RBL SK5009 Advanced Artificial Intelligence
Pumpkin Seeds Classification
with Machine Learning

Ahmad Zaini Zahrandika (20923302)

Purpose

- Create machine learning models with Support Vector Machine (SVM), k-Nearest Neighbors (k-NN), Random Forest (RF), and Logistic Regression (LR) for pumpkin seeds (*Cucurbita pepo L.*) classification.

Machine Learning Models

- Support Vector Machine (SVM)

Support vector machines predicted a suitable hyperplane function to statistically separate two classes on the multidimensional plane.

- k-Nearest Neighbors (k-NN)

The k-NN or k-nearest neighbor algorithm determined the nearest k points in the same space with each data in the training set, usually by considering the Euclidean distance.

Machine Learning Models

- Random Forest (RF)

The random forest classifier classified many random samples, which were sampled independently of the input vector by considering the combination of predictors that received the highest vote from all the tree estimators.

- Logistic Regression (LR)

Logistic regression is a statistical method used to analyze a dataset with independent variables to determine an outcome. It constructs a dividing hyper-plane between two data sets and provides a functional form and parameter vector to express the probability of a certain outcome given the input variables.

- k-Fold Cross-Validation

A statistical method used to assess the performance and generalizability of machine learning models. It helps prevent issues like overfitting and provides a more reliable estimate of a model's performance on unseen data. The dataset was divided into k parts, and the k-1 part was used as train data. This process was repeated k times, and the average of the accuracy value found in each iteration was accepted as the performance of cross-validation.

- Feature scaling

No	Scaler	Formula	Explanation
1	StandardScaler	$z = rac{x-\mu}{\sigma}$	This scaler standardizes the data by removing the mean and scaling it to unit variance.
2	MinMaxScaler	$x' = rac{x - x_{\min}}{x_{\max} - x_{\min}}$	This scaler scales the data to a specific range, typically [0, 1].

- Confusion matrix

A confusion matrix is a table used to evaluate the performance of a classification model by comparing the predicted outcomes to the actual outcomes. It is particularly useful for assessing models on imbalanced datasets or when multiple classes are involved.

	Predicted Positive	Predicted Negative
Actual Positive	True Positive (tp)	False Negative (fn)
Actual Negative	False Positive (fp)	True Negative (tn)

- Performance criteria table

No	Performance Measure	Formula	Evaluation
1	Accuracy	$\tfrac{tp+tn}{tp+fp+tn+fn}$	The ratio of correct estimates to the total is the number of samples evaluated.
2	Precision	$rac{tp}{tp+fp}$	It is used to measure the positive patterns that are correctly predicted from the total predictive forms in a positive class.
3	Recall	$rac{tp}{tp+fn}$	Used to measure the proportion of correctly classified positive patterns.
4	F-score	$rac{2tp}{2tp+fp+fn}$	Represents the harmonic mean between Recall and Precision values.

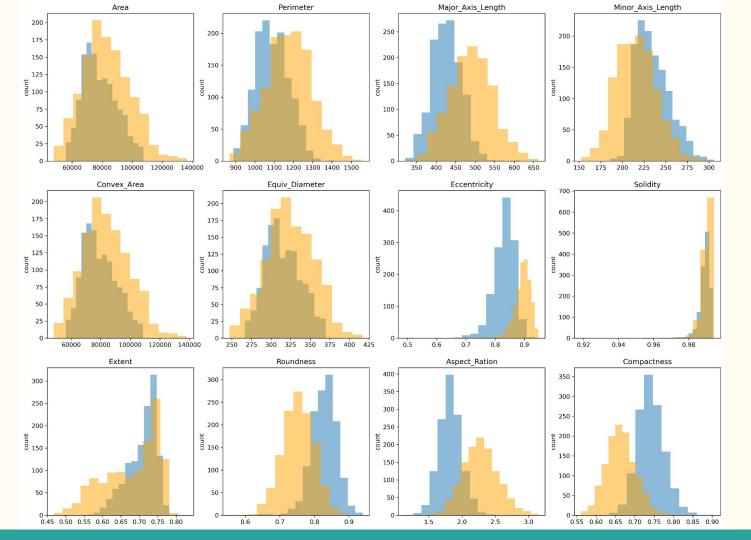
Steps

- 1. Check if there are empty values in the dataset,
- 2. Check feature types. Convert to numerical if the feature is categorical,
- 3. Split the dataset to X (data) and y (target),
- 4. Split X, y: 70% training, 30% test,
- 5. Perform hyperparameter tuning to find the best parameter for each model.
- 6. Calculate the performance of each model.

