Application and Comparison of Machine Learning Techniques in Business

by

Shukhrat Khuseynov

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Application and Comparison of Machine Learning Techniques in Business

Koç University

Graduate School of Sciences and Engineering

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Shukhrat Khuseynov

and have found that it is complete and satisfactory in all respects, and that any and all revisions required by the final examining committee have been made.

Con	nmittee Members:
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	Asst. Prof. David Carlson (Advisor)
-	Asst. Prof. Jeffrey Ziegler
-	Assoc. Prof. Özden Gür Ali
_	
Date	e:

ABSTRACT

Application and Comparison of Machine Learning Techniques in Business Shukhrat Khuseynov Master of Science in Data Science August 1, 2021

Data science methodology and, particularly, machine learning techniques, are being widely used in many different fields today. There is the same trend in the private sector, where using machine learning is highly advantageous. The prediction of key variables in business given multiple input parameters is important and may affect the firm's profitability. This research aims to investigate several problems in business, being churn estimation, housing price prediction and sentiment analysis, which require a certain data framework and an algorithmic approach. The applied algorithms are compared for each case. The regarded cases involve an application of classification, regression, and text analysis types of machine learning models, demonstrating their diversity and useful application in the industry.

ÖZETÇE

Makine Öğrenmesi Tekniklerinin İşletmede Uygulanması ve Karşılaştırılması Shukhrat Khuseynov Veri Bilimleri, Yüksek Lisans 1 Ağustos 2021

Veri bilimi metodolojisi ve özellikle makine öğrenmesi teknikleri günümüzde birçok farklı alanda yaygın olarak kullanılmaktadır. Özel sektörde de kullanımı oldukça avantajlı olan makine öğrenmesini uygulama konusunda aynı eğilim bulunmaktadır. Birden fazla girdi parametresinin olduğu bilinen işletmedeki kilit değişkenlerin tahmini önemlidir ve bu, firmanın kar düzeyini etkileyebilmektedir. Bu araştırma, belirli bir veri çerçevesi ve algoritmik bir yaklaşım gerektiren iş dünyasındaki çeşitli sorunları; bu kapsamda müşteri kaybı tahmini, konut fiyat tahmini ve duygu analizini araştırmayı amaçlamaktadır. Uygulanan algoritmalar her bir vaka için karşılaştırılmaktadır. Ele alınan vakalar, makine öğrenmesi modellerinin sınıflandırma, regresyon ve metin analizi türlerinin birer uygulamasını içermekte ve bunların endüstrideki çeşitliliğiyle birlikte faydalı uygulamalarını da göstermektedir.

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ABBREVIATIONS

AUROC Area Under the ROC curve

BoW Bag-of-Words

KNN K Nearest Neighbors

OLS Ordinary Least Squares

RMSE Root Mean Square Error

ROC Receiver Operating Characteristic

SVM Support Vector Machines

TF-IDF Term Frequency – Inverse Document Frequency

XGBoost Extreme Gradient Booster

Introduction 1

Chapter 1:

INTRODUCTION

The application of data science and machine learning techniques is very important nowadays since it yields efficiencies and high profits in return, especially in the private sector. Other than direct business applications, artificial intelligence and machine learning can be observed in e-commerce, chatbot applications, banking, healthcare, cyber security, and commercial applications [Rao & Bhattacharyya, 2019]. The study also notes that this methodology brings gradually more and more security and profits to the industry, reducing the unneeded costs. Some of the recent examples for innovative business models include car-sharing and house-sharing services, and even manufacturing industry with Internet of things (IoT), utilizing big data and machine learning algorithms for better productivity [Jeong, 2018]. All different business tasks that use data science methodology generate vast amounts of data, sometimes even in real time, which then are taken as an input to a certain machine learning algorithm or a framework of algorithms to make some decision, prediction or detection. There is a wide range of learning models, depending on a type of problem. The main traditional models are regression analyses, Bayesian methods, random forests, and support vector machines. The more recent approaches also comprise deep learning algorithms, which outperform the classical models in some cases, especially for larger amounts of data [Kraus, Feuerriegel, & Oztekin, 2020]. Consequently, the comparison of algorithms for several business cases is worthwhile, emphasizing the diversity of machine learning algorithms and their beneficial application in the industry. Thus, this paper aims to compare a range of algorithms for three distinct business problems with a specific setup. The first chapter discusses the churn estimation, which involves predicting whether clients will stop doing business with a particular company, representing a binary classification task. Then, the second chapter describes the prediction of housing prices, given the key details of the apartments, using various regression models. Finally, the third chapter involves a sentiment analysis of e-commerce reviews, first implementing the text analysis, then the classification task to find a relationship between the reviews and the recommendations with the ratings given.

