Qualitative Bankruptcy Prediction

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Executive Summary

The data in this study is the qualitative bankruptcy data set from UCI [1]. We are to classify companies and predict bankruptcy or non-bankruptcy from qualitative features assessed by experts [2]. This study was inspired by the predictive system [3] where the algorithms used are Logistic Regression, Naive Bayes, Neural Networks, Random Forest and RBF Support Vector Machine. We then added kNN, Decision Trees and Gradient Boosting classifiers to potentially increase the test accuracy. In the previous paper, a 70% - 30% train test split was used. The test accuracy achieved is between 97.2% to a maximum of 99.6%. In this paper, we are able get a test accuracy of 97.5% up to 100%. The latter was achieved through Gradient Boosting Method. The top predictors of accuracy were Competitiveness and Credibility.

Data Description

There are 250 instances in the data with 6 attributes corresponding to the qualitative parameters in bankruptcy.

Feature	Description	Legend		
Industrial Risk	Government policies and international agreements, competition, sensitivity to changes in macroeconomic factors and competitive power			
Management risk	Ability and competence of the management, growth process and business performance, short term andlong term business planning	Positive (P), Average (A), Negative (N)		
Financial Flexibility	Forms of financing			
Credibility	Credibility Credit history, reliability of information and relationship with financial institutes			
Competitiveness	Competitiveness Market position, level of core capacities and differentiated strategy			
Operating Risk Stability and diversity of procurement, transaction, efficiency of production, sales diversification, collection of A/R, effectiveness sale network		Bankruptcy (B), Non-Bankruptcy (NB)		

Data Pre-Processing

We now look at the data and label the headers. Then we looked into the numbers of bankruptcy or non-bankruptcy to check whether undersampling (or oversampling) needs to be done. The data has a 40% - 60% proportion of bankruptcy and non-bankruptcy respectively.

We then proceeded to use get_dummies into the target column to convert the categorical variable. Drop_first was used so that we have ones for NB and zeros for B and remove the redundant feature.

For the other features in the data set, it was replaced with 0 for Negative, 0.5 for Average and 1 for Positive. This is similar to the concept of MinMaxScaler.

In [40]: import pandas as pd import numpy as np import matplotlib.pyplot as plt from collections import Counter import seaborn as sns

In [2]: df = pd.read_csv('Qualitative_Bankruptcy.data.txt', header=None)

In [3]: df.head()

Out[3]:

	0	1	2	3	4	5	6
0	Р	Ρ	Α	Α	Α	Ρ	NB
1	Ζ	Z	Α	Α	Α	Ζ	NB
2	Α	Α	Α	Α	Α	Α	NB
3	Р	Р	Р	Р	Р	Р	NB
4	Ν	Ζ	Ρ	Ρ	Р	Ν	NB

Out[6]:

	Industrial Risk	Management Risk	Financial Flexibility	Credibility	Competitiveness	Operating Risk	Target
0	Р	Р	Α	Α	Α	Р	NB
1	N	N	Α	Α	Α	Ν	NB
2	А	А	А	А	Α	Α	NB
3	Р	Р	Р	Р	Р	Р	NB
4	N	N	Р	Р	Р	N	NB

In [7]: df['Target'].value_counts()

Out[7]: NB 143 B 107

Name: Target, dtype: int64

Out[8]:

	Industrial Risk	Management Risk	Financial Flexibility	Credibility	Competitiveness	Operating Risk	Target
0	Р	Р	А	Α	А	Р	1
1	N	N	Α	Α	А	N	1
2	А	А	Α	Α	А	Α	1
3	Р	Р	Р	Р	Р	Р	1
4	N	N	Р	Р	Р	N	1

```
In [10]: df.replace('N',0,axis=1, inplace=True)
    df.replace('A',0.5,axis=1, inplace=True)
    df.replace('P',1,axis=1, inplace=True)

    df.head()
```

/anaconda3/lib/python3.6/site-packages/pandas/core/generic.py:4480:
UserWarning: the "axis" argument is deprecated and will be removed i
nv0.13; this argument has no effect
warnings.warn('the "axis" argument is deprecated '

