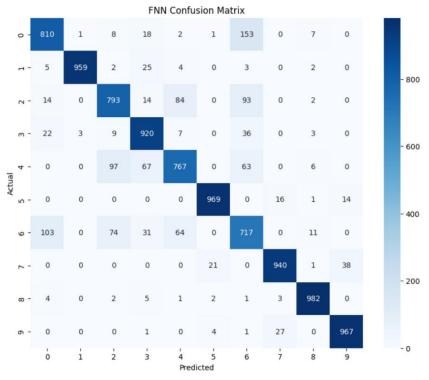
```
# Optional: Visualize predictions
print("\n--- Sample CNN Predictions ---")
class names mnist = [str(i) \text{ for } i \text{ in } range(10)]
plt.figure(figsize=(10, 10))
for i in range(25):
  plt.subplot(5, 5, i + 1)
  plt.xticks([])
  plt.yticks([])
  plt.grid(False)
  plt.imshow(x test cnn[i].reshape(28, 28), cmap=plt.cm.binary)
  true label = y test cnn[i]
  predicted label = y pred cnn[i]
  color = 'green' if true label == predicted label else 'red'
  plt.xlabel(f"True:
                                                       {class names mnist[true label]}\nPred:
{class_names_mnist[predicted_label]}", color=color)
plt.suptitle("Sample CNN Predictions (Green: Correct, Red: Incorrect)", y=1.02, fontsize=16)
plt.tight layout(rect=[0, 0, 1, 0.98])
plt.show()
OUTPUT:
```

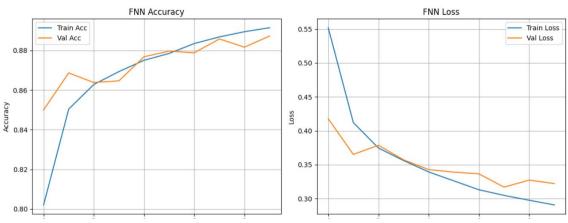
FNN Test Loss: 0.3404 FNN Test Accuracy: 0.8824

--- FNN Classification Report ---

	precision	recall	f1-score	support
0	0.85	0.81	0.83	1000
1	1.00	0.96	0.98	1000
2	0.81	0.79	0.80	1000
3	0.85	0.92	0.88	1000
4	0.83	0.77	0.80	1000
5	0.97	0.97	0.97	1000
6	0.67	0.72	0.69	1000
7	0.95	0.94	0.95	1000
8	0.97	0.98	0.97	1000
9	0.95	0.97	0.96	1000
accuracy			0.88	10000
macro avg	0.88	0.88	0.88	10000
weighted avg	0.88	0.88	0.88	10000

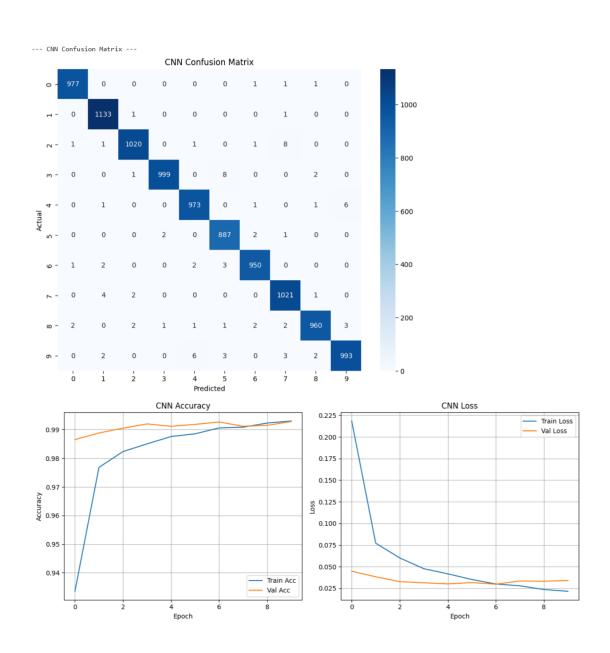




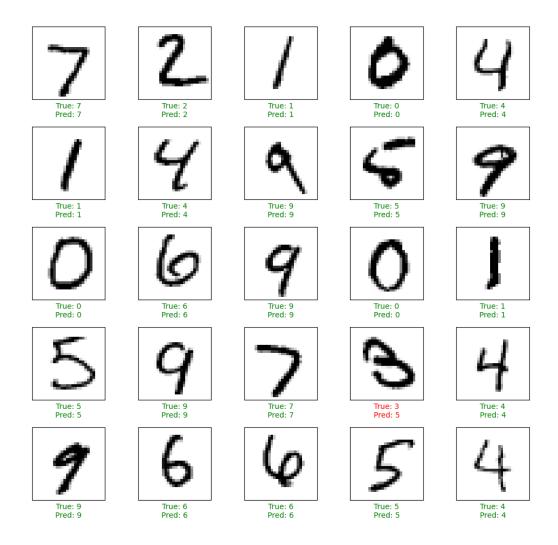


CNN Test Loss: 0.0285 CNN Test Accuracy: 0.9913

CNN	Class	ification R	eport		
		precision	recall	f1-score	support
	0	1.00	1.00	1.00	980
	1	0.99	1.00	0.99	1135
	2	0.99	0.99	0.99	1032
	_				
	3	1.00	0.99	0.99	1010
	4	0.99	0.99	0.99	982
	5	0.98	0.99	0.99	892
	6	0.99	0.99	0.99	958
	7	0.98	0.99	0.99	1028
	8	0.99	0.99	0.99	974
	9	0.99	0.98	0.99	1009
accı	ıracy			0.99	10000
macro	avg	0.99	0.99	0.99	10000
weighted	avg	0.99	0.99	0.99	10000



CNN Predictions (Green = Correct, Red = Incorrect)



Exp No: 7	Generative Models with GANs: Creating and Training a Generative Adversarial Network
	Creating and Training a Generative Adversarial Network

Aim:

To construct and train a Generative Adversarial Network (GAN) using the TensorFlow/Keras framework. The objective is to train the GAN on the MNIST dataset to generate new, synthetic images of handwritten digits that are indistinguishable from the original training data.