Chapter 2:

CHURN ESTIMATION

One of the common business problems in many companies is to find out which customers will potentially be using their products or services continuously in a long-term perspective, becoming loyal clients, and which ones will possibly churn. It is very important to identify these types of customers to manage the time and resources more efficiently, also to try to take precautionary measures to keep those who have the tendency to leave. The churn estimation is based on the machine learning algorithms of binary classification predicting whether a client will stop doing business with the company or not, given multiple parameters. The target variable, the churn of customers, usually represents a sparse vector due to the fact that only some customers tend to leave. The most common fields of churn estimation problem are telecom industry, insurance companies, banking sector, financial and subscription services [Siemes, 2016].

The data are obtained from a mobile company and anonymously shared on Kaggle¹. The dataset consists of 66469 entries depicting the customers from a few months of 2013. There are 66 variables related to different measurements of calls, messages, durations and other information extracted via data mining techniques. The variables include the id, which is dropped before modeling since it is different for each customer and, therefore, useless in predictions, and the churn itself, which is the target variable in the models. The churn distribution can be observed in the data. Apparently, 79.1% of the customers are retained and 20.9% are exited customers. Such a distribution implies that guessing all customers are retained in the business regardless of any variable would result in almost 80% of accuracy score, which is the baseline to beat. The underrepresented group of exited clients has to be dealt with more carefully. For the same reason different measurements along with the accuracy score, such as the AUROC or, in other words, the area under the Receiver Operating Characteristic (ROC) curve, which constitutes a tradeoff between the true positive (TP) and the false positive (FP) rates, have to be used since the accuracy score can be misleading. The AUROC measures the overall performance of a classifier given any possible threshold for the probability between predicting one class or the other [Mand'ák & Hančlová, 2019].

¹ Can be retrieved from www.kaggle.com/dimitaryanev/mobilechurndataxlsx

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Before the application of various machine learning models, the dataset is normalized using the min-max scaler, which normalizes by only keeping the relative distances between the data points. It is done to use the algorithms more efficiently and faster, because normalization allows some algorithms to approach the solution faster. Furthermore, after deciding the best parameters for each model, where it is required, in a grid search of 5-fold cross validation, maximizing the area under ROC (AUROC), 80% of the data are used for model training and the remaining 20% are for model testing purposes. The ROC curves of all models are summarized at the end.

The first model is the Naïve Bayes classifier. The Gaussian Naïve Bayes algorithm is used for the binary classification in this manner:

$$P(y \mid x) = \frac{P(y)P(x \mid y)}{P(x)} = \frac{P(y)\prod_{i=1}^{N} P(x_i \mid y)}{P(x)} \propto$$

$$\propto P(y)\prod_{i=1}^{N} P(x_i \mid y)$$
(2.1)

given that
$$P(x_i | y) = \frac{1}{\sqrt{2\pi\sigma_y^2}} \exp(-\frac{(x_i - \mu_y)^2}{2\sigma_y^2}),$$

where y and x are the target, which is churn, and the features, respectively, with the mean and the variance of y being obtained from maximum likelihood estimation [Scikit-learn, 1.9.1].

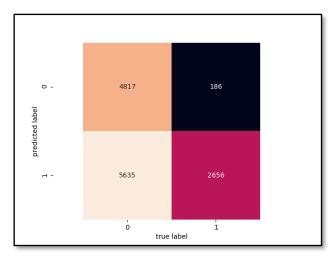


Figure 2.1: Confusion matrix for Naïve Bayes classifier.

