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1 Glass Classification

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1.0.1 Context and Content

Context

This is a Glass Identification Data Set from UCI. It contains 9 attributes.

The response is glass type(discrete 7 values)

Kaggle Link

UCI dataset Link

Content

Attribute Information

- RI: refractive index
- Na: Sodium (unit measurement: weight percent in corresponding oxide, as are attributes 4-10)
- Mg: Magnesium
- Al: Aluminum
- Si: Silicon
- K: Potassium
- Ca: Calcium
- Ba: Barium
- Fe: Iron
- Type of glass: (class attribute)
 - 1 building_windows_float_processed

```
- 2 building_windows_non_float_processed
```

- 3 vehicle_windows_float_processed
- 4 vehicle_windows_non_float_processed (none in this database)
- 5 containers
- 6 tableware
- 7 headlamps

1.0.2 Data Loading and Analysis

```
[1]: # importing essential libraries
import numpy as np
import pandas as pd
import matplotlib.pyplot as plt
import seaborn as sns
```

```
[74]: # loading the dataset

df = pd.read_csv("./glass.csv")

df.head()
```

```
[74]:
                               Al
                                      Si
                                            K
                                                 Ca
             RΙ
                   Na
                         Mg
                                                      Ba
                                                           Fe
                                                              Туре
     0 1.52101 13.64
                      4.49 1.10 71.78
                                         0.06
                                               8.75
                                                     0.0
                                                          0.0
                                                                 1
     1 1.51761 13.89 3.60
                            1.36 72.73
                                               7.83
                                                          0.0
                                         0.48
                                                     0.0
                                                                 1
     2 1.51618 13.53 3.55
                            1.54 72.99
                                         0.39
                                               7.78
                                                     0.0
                                                          0.0
                                                                 1
     3 1.51766 13.21 3.69
                            1.29 72.61
                                         0.57
                                               8.22
                                                     0.0
                                                          0.0
                                                                 1
     4 1.51742 13.27 3.62 1.24 73.08 0.55 8.07
                                                     0.0
                                                          0.0
                                                                 1
```

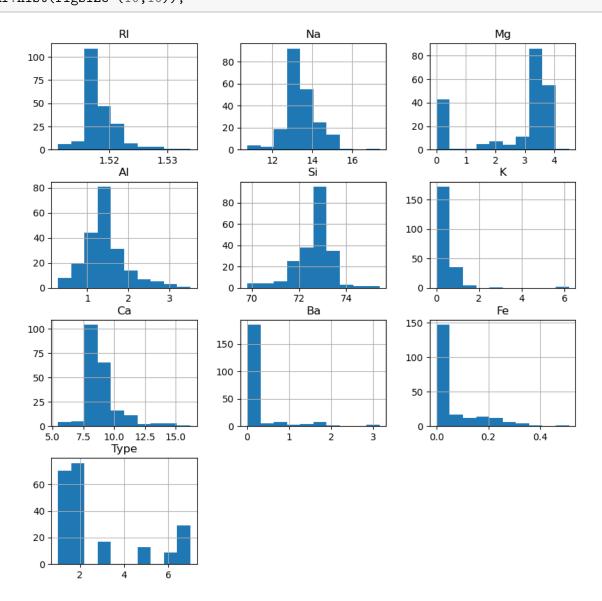
```
[3]: #getting information about the data df.info()
```

<class 'pandas.core.frame.DataFrame'>
RangeIndex: 214 entries, 0 to 213
Data columns (total 10 columns):