Algorithm:

Generative Adversarial Networks (GANs)

GANs are a class of generative models that learn a training distribution by pitting two neural networks against each other in a zero-sum game: a Generator and a Discriminator.

- **1. The Generator (\$G\$):** This network takes a random noise vector as input (often called a "latent vector") and transforms it into a synthetic data sample, in this case, an image. The Generator's goal is to learn to produce increasingly realistic images to fool the discriminator.
- **2.** The Discriminator (\$D\$): This is a binary classifier network. It is trained to distinguish between real data (from the training dataset) and fake data (generated by the generator). Its goal is

to get better at identifying which images are real and which are fake.

3. The Adversarial Process:

Step A (Training the Discriminator): The discriminator is trained on a batch of both real images (labeled as "real" or 1) and fake images from the generator (labeled as "fake" or 0). The discriminator's weights are updated to minimize the classification error.

Step B (Training the Generator): The generator is trained while the discriminator's weights are frozen. The generator creates fake images and feeds them to the discriminator. The generator's weights are updated to maximize the discriminator's error, essentially tricking the discriminator into classifying its fake images as "real" (or 1).

This iterative process continues, with both networks improving, until the generator can produce images so realistic that the discriminator can no longer reliably tell the difference between real and

fake.

CODE:

Import necessary libraries

```
import numpy as np
import matplotlib.pyplot as plt
import tensorflow as tf
from tensorflow import keras
from tensorflow.keras import layers
from tensorflow.keras.datasets import mnist
import os
# Suppress TensorFlow warnings for cleaner output
tf.keras.utils.disable interactive logging()
# --- Part 1: Dataset Loading and Preprocessing ---
print("--- Part 1: Loading and Preprocessing the MNIST Dataset ---")
(x_{train}, ), (, ) = mnist.load_data()
x train = x train.reshape(x train.shape[0], 28, 28, 1).astype('float32')
x train = (x \text{ train} - 127.5) / 127.5 \# \text{Normalize to } [-1, 1]
print(f"Normalized training data shape: {x train.shape}")
print("Example of a normalized pixel value:", x train[0, 0, 0, 0])
# --- Part 2: Building the Generator and Discriminator Models ---
print("\n--- Part 2: Building the GAN Components ---")
latent dim = 100
# Generator
def build generator():
  model = keras. Sequential(name="generator")
  model.add(layers.Dense(7 * 7 * 256, use bias=False, input shape=(latent dim,)))
  model.add(layers.BatchNormalization())
  model.add(layers.LeakyReLU())
  model.add(layers.Reshape((7, 7, 256)))
  model.add(layers.Conv2DTranspose(128,
                                              (5,
                                                  5),
                                                         strides=(1, 1),
                                                                            padding='same',
use bias=False))
  model.add(layers.BatchNormalization())
  model.add(layers.LeakyReLU())
  model.add(layers.Conv2DTranspose(64,
                                             (5,
                                                   5),
                                                         strides=(2,
                                                                      2),
                                                                            padding='same',
use bias=False))
```

```
model.add(layers.BatchNormalization())
  model.add(layers.LeakyReLU())
  model.add(layers.Conv2DTranspose(1, (5, 5), strides=(2, 2), padding='same',
                      use bias=False, activation='tanh'))
  return model
generator = build generator()
print("\n--- Generator Model Summary ---")
generator.summary()
# Discriminator
def build discriminator():
  model = keras.Sequential(name="discriminator")
  model.add(layers.Conv2D(64, (5, 5), strides=(2, 2), padding='same', input shape=[28, 28,
1]))
  model.add(layers.LeakyReLU())
  model.add(layers.Dropout(0.3))
  model.add(layers.Conv2D(128, (5, 5), strides=(2, 2), padding='same'))
  model.add(layers.LeakyReLU())
  model.add(layers.Dropout(0.3))
  model.add(layers.Flatten())
  model.add(layers.Dense(1, activation='sigmoid'))
  return model
discriminator = build discriminator()
print("\n--- Discriminator Model Summary ---")
discriminator.summary()
# --- Part 3: Training Setup ---
cross entropy = keras.losses.BinaryCrossentropy(from logits=False)
def discriminator loss(real output, fake output):
  real loss = cross entropy(tf.ones like(real output), real output)
  fake loss = cross entropy(tf.zeros like(fake output), fake output)
  return real loss + fake loss
def generator loss(fake output):
  return cross entropy(tf.ones like(fake output), fake output)
generator optimizer = tf.keras.optimizers.Adam(learning rate=1e-4)
```

```
discriminator optimizer = tf.keras.optimizers.Adam(learning rate=1e-4)
@tf.function
def train step(images, latent dim=latent dim):
  noise = tf.random.normal([batch size, latent dim])
  with tf.GradientTape() as gen tape, tf.GradientTape() as disc tape:
    generated images = generator(noise, training=True)
    real output = discriminator(images, training=True)
    fake output = discriminator(generated images, training=True)
    gen loss = generator loss(fake output)
    disc loss = discriminator loss(real output, fake output)
  gradients of generator = gen tape.gradient(gen loss, generator.trainable variables)
  gradients of discriminator
                                                                disc tape.gradient(disc loss,
discriminator.trainable variables)
  generator optimizer.apply gradients(zip(gradients of generator,
generator.trainable variables))
  discriminator optimizer.apply gradients(zip(gradients of discriminator,
discriminator.trainable variables))
  return gen loss, disc loss
def generate and save images(model, epoch, test input):
  predictions = model(test input, training=False)
  predictions rescaled = (predictions * 0.5) + 0.5 # Scale back to [0, 1]
  fig = plt.figure(figsize=(4, 4))
  for i in range(predictions.shape[0]):
    plt.subplot(4, 4, i + 1)
    plt.imshow(predictions rescaled[i, :, :, 0], cmap='gray')
    plt.axis('off')
  plt.suptitle(f"Epoch {epoch}", fontsize=16)
  if not os.path.exists('images'):
    os.makedirs('images')
  plt.savefig(fimages/image at epoch {epoch:04d}.png')
  plt.show()
# Training parameters
EPOCHS = 200
batch size = 256
num examples to generate = 16
seed = tf.random.normal([num examples to generate, latent dim])
```

```
train dataset
tf.data.Dataset.from tensor slices(x train).shuffle(x train.shape[0]).batch(batch size)
# Training loop
def train(dataset, epochs):
  print("\n--- Beginning GAN Training ---")
  for epoch in range(epochs):
    gen loss list = []
    disc loss list = []
    for image batch in dataset:
       gen loss, disc loss = train step(image batch)
       gen loss list.append(gen loss.numpy())
       disc loss list.append(disc loss.numpy())
    avg gen loss = np.mean(gen loss list)
    avg disc loss = np.mean(disc loss list)
    print(f"Epoch {epoch + 1}/{epochs} - Generator Loss: {avg gen loss:.4f}, Discriminator
Loss: {avg disc loss:.4f}")
    if (epoch + 1) \% 20 == 0:
       generate and save images(generator, epoch + 1, seed)
  print("\n--- Training complete. Generating final images. ---")
  generate_and_save_images(generator, epochs, seed)
# Run training
train(train dataset, EPOCHS)
```

