5- class Sentiment Analysis

Using

Support Vector Machine,
Decision Tree,
Random Forest,
Adaboosting,
Gradient Boosting

Submitted by Shristi Maskay

Abstract

The data set we have is a categorical data, consisting of Phrase and sentiment levels 0-5 as very sad, sad, neutral, happy, very happy. The task is to use various classification model to classify the dataset into categories and to used different param_grids and find the best parameters and best model that can be used

Required Libraries used

import numpy as np import pandas as pd import matplotlib.pyplot as plt import seaborn as sns %matplotlib inline

1. Import Dataset

Using Pandas library to import the dataset and store the data in dataframe and then performing replace action to replace NaN data with 0 in the dataframe

```
df = pd.read_csv('/home/shristi/Downloads/SVM/sentiment_5_class - sentiment_5_class.csv')
df.head()

df.replace('NaN', 0)|
```

Sentiment	Phrase	
3	injects just enough freshness into the proceed	0
2	that	1
0	never plays as dramatic even when dramatic thi	2
0	None of this is very original, and it is n't	3
3	, Madonna gives her best performance since Abe	4
***	***	
3	to balance pointed , often incisive satire and	18384
4	have to be a most hard-hearted person not to b	18385
3	could young romantics out on a date	18386
3	could be this good	18387
0	such a dungpile	18388

18389 rows × 2 columns

2. Data exploration: Data analysis and visualization

The data set we have is a categorical data, consisting of columns: Phrase and sentiment. Sentiment column shows the level of sentiments range from 0-5 as very sad, sad, neutral, happy and very happy as per the contents in the phrases.

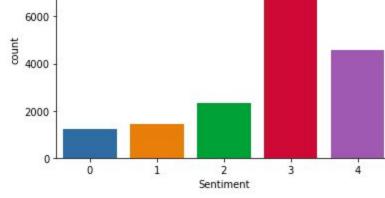
The below command helps to get the columns in the dataframe

```
1 df.columns
Index(['Phrase', 'Sentiment'], dtype='object')
```

Now creating a countplot for Sentiment data got the graph of the phrase categorized into 5 classes. We can state that existing data seems to be imbalanced set as we have data of happy to be quite high in number than data of sad, very sad and neutral

```
1 sns.countplot(x="Sentiment", data=df)
<matplotlib.axes._subplots.AxesSubplot at 0x7f3b42eb9828>

8000 - 6000 - E
```



3. Data Split

To get the model for proper classification now we need to split the data into train set and test set and used the train set to fit the model and use the test set to evaluate the model prediction and determine the accuracy of the model

```
X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.2, random_state=1, stratify=y
```

Splitting the data to train and test set considering 80% of data to train and 20% to test and using the data from the Phrase column need to predict the categories on which the phrase lies.

4. Features: Extraction and normalization

Since the data in Phrase Column is in a text form so, converting it to a matrix of token counts. Initially creating an a-priori dictionary using the words in phrases and then using analyzer for feature selection then the number of features will be equal to the vocabulary size found by analyzing the data.

Using Count Feature Vectorization

It is used to create vectors having dimensionality equal to the size of our vocabulary, and if the phrase consists of words available in vocab we put 1 in dimension and if not found then 0 for each and every word on the phrase. The result thus will be very large vector which will help to give accurate count

```
def count_feature_vectorization(X_train, X_test):
    c_vectorizer = CountVectorizer()
    c_vectorizer.fit(X_train) #only use training data set to fit model
    c_vectorizer.get_feature_names() #vocab list
    print(c_vectorizer)

#to vecorize the train and test data i.e encoding the phrases as per the dictionary created
    c_train_v = c_vectorizer.transform(X_train)
    c_test_v = c_vectorizer.transform(X_test)
    return c_train_v, c_test_v, c_vectorizer

c_train_v, c_test_v, c_vectorizer = count_feature_vectorization(X_train, X_test)
    c_train_v.toarray()
    c_test_v.toarray()
```