Out[10]: ____

	Industrial Risk	Management Risk		Credibility	Competitiveness	Operating Risk	Target
0	1.0	1.0	0.5	0.5	0.5	1.0	1
1	0.0	0.0	0.5	0.5	0.5	0.0	1
2	0.5	0.5	0.5	0.5	0.5	0.5	1
3	1.0	1.0	1.0	1.0	1.0	1.0	1
4	0.0	0.0	1.0	1.0	1.0	0.0	1

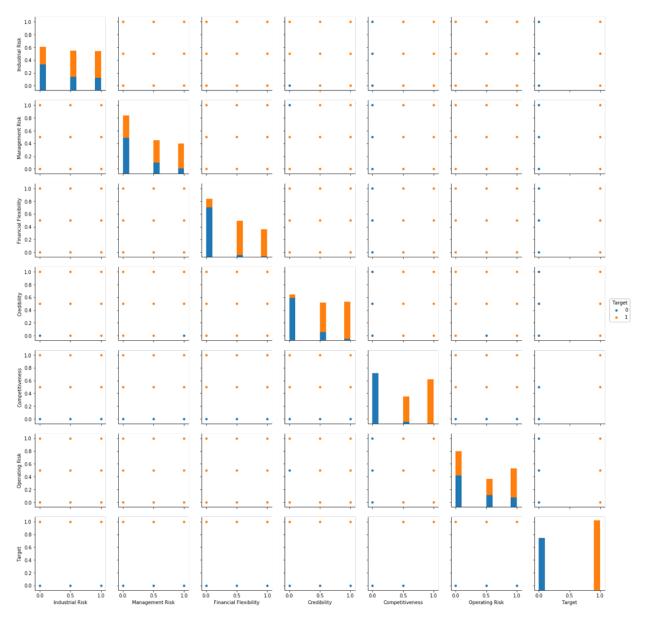
Exploratory Data Analysis

We now looked into the pairplot of the features to have a visual check of possible predictors of the target.

We can see below that Competitiveness can somehow be clearly separate the Target. The same is true for Credibility and Financial Flexibility. We'll park this for now to see if we can use the following information for feature selection later on.

In [41]: sns.pairplot(df, hue="Target")

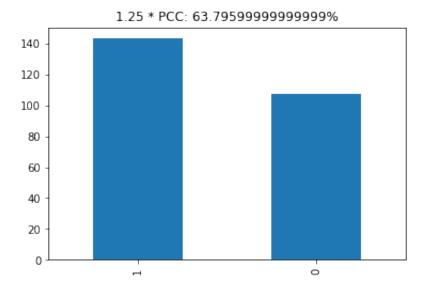
Out[41]: <seaborn.axisgrid.PairGrid at 0x1a1d8e2e48>



We now achieved a score to beat of 63.795%. This is 1.25 times the Proportional Chance Criterion or the accuracy we need to beat inorder to say that what we have in our output is indeed a good predictor.

```
In [42]: state_counts = Counter(df['Target'])
    df_state = pd.DataFrame.from_dict(state_counts, orient='index')
    df_state.plot(kind='bar', colormap='tab10')

pcc1 = (df_state[0]/df_state[0].sum())**2
    score_to_beat = 1.25*100*pcc1.sum()
    plt.title("1.25 * PCC: {}%".format(score_to_beat));
    plt.legend().set_visible(False)
```



Model

We use the machine learning models in the previous paper and added k-Nearest Neighbor, Decision Trees and Gradient Boosting classifier. We will go over the added models briefly.

k-Nearest Neighbor is one of the simplest models in machine learning. It is non-parametric and instance based learning algorithm, meaning it does not explicitly learn the model

Decision Trees is an algorithm where the data does not need scaling. Essentially, they learn a hierarchy of if-else questions leading to a decision.

