```
In [ ]: import numpy as np
        import matplotlib.pyplot as plt
        import cv2
        import os
        from sklearn.preprocessing import StandardScaler
        from sklearn.metrics import accuracy score, f1 score
In [ ]: # mapping labels to 0 to 3 for all the four classes
        labmap = {0: "n02089078-black-and-tan coonhound"
                  ,1: "n02091831-Saluki"
                  ,2:"n02092002-Scottish deerhound"
                  ,3:"n02095314-wire-haired fox terrier"}
        # giving the paths of the cropped images
        paths = [r'../DataSet/ProcessedDatasets/n02089078-black-and-tan coonhound/'
                 ,r'../DataSet/ProcessedDatasets/n02091831-Saluki/'
                 ,r'../DataSet/ProcessedDatasets/n02092002-Scottish deerhound/'
                 ,r'../DataSet/ProcessedDatasets/n02095314-wire-haired fox terrier/']
        data set = [] # empty array to save the dataset
        labels = [] # empty array to save the labels
        # 1. Loop to convert the images to gray scale pixel intensity histograms and save the labels and the dataset
        for i in paths:
            for dog in os.listdir(i):
                img = cv2.imread(i + dog,cv2.IMREAD GRAYSCALE)
                hist = cv2.calcHist(img, [0], None, [256], [0, 256])
                data set.append(hist)
                labels.append(paths.index(i))
        # 2. standardize the data set
        standard dset = StandardScaler().fit transform(np.array(data set)[:,:,0])
        # convert the python list to a numpy array for appending it to the data
        labels = np.array(labels)
        # final data that consists of the normalized data and labels.
        final data = np.column stack((standard dset, labels))
```

```
In [ ]: from sklearn.model_selection import train_test_split
        # referred from https://scikit-learn.org/stable/modules/generated/sklearn.model selection.train test split.html
        X_train, X_test, Y_train, Y_test = [], [], [], []
        # 3. Loop to split training and testing data to 80/20 percent
        for i in labmap.keys():
            class_indices = np.where(labels == i)[0]
            X class = standard dset[class indices]
            y class = labels[class indices]
            X train class, X test class, Y train class, Y test class = train test split(X class, y class, test size=0.2)
            X_train.extend(X_train_class)
            X test.extend(X test class)
            Y train.extend(Y train class)
            Y test.extend(Y test class)
        X train = np.array(X train)
        X test = np.array(X test)
        Y train = np.array(Y train)
        Y test = np.array(Y test)
```

# Standard 5-Fold cross-validation

```
In []: from sklearn.model_selection import cross_val_score, KFold,StratifiedKFold
    from sklearn.neighbors import KNeighborsClassifier
    # referred from https://scikit-learn.org/stable/modules/generated/sklearn.model_selection.KFold.html

# list of k values to iterate through
    k_values = [1, 3, 5, 7, 10, 20]

# Lists of validation and training errors for standard KFold cross-validation
    standard_validation_errors = []
    standard_training_errors = []

num_folds = 5

# Initialize a KFold cross-validator
    kf = KFold(n_splits=num_folds, shuffle=True, random_state=57)
```

```
for k in k_values:
   # creating a k-NN classifier with the current k value
   knn_standard = KNeighborsClassifier(n_neighbors=k)
   # Initialize error variables for this k value
   validation errors k = []
   training_errors_k = []
   for train index, validation index in kf.split(X train):
       # Split the data into training and validation sets
       X train fold, X validation fold = X train[train index], X train[validation index]
       Y train fold, Y validation fold = Y train[train index], Y train[validation index]
       # Fit the k-NN classifier on the training data
       knn standard.fit(X train fold, Y train fold)
       # Predict on the training and validation sets
       Y pred train = knn standard.predict(X train fold)
       Y pred validation = knn standard.predict(X validation fold)
       # Calculate training and validation errors for this fold
       training errors k.append(1 - (Y pred train == Y train fold).mean())
       validation errors k.append(1 - (Y pred validation == Y validation fold).mean())
   # Calculate the mean errors across all folds for this k value
   mean training error = np.mean(training errors k)
   mean validation error = np.mean(validation errors k)
   # Append errors to the respective lists
   standard training errors.append(mean training error)
   standard validation errors.append(mean validation error)
```