Using Tf-IDF Feature Vectorization

This is used to transform a count matrix to tf-idf representation i.e term frequency times inverse document frequency. The purpose of using Tf-idf is to scale down the impact of tokens that occur very frequently in given data.

```
from sklearn.feature extraction.text import TfidfVectorizer
def Tfidf_feature_vectorization(X_train, X_test):
    t vectorizer = TfidfVectorizer()
    t_vectorizer.fit(X_train) #only use training data set to fit model t_vectorizer.get_feature_names() #vocab list
    print(t vectorizer)
    #to vecorize the train and test data i.e encoding the phrases as per the dictionary created
    t train v =t vectorizer.transform(X train)
    t_test_v = t_vectorizer.transform(X_test)
    return t train v, t test v, t vectorizer
t_train_v, t_test_v, t_vectorizer = Tfidf_feature_vectorization(X_train, X_test)
```

```
t_train_v.toarray()
t_test_v.toarray()
```

```
TfidfVectorizer(analyzer='word', binary=False, decode error='strict',
                dtype=<class 'numpy.float64'>, encoding='utf-8',
                input='content', lowercase=True, max_df=1.0, max_features=None,
                min_df=1, ngram_range=(1, 1), norm='l2', preprocessor=None,
                smooth_idf=True, stop_words=None, strip_accents=None,
                sublinear tf=False, token pattern='(?u)\\b\\w\\w+\\b',
                tokenizer=None, use idf=True, vocabulary=None)
```

5. Model Building

SVM (Support Vector Machine)

"Support Vector Machine" (SVM) is a supervised machine learning algorithm that can be used for both classification or regression. However, it is mostly used in classification problems. It uses sets of training points in the function thus is memory efficient and has different kernel functions can be used in decision functions. Sometimes SVM might lead to overfitting when features are greater than number of samples. So in this case 5fold cross validation might be used which might be expensive.

To determine the best parameters for SVM model, the parameters used in model training are:

```
grid param = [{'kernel': ['rbf'], 'gamma': [1e-3, 1e-4, 1e-2],
              'C': [1, 10, 100, 1000]},
              {'kernel': ['linear'], 'C': [1, 10, 100, 1000]}]
```

And considering the score from the fitted model we get the best parameters that can be used in classification for the specific model and develop the model. Here,

Best parametes:

Score:

Library: from sklearn.svm import SVC

```
def svm model(X train v, y train):
   model = SVC(random state = 1)
   model.fit(X_train_v, y_train)
   return model
```

Now fitting the model using vectorized train data we get a model and now using the model

Model evaluation using the test set.

```
def prediction(model, feature_data):
    y_pred = model.predict(feature_data)
    return y_pred
```

Using model generated using tf-idf vectorizer train data and count vectorized data we predicted the output and generated a metric report

from sklearn import metrics print(metrics.classification_report(y_test, y_pred)

	precision	recall	f1-score	support	
0	0.79	0.45	0.57	247	
1	0.62	0.30	0.40	291	
2	0.78	0.20	0.32	469	
3	0.60	0.91	0.73	1759	
4	0.76	0.52	0.62	912	
accuracy			0.64	3678	
macro avg	0.71	0.48	0.53	3678	
ghted avg	0.68	0.64	0.61	3678	
Classifica	precision		f1-score	support	raıı
0	0.81	0.55	0.65	247	
1	0.64	0.44	0.52	291	
-	0.74	0.52	0.61	469	
2	0.68	0.89	0.77	1759	
3	0.00		0 66	912	
	0.78	0.58	0.66		
3		0.58	0.71	3678	
3 4		0.58	0.71	3678 3678	
3 4 accuracy	0.78		0.71 0.64		

This shows that the model generated using TF-IDF is more accurate than Count vectorization

Model Selection

For Model selection, GridSearchCV library is used for hyper parameter tuning process to determine optimal values for the given model as the model depends on the specified