Gradient Boosting Method follows a weaker decision tree then learns from it, that's why it is one of the most powerful models.

```
In [33]: X = df.drop('Target', axis=1)
y = df['Target']
```

```
In [35]: #from sklearn.grid search import GridSearchCV
         from sklearn.model selection import GridSearchCV, train test split
         from sklearn.neighbors import KNeighborsClassifier
         from sklearn.linear model import LogisticRegression
         from sklearn.svm import LinearSVC, SVC
         from sklearn.tree import DecisionTreeClassifier
         from sklearn.ensemble import RandomForestClassifier, GradientBoostingC
         lassifier
         from sklearn.naive bayes import GaussianNB
         def knn clf(X, y, n neighbor range=np.arange(1, 20)):
             param grids = {'n neighbors' : n neighbor range}
             clf = KNeighborsClassifier()
             gs clf = GridSearchCV(clf, param grids, n jobs=-1, cv=50).fit(X, y
         )
             return pd.DataFrame(
                 {'Params' : gs clf.cv results ['params'],
                   'Test scores' : gs clf.cv results ['mean test score']}
                                 ).sort values(by='Test scores',
                                               ascending=False).iloc[0,:].tolist
         ()
         def lin_clf(X, y, kind, C_range=[ 1e-4, 1e-3,0.1, 0.2,0.4,
                                     0.75, 1, 1.5, 3, 5, 10, 15, 20, 100, 300,
         1000]):
             param_grids = {'penalty' : ('l1', 'l2'), 'C' : C range}
             if kind == 'logistic':
                 clf = LogisticRegression(dual=False)
             if kind == 'lin svc':
                 clf = LinearSVC(dual=False)
             gs_clf = GridSearchCV(clf, param_grids, n_jobs=-1, cv=50).fit(X, y
         )
             return pd.DataFrame(
                 {'Params' : gs clf.cv results ['params'],
                   'Test scores' : gs clf.cv results ['mean test score']}
                                 ).sort values(by='Test scores',
                                               ascending=False).iloc[0,:].tolist
         ()
         def non_lin_svm_clf(X, y, C_range=[ 1e-4, 1e-3,0.1, 0.2,0.4,
                                      0.75, 1, 1.5, 3, 5, 10, 15, 20, 100, 300,
         1000],
                         gamma range=np.arange(0.1,2.1,0.1), degree range=[2,3,
         4]):
             param_grids = [{'kernel' : ['rbf'], 'C' : C_range,
```

```
'gamma' : gamma range},
                   {'kernel' : ['poly'], 'C' : C range,
                   'gamma' : gamma range, 'degree' : degree range}]
    clf = SVC()
    gs clf = GridSearchCV(clf, param grids, n jobs=-1, cv=50).fit(X, y
)
    return pd.DataFrame(
        {'Params' : gs clf.cv results ['params'],
         'Test scores' : gs clf.cv results ['mean test score']}
                       ).sort values(by='Test scores',
                                      ascending=False).iloc[0,:].tolist
()
def trees clf(X, y, kind, max depth range=np.arange(2, 8),
              learning rate range=np.arange(0.1, 2.1, 0.1),
              max feat gb=[3,4,5]):
    if kind == 'des tree':
        param grids = {'max depth' : max depth range}
        clf = DecisionTreeClassifier()
    if kind == 'rand for':
        mfeat = int(np.sqrt(X.shape[1]))
#
          param grids = {'max depth' : max depth range,
#
                          'max features' : np.arange(mfeat - 2, mfeat)}
        param grids = {'max depth' : max depth range,
                       'max features' : [3,4]}
        clf = RandomForestClassifier(n estimators=100, n jobs=-1)
    if kind == 'grad boost':
        param_grids = {'learning_rate' : learning_rate_range,
                        'max depth' : max depth range,
                       'max features' : max feat gb}
        clf = GradientBoostingClassifier(n estimators=100)
    gs clf = GridSearchCV(clf, param grids, n jobs=-1, cv=50).fit(X, y
)
    return pd.DataFrame(
        {'Params' : gs clf.cv results ['params'],
         'Test scores' : gs clf.cv_results_['mean_test_score']}
                       ).sort values(by='Test scores',
                                      ascending=False).iloc[0,:].tolist
()
def nb clf(X, y):
    clf = GaussianNB()
    test acc = []
    for i in range(100):
        X_train, X_test, y_train, y_test = train_test_split(X, y,
                                                             test size=
0.25,
                                                             random sta
te=i)
        clf.fit(X train, y train)
```

```
test_acc.append(clf.score(X_test,y_test))
return np.mean(test acc)
```

```
In [36]:
         knn res = knn clf(X, y,
                      n neighbor range=np.arange(1, 20))
         #print('knn Done!')
         log res = lin clf(X, y,
                      'logistic', C range=[ 1e-4, 1e-3,0.1, 0.2,0.4,
                                      0.75, 1, 1.5, 3, 5, 10, 15, 20, 100, 300,
         10001)
         #print('logistic reg Done!')
         \lim svc res = \lim clf(X, y,
                      'lin svc', C range=[ 1e-4, 1e-3,0.1, 0.2,0.4,
                                      0.75, 1, 1.5, 3, 5, 10, 15, 20, 100, 300,
         1000])
         #print('lin svc Done!')
         non_lin_svc_res = non_lin_svm_clf(X, y,
                          C range=[ 1e-4, 1e-3, 0.1, 0.2, 0.4,
                                      0.75, 1, 1.5, 3, 5, 10, 15, 20, 100, 300,
         1000],
                          gamma range=np.arange(0.1,2.1,0.1), degree range=[2,3,
         41)
         #print('nonlin svc Done!')
         des t res = trees clf(X, y,
                        'des tree', max depth range=np.arange(2, 8))
         #print('des trees Done!')
         rand f res = trees clf(X, y,
                        'rand for', max depth range=np.arange(2, 5))
         #print('rand forest Done!')
         grad b ref = trees clf(X, y,
                        'grad boost', max depth range=np.arange(2, 5),
                       learning rate range=np.arange(0.1, 0.5, 0.1),
                       \max feat gb=[3,4])
         #print('grad boost Done!')
         nb res = nb clf(X, y)
         #print('bayes Done!')
```