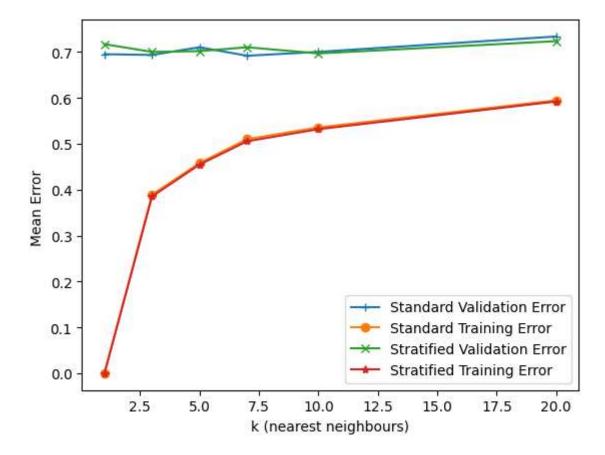
## Stratified 5-Fold cross-validation

```
import math
# Lists of validation and training errors for stratified 5-fold cross-validation
stratified_validation_errors = []
stratified_training_errors = []
```

```
best k values = []
 # reffered from https://scikit-learn.org/stable/modules/generated/sklearn.model selection.StratifiedKFold.html#sklear
 kfs = StratifiedKFold(n splits=num folds,shuffle=True,random state=57)
 for k in k values:
     validation_errors = []
     training_errors = []
     # loop to split the train dataset into stratified 5 fold validation sets
     for train index, val index in kfs.split(X train, Y train):
         X_train_fold, X_val_fold = X_train[train_index], X_train[val_index] # Getting training data and validation de
         Y train fold, Y val fold = Y train[train index], Y train[val index] # Getting training labels and validation
         knn stratified = KNeighborsClassifier(n neighbors=k)
         knn_stratified.fit(X_train_fold, Y_train_fold)
         y pred train = knn stratified.predict(X train fold)
         y pred val = knn stratified.predict(X val fold)
         validation errors.append(1- (y pred val == Y val fold).mean())
         training errors.append(1- (y pred train == Y train fold).mean())
     mean validation error = np.mean(validation errors)
     mean training error = np.mean(training errors)
     stratified validation errors.append(mean validation error)
     stratified training errors.append(mean training error)
     best k = k values[np.argmin(stratified validation errors)]
     best k values.append(best k)
 print(f'validation errors = {stratified validation errors}')
 print(f'training errors = {stratified training errors}')
validation errors = [0.7170028011204482, 0.7001120448179272, 0.7017787114845939, 0.7102100840336135, 0.6967787114845
938, 0.7235714285714285]
training errors = [0.0, 0.38526266852626684, 0.45476873415611874, 0.5054559967720149, 0.5318316184661807, 0.59213354
03454295]
```