hyperparameter values. Cross validation process is performed in order to determine the best score and parameters for the model.

```
from sklearn.model selection import GridSearchCV
  1 grid param = {'kernel': ('linear', 'rbf'), 'C':(1,10, 0.5)}
  2 grid param
 {'kernel': ('linear', 'rbf'), 'C': (1, 10, 0.5)}
    grid param = [{'kernel': ['rbf'], 'gamma': [1e-3, 1e-4, 1e-2],
                          'C': [1, 10, 100, 1000]},
  2
                         {'kernel': ['linear'], 'C': [1, 10, 100, 1000]}]
  3
  4 grid param
 [{'kernel': ['rbf'], 'qamma': [0.001, 0.0001, 0.01], 'C': [1, 10, 100, 1000]},
 {'kernel': ['linear'], 'C': [1, 10, 100, 1000]}]
  1 from sklearn.metrics import make scorer, fl score
  2 scorer = make scorer(f1 score, average='micro')
 #using tfidf vectorised train data
 scorer = make scorer(f1 score, average='micro')
 clf tfidf = GridSearchCV(SVC(), grid param,scoring=scorer)
 clf tfidf.fit(t train v, y train)
 print(clf tfidf)
 print(clf tfidf.best score , clf tfidf.best params )
Performing gridsearch with kernel parameters and getting the best_score result we got,
 GridSearchCV(cv=None, error score=nan,
              estimator=SVC(C=1.0, break ties=False, cache size=200,
                            class weight=None, coef0=0.0,
                            decision function shape='ovr', degree=3,
                            gamma='scale', kernel='rbf', max iter=-1,
                            probability=False, random state=None, shrinking=True,
                            tol=0.001, verbose=False),
              iid='deprecated', n jobs=None,
              param grid=[{'C': [1, 10, 100, 1000],
                           'gamma': [0.001, 0.0001, 0.01], 'kernel': ['rbf']},
                          {'C': [1, 10, 100, 1000], 'kernel': ['linear']}],
              pre_dispatch='2*n jobs', refit=True, return train score=False,
              scoring=make scorer(f1 score, average=micro), verbose=0)
 0.6851333505653415 {'C': 100, 'gamma': 0.01, 'kernel': 'rbf'}
Best score = 0.6851
Best parameters = C:100, gamma:0.01, kernel: rbf
```

Now using the best parameters obtained from the gridsearch

Final Model using best parameter for SVM

```
model = SVC(random_state=1, kernel="rbf", C=100.0, gamma=0.01)
model.fit(t_train_v, y_train)
y_pred = model.predict(t_test_v)
print(metrics.classification_report(y_test, y_pred))
#FInal result obtained
```

	precision	recall	f1-score	support
0	0.73	0.66	0.70	247
1	0.60	0.53	0.56	291
2	0.68	0.58	0.63	469
3	0.72	0.84	0.78	1759
4	0.72	0.61	0.66	912
accuracy			0.71	3678
macro avg	0.69	0.64	0.66	3678
weighted avg	0.71	0.71	0.71	3678

Classification report generated for final model

5.2. Decision Tree

Decision tree is another non-linear model formed from a combination of linear boundaries. It is simply a tree formed by a series of yes/no questions asked about data leading to a predicted class. This is an interpretable model because it makes classifications much like we do: we ask a sequence of queries about the available data we have until we arrive at a decision.

Attributes in decision tree model:

__init__(self, *, criterion='gini', splitter='best', max_depth=None, min_samples_split=2, min_samples_leaf=1, min_weight_fraction_leaf=0.0, max_features=None, random_state=None, max_leaf_nodes=None, min_impurity_decrease=0.0, min_impurity_split=None, class_weight=None, presort='deprecated', ccp_alpha=0.0)

Parameters used in model training:

```
grid_params = {
    'max_depth': (5, 10, 20, 50, 100, 500),
    'min_samples_split': (2, 3, 4, 6, 8, 16, 20)
}
```

```
GridSearchCV(cv=3, error score=nan,
             estimator=DecisionTreeClassifier(ccp alpha=0.0, class weight=None,
                                              criterion='gini', max depth=None,
                                              max features=None,
                                              max leaf nodes=None,
                                              min impurity decrease=0.0,
                                              min_impurity_split=None,
                                              min samples leaf=1,
                                              min samples split=2,
                                              min weight fraction leaf=0.0,
                                              presort='deprecated',
                                              random state=None,
                                              splitter='best'),
             iid='deprecated', n_jobs=None,
             param grid={'max depth': (5, 10, 20, 50, 100, 500),
                         'min samples split': (2, 3, 4, 6, 8, 16, 20)},
             pre dispatch='2*n jobs', refit=True, return train score=False,
             scoring=make scorer(f1 score, average=micro), verbose=0)
```

