```
import numpy as np
import matplotlib.pyplot as plt
import pandas as pd
from google.colab import drive
import seaborn as sns
from itertools import product
def euclidean_distance(vector1, vector2):
    Calculate the Euclidean distance between two vectors.
    Parameters:
    vector1 (numpy array or list): The first vector.
    vector2 (numpy array or list): The second vector.
    Returns:
    float: The Euclidean distance between the two vectors.
    vector1 = np.array(vector1)
    vector2 = np.array(vector2)
    if vector1.shape != vector2.shape:
       raise ValueError("Vectors must have the same length")
    return np.linalg.norm(vector1 - vector2)
def manhattan_distance(vector1, vector2):
    Calculate the Manhattan distance between two vectors.
    Parameters:
    vector1 (numpy array or list): The first vector.
    vector2 (numpy array or list): The second vector.
    Returns:
    float: The Manhattan distance between the two vectors.
    vector1 = np.array(vector1)
    vector2 = np.array(vector2)
    if vector1.shape != vector2.shape:
       raise ValueError("Vectors must have the same length")
    return np.sum(np.abs(vector1 - vector2))
def accuracy_and_generalization_error(true_labels, predicted_labels):
    Calculate accuracy and generalization error between two vectors of labels.
    Parameters:
    true_labels (numpy array or list): The true labels.
    predicted_labels (numpy array or list): The predicted labels.
    tuple: (accuracy, generalization_error)
    true_labels = np.array(true_labels)
    predicted_labels = np.array(predicted_labels)
    if true_labels.shape != predicted_labels.shape:
       raise ValueError("Vectors must have the same length")
    correct_predictions = np.sum(true_labels == predicted_labels)
    total_predictions = len(true_labels)
    accuracy = correct_predictions / total_predictions
    generalization_error = 1 - accuracy
    return accuracy, generalization_error
def precision(true_labels, predicted_labels):
    Calculate precision.
```

```
Parameters:
   true_labels (numpy array or list): The true labels.
   predicted_labels (numpy array or list): The predicted labels.
   Returns:
   float: Precision score.
   true_labels = np.array(true_labels)
   predicted_labels = np.array(predicted_labels)
   true_positives = np.sum((true_labels == 1) & (predicted_labels == 1))
    false_positives = np.sum((true_labels == 0) & (predicted_labels == 1))
    return true_positives / (true_positives + false_positives + 1e-10) # Add small epsilon to avoid division by zero
def recall(true_labels, predicted_labels):
   Calculate recall.
   Parameters:
   true_labels (numpy array or list): The true labels.
   predicted_labels (numpy array or list): The predicted labels.
   Returns:
   float: Recall score.
   true_labels = np.array(true_labels)
   predicted_labels = np.array(predicted_labels)
   true_positives = np.sum((true_labels == 1) & (predicted_labels == 1))
   false_negatives = np.sum((true_labels == 1) & (predicted_labels == 0))
   return true_positives / (true_positives + false_negatives + 1e-10) # Add small epsilon to avoid division by zero
def f1_score(true_labels, predicted_labels):
   Calculate F1 score.
   Parameters:
   true_labels (numpy array or list): The true labels.
   predicted_labels (numpy array or list): The predicted labels.
   Returns:
   float: F1 score.
   prec = precision(true_labels, predicted_labels)
   rec = recall(true_labels, predicted_labels)
   return 2 * (prec * rec) / (prec + rec + 1e-10) # Add small epsilon to avoid division by zero
def confusion_matrix(true_labels, predicted_labels):
   Compute the confusion matrix.
   true_labels (numpy array or list): The true labels.
   predicted_labels (numpy array or list): The predicted labels.
   Returns:
   numpy array: 2x2 confusion matrix.
   true_labels = np.array(true_labels)
   predicted_labels = np.array(predicted_labels)
   tn = np.sum((true_labels == 0) & (predicted_labels == 0))
   fp = np.sum((true_labels == 0) & (predicted_labels == 1))
   fn = np.sum((true_labels == 1) & (predicted_labels == 0))
   tp = np.sum((true_labels == 1) & (predicted_labels == 1))
   return np.array([[tn, fp], [fn, tp]])
def roc_curve(true_labels, predicted_scores):
   Generate the ROC curve.
```

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true_labels (numpy array or list): The true labels.
    predicted_scores (numpy array or list): The predicted scores (probabilities).