#### Plotting the validation and training errors for both the cross-validations

```
In [ ]: import matplotlib.pyplot as plt
        # Plot the validation and training error curves for standard 5-fold cross-validation
        plt.plot(k_values, standard_validation_errors, label='Standard Validation Error', marker='+')
        plt.plot(k values, standard training errors, label='Standard Training Error', marker='o')
        # Plot the validation and training error curves for stratified 5-fold cross-validation
        plt.plot(k values, stratified validation errors, label='Stratified Validation Error', marker='x')
        plt.plot(k values, stratified training errors, label='Stratified Training Error', marker='*')
        # Label the axes and add a legend
        plt.xlabel('k (nearest neighbours)')
        plt.ylabel('Mean Error ')
        plt.legend()
        # Find the best k for each curve
        best k standard validation = k values[standard validation errors.index(min(standard validation errors))]
        best k standard training = k values[standard training errors.index(min(standard training errors))]
        best k stratified validation = k values[stratified validation errors.index(min(stratified validation errors))]
        best k stratified training = k values[stratified training errors.index(min(stratified training errors))]
        # Print the best k values for each curves
        print(f'K with the least Standard Validation Error: {best k standard validation} with {min(standard validation errors
        print(f'K with the least Standard Training Error: {best k standard training}' )
        print(f'K with the least Stratified Validation Error: {best k stratified validation} with with {min(stratified valid
        print(f'K with the least Stratified Training Error: {best k stratified training}')
        # Show the plot
        plt.show()
       K with the least Standard Validation Error: 7 with 0.6917366946778711 validation error
       K with the least Standard Training Error: 1
       K with the least Stratified Validation Error: 10 wit h with 0.6967787114845938 validation error
       K with the least Stratified Training Error: 1
```



### **Comments**

5 nearest neighbours has the least mean validation error for stratified cross fold validations 7 nearest neighbours has the least mean validation error for standard cross fold validations 1 nearest neighbours has the least mean training error for both standard and stratified cross fold validations 1 nearest neighbours has the least mean training error for both standard and stratified cross fold validations From the above curves The training error increases gradually as the number of neighbours increases. The model complexity depends on the number of number of zones or clusters that the model can make. If the number of nearest neighbours is 1 then for each datapoint the model tries to draw a boundary to it. Hence for a less k value the model is much complex and over fit. If the k value is more the model has big clustering zones and the model is very simple and the is under fit as it a lot of data points from other class may also fit into the current class. The best spot is somewhere between to have a less complex model but

with better fit.

From the give nK values the table can say whether the model is complex or simple and overfit or under fit.

K neighbours	complexity	overfit/underfit
1	more complexity	overfit
3	less complexity	close to best fit
5	less complexity	best fit
7	less complexity	close to best fit
10	simple	underfit
20	simple	underfit

```
In [ ]: knn_s = KNeighborsClassifier(n_neighbors=best_k_stratified_validation)

# Train the k-NN classifier on the entire training dataset
knn_s.fit(X_train, Y_train)

# Evaluate the model on the test dataset and calculate the test error
Y_pred_test = knn_s.predict(X_test)
test_error = 1 - (Y_pred_test == Y_test).mean() # Test error in terms of misclassification rate

print(f'Test Error with k={best_k_stratified_validation}: {test_error * 100:.2f}% ')
```