OUTPUT:

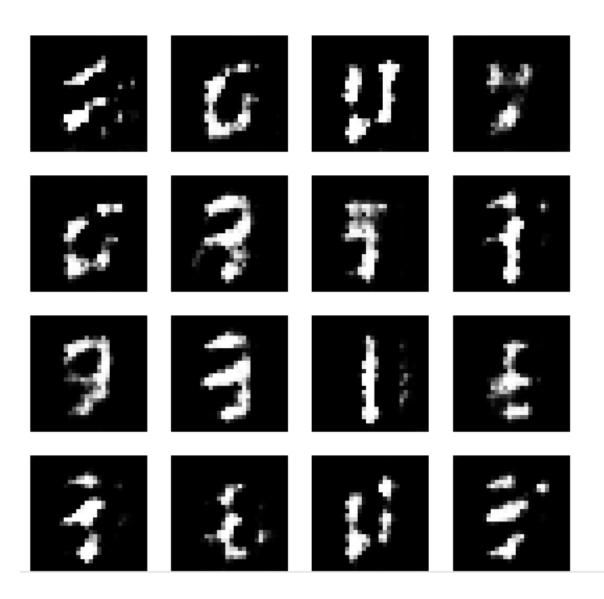
--- Part 1: Loading and Preprocessing the MNIST Dataset --- Normalized training data shape: (60000, 28, 28, 1)

Example normalized pixel value: -1.0

--- Beginning GAN Training ---

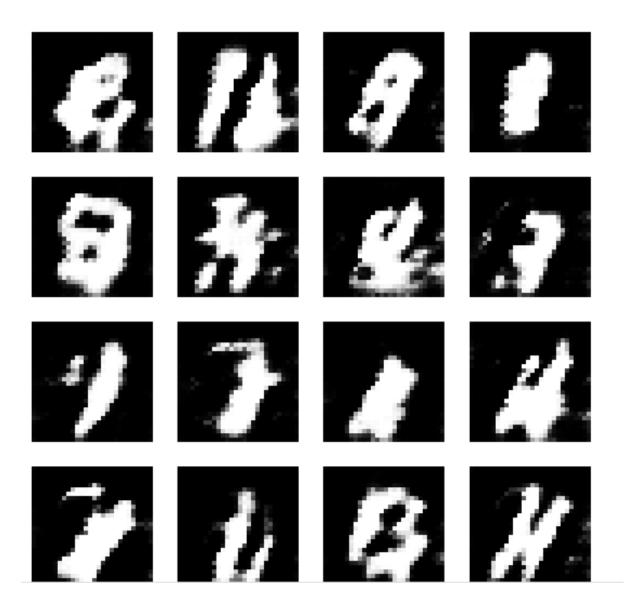
Epoch 1/20 - Generator Loss: 0.7877, Discriminator Loss: 1.0228 Epoch 2/20 - Generator Loss: 0.8148, Discriminator Loss: 1.2225 Epoch 3/20 - Generator Loss: 0.8448, Discriminator Loss: 1.3034 Epoch 4/20 - Generator Loss: 0.8534, Discriminator Loss: 1.2366 Epoch 5/20 - Generator Loss: 0.8372, Discriminator Loss: 1.2497

Epoch 5



Epoch 6/20 - Generator Loss: 0.8516, Discriminator Loss: 1.2705 Epoch 7/20 - Generator Loss: 0.8888, Discriminator Loss: 1.3028 Epoch 8/20 - Generator Loss: 0.8739, Discriminator Loss: 1.2512 Epoch 9/20 - Generator Loss: 0.8691, Discriminator Loss: 1.3130 Epoch 10/20 - Generator Loss: 0.8862, Discriminator Loss: 1.2320