Information in Dataset

```
# Check the features of the data. There are 13 of them.
# The only feature with categorical type is "Class".
df.info()
<class 'pandas.core.frame.DataFrame'>
RangeIndex: 2500 entries, 0 to 2499
Data columns (total 13 columns):
     Column
                        Non-Null Count Dtype
                                        int64
     Area
                        2500 non-null
     Perimeter
                        2500 non-null
                                        float64
     Major_Axis_Length
                        2500 non-null
                                       float64
                                       float64
    Minor_Axis_Length
                        2500 non-null
     Convex Area
                        2500 non-null
                                        int64
     Equiv_Diameter
                        2500 non-null
                                        float64
     Eccentricity
                        2500 non-null
                                        float64
                                        float64
     Solidity
                        2500 non-null
     Extent
                        2500 non-null
                                        float64
                                        float64
     Roundness
                        2500 non-null
     Aspect_Ration
                        2500 non-null
                                        float64
    Compactness
                        2500 non-null
                                        float64
 12 Class
                        2500 non-null
                                        object
dtypes: float64(10), int64(2), object(1)
memory usage: 254.0+ KB
```

No empty values in the dataset.



Support Vector Machine (SVM)

```
# Define a pipeline with a placeholder for the scaler and the SVC model
pipeline = Pipeline([
    ("scaler", "passthrough"), # Placeholder for scaler
    ("svc", SVC())
1)
# Define the parameter grid for GridSearchCV
param grid = {
    "scaler": [StandardScaler(), MinMaxScaler()],
    "svc C": [0.1, 1, 10, 100],
    "svc kernel": ["linear", "rbf", "sigmoid"],
    "svc gamma": ["auto", 1/100]
# Perform GridSearchCV with 5-fold cross-validation on the training set
grid_search = GridSearchCV(pipeline, param_grid, cv=5)
grid search.fit(X train, y train)
```

Support Vector Machine (SVM)

Best Parameters: {'scaler': StandardScaler(), 'svc_C': 1, 'svc_gamma': 'auto', 'svc_kernel': 'rbf'} Best Cross-Validation Score: 0.8891428571428571

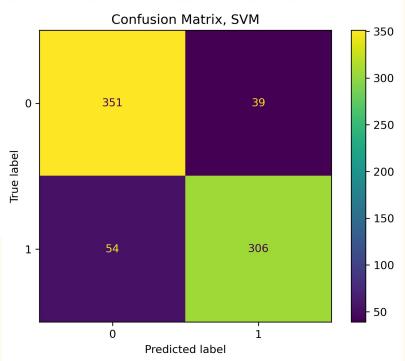
Accuracy: 0.8760

Classification Report:

		precision	recall	f1-score	support
	0	0.8667	0.9000	0.8830	390
	1	0.8870	0.8500	0.8681	360
accur	racy			0.8760	750
macro	avg	0.8768	0.8750	0.8756	750
weighted	avg	0.8764	0.8760	0.8759	750

0: Çerçevelik

1: Ürgüp Sivrisi



k-Nearest Neighbors (k-NN)

```
# Define a pipeline with a placeholder for the scaler and the k-NN model
pipeline = Pipeline([
    ("scaler", "passthrough"), # Placeholder, to be replaced during grid search
    ("knn", KNeighborsClassifier())
])
# Define the parameter arid for GridSearchCV
param grid = {
    "scaler": [StandardScaler(), MinMaxScaler()],
    "knn n neighbors": [5, 10, 15, 20],
    "knn__metric": ["euclidean", "cosine", "manhattan"]
# Perform GridSearchCV with 5-fold cross-validation on the training set
grid search = GridSearchCV(pipeline, param grid, cv=5)
grid search.fit(X train, y train)
```

k-Nearest Neighbors (k-NN)

Best Parameters: {'knn_metric': 'manhattan', 'knn_n_neighbors': 20, 'scaler': StandardScaler()}

Best Cross-Validation Score: 0.885142857142857

Accuracy: 0.8733

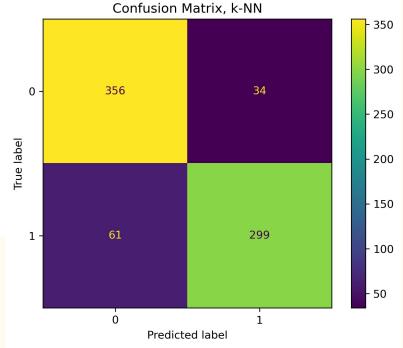
Confusion Matrix, k-NN

Classification Report:

		precision	recall	f1-score	support
	0	0.8537	0.9128	0.8823	390
	1	0.8979	0.8306	0.8629	360
accui	racy			0.8733	750
macro	avg	0.8758	0.8717	0.8726	750
weighted	avg	0.8749	0.8733	0.8730	750