    Returns:
    tuple: (fpr, tpr, thresholds)
    true_labels = np.array(true_labels)
    predicted_scores = np.array(predicted_scores)
    thresholds = np.sort(np.unique(predicted_scores))[::-1]
    for threshold in thresholds:
        predicted_labels = (predicted_scores >= threshold).astype(int)
        tp = np.sum((true_labels == 1) & (predicted_labels == 1))
        fn = np.sum((true_labels == 1) & (predicted_labels == 0))
        fp = np.sum((true_labels == 0) & (predicted_labels == 1))
        tn = np.sum((true_labels == 0) & (predicted_labels == 0))
        tpr.append(tp / (tp + fn + 1e-10))
        fpr.append(fp / (fp + tn + 1e-10))
    plt.plot(fpr, tpr, marker='o')
    plt.xlabel('False Positive Rate (FPR)')
    plt.ylabel('True Positive Rate (TPR)')
    plt.title('ROC Curve')
    plt.show()
    return fpr, tpr, thresholds
def auc_roc(fpr, tpr):
    Compute the area under the ROC curve (AUC).
    Parameters:
    fpr (list or numpy array): False Positive Rates.
    tpr (list or numpy array): True Positive Rates.
    Returns:
    float: AUC score.
    return np.trapz(tpr, fpr)
def precision_recall_curve(true_labels, predicted_scores):
    Generate the precision-recall curve.
    Parameters:
    true_labels (numpy array or list): The true labels.
    predicted_scores (numpy array or list): The predicted scores (probabilities).
    Returns:
    tuple: (precision, recall, thresholds)
    true_labels = np.array(true_labels)
    predicted_scores = np.array(predicted_scores)
    thresholds = np.sort(np.unique(predicted_scores))[::-1]
    precision_list = []
    recall_list = []
    for threshold in thresholds:
        predicted_labels = (predicted_scores >= threshold).astype(int)
        prec = precision(true_labels, predicted_labels)
       rec = recall(true_labels, predicted_labels)
        precision_list.append(prec)
       recall_list.append(rec)
    plt.plot(recall_list, precision_list, marker='o')
    plt.xlabel('Recall')
    plt.ylabel('Precision')
    plt.title('Precision-Recall Curve')
    plt.show()
```

```
class KNN_Classifier:
    def __init__(self):
        Initialize the KNN classifier.
        self.X_train = None
        self.Y_train = None
        self.n_neighbors = None
        self.weights = None
    def fit(self, X, Y, n_neighbors=5, weights='uniform', **kwargs):
        Fit the KNN model by storing the training data and parameters.
        self.X_train = X
        self.Y_train = Y
        self.n_neighbors = n_neighbors
        self.weights = weights
    def _compute_distance(self, x1, x2):
    """
        Compute the Euclidean distance between two vectors.
        return np.sqrt(np.sum((x1 - x2) ** 2))
    def predict(self, X):
        Predict the labels for the input data.
        predictions = []
        for x in X:
            # Compute distances between the current sample and all training samples
            distances = np.array([self._compute_distance(x, x_train) for x_train in self.X_train])
            # Get the indices of the k-nearest neighbors
            nearest_indices = np.argsort(distances)[:self.n_neighbors]
            # Get the labels of the k-nearest neighbors
            nearest_labels = self.Y_train[nearest_indices]
            # Apply weights if necessary
            if self.weights == 'uniform':
                # Uniform weights: all neighbors have equal influence
                prediction = np.bincount(nearest_labels).argmax()
            elif self.weights == 'distance':
                # Distance weights: closer neighbors have more influence
                nearest_distances = distances[nearest_indices]
                weights = 1 / (nearest_distances + 1e-10) # Add small epsilon to avoid division by zero
                weighted_votes = np.zeros(np.max(self.Y_train) + 1)
                for label, weight in zip(nearest_labels, weights):
                    weighted_votes[label] += weight
                prediction = weighted_votes.argmax()
                raise ValueError("Invalid weight parameter. Use 'uniform' or 'distance'.")