Epoch 10



```
Epoch 11/20 - Generator Loss: 0.9361, Discriminator Loss: 1.2244

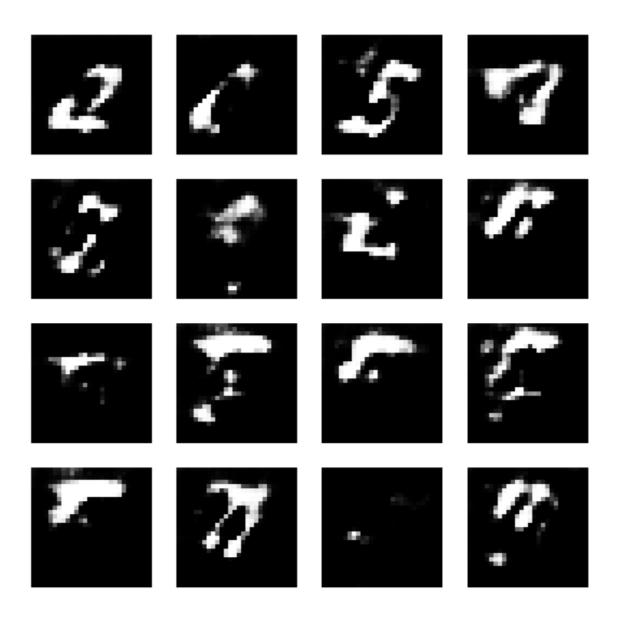
Epoch 12/20 - Generator Loss: 0.9946, Discriminator Loss: 1.1719

Epoch 13/20 - Generator Loss: 0.9948, Discriminator Loss: 1.1944

Epoch 14/20 - Generator Loss: 0.9786, Discriminator Loss: 1.1809

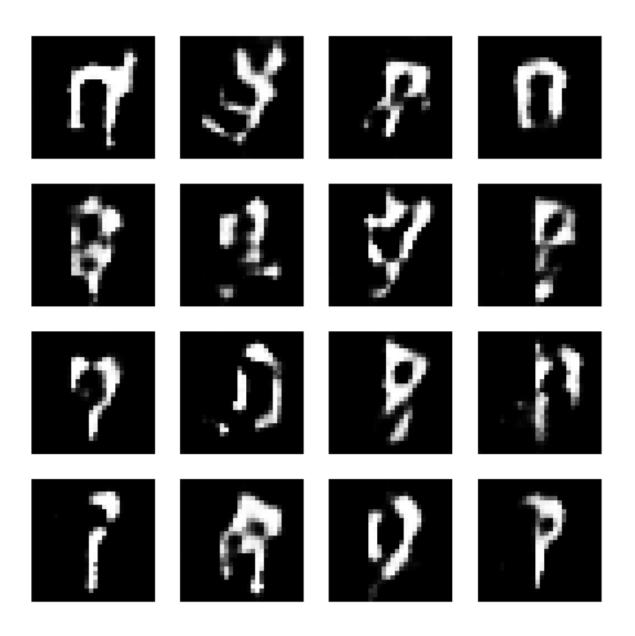
Epoch 15/20 - Generator Loss: 1.0420, Discriminator Loss: 1.1079
```

Epoch 15

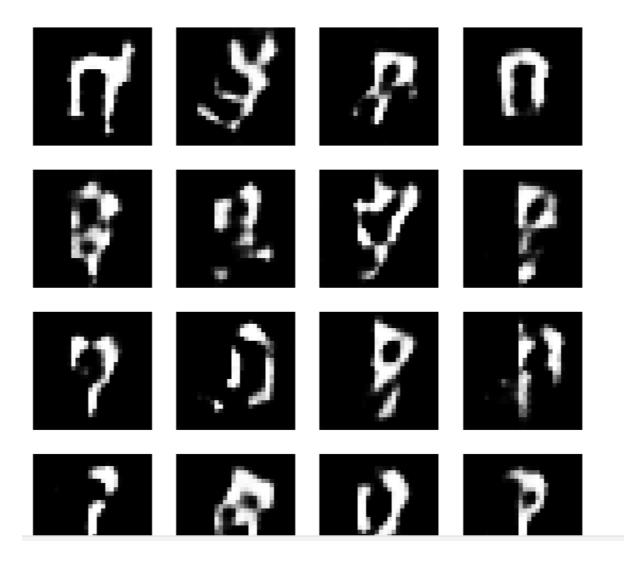


Epoch 16/20 - Generator Loss: 1.2020, Discriminator Loss: 1.0483 Epoch 17/20 - Generator Loss: 1.2648, Discriminator Loss: 1.0605 Epoch 18/20 - Generator Loss: 1.1657, Discriminator Loss: 1.0404 Epoch 19/20 - Generator Loss: 1.1644, Discriminator Loss: 1.0897 Epoch 20/20 - Generator Loss: 1.1770, Discriminator Loss: 1.0938

Epoch 20



Epoch 20



Exp No: 8	Model Evaluation and Improvement: Hyperparameter Tuning with Grid Search and Cross-Validation

Aim:

To demonstrate key techniques for model evaluation and improvement:

- 1. Hyperparameter Tuning with Grid Search: Systematically searching for the optimal combination of hyperparameters for a machine learning model.
- **2.** Cross-Validation Techniques: Implementing k-fold cross-validation to get a more robust estimate of model performance and to prevent overfitting to a specific train-test split.

Algorithm:

1. Hyperparameter Tuning with Grid Search

Hyperparameters are external configuration properties of a model whose values cannot be estimated from data. Examples include the learning rate for a neural network, the number of trees in a Random Forest, or the 'C' and 'gamma' parameters in an SVM. Tuning these parameters is crucial for optimal model performance.

Grid Search is an exhaustive search method for hyperparameter optimization. **Steps:**

- 1. Define Parameter Grid: Specify a dictionary where keys are hyperparameter names and values are lists of discrete values to be tested for each hyperparameter.
- 2. Instantiate Model: Choose a machine learning model.
- 3. Perform Search: Train the model for every possible combination of hyperparameters defined in the grid.
- 4. Evaluate: For each combination, evaluate the model's performance using a specified scoring

metric (e.g., accuracy, F1-score) and often in conjunction with cross-validation.

5. Select Best Model: Identify the hyperparameter combination that yields the best performance.

2. Cross-Validation Techniques

Cross-validation is a resampling procedure used to evaluate machine learning models on a limited data sample. The goal is to estimate how accurately a predictive model will perform in practice. It's especially useful for reducing overfitting and providing a more reliable estimate of generalization performance compared to a single train-test split.

k-Fold Cross-Validation:

Steps:

- 1. Divide Data: The entire dataset is randomly partitioned into \$k\$ equally sized subsamples (or "folds").
- 2. Iterate \$k\$ Times:

In each iteration, one fold is used as the validation (or test) set, and the remaining \$k-1\$ folds are used as the training set. The model is trained on the training set and evaluated on the validation set.