Best_parameters: {'max_depth': 500, 'min_samples_split': 6

Best Score: 0.5981915944749566}

Final Model using best parameter for Decision Tree

```
import sklearn.metrics as metrics
model =DecisionTreeClassifier(random_state=1,max_depth=500,min_samples_split=6)
model.fit(c_train_v, y_train)
preds = model.predict(c_test_v)
print(metrics.classification_report(y_test, preds))
```

		precision	recall	f1-score	support
	0	0.59	0.53	0.56	247
	1	0.54	0.33	0.41	291
	2	0.51	0.29	0.37	469
	3	0.63	0.82	0.71	1759
	4	0.69	0.55	0.61	912
accur	асу			0.63	3678
macro	avg	0.59	0.50	0.53	3678
weighted	avg	0.62	0.63	0.61	3678

5.3. Random Forest

It technically is an ensemble method of decision trees generated on a randomly split dataset. This collection of decision tree classifiers is also known as the forest. The individual decision trees are generated using an attribute selection indicator such as information gain, gain ratio, and Gini index for each attribute. Each tree depends on an independent random sample. In a classification problem, each tree votes and the most popular class is chosen as the final result. In the case of regression, the average of all the tree outputs is considered as the final result. It is simpler and more powerful compared to the other non-linear classification algorithms. It is considered as an accurate and robust method because of the use of a number of decision trees.

Steps:

- 1. Select random samples from a given dataset.
- 2. Construct a decision tree for each sample and get a prediction result from trees.
- 3. Perform a vote for each predicted result.
- 4. Select the prediction result with the most votes as the final prediction.

```
Parameters used in model training:
```

```
grid params = {
        'max depth': (3,10, 13, 50),
        'min_samples_split': (2, 4, 8, 10),
        'n estimators':(10, 20),
      }
: scorer = make scorer(f1 score, average='micro')
  clf=GridSearchCV(RandomForestClassifier(),grid params,scoring=scorer,cv=3)
  clf.fit(c train v, y train)
: GridSearchCV(cv=3, error score=nan,
               estimator=RandomForestClassifier(bootstrap=True, ccp alpha=0.0,
                                                 class weight=None,
                                                  criterion='gini', max depth=None,
                                                 max features='auto',
                                                 max leaf nodes=None,
                                                 max samples=None,
                                                 min impurity decrease=0.0,
                                                 min_impurity_split=None,
                                                 min samples leaf=1,
                                                 min samples split=2,
                                                 min weight fraction leaf=0.0,
                                                 n estimators=100, n jobs=None,
                                                 oob score=False,
                                                  random state=None, verbose=0,
                                                 warm start=False),
               iid='deprecated', n jobs=None,
               param grid={'max depth': (3, 10, 13, 50),
                            'min samples split': (2, 4, 8, 10),
                            'n estimators': (10, 20)},
```

Best_parameters: {'max_depth': 50, 'min_samples_split': 2, 'n_estimators': 20}

Best Score: 0.5466657283990215

Final Model Score

```
from sklearn.model_selection import cross_val_score
forest_scores = cross_val_score(model, c_train_v, y_train, cv=10)
forest_scores.mean()
```

pre_dispatch='2*n_jobs', refit=True, return_train_score=False, scoring=make scorer(f1 score, average=micro), verbose=0)

: 0.5574050760352319

5.4. Adaboosting

Adaboosting depends on classification problems and aims in converting weak classifiers to strong. It is a weighted combination of M weak classifiers. In the process of iteration if the accuracy is higher than 50% the weight is positive and more accurate the classifier, larger the weight while less than 50% accuracy gives negative weight.

```
Parameter used for model training:
      grid params = {
        'learning rate': (0.001,0.1, 0.5,0.8),
        'n estimators':(10,50,100),
      }
scorer = make scorer(f1 score, average='micro')
clf=GridSearchCV(AdaBoostClassifier(),grid params,scoring=scorer,cv=3)
clf.fit(c train v, y train)
GridSearchCV(cv=3, error score=nan,
              estimator=AdaBoostClassifier(algorithm='SAMME.R',
                                            base estimator=None,
                                             learning rate=1.0, n estimators=50,
                                             random state=None),
              iid='deprecated', n jobs=None,
              param grid={'learning rate': (0.001, 0.1, 0.5, 0.8),
                           'n estimators': (10, 50, 100)},
              pre dispatch='2*n jobs', refit=True, return train score=False,
              scoring=make scorer(fl score, average=micro), verbose=0)
```