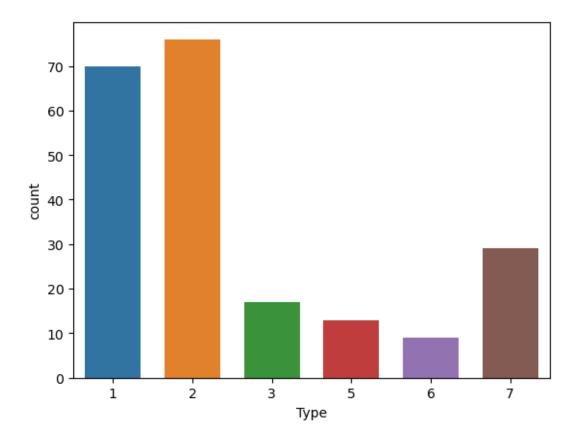
#	Column	Non-Null Count	Dtype
0	RI	214 non-null	float64
1	Na	214 non-null	float64
2	Mg	214 non-null	float64
3	Al	214 non-null	float64
4	Si	214 non-null	float64
5	K	214 non-null	float64
6	Ca	214 non-null	float64
7	Ba	214 non-null	float64
8	Fe	214 non-null	float64
9	Type	214 non-null	int64

dtypes: float64(9), int64(1)
memory usage: 16.8 KB

[4]: # understanding the correlation between the data df.hist(figsize=(10,10));



[5]: # analyzing the number and types of outputs
sns.countplot(x=df["Type"], width=0.7);



```
[6]: # finding the total number of possible outcomes
df.Type.unique(), len(df.Type.unique())
```

[6]: (array([1, 2, 3, 5, 6, 7]), 6)

Inference

As the data outcomes are discrete, we would use a multiclass classification algorithm such as non-binary multiclass classification.

For this data, I will use Random Forest Classifier Algorithm.

1.0.3 Model Creation and Training

```
[7]: # splitting the data into labels and outcomes
X = df.iloc[:, :-1]
y = df.iloc[:, -1]
X.shape , y.shape
```