- 3. Aggregate Results: The performance metric (e.g., accuracy) from each of the \$k\$ iterations is collected.
- 4. Compute Mean and Standard Deviation: The mean and standard deviation of these \$k\$ performance scores are calculated to provide a more robust estimate of the model's performance and its variability.

CODE:

Import necessary libraries
import numpy as np
import pandas as pd
import matplotlib.pyplot as plt
import seaborn as sns
from sklearn.datasets import load_iris # A classic dataset for classification
from sklearn.model_selection import train_test_split, KFold, cross_val_score, GridSearchCV
from sklearn.svm import SVC # Support Vector Classifier, a common model for tuning
from sklearn.metrics import accuracy_score, classification_report, confusion_matrix
from sklearn.preprocessing import StandardScaler

--- Part 1: Hyperparameter Tuning with Grid Search ---

```
print("--- Part 1: Hyperparameter Tuning with Grid Search ---")

# 1. Load a Dataset (Iris Dataset for classification)

# The Iris dataset is a classic and simple dataset for classification tasks.

# It contains measurements of iris flowers (sepal length, sepal width, petal length, petal width)

# and their corresponding species (Setosa, Versicolor, Virginica).

iris = load_iris()

X = iris.data

y = iris.target

feature_names = iris.feature_names

target_names = iris.target_names
```

```
print(f"\nDataset Features (X) shape: {X.shape}")
print(f"Dataset Labels (y) shape: {y.shape}")
print(f"Feature Names: {feature names}")
print(f"Target Names: {target names}")
# 2. Split Data into Training and Testing Sets
# It's crucial to split the data before scaling to prevent data leakage.
# The test set will be used for final model evaluation, after tuning.
X train, X test, y train, y test = train test split(X, y, test size=0.3, random state=42,
stratify=y)
print(f"\nTraining set size: {X train.shape[0]} samples")
print(f"Test set size: {X test.shape[0]} samples")
# 3. Standardize Features
# Scaling features is important for SVMs as they are sensitive to feature scales.
# Fit scaler only on training data to prevent data leakage.
scaler = StandardScaler()
X train scaled = scaler.fit transform(X train)
X test scaled = scaler.transform(X test)
print("\nFeatures standardized.")
# 4. Define the Model and Hyperparameter Grid
# We'll use a Support Vector Classifier (SVC) as our model.
# Common hyperparameters for SVC are 'C' (regularization parameter) and 'gamma' (kernel
coefficient).
# 'kernel' also can be tuned (e.g., 'linear', 'rbf').
# Define the parameter grid for Grid Search
param grid = {
  'C': [0.1, 1, 10, 100],
                              # Regularization parameter
  'gamma': [1, 0.1, 0.01, 0.001], # Kernel coefficient for 'rbf', 'poly' and 'sigmoid'
  'kernel': ['rbf', 'linear']
                             # Type of kernel function
}
print("\nHyperparameter grid defined:")
for param, values in param grid.items():
  print(f" {param}: {values}")
```

```
# 5. Perform Grid Search with Cross-Validation
# GridSearchCV automatically performs k-fold cross-validation for each combination.
# cv=5 means 5-fold cross-validation.
# scoring='accuracy' means we want to optimize for accuracy.
grid search = GridSearchCV(SVC(), param grid, cv=5, scoring='accuracy', verbose=1,
n jobs=-1
print("\nStarting Grid Search with 5-fold Cross-Validation...")
# Fit GridSearchCV on the scaled training data
grid search.fit(X train scaled, y train)
print("\nGrid Search completed.")
# 6. Get the Best Parameters and Best Score
print(f"\nBest hyperparameters found: {grid search.best params }")
print(f"Best cross-validation accuracy: {grid search.best score :.4f}")
# 7. Evaluate the Best Model on the Test Set
# The best estimator attribute provides the model trained with the best parameters.
best model = grid search.best estimator
y pred tuned = best model.predict(X test scaled)
test accuracy tuned = accuracy_score(y_test, y_pred_tuned)
print(f"\nTest set accuracy with tuned model: {test accuracy tuned:.4f}")
print("\n--- Classification Report for Tuned Model ---")
print(classification report(y test, y pred tuned, target names=target names))
print("\n--- Confusion Matrix for Tuned Model ---")
cm tuned = confusion matrix(y test, y pred tuned)
plt.figure(figsize=(8, 6))
sns.heatmap(cm tuned, annot=True, fmt='d', cmap='Blues', xticklabels=target names,
yticklabels=target names)
plt.title('Confusion Matrix (Tuned SVM)')
plt.xlabel('Predicted Label')
plt.ylabel('True Label')
plt.show()
# Visualize Grid Search results (optional, but good for understanding)
```