            predictions.append(prediction)
        return np.array(predictions)
from sklearn.neighbors import KNeighborsClassifier
# Example data
X_train = np.array([
    [1, 1], # Class 0
   [1, 2], # Class 0
[2, 1], # Class 0
[2, 2], # Class 0
    [5, 5], # Class 1
    [5, 6], # Class 1
   [6, 6], # Class 1
[3, 3], # Class 2
    [3, 4], # Class 2
```

```
4, 3], # Class 2
    [4, 4]
            # Class 2
Y_{train} = np.array([0, 0, 0, 0, 1, 1, 1, 1, 2, 2, 2, 2])
X_test = np.array([
    [1.5, 1.5], # Close to Class 0
    [5.5, 5.5], # Close to Class 1
    [3.5, 3.5], # Close to Class 2
    [2.5, 5.0] # Equidistant to Class 0 and Class 1
# Custom KNN implementation
knn_custom = KNN_Classifier()
knn\_custom.fit(X\_train, Y\_train, n\_neighbors=3, weights='distance')
predictions custom = knn custom.predict(X test)
# scikit-learn KNN implementation
knn_sklearn = KNeighborsClassifier(n_neighbors=3, weights='distance')
knn_sklearn.fit(X_train, Y_train)
predictions_sklearn = knn_sklearn.predict(X_test)
# Compare predictions
print("Custom KNN Predictions:", predictions_custom)
print("scikit-learn Predictions:", predictions_sklearn)
print("Predictions Match:", np.array_equal(predictions_custom, predictions_sklearn))
→ Custom KNN Predictions: [0 1 2 2]
     scikit-learn Predictions: [0 1 2 2]
     Predictions Match: True
drive.mount('/content/drive')
df = pd.read_csv('/content/drive/MyDrive/winequality-white.csv', delimiter=';')
df.head()
🔂 Drive already mounted at /content/drive; to attempt to forcibly remount, call drive.mount("/content/drive", force_remount=True).
                                                                    free sulfur total sulfur
                                                        chlorides
      0
                                                                                                                                       6
               7.0
                            0.27
                                      0.36
                                                  20.7
                                                            0.045
                                                                           45 0
                                                                                         170.0
                                                                                                1.0010 3.00
                                                                                                                   0.45
                                                                                                                             8.8
      2
                                      0.40
                                                   6.9
                                                                                                                                       6
                8.1
                            0.28
                                                            0.050
                                                                           30.0
                                                                                                0.9951 3.26
                                                                                                                   0.44
                                                                                                                            10.1
                                                                                         97 0
      4
                7.2
                            0.23
                                      0.32
                                                   8.5
                                                            0.058
                                                                           47.0
                                                                                         186.0
                                                                                                0.9956 3.19
                                                                                                                   0.40
                                                                                                                             9.9
                                                                                                                                       6
 Next steps: ( Generate code with df
                                    View recommended plots