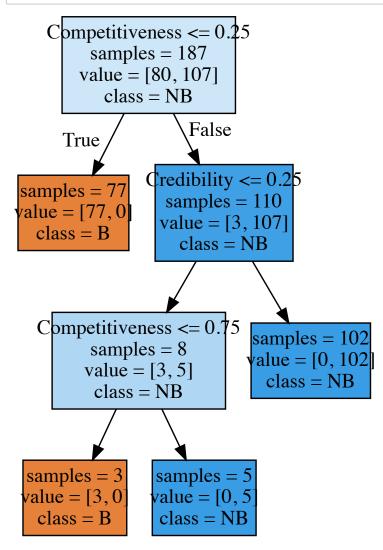
Out[46]:

	Test Accuracy	Best Parameters
k-Nearest Neighbors	0.996000	{'n_neighbors': 1}
Logistic Regression	0.996000	{'C': 3, 'penalty': 'I1'}
Linear SVC	0.996000	{'C': 0.75, 'penalty': 'I1'}
Non-linear SVC	0.996000	{'C': 1000, 'gamma': 1.4000000000000001, 'kern
Decision Trees	0.992000	{'max_depth': 3}
Random Forest	0.996000	{'max_depth': 3, 'max_features': 4}
Gradient Boosting	1.000000	{'learning_rate': 0.1, 'max_depth': 2, 'max_fe
Naive Bayes	0.975556	NA

```
In [49]: grad_b_ref[0]
Out[49]: {'learning_rate': 0.1, 'max_depth': 2, 'max_features': 3}
```

We now look into the top predictors of the data. Decision tree is not the model with highest accuracy but we use it for interpretability of results. We can see from the Decision Tree below that we have Competitiveness and Credibitlity as the top predictors.

Out[47]:



Conclusion

Using all the features in the dataset, we are able to predict if a company is going under. Using the Gradient Boosting Method, we are able to predict with a test accuracy of up to 100% with a learning rate of 0.1, max_depth of 2 and max_features of 3. We are able to create a better algorithm than the previous paper [3].

We are also able to show in this paper that competitiveness, which refers to differentiated stategy, and credibility, in terms of credit score, are the top predictors. Companies need to highlight these factors and make sure that they are always positive so as not to go bankrupt and prevent financial crisis.

References

[1] https://archive.ics.uci.edu/ml/datasets/qualitative_bankruptcy (https://archive.ics.uci.edu/ml/datasets/qualitative_bankruptcy)

[2]

http://neuron.csie.ntust.edu.tw/homework/95/neuron/homework/Homework%233/M9509007/paper/5.htm (http://neuron.csie.ntust.edu.tw/homework/95/neuron/homework/Homework%233/M9509007/paper/5.htm) [3] https://arxiv.org/pdf/1502.03601.pdf (https://arxiv.org/pdf/1502.03601.pdf)