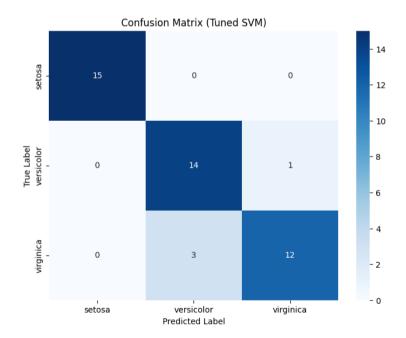
```
# Convert results to a DataFrame for easier analysis
results df = pd.DataFrame(grid search.cv results)
print("\n--- Top 5 Grid Search Results ---")
print(results df[['param C',
                                 'param gamma',
                                                       'param kernel',
                                                                           'mean test score',
'rank test score']].sort values(by='rank test score').head())
# --- Part 2: Cross-Validation Techniques (k-fold) ---
print("\n--- Part 2: Cross-Validation Techniques (k-fold) ---")
# We will demonstrate k-fold cross-validation on a simple SVM without explicit tuning for
clarity,
# to focus solely on the CV process.
# 1. Instantiate a Model (using default or chosen parameters)
model cv = SVC(random state=42) # Using default parameters for simplicity
# 2. Define k-fold Cross-Validation Strategy
# We'll use 5-fold cross-validation.
# KFold ensures that each fold is distinct.
# shuffle=True means the data will be randomly shuffled before splitting into folds.
# random state for reproducibility.
k 	ext{ folds} = 5
kf = KFold(n splits=k folds, shuffle=True, random state=42)
print(f"\nPerforming {k folds}-fold cross-validation...")
# 3. Perform Cross-Validation and Get Scores
# cross val score performs the KFold splitting, training, and evaluation automatically.
# It returns an array of scores, one for each fold.
cv scores = cross val score(model cv, X train scaled, y train, cv=kf, scoring='accuracy')
print(f"\nCross-validation scores for each fold: {cv scores}")
print(f"Mean cross-validation accuracy: {np.mean(cv scores):.4f}")
print(f'Standard deviation of cross-validation accuracy: {np.std(cv scores):.4f}")
# 4. Visualize Cross-Validation Scores
plt.figure(figsize=(8, 5))
plt.bar(range(1, k folds + 1), cv scores, color='skyblue')
plt.axhline(y=np.mean(cv scores),
                                      color='r',
                                                  linestyle='--',
                                                                  label=f'Mean
                                                                                    Accuracy
```

```
({np.mean(cv_scores):.4f})')
plt.title(f'{k_folds}-Fold Cross-Validation Accuracy Scores')
plt.xlabel('Fold Number')
plt.ylabel('Accuracy')
plt.ylim(0.8, 1.0) # Set y-axis limits for better visualization
plt.legend()
plt.grid(axis='y', linestyle='--')
plt.show()

# 5. Discuss why CV is useful
print("\n--- Why is Cross-Validation Important? ---")
print("1. More Reliable Performance Estimate: Reduces bias from a single train-test split.")
print("2. Better Generalization: Helps ensure the model performs well on unseen data.")
print("3. Efficient Data Usage: All data points are used for both training and validation across different folds.")
print("4. Detects Overfitting/Underfitting: Variability in scores can indicate instability.")
```

OUTPUT:

```
--- Part 1: Hyperparameter Tuning with Grid Search ---
Dataset Features (X) shape: (150, 4)
Dataset Labels (y) shape: (150,)
Feature Names: ['sepal length (cm)', 'sepal width (cm)', 'petal length (cm)', 'petal width (cm)']
Target Names: ['setosa' 'versicolor' 'virginica']
Training set size: 105 samples
Test set size: 45 samples
Features standardized.
Hyperparameter grid defined:
  C: [0.1, 1, 10, 100]
  gamma: [1, 0.1, 0.01, 0.001]
kernel: ['rbf', 'linear']
Starting Grid Search with 5-fold Cross-Validation...
Fitting 5 folds for each of 32 candidates, totalling 160 fits
Grid Search completed.
Best hyperparameters found: {'C': 1, 'gamma': 0.1, 'kernel': 'rbf'}
Best cross-validation accuracy: 0.9810
Test set accuracy with tuned model: 0.9111
--- Classification Report for Tuned Model ---
              precision recall f1-score support
      setosa
                    1.00
                           1.00
                                         1.00
                 0.82 0.93
0.92 0.80
  versicolor
                                        0.88
                                                      15
                             0.80
   virginica
                                         0.86
                                         0.91
                                                      45
   accuracy
macro avg 0.92 0.91
weighted avg 0.92 0.91
                                         0.91
                                                      45
                                         0.91
                                                      45
```

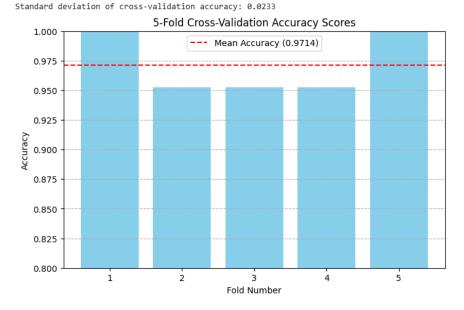


```
--- Top 5 Grid Search Results ---
    .
param_C param_gamma param_kernel mean_test_score rank_test_score
10
       1.0
                   0.100
                                 rbf
                                              0.980952
27
      100.0
                   0.100
                               linear
                                              0.980952
31
      100.0
                   0.001
                               linear
                                              0.980952
29
      100.0
                   0.010
                               linear
                                              0.980952
                                                                      1
                                              0.980952
25
      100.0
                   1.000
                               linear
```

--- Part 2: Cross-Validation Techniques (k-fold) ---

Performing 5-fold cross-validation...

Cross-validation scores for each fold: [1. 0.95238095 0.95238095 0.95238095 1. Mean cross-validation accuracy: 0.9714



- --- Why is Cross-Validation Important? ---
- 1. More Reliable Performance Estimate: Reduces bias from a single train-test split.
- 2. Better Generalization: Helps ensure the model performs well on unseen data.
- 3. Efficient Data Usage: All data points are used for both training and validation across different folds.
- 4. Detects Overfitting/Underfitting: Variability in scores can indicate instability.

Exp No: 9	MINI PROJECT

ABSTRACT:

Customer segmentation is a vital strategy in modern retail analytics, helping businesses understand and target their customers effectively. This project, Customer Segmentation for Retail Using K-Means and PCA, aims to group customers based on purchasing patterns and demographic attributes to support data-driven marketing decisions. By identifying segments such as high-income high-spenders or budget-conscious customers, retailers can design personalized offers and improve overall customer satisfaction.

The project employs the K-Means clustering algorithm to discover natural customer groupings and uses Principal Component Analysis (PCA) for dimensionality reduction and visual interpretation of clusters. The dataset is preprocessed by handling missing values, encoding categorical data, and standardizing numerical features. The Elbow Method and Silhouette Score are applied to determine the optimal number of clusters, ensuring meaningful segmentation and strong cluster separation.