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Random Forest (RF)

```
# Define a pipeline with a placeholder for the scaler and the RF model
pipeline = Pipeline([
    ("scaler", "passthrough"), # Placeholder, to be replaced during grid search
    ("rf", RandomForestClassifier())
1)
# Define the parameter grid for GridSearchCV
param grid = {
    "scaler": [StandardScaler(), MinMaxScaler()],
    "rf__criterion" : ["gini", "entropy", "log_loss"],
    "rf n estimators": [10, 20, 50, 75, 100]
# Perform GridSearchCV with 5-fold cross-validation on the training set
grid_search = GridSearchCV(pipeline, param_grid, cv=5)
grid_search.fit(X_train, y_train)
```

Random Forest (RF)

Best Parameters: {'rf__criterion': 'gini', 'rf__n_estimators': 100, 'scaler': MinMaxScaler()}

Best Cross-Validation Score: 0.8942857142857144

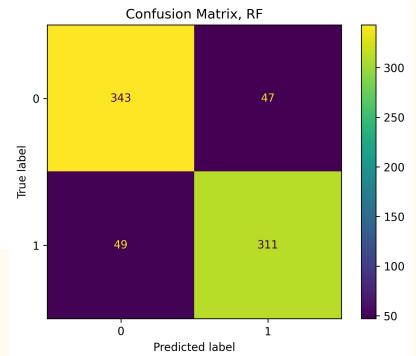
Accuracy: 0.8720

Classification Report:

	precision	recall	f1-score	support
0	0.8750	0.8795	0.8772	390
1	0.8687	0.8639	0.8663	360
accuracy			0.8720	750
macro avg	0.8719	0.8717	0.8718	750
weighted avg	0.8720	0.8720	0.8720	750

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Logistic Regression (LR)

```
# Define a pipeline with a placeholder for the scaler and the LR model
pipeline = Pipeline([
    ("scaler", "passthrough"), # Placeholder, to be replaced during grid search
    ("lr", LogisticRegression())
# Define the parameter grid for GridSearchCV
param grid = {
    "scaler": [StandardScaler(), MinMaxScaler()],
    "lr C": [0.01, 0.1, 1, 10, 100],
    "lr__solver": ["liblinear", "saga", "lbfgs", "newton-cg"],
    "lr max iter": [1000, 5000, 10000]
# Perform GridSearchCV with 5-fold cross-validation on the training set
grid search = GridSearchCV(pipeline, param grid, cv=5)
grid search.fit(X train, y train)
```

Logistic Regression (LR)

Best Parameters: {'lr_C': 100, 'lr_max_iter': 5000, 'lr_solver': 'saga', 'scaler': StandardScaler()} Best Cross-Validation Score: 0.8868571428571428 Confusion Matrix, LR Accuracy: 0.8733 - 350 Classification Report: - 300 precision recall f1-score support 352 38 0 - 250 0.8606 0.9026 0.8811 390 True label 1 0.8886 0.8417 360 0.8645 - 200 0.8733 750 accuracy - 150 0.8721 0.8728 macro avg 0.8746 750 weighted avg 0.8740 0.8733 0.8731 750 1 -57 303 - 100 0: Çerçevelik 1: Ürgüp Sivrisi

Predicted label

Comparison

- In this RBL, SVM is slightly more accurate than k-NN, RF, and LR (0.8760 vs 0.8733, 0.8720, 0.8733).
- The result may differ if there are more variables and values in hyperparameter tuning.

Comparison (with the reference)

	Accuracy (RBL)	Accuracy (Reference)
SVM	0.8760	0.8864
k-NN	0.8733	0.8764
RF	0.8720	0.8756
LR	0.8733	0.8792

Comparison (with the reference)

- There are some differences in methods.
- In the reference, 10-fold cross-validation was used in the whole dataset (without splitting it to training and test).
- SVM: The sigmoid hyperplane function was determined, and the gamma value was accepted as "1/feature number".
- k-NN: Used Euclidean distance, k value was considered as five.

Comparison (with the reference)

- RF: The number of trees in the forest was 100. Information acquisition was calculated according to the entropy.
- LR: The Newton method was used for the optimization during classification.
- The reference didn't mention the scaler that was used; StandardScaler, MinMaxScaler, or none.

References

- https://www.kaggle.com/datasets/muratkokludataset/pumpkin-seeds-dataset
- https://link.springer.com/article/10.1007/s10722-021-01226-0
- https://link.springer.com/content/pdf/10.1007/s10722-021-01226-0.pdf
- https://www.muratkoklu.com/datasets/