                                                                 New interactive sheet
# Convert quality to binary target
df['target'] = df['quality'].apply(lambda x: 1 if x > 5 else 0)
# Drop the original quality column
df.drop(columns=['quality'], inplace=True)
print(df['target'].value_counts())
₹
    target
          3258
         1640
     Name: count, dtype: int64
# Summary statistics
summary = df.describe().transpose()
summary[['mean', 'std', '25%', '50%', '75%']]
print(summary)
₹
                                                                           25% \
                                                      std
                                                               min
     fixed acidity
                           4898.0
                                      6.854788
                                                 0.843868
                                                           3.80000
                                                                      6.300000
     volatile acidity
                           4898.0
                                     0.278241
                                                 0.100795
                                                           0.08000
                                                                      0.210000
     citric acid
                           4898.0
                                     0.334192
                                                 0.121020 0.00000
                                                                      0.270000
     residual sugar
                           4898.0
                                     6.391415
                                                 5.072058 0.60000
                                                                      1.700000
                                     0.045772
                                                 0.021848 0.00900
                                                                      0.036000
     chlorides
                           4898.0
```

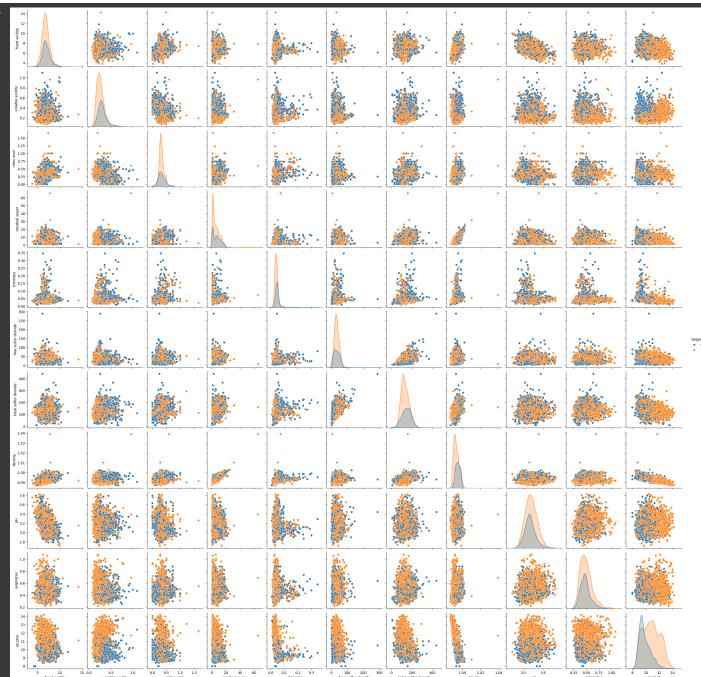
```
free sulfur dioxide 4898.0 35.308085 17.007137 2.00000 23.000000 total sulfur dioxide 4898.0 138.360657 42.498065 9.00000 108.000000
density
                4898.0 0.994027 0.002991 0.98711
                                                                    0.991723
                                 3.188267 0.151001 2.72000
0.489847 0.114126 0.22000
                       4898.0
                                                                     3.090000
                    4898.0
                                                                     0.410000
sulphates
alcohol
                     4898.0 10.514267 1.230621 8.00000
                                                                     9.500000
                       4898.0 0.665169 0.471979 0.00000
                                                                     0.000000
target
                      6.80000
                                      7.3000
fixed acidity
                                                14.20000
volatile acidity
                                      0.3200
                                                1.10000
citric acid
                        0.32000
                                      0.3900
                                                1.66000
residual sugar
                         5.20000
                                      9.9000
                                                65.80000
                        0.04300 0.0500 0.34600
34.00000 46.0000 289.00000
chlorides
free sulfur dioxide 34.00000 46.0000 289.00000 total sulfur dioxide 134.00000 167.0000 440.00000
density
                        0.99374
                                    0.9961
                                                 1.03898
                          3.18000
                                      3.2800
                                                 3.82000
                                    0.5500
sulphates
                                                 1.08000
                         0.47000
                                   11.4000
alcohol
                        10.40000
                                               14.20000
                         1.00000
                                     1.0000
                                                 1.00000
target
```

Shuffle the dataset df = df.sample(frac=1, random_state=42).reset_index(drop=True) print(df.head())