[7]: ((214, 9), (214,))

```
[8]: # splitting the data into train and test set with test size of 0.2 as the
      →availability of data is less
     from sklearn.model_selection import train_test_split
     →random state=42)
 [9]: # importing necessary libraries for model creation
     from sklearn.pipeline import Pipeline
     from sklearn.model_selection import GridSearchCV
     from sklearn.ensemble import RandomForestClassifier
     from sklearn.preprocessing import StandardScaler
[10]: # creating the pipeline for the model and getting the pipeline parameters
     pipe = Pipeline([
         ("scaler", StandardScaler()),
         ("randomForest", RandomForestClassifier())
     pipe.get_params()
[10]: {'memory': None,
      'steps': [('scaler', StandardScaler()),
       ('randomForest', RandomForestClassifier())],
      'verbose': False,
      'scaler': StandardScaler(),
      'randomForest': RandomForestClassifier(),
      'scaler__copy': True,
      'scaler__with_mean': True,
      'scaler__with_std': True,
      'randomForest_bootstrap': True,
      'randomForest__ccp_alpha': 0.0,
      'randomForest class weight': None,
      'randomForest__criterion': 'gini',
      'randomForest__max_depth': None,
      'randomForest__max_features': 'sqrt',
      'randomForest__max_leaf_nodes': None,
      'randomForest__max_samples': None,
      'randomForest__min_impurity_decrease': 0.0,
      'randomForest__min_samples_leaf': 1,
      'randomForest__min_samples_split': 2,
      'randomForest__min_weight_fraction_leaf': 0.0,
      'randomForest_n_estimators': 100,
      'randomForest__n_jobs': None,
      'randomForest__oob_score': False,
      'randomForest__random_state': None,
      'randomForest verbose': 0,
      'randomForest__warm_start': False}
```

```
[11]: # creating a cross validation estimator grid with estimators and maximum depth.
       ⇔as grid parameters
      estimator = GridSearchCV(estimator=pipe, param_grid={
          'randomForest__n_estimators':[i for i in range(50,150,5)],
          'randomForest__max_depth':[i for i in range(1,11)]
      }, scoring='accuracy', cv=3)
[52]: # fitting the estimator
      estimator.fit(X_train, y_train)
[52]: GridSearchCV(cv=3,
                   estimator=Pipeline(steps=[('scaler', StandardScaler()),
                                             ('randomForest',
                                              RandomForestClassifier())]),
                   param_grid={'randomForest__max_depth': [1, 2, 3, 4, 5, 6, 7, 8, 9,
                               'randomForest__n_estimators': [50, 55, 60, 65, 70, 75,
                                                              80, 85, 90, 95, 100,
                                                               105, 110, 115, 120, 125,
                                                               130, 135, 140, 145]},
                   scoring='accuracy')
[53]: # storing the results of the estimator in a dataframe
      result_df = pd.DataFrame(estimator.cv_results_)
[54]: # getting the parameters for the best optimum model
      print(estimator.best_index_)
      print(estimator.best_params_)
      print(estimator.best score )
     125
     {'randomForest__max_depth': 7, 'randomForest__n_estimators': 75}
     0.7543859649122807
[55]: # getting the details about best optimum model from result database
      result_df.iloc[150]
[55]: mean_fit_time
     0.122424
      std_fit_time
      0.002658
     mean_score_time
     0.010924
     std_score_time
      0.001117
     param_randomForest__max_depth
```

```
8
      param_randomForest_n_estimators
      100
                                          {'randomForest_max_depth': 8,
      params
      'randomForest__...
      split0_test_score
      0.578947
      split1_test_score
      0.842105
      split2_test_score
      0.701754
     mean_test_score
      0.707602
      std_test_score
      0.107513
      rank_test_score
      93
      Name: 150, dtype: object
[56]: |# creating an improvised estimator with parameters close to best estimate of
       ⇔previous estimator
      est_depth = estimator.best_params_['randomForest__max_depth']
      est_estimators = estimator.best_params_['randomForest__n_estimators']
      delta1 = 5
      delta2 = 2
      estimator_improvised = GridSearchCV(estimator=pipe, param_grid={
          'randomForest__n_estimators':[i for i in_
       →range(est_estimators-delta1,est_estimators+delta1+1)],
          'randomForest__max_depth':[i for i in__
       →range(est_depth-delta2,est_depth+delta2+1)],
      }, scoring='accuracy', cv=3)
[57]: # fitting the improvised estimator
      estimator_improvised.fit(X_train, y_train)
[57]: GridSearchCV(cv=3,
                   estimator=Pipeline(steps=[('scaler', StandardScaler()),
                                              ('randomForest',
                                              RandomForestClassifier())]),
                   param_grid={'randomForest__max_depth': [5, 6, 7, 8, 9],
                                'randomForest__n_estimators': [70, 71, 72, 73, 74, 75,
                                                              76, 77, 78, 79, 80]},
                   scoring='accuracy')
[58]: # storing the results of improvised estimator in a dataframe
      result2_df = pd.DataFrame(estimator.cv_results_)
```

```
[59]: # getting the parameters for the best optimum model from the improvised
       \hookrightarrow estimator
      print(estimator_improvised.best_index_)
      print(estimator_improvised.best_params_)
      print(estimator_improvised.best_score_)
     27
     {'randomForest_max_depth': 7, 'randomForest_n_estimators': 75}
     0.7485380116959064
[60]: # getting the details about best optimum model from improvised result database
      result2_df.iloc[181]
[60]: mean_fit_time
     0.071356
      std_fit_time
      0.000997
      mean_score_time
      0.008694
      std_score_time
      0.000641
      param_randomForest__max_depth
     param_randomForest__n_estimators
      55
                                           {'randomForest__max_depth': 10,
      params
      'randomForest_...
      split0_test_score
      0.596491
      split1_test_score
      0.894737
      split2_test_score
      0.684211
     mean_test_score
      0.725146
      std_test_score
      0.125152
      rank_test_score
      30
      Name: 181, dtype: object
[67]: # getting the top 2 estimators
      model1 = estimator.best_estimator_
      model2 = estimator_improvised.best_estimator_
```