An interactive Streamlit application is developed to make the model accessible and user-friendly. The app allows users to upload data, adjust clustering parameters, visualize customer clusters in real time, and predict the segment of new customers. This project demonstrates how unsupervised learning can transform raw retail data into actionable business insights, enabling smarter marketing strategies and more efficient customer relationship management.

CODE:

```
import streamlit as st
import pandas as pd
import numpy as np
import plotly.express as px
from sklearn.cluster import KMeans
from sklearn.preprocessing import StandardScaler, OneHotEncoder
from sklearn.impute import SimpleImputer
from sklearn.decomposition import PCA
from sklearn.metrics import silhouette score
from sklearn.pipeline import Pipeline
from sklearn.compose import ColumnTransformer
import io
# --- App Configuration ---
st.set page config(
  page title="Retail Customer Segmentation",
  page icon="ba",
```

```
layout="wide"
)

# --- Abstract ---
APP ABSTRACT = """
```

Customer segmentation is a vital strategy in modern retail analytics, helping businesses understand and target their customers effectively. This project, Customer Segmentation for Retail Using K-Means and PCA, aims to group customers based on purchasing patterns and demographic attributes to support datadriven marketing decisions. By identifying segments such as high-income high-spenders or budget-conscious customers, retailers can design personalized offers and improve overall customer satisfaction.

The project employs the K-Means clustering algorithm to discover natural customer groupings and uses Principal Component Analysis (PCA) for dimensionality reduction and visual interpretation of clusters. The dataset is preprocessed by handling missing values, encoding categorical data, and standardizing numerical features. The Elbow Method and Silhouette Score are applied to determine the optimal number of clusters, ensuring meaningful segmentation and strong cluster separation.

An interactive Streamlit application is developed to make the model accessible and user-friendly. The app allows users to upload data, adjust clustering parameters, visualize customer clusters in real time, and predict the segment of new customers. This project demonstrates how unsupervised learning can transform raw retail data into actionable business insights, enabling smarter marketing strategies and more efficient customer relationship management.

```
# --- Helper Functions ---
@st.cache data
def load default data():
  """Loads the default Mall Customers dataset."""
  csv data = """CustomerID, Gender, Age, Annual Income (k$), Spending Score (1-100)
1,Male,19,15,39
2, Male, 21, 15, 81
3,Female,20,16,6
4,Female,23,16,77
5, Female, 31, 17, 40
6,Female,22,17,76
7,Female,35,18,6
8,Female,23,18,94
9,Male,64,19,3
10, Female, 30, 19, 72
11, Male, 67, 19, 14
12,Female,35,19,99
13,Female,58,20,15
14, Female, 24, 20, 77
```

```
15, Male, 37, 20, 13
16,Male,22,20,79
17,Female,35,21,35
18, Male, 20, 21, 66
19, Male, 52, 23, 29
20,Female,35,23,98
21,Male,35,24,35
22,Male,25,24,73
23,Female,46,25,5
24,Male,31,25,73
25,Female,54,28,14
26,Male,29,28,82
27,Female,45,28,32
28,Male,35,28,61
29,Female,40,29,31
30,Female,23,29,87
31, Male, 60, 30, 4
32,Female,21,30,73
33,Male,53,33,4
34, Female, 18, 33, 92
35, Female, 49, 33, 14
36,Female,21,33,81
37, Female, 42, 34, 17
38, Female, 30, 34, 73
39,Female,36,37,26
40,Female,65,38,35
41,Female,20,38,92
42, Male, 48, 39, 36
43,Female,24,39,65
44,Female,31,39,8
45,Female,49,39,28
46,Female,24,39,68
47,Female,50,40,55
48, Female, 27, 40, 47
49,Female,29,40,42
50, Female, 31, 40, 42
  return pd.read csv(io.StringIO(csv data))
@st.cache resource
def create preprocessing pipeline(df):
  """Creates a preprocessing pipeline based on dataframe columns."""
```