9 1 2 3 4	fixed acidity 6.0 5.4 7.1 7.3 6.5	volatile ac	idity 0.29 0.53 0.25 0.28 0.32		41 16 39 35	10	0.8 0. 2.7 0. 2.1 0. 1.6 0.	.des 048 036 036 054 044
0 1 2 3 4	free sulfur did	oxide total 55.0 34.0 30.0 31.0 27.0	sulfur	dioxide 149.0 128.0 124.0 148.0 91.0	density 0.99370 0.98856 0.99080 0.99178 0.99184	pH 3.09 3.20 3.28 3.18 3.28	sulphates 0.59 0.53 0.43 0.47 0.60	

alcohol target 0 10.966667 1 13.200000 2 12.200000 3 10.700000 4 12.000000

sns.pairplot(df, hue='target') plt.show()



```
tuple: (X_train, X_test, y_train, y_test)
    n_samples = features.shape[0]
    n_test = int(n_samples * t)
    # Split the data
    X_train, X_test = features[:-n_test], features[-n_test:]
    y_train, y_test = target[:-n_test], target[-n_test:]
    return X_train, X_test, y_train, y_test
# Example usage
features = df.drop(columns=['target']).values
target = df['target'].values
X_train, X_test, y_train, y_test = partition(features, target, t=0.2)
print("Training set size:", X_train.shape)
print("Test set size:", X_test.shape)
→ Training set size: (3919, 10)
     Test set size: (979, 10)
# Initialize and fit the KNN classifier
knn = KNN_Classifier()
knn.fit(X_train, y_train, n_neighbors=5, weights='uniform')
# Make predictions
predictions = knn.predict(X_test)
# Calculate accuracy and F1 score
accuracy = np.mean(predictions == y_test)
f1 = f1_score(y_test, predictions)
print("Accuracy (Unscaled Data):", accuracy)
print("F1 Score (Unscaled Data):", f1)
Accuracy (Unscaled Data): 0.7088866189989785
     F1 Score (Unscaled Data): 0.78490566032732
# Standardize the training data
mean = X_train.mean(axis=0)
std = X_train.std(axis=0)
X_train_scaled = (X_train - mean) / std
# Standardize the test data using the training mean and std
X_test_scaled = (X_test - mean) / std
# Initialize and fit the KNN classifier on standardized data
knn_scaled = KNN_Classifier()
knn\_scaled.fit(X\_train\_scaled, y\_train, n\_neighbors=5, weights='uniform')
# Make predictions
predictions_scaled = knn_scaled.predict(X_test_scaled)
# Calculate accuracy and F1 score
accuracy_scaled = np.mean(predictions_scaled == y_test)
f1_scaled = f1_score(y_test, predictions_scaled)
print("Accuracy (Scaled Data):", accuracy_scaled)
print("F1 Score (Scaled Data):", f1_scaled)
Accuracy (Scaled Data): 0.7354443309499489
     F1 Score (Scaled Data): 0.7993803252790892
print("Comparison:")
print(f"Unscaled \ Data \ - \ Accuracy: \ \{accuracy\}, \ F1 \ Score: \ \{f1\}")
print(f"Scaled Data - Accuracy: {accuracy_scaled}, F1 Score: {f1_scaled}")
→ Comparison:
     Unscaled Data - Accuracy: 0.7088866189989785, F1 Score: 0.78490566032732
     Scaled Data - Accuracy: 0.7354443309499489, F1 Score: 0.7993803252790892
```

```
# Initialize and fit the KNN classifier with distance weights
knn distance = KNN Classifier()
knn_distance.fit(X_train_scaled, y_train, n_neighbors=5, weights='distance')
# Make predictions
predictions_distance = knn_distance.predict(X_test_scaled)
# Calculate accuracy and F1 score
accuracy_distance = np.mean(predictions_distance == y_test)
f1_distance = f1_score(y_test, predictions_distance)
print("Accuracy (Distance Weights):", accuracy_distance)
print("F1 Score (Distance Weights):", f1_distance)
Accuracy (Distance Weights): 0.7926455566905005
     F1 Score (Distance Weights): 0.8427575522349305
# Hyperparameter combinations
k_{values} = [1, 5, 9, 11]
distance_metrics = ['Euclidean', 'Manhattan']
weight_schemes = ['uniform', 'distance']
# Function to evaluate the model
def evaluate_knn(X_train, y_train, X_test, y_test, k, distance, weights):
    Evaluate the KNN model with given hyperparameters.
    Parameters:
    X_train (ndarray): Training feature matrix.
    y_train (ndarray): Training target vector.
    X_test (ndarray): Test feature matrix.
    y_test (ndarray): Test target vector.
    k (int): Number of neighbors.
    distance (str): Distance metric ('Euclidean' or 'Manhattan').
    weights (str): Weighting scheme ('uniform' or 'distance').
    Returns:
    dict: Performance metrics (accuracy, precision, recall, F1 score, confusion matrix).
    # Initialize the KNN classifier
    knn = KNN_Classifier()
    # Fit the model
    knn.fit(X_train, y_train, n_neighbors=k, weights=weights)
    # Make predictions
    predictions = knn.predict(X_test)
    # Compute performance metrics
    acc = np.mean(predictions == y_test)
    prec = precision(y_test, predictions)
    rec = recall(y_test, predictions)
    f1 = f1_score(y_test, predictions)
    cm = confusion_matrix(y_test, predictions)