Test Error with k=10: 68.87%

```
In []: import numpy as np
    import matplotlib.pyplot as plt
    from sklearn.model_selection import StratifiedKFold
    from sklearn.metrics import confusion_matrix
    from sklearn.naive_bayes import GaussianNB
    from sklearn.neural_network import MLPClassifier
    from sklearn.ensemble import RandomForestClassifier
    # reffered from https://scikit-learn.org/stable/modules/generated/sklearn.metrics.confusion_matrix.html

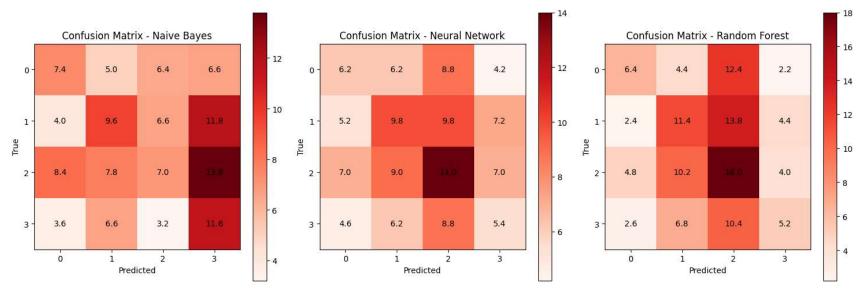
class_labels = np.unique(Y_train)

naive_bayes = GaussianNB()
```

```
neural network = MLPClassifier(hidden layer sizes=(10, 10, 10))
random_forest = RandomForestClassifier()
models = [naive bayes, neural network, random forest]
model names = ['Naive Bayes', 'Neural Network', 'Random Forest']
skf = StratifiedKFold(n splits=5, shuffle=True, random state=57)
confusion matrices = []
confusion matrices test = []
validation accuracies = []
for model, model name in zip(models, model names):
    fold matrices = []
    validation accuracy model = []
    for train_index, tv_index in skf.split(X_train, Y_train):
        X train fold, X tv fold = X train[train index], X train[tv index]
        Y train fold, y tv fold = Y train[train index], Y train[tv index]
        model.fit(X train fold, Y train fold)
        Y val pred = model.predict(X tv fold) # predictions based on validation set
        fold matrix = confusion matrix(y tv fold, Y val pred, labels=class labels) # confusion matrix based on test s
        validation accuracy model.append(accuracy score(y tv fold , Y val pred)) # to find validation accuracy
        fold matrices.append(fold matrix)
    validation accuracies.append(np.array(validation accuracy model).mean())
    confusion matrices.append(fold matrices)
best model = ""
best accuracy = 0
plt.figure(figsize=(15, 5))
for i, model name in enumerate(model names):
    plt.subplot(1, 3, i + 1)
    average matrix = np.mean(confusion matrices[i], axis=0)
    plt.imshow(average matrix, cmap=plt.cm.Reds)
    plt.title(f'Confusion Matrix - {model name}')
    plt.colorbar()
    tick marks = np.arange(len(class labels))
    plt.xticks(tick marks, class labels)
    plt.yticks(tick marks, class labels)
    plt.xlabel('Predicted')
    plt.ylabel('True')
```

```
for i in range(len(class_labels)):
       for j in range(len(class_labels)):
           plt.text(j, i, f'{average_matrix[i, j]}', ha='center', va='center', color='black')
    print(f'Confusion Matrix - {model name}')
    print(average matrix)
    sum = 0
   for i in range(len(average matrix)):
       for j in range(len(average matrix[0])):
           if i==j:
                sum+=average_matrix[i,j]
   print(f'{sum} Average correct prediction for {model name} on stratified 5 fold on test dataset.')
   if(sum>best accuracy):
        best accuracy = sum
        best_model = model_name
plt.tight layout()
plt.show()
print(f"Best model according to confusion matrices is {best model} with average correct prediction of {best accuracy
print(f"Best model according to validation accuracy is {model names[validation accuracies.index(max(validation accuracy
```

```
c:\Users\kaasa\AppData\Local\Programs\Python\Python310\lib\site-packages\sklearn\neural network\ multilayer perceptro
n.py:691: ConvergenceWarning: Stochastic Optimizer: Maximum iterations (200) reached and the optimization hasn't conv
erged yet.
  warnings.warn(
c:\Users\kaasa\AppData\Local\Programs\Python\Python310\lib\site-packages\sklearn\neural_network\_multilayer_perceptro
n.py:691: ConvergenceWarning: Stochastic Optimizer: Maximum iterations (200) reached and the optimization hasn't conv
erged yet.
  warnings.warn(
c:\Users\kaasa\AppData\Local\Programs\Python\Python310\lib\site-packages\sklearn\neural_network\_multilayer_perceptro
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 warnings.warn(
c:\Users\kaasa\AppData\Local\Programs\Python\Python310\lib\site-packages\sklearn\neural_network\_multilayer_perceptro
n.py:691: ConvergenceWarning: Stochastic Optimizer: Maximum iterations (200) reached and the optimization hasn't conv
erged yet.
 warnings.warn(
c:\Users\kaasa\AppData\Local\Programs\Python\Python310\lib\site-packages\sklearn\neural_network\_multilayer_perceptro
n.py:691: ConvergenceWarning: Stochastic Optimizer: Maximum iterations (200) reached and the optimization hasn't conv
erged yet.
 warnings.warn(
Confusion Matrix - Naive Bayes
[ [ 7.4 5. 6.4 6.6 ]
[ 4. 9.6 6.6 11.8]
[ 8.4 7.8 7. 13.8]
[ 3.6 6.6 3.2 11.6]]
35.6 Average correct prediction for Naive Bayes on stratified 5 fold on test dataset.
Confusion Matrix - Neural Network
[[ 6.2 6.2 8.8 4.2]
[5.2 9.8 9.8 7.2]
[7. 9. 14. 7.]
[ 4.6 6.2 8.8 5.4]]
35.4 Average correct prediction for Neural Network on stratified 5 fold on test dataset.
Confusion Matrix - Random Forest
[[ 6.4 4.4 12.4 2.2]
[ 2.4 11.4 13.8 4.4]
[ 4.8 10.2 18. 4. ]
[ 2.6 6.8 10.4 5.2]]
41.0 Average correct prediction for Random Forest on stratified 5 fold on test dataset.
```