Identify numeric and categorical features (exclude CustomerID)

```
numeric features = df.select dtypes(include=np.number).columns.tolist()
  if 'CustomerID' in numeric features:
     numeric features.remove('CustomerID')
  categorical features = df.select dtypes(include=['object', 'category']).columns.tolist()
  # Create transformers
  numeric transformer = Pipeline(steps=[
     ('imputer', SimpleImputer(strategy='median')),
     ('scaler', StandardScaler())
  1)
  categorical transformer = Pipeline(steps=[
     ('imputer', SimpleImputer(strategy='most frequent')),
     ('onehot', OneHotEncoder(handle unknown='ignore'))
  ])
  # Create the full preprocessor
  preprocessor = ColumnTransformer(
     transformers=[
       ('num', numeric transformer, numeric features),
       ('cat', categorical transformer, categorical features)
     ],
     remainder='passthrough'
  )
  return preprocessor, numeric features, categorical features
@st.cache data
def calculate cluster metrics( processed data):
  """Calculates inertia and silhouette scores for a range of k."""
  inertia = []
  silhouette scores = []
  k range = range(2, 11)
  for k in k range:
     kmeans = KMeans(n clusters=k, init='k-means++', n init=10, random state=42)
     kmeans.fit( processed data)
     inertia.append(kmeans.inertia)
     score = silhouette score(processed data, kmeans.labels)
     silhouette scores.append(score)
  metrics df = pd.DataFrame({
     'K': k range,
```

```
'Inertia (Elbow)': inertia,
     'Silhouette Score': silhouette scores
  })
  return metrics df
# --- Sidebar Controls ---
st.sidebar.title(" Segmentation Controls")
st.sidebar.markdown("Configure the analysis parameters.")
uploaded file = st.sidebar.file uploader("Upload your own CSV", type=["csv"])
if uploaded file is not None:
  df = pd.read csv(uploaded file)
else:
  st.sidebar.info("Using default 'Mall Customers' dataset. Upload a CSV to use your own data.")
  df = load default data()
# Drop CustomerID if it exists, as it's just an identifier
if 'CustomerID' in df.columns:
  df analysis = df.drop(columns=['CustomerID'])
else:
  df analysis = df.copy()
# Create pipeline and process data
try:
  preprocessor, numeric cols, cat cols = create preprocessing pipeline(df analysis)
  processed data = preprocessor.fit transform(df analysis)
  st.sidebar.markdown("---")
  st.sidebar.header("Model Parameters")
  k slider = st.sidebar.slider(
     "Select Number of Clusters (K)",
     min value=2,
     max value=10,
     value=5,
     step=1,
     help="Use the 'Optimal K Analysis' tab to find the best value for K."
  # --- Main Application ---
  st.title("Retail Customer Segmentation Dashboard")
  st.info(APP ABSTRACT)
  tab1, tab2, tab3, tab4 = st.tabs([
```

```
"Data Overview".
     "Optimal K Analysis",
     "Segmentation Results",
     "Predict New Customer"
  1)
  # --- Tab 1: Data Overview ---
  with tab1:
     st.header("Data Preview")
     st.dataframe(df.head(), width='stretch')
     st.header("Data Statistics")
     st.dataframe(df.describe(), width='stretch')
     st.header("Features Identified")
     st.markdown(f"*Numeric Features:* {', '.join(numeric cols)}")
     st.markdown(f"*Categorical Features:* {', '.join(cat cols)}")
  # --- Tab 2: Optimal K Analysis ---
  with tab2:
     st.header("Finding the Optimal Number of Clusters (K)")
     st.markdown("""
     We use two methods to find the best 'K':
     1. *Elbow Method:* Look for the "elbow" point where the inertia (sum of squared distances) stops
decreasing rapidly.
     2. *Silhouette Score:* Measures how similar an object is to its own cluster compared to other
clusters. A higher score (closer to 1) is better.
     ("""
     with st.spinner("Calculating cluster metrics for K=2 to 10..."):
       metrics df = calculate cluster metrics(processed data)
     col1, col2 = st.columns(2)
     with col1:
       st.subheader("Elbow Method (Inertia)")
       fig elbow = px.line(metrics df, x='K', y='Inertia (Elbow)', title='Elbow Method',
markers=True)
       st.plotly chart(fig elbow, use container width=True) # Note: Keeping use container width as
'width' is not a direct replacement for plotly
     with col2:
       st.subheader("Silhouette Score")
       fig sil = px.line(metrics df, x='K', y='Silhouette Score', title='Silhouette Score', markers=True)
```

```
st.plotly_chart(fig_sil, use_container_width=True) # Note: Keeping use_container_width as 'width' is not a direct replacement for plotly
```

```
# --- Tab 3: Segmentation Results ---
  with tab3:
     st.header(f''Customer Segments (K={k slider})")
     # Run K-Means with the selected K
     kmeans model = KMeans(n clusters=k slider, init='k-means+++', n init=10, random state=42)
     kmeans model.fit(processed data)
     # Add cluster labels to the original dataframe
     df results = df analysis.copy()
     df results['Cluster'] = kmeans model.labels .astype(str) # Convert to string for Plotly
     # Run PCA for visualization
     pca = PCA(n components=2, random state=42)
     pca features = pca.fit transform(processed data)
     df results['PC1'] = pca features[:, 0]
     df results['PC2'] = pca features[:, 1]
     # Create PCA Scatter Plot
     st.subheader("Customer Segments (PCA Visualization)")
     hover cols = numeric cols + cat cols + ['Cluster']
     fig pca = px.scatter(
       df results,
       x='PC1',
       y='PC2'
       color='Cluster',
       title=f"2D PCA Visualization of {k slider} Clusters",
       hover data=[col for col in hover cols if col in df results.columns],
       color discrete sequence=px.colors.qualitative.Plotly
     st.plotly chart(fig pca, use container width=True) # Note: Keeping use container width as
'width' is not a direct replacement for plotly
     # Show Cluster Profiles
     st.subheader("Cluster Profiles (Mean Values)")
     st.markdown("This table shows the average values for each numeric feature by cluster.")
     cluster profile = df results.groupby('Cluster')[numeric cols].mean().reset index()
     st.dataframe(cluster profile, width='stretch')
  # --- Tab 4: Predict New Customer ---
```

```
with tab4:
  st.header("Predict New Customer Segment")
  st.markdown("Enter the details of a new customer to predict their segment.")
  # Use a form for input
  with st.form(key="prediction form"):
     input data = \{\}
    # Create input fields dynamically
     for col in numeric cols:
       input data[col] = st.number input(
         f"Enter {col}",
         value=float(df analysis[col].median())
       )
     for col in cat cols:
       unique vals = df analysis[col].unique().tolist()
       input data[col] = st.selectbox(
         f"Select {col}",
         options=unique vals,
         index=0
       )
    submit button = st.form submit button(label="Predict Segment")
  if submit button:
    # Create a DataFrame from the input
    new customer df = pd.DataFrame([input data])
    # Preprocess the new data
    try:
       new data processed = preprocessor.transform(new customer df)
       # Predict the cluster
       prediction = kmeans model.predict(new data processed)
       st.success(f"This customer belongs to *Cluster {prediction[0]}*")
       # Show profile of the predicted cluster
       st.subheader(f"Profile of Predicted Cluster {prediction[0]}")
       st.dataframe(cluster profile[cluster profile['Cluster'] == str(prediction[0])], width='stretch')
    except Exception as e:
       st.error(f"An error occurred during prediction: {e}")
```

except Exception as e:

st.error(f"An error occurred during data processing: {e}")

st.error("Please check your uploaded CSV file. It may have an incorrect format, missing columns, or data types that don't match the expected schema (numeric and categorical).")

Output:

