EXP NO: 4) B

#### **ENSEMBLE METHODS: RANDOM FOREST**

DATE: 21/08/2025

#### AIM:

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

#### **ALGORITHM:**

**STEP 1**: Import libraries.

**STEP 2**: Load data (use same dataset to compare with SVM).

**STEP 3**: Train/Test split with stratification.

STEP 4: (Optional) Preprocess: Random Forests don't require scaling; we'll use raw features.

STEP 5: Model: RandomForestClassifier.

**STEP 6**: Hyperparameter tuning: Grid search over n\_estimators, max\_depth, min\_samples\_split, min\_samples leaf.

**STEP 7**: Train the best model on training data.

STEP 8: Evaluate with Accuracy, Precision, Recall, F1, Confusion Matrix, ROC-AUC.

**STEP 9**: Interpretation: Plot top feature importances.

### **CODE:**

```
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()
```

# **OUTPUT:**

```
Best Parameters (CV): {'max_depth': None, 'min_samples_leaf': 1, 'min_samples_split': 2, 'n_estimators': 100}
```

=== Random Forest - Test Metrics ===

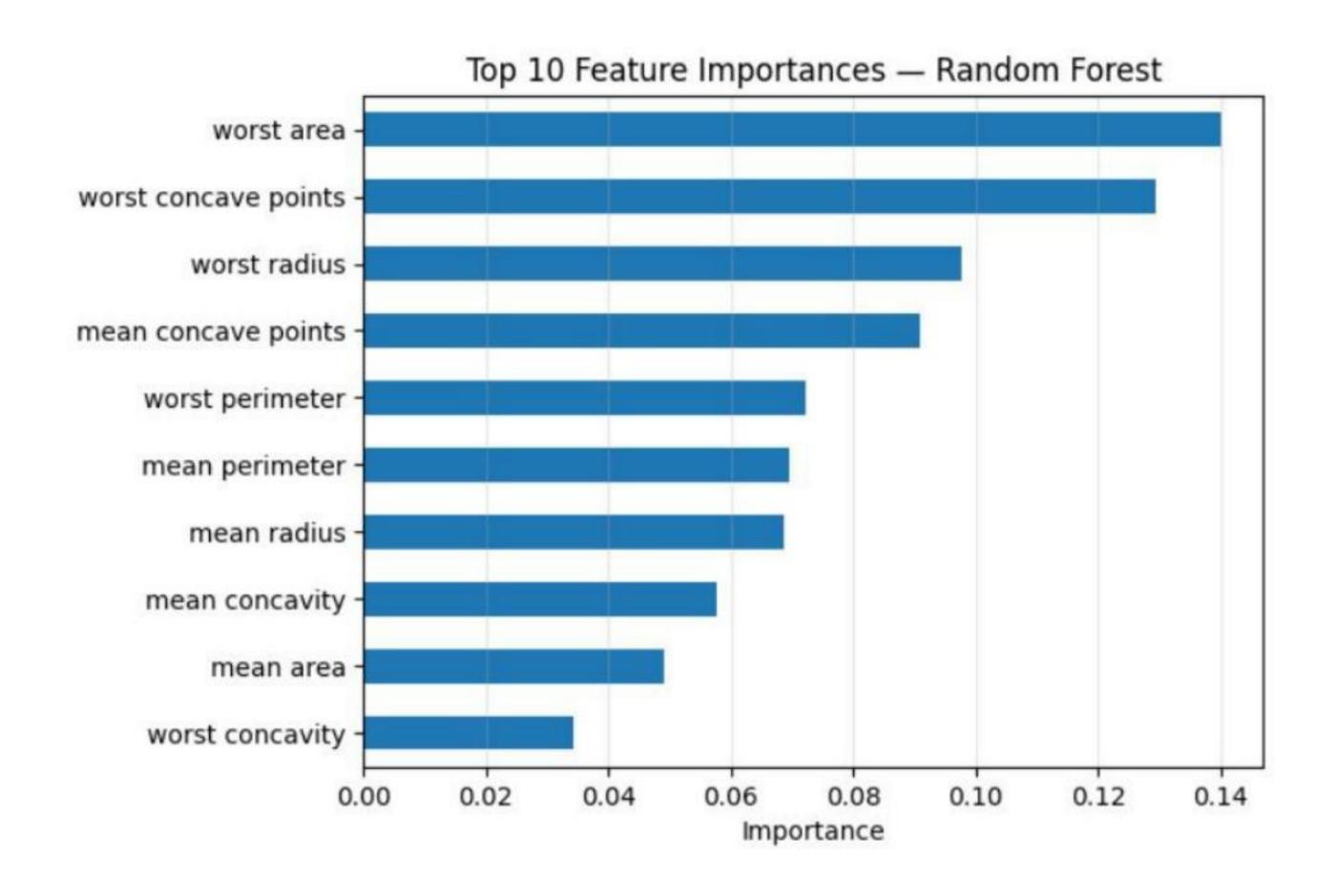
Accuracy: 0.9561
Precision: 0.9589
Recall: 0.9722
F1-Score: 0.9655
ROC-AUC: 0.9937

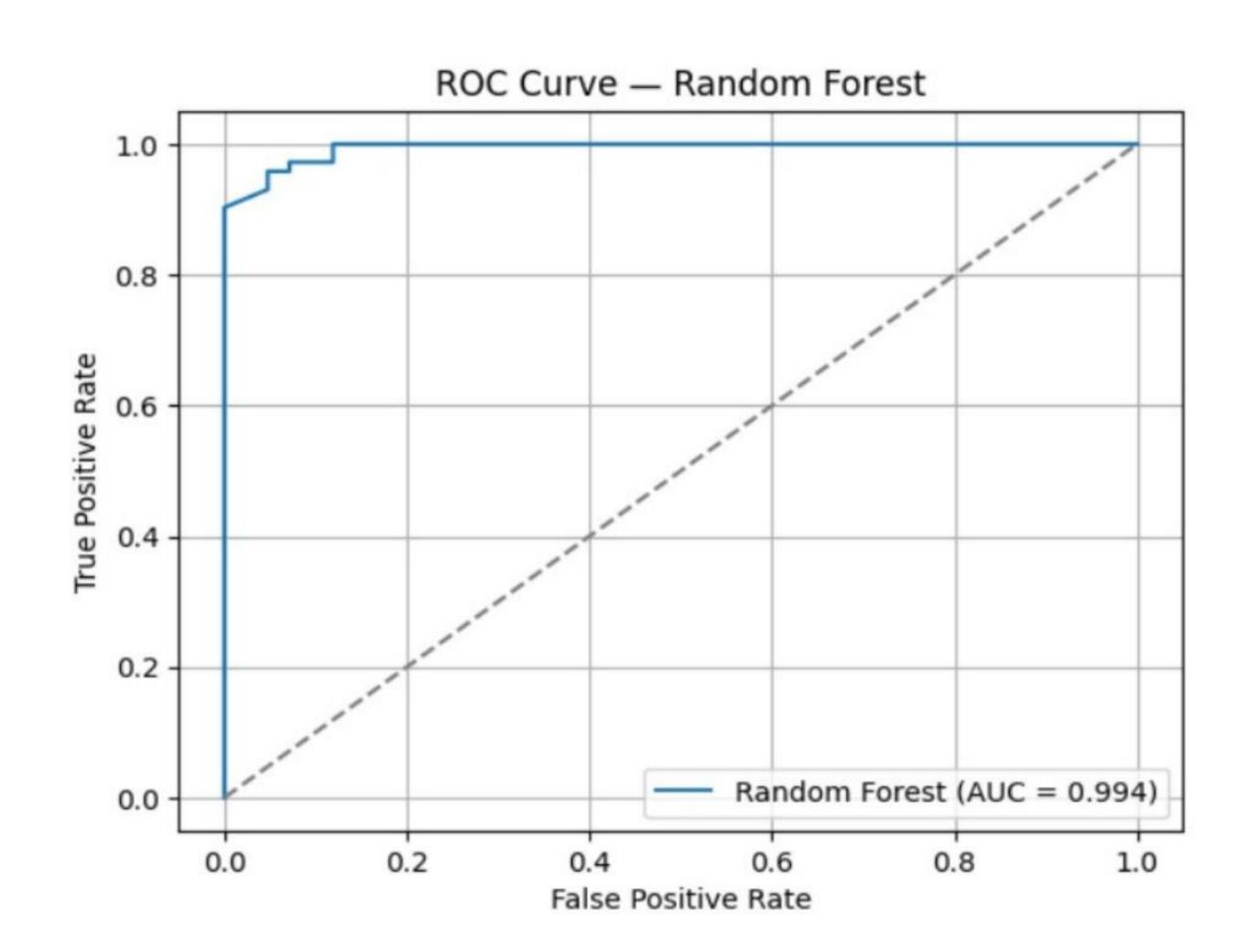
### Confusion Matrix:

[[39 3] [ 2 70]]

## Classification Report:

	precision	recall	f1-score	support
0	0.95	0.93	0.94	42
1	0.96	0.97	0.97	72
accuracy			0.96	114
macro avg	0.96	0.95	0.95	114
weighted avg	0.96	0.96	0.96	114





# **RESULT:**

Thus, the execution successfully implemented a **Random Forest classifier**, tuned key **hyperparameters**, evaluated its performance using appropriate metrics, and analyzed **feature importance** to interpret the model's decision-making process effectively.