    return {
        'distance': distance,
        'weights': weights,
        'accuracy': acc,
        'precision': prec,
        'recall': rec,
        'f1_score': f1,
        'confusion_matrix': cm
# Evaluate all combinations
results = []
for k, distance, weights in product(k_values, distance_metrics, weight_schemes):
    result = evaluate_knn(X_train_scaled, y_train, X_test_scaled, y_test, k, distance, weights)
    results.append(result)
# Display results
for result in results:
    print(f"k=\{result['k']\},\ Distance=\{result['distance']\},\ Weights=\{result['weights']\}")
    print(f"Accuracy: {result['accuracy']:.4f}, Precision: {result['precision']:.4f}, Recall: {result['recall']:.4f}, F1 Score: {result['f1_
```

```
print("Confusion Matrix:")
    print(result['confusion_matrix'])
    print("-" * 50)
→ k=1, Distance=Euclidean, Weights=uniform
     Accuracy: 0.7783, Precision: 0.8341, Recall: 0.8223, F1 Score: 0.8282
     Confusion Matrix:
     [[239 104]
      [113 523]]
     k=1, Distance=Euclidean, Weights=distance
     Accuracy: 0.7783, Precision: 0.8341, Recall: 0.8223, F1 Score: 0.8282
     Confusion Matrix:
     [[239 104]
      [113 523]]
     k=1, Distance=Manhattan, Weights=uniform
     Accuracy: 0.7783, Precision: 0.8341, Recall: 0.8223, F1 Score: 0.8282
     Confusion Matrix:
     [[239 104]
     k=1, Distance=Manhattan, Weights=distance
     Accuracy: 0.7783, Precision: 0.8341, Recall: 0.8223, F1 Score: 0.8282
     Confusion Matrix:
     [[239 104]
      [113 523]]
     k=5, Distance=Euclidean, Weights=uniform
     Accuracy: 0.7354, Precision: 0.7878, Recall: 0.8113, F1 Score: 0.7994
     Confusion Matrix:
     [[204 139]
     [120 516]]
     k=5, Distance=Euclidean, Weights=distance
     Accuracy: 0.7926, Precision: 0.8305, Recall: 0.8553, F1 Score: 0.8428
     [[232 111]
     [ 92 544]]
     k=5, Distance=Manhattan, Weights=uniform
     Accuracy: 0.7354, Precision: 0.7878, Recall: 0.8113, F1 Score: 0.7994
     Confusion Matrix:
     [[204 139]
      [120 516]]
     k=5, Distance=Manhattan, Weights=distance
     Accuracy: 0.7926, Precision: 0.8305, Recall: 0.8553, F1 Score: 0.8428
     Confusion Matrix:
     [[232 111]
     [ 92 544]]
     k=9, Distance=Euclidean, Weights=uniform
     Accuracy: 0.7426, Precision: 0.7815, Recall: 0.8381, F1 Score: 0.8088
     Confusion Matrix:
     [[194 149]
      [103 533]]
     k=9, Distance=Euclidean, Weights=distance
     Accuracy: 0.8029, Precision: 0.8291, Recall: 0.8774, F1 Score: 0.8526
     Confusion Matrix:
     [[228 115]
k=1, Distance=Euclidean, Weights=uniform
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[103 533]]
k=9, Distance=Euclidean, Weights=distance
Accuracy: 0.8029, Precision: 0.8291, Recall: 0.8774, F1 Score: 0.8526
Confusion Matrix:
[[228 115]
[ 78 558]]
k=9, Distance=Manhattan, Weights=uniform
Accuracy: 0.7426, Precision: 0.7815, Recall: 0.8381, F1 Score: 0.8088
Confusion Matrix:
[[194 149]
[103 533]]
k=9, Distance=Manhattan, Weights=distance
Accuracy: 0.8029, Precision: 0.8291, Recall: 0.8774, F1 Score: 0.8526
Confusion Matrix:
[[228 115]
[ 78 558]]
k=11, Distance=Euclidean, Weights=uniform
Accuracy: 0.7395, Precision: 0.7781, Recall: 0.8381, F1 Score: 0.8070
Confusion Matrix:
[[191 152]
[103 533]]
k=11, Distance=Euclidean, Weights=distance
Accuracy: 0.8039, Precision: 0.8265, Recall: 0.8836, F1 Score: 0.8541
Confusion Matrix:
[[225 118]
[ 74 562]]
k=11, Distance=Manhattan, Weights=uniform
Accuracy: 0.7395, Precision: 0.7781, Recall: 0.8381, F1 Score: 0.8070
Confusion Matrix:
[[191 152]
[103 533]]
k=11, Distance=Manhattan, Weights=distance
Accuracy: 0.8039, Precision: 0.8265, Recall: 0.8836, F1 Score: 0.8541
Confusion Matrix:
[[225 118]
[ 74 562]]
```

1.1

The best model is the KNN classifier with k=11, Distance=Manhattan, and Weights=distance. This model achieves the highest accuracy (0.8039), recall (0.8836), and F1 score (0.8541), indicating strong overall performance. The high recall means it excels at identifying true positives (good wines), which is crucial for minimizing false negatives. Additionally, the F1 score reflects a good balance between precision and recall, ensuring reliable predictions. The use of distance weighting further enhances performance by giving more importance to closer neighbors, while the larger k value (k=11) helps smooth out noise and improve generalization. Overall, this model strikes the best balance between accuracy, precision, and recall, making it the optimal choice for this dataset.