Best model according to confusion matrices is Random Forest with average correct prediction of 41.0

Best model according to validation accuracy is Random Forest with average validation accuracy of 0.3434733893557423

```
In [ ]: y true = Y test
        models[0].fit(X train, Y train)
        models[1].fit(X train,Y train)
        models[2].fit(X train,Y train)
        y pred naive bayes = models[0].predict(X test)
        y pred neural network = models[1].predict(X test)
        y pred random forest = models[2].predict(X test)
        accuracy naive bayes = accuracy score(y true, y pred naive bayes)
        accuracy neural network = accuracy score(y true, y pred neural network)
        accuracy random forest = accuracy score(y true, y pred random forest)
        f1 naive bayes = f1 score(y true, y pred naive bayes, average='weighted')
        f1 neural network = f1 score(y true, y pred neural network, average='weighted')
        f1 random forest = f1 score(y true, y pred random forest, average='weighted')
        print(f"Test Accuracy - Naive Bayes: {accuracy naive bayes * 100:.2f}%")
        print(f"Test Accuracy - Neural Network: {accuracy neural network * 100:.2f}%")
        print(f"Test Accuracy - Random Forest: {accuracy random forest * 100:.2f}%")
```

```
print(f"F-Measure - Naive Bayes: {f1_naive_bayes:.4f}")
print(f"F-Measure - Neural Network: {f1_neural_network:.4f}")
print(f"F-Measure - Random Forest: {f1_random_forest:.4f}")

c:\Users\kaasa\AppData\Local\Programs\Python\Python310\lib\site-packages\sklearn\neural_network\_multilayer_perceptro
n.py:691: ConvergenceWarning: Stochastic Optimizer: Maximum iterations (200) reached and the optimization hasn't conv
erged yet.
   warnings.warn(
Test Accuracy - Naive Bayes: 27.81%
Test Accuracy - Neural Network: 24.50%
```

Test Accuracy - Naive Bayes: 27.81%
Test Accuracy - Neural Network: 24.50%
Test Accuracy - Random Forest: 32.45%
F-Measure - Naive Bayes: 0.2650
F-Measure - Neural Network: 0.2437
F-Measure - Random Forest: 0.3100

Random forest is the best classification according to test Accuracy. Random forest is the best classification according to F-measure

## **References:**

- 1. https://scikit-learn.org/stable/modules/generated/sklearn.metrics.confusion\_matrix.html
- 2. https://scikit-

 $learn. org/stable/modules/generated/sklearn. model\_selection. Stratified KFold. html \#sklearn. model\_selection. Stratified KFold. split the sklearn was also become a support of the sklear was also become a sklear was also become a support of the sklear was also become a skl$ 

- 3. https://scikit-learn.org/stable/modules/generated/sklearn.model\_selection.train\_test\_split.html
- 4. https://scikit-learn.org/stable/modules/generated/sklearn.model\_selection.KFold.html