1.0.4 Model Evaluation

```
[68]: # comparing the top two models
     from sklearn.metrics import accuracy_score
     from sklearn.metrics import confusion_matrix
     preds1 = model1.predict(X_test)
     preds_train1 = model1.predict(X_train)
     accuracy_test1 = accuracy_score(y_true=y_test, y_pred=preds1)
     accuracy_train1 = accuracy_score(y_true=y_train, y_pred=preds_train1)
     cmatrix1 = confusion_matrix(y_true=y_test, y_pred=preds1)
     print(f'Model 1 Metrics')
     print(f'accuracy (test)
                                : {accuracy_test1}\n\tin precentage :
      print(f'accuracy (train) : {accuracy_train1}\n\tin precentage :
      print(f'confusion matrix (for 6 outcomes):\n{cmatrix1}')
     preds2 = model2.predict(X_test)
     preds_train2 = model2.predict(X_train)
     accuracy_test2 = accuracy_score(y_true=y_test, y_pred=preds2)
     accuracy_train2 = accuracy_score(y_true=y_train, y_pred=preds_train2)
     cmatrix2 = confusion_matrix(y_true=y_test, y_pred=preds2)
     print(f'\nModel 2 Metrics')
                                : {accuracy_test2}\n\tin precentage :
     print(f'accuracy (test)
      print(f'accuracy (train)
                              : {accuracy_train2}\n\tin precentage :
      →\t{accuracy_train2*100} %\n')
     print(f'confusion matrix (for 6 outcomes):\n{cmatrix2}')
    Model 1 Metrics
                         : 0.9069767441860465
    accuracy (test)
            in precentage: 90.69767441860465 %
    accuracy (train)
                        : 0.9766081871345029
            in precentage: 97.6608187134503 %
    confusion matrix (for 6 outcomes):
     [[11 0 0 0 0 0]
     [311 0 0 0 0]
     [1 0 2 0 0 0]
     [0 0 0 4 0 0]
      [000030]
     [0 0 0 0 0 8]
    Model 2 Metrics
    accuracy (test)
                       : 0.813953488372093
```

```
in precentage: 81.3953488372093 %
                        : 0.9766081871345029
    accuracy (train)
            in precentage: 97.6608187134503 %
    confusion matrix (for 6 outcomes):
     [[11 0 0 0
                     0]
      [ 3 10 0 0
                  0
                     17
     [1 1 1 0 0 0]
     [0 2 0 2 0 0]
      [0 0 0 0 3 0]
      [000008]]
[75]: print(f'Test delta : {(accuracy_test1-accuracy_test2)*100}%')
     print(f'Train delta : {(accuracy_train1-accuracy_train2)*100}%')
```

Test delta : 9.302325581395344%

Train delta: 0.0%

We can see the metrics as:

- Model1
 - Test Accuracy : 90.69%
 - Train Accuracy: 97.66%
- Model2
 - Test Accuracy : 81.39%
 - Train Accuracy: 97.66%

We can see that *model1* perfoms better on train set than test set by a margin of **9.30%** and *model1* perfoms better on test set than train set by a margin of **0.00%**.

Hence, from this observation, we can say that model outperfoms model 2.

Therefore model1 is the most optimum model.

1.0.5 Exporting the Model

```
[71]: # exporting the model in .joblib format
    from joblib import dump
    dump(model1, "glass_pred.joblib")

[71]: ['glass_pred.joblib']

[73]: # loading the model and finding its accuracy
    from joblib import load
    from sklearn.metrics import accuracy_score
```

```
load_model = load("./glass_pred.joblib")

pred_load = load_model.predict(X_test)
accuracy_load_model = accuracy_score(y_true=y_test,y_pred=pred_load)
print(f'Accuracy of the model loaded : {accuracy_load_model*100} %')
```

Accuracy of the model loaded : 90.69767441860465 %

1.0.6 Conclusion

Hence, a Random Forest Classifier algorithm is implemented on the given data and a Machine Learning model with 90.698% accuracy is obtained.