Best_parametes: {'learning_rate': 0.8, 'n_estimators': 100}

F1_score: 0.512677440441354

Final Model using Best parameter

	0	0.84	0.09	0.15	247
	1	0.33	0.00	0.01	291
	2	0.20	0.00	0.00	469
	3	0.49	1.00	0.65	1759
	4	0.54	0.02	0.03	912
accurac	У			0.49	3678
macro av	g	0.48	0.22	0.17	3678
weighted av	g	0.47	0.49	0.33	3678

5.5. Gradient Boosting

This model is used for regression and classification models to predict models in the form of an ensemble of weak prediction models. It is similar to other boosting models allowing optimization by using differential loss function. It helps in combination of weak learners into one strong learner in iterative fashion. It is used with decision trees. It's performance depends on the size of trees, and number of leaves. In cases when training sets are too closed it can lead to overfitting and that can be reduced with regularization technique

```
Parameter used for model training:
      grid params = {
        'min_samples_split': (2, 4, 6, 8, 10, 20),
        'n estimators':(50,100),
      }
scorer = make scorer(f1 score, average='micro')
clf=GridSearchCV(GradientBoostingClassifier(),grid params,scoring=scorer,cv=3)
clf.fit(c train v, y train)
GridSearchCV(cv=3, error score=nan,
              estimator=GradientBoostingClassifier(ccp alpha=0.0,
                                                     criterion='friedman mse',
                                                     init=None, learning rate=0.1,
                                                     loss='deviance', max depth=3,
                                                    max features=None,
                                                    max leaf nodes=None,
                                                    min impurity decrease=0.0,
                                                    min impurity split=None,
                                                    min samples leaf=1,
                                                    min samples split=2,
                                                    min weight fraction leaf=0.0,
                                                    n estimators=100,
                                                     n iter no change=None,
                                                    presort='deprecated',
                                                     random state=None,
                                                     subsample=1.0, tol=0.0001,
                                                    validation fraction=0.1,
                                                    verbose=0, warm start=False),
              iid='deprecated', n jobs=None,
              param grid={'min samples split': (2, 4, 6, 8, 10, 20),
                           'n estimators': (50, 100)},
              pre dispatch='2*n jobs', refit=True, return train score=False,
              scoring=make scorer(f1 score, average=micro), verbose=0)
```

Best_parametes: {'min_samples_split': 2, 'n_estimators': 100}

Best F1 score: 0.5529199171928897

Final Model using Best parameter

```
model =GradientBoostingClassifier(min_samples_split=2, n_estimators=100)
model.fit(c_train_v, y_train)
preds = model.predict(c_test_v)
print(metrics.classification_report(y_test, preds))
```

	precision	recall	f1-score	support
0	0.84	0.23	0.37	247
1	0.76	0.11	0.19	291
2	0.78	0.06	0.11	469
3	0.53	0.94	0.68	1759
4	0.68	0.29	0.40	912
accuracy			0.55	3678
macro avg	0.72	0.33	0.35	3678
weighted avg	0.64	0.55	0.48	3678

Model Comparison

Model	Parameter	F1_score (training)	F1_score final
SVM	C:100, gamma:0.01, kernel: rbf	0.6851	0.75
Decision Tree	'max_depth': 500, 'min_samples_split': 6	0.5819	0.63
Random Forest	{'max_depth': 50, 'min_samples_split': 2, 'n_estimators': 20}	0.54666	0.5574
Adaboosting	{'learning_rate': 0.8, 'n_estimators': 100}	0.512677	0.49
Gradient Boosting	{'min_samples_split': 2, 'n_estimators': 100}	0.5529	0.55

From the 5 model comparison SVM is found to be best model for the categorical data classification

Code link https://github.com/Kristiee/SVM_sentiment

SVM https://github.com/Kristiee/SVM_sentiment/blob/master/SVM-sentiment-final.ipynb
Other 4 models (Decision tree, Random forest, adaboosting, gradient boosting)
https://github.com/Kristiee/SVM_sentiment/blob/master/sentiment_analysis_other_models.ipynb