It is among the simplest and fastest models for classification, using a simple formula with a naïve assumption of no covariance between the variables. It sometimes even outperforms the more complicated models, especially when the dataset is relatively small. Here, the calculated accuracy score is 56.21%, which is quite low but better than random, and AUROC is 69.77%, which is higher due to better accuracy for the underrepresented group, as seen in the confusion matrix represented in Figure 2.1.

The second model is Logistic Regression, which is classical for this type of problem. It uses a sigmoid-shaped activation function to produce binary output, either zero or one. The model is as follows:

$$\log\left(\frac{P(y\mid x)}{1 - P(y\mid x)}\right) = w_0 + w_1 x_1 + w_2 x_2 + \dots + w_N x_N, \quad (2.2)$$

which I choose to solve with Stochastic Average Gradient descent (SAG), since it is faster for large datasets [Scikit-learn, 1.1.11]. The accuracy score is 86.45% and the AUROC score is 78.18%. Its confusion matrix is summarized in Figure 2.2 below.

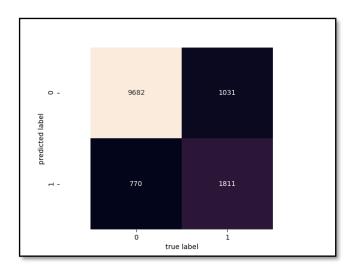


Figure 2.2: Confusion matrix for Logistic Regression classifier.

The next model is K Nearest Neighbors (KNN) classifier. It is a nonparametric method, using only the distances and closest neighbors to decide on prediction [Scikitlearn, 1.6.2]. The number of neighbors is chosen to be 50, although more neighbors would improve the result; however, the change is not so significant, as tests show. The observed accuracy score is 87.33% and AUROC is 79.6%. Figure 2.3 demonstrates the confusion matrix.

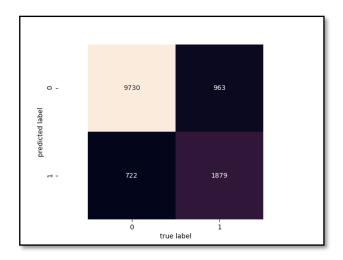


Figure 2.3: Confusion matrix for KNN classifier.

The fourth model is Random Forest classifier, which is an ensemble of decision trees. A decision tree left alone is a very weak nonparametric model; however, as a combination of many trees using the boosting aggregation, it is a powerful tool. The algorithm takes the average of probabilistic prediction made by each decision tree in a sample with replacement to find the classes [Scikit-learn, 1.11.2.1]. Similarly, the number of estimators is chosen as 50 and not more because the improvement is insignificant and time consuming, as tests demonstrate. The accuracy and AUROC scores are 86.93% and 78.43%, respectively. The confusion matrix is given in Figure 2.4 below.

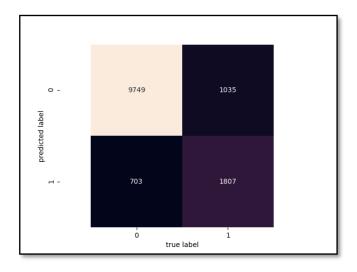


Figure 2.4: Confusion matrix for Random Forest classifier.

The fifth model is Support Vector Machines (SVM) classifier. It is a strong model, based on several optimizations, which require some time for computations. The algorithm involves primal and dual optimization problems with the following specifications:

* Primal:

$$\min_{w,b,\varsigma} \frac{1}{2} w^{T} w + C \sum_{i=1}^{n} \varsigma_{i}$$
subject to $y_{i}(w^{T} \phi(x_{i}) + b) \ge 1 - \varsigma_{i}$,
 $\varsigma_{i} \ge 0$, $i = 1, ..., n$
* Dual:

$$\min_{\alpha} \frac{1}{2} \alpha^{T} Q \alpha - e^{T} \alpha$$
subject to $y^{T} \alpha = 0$, $0 \le \alpha_{i} \le C$, $i = 1, ..., n$
[$Q_{i,i} = y_{i} y_{i} K(x_{i}, x_{i}) & K(x_{i}, x_{i}) = \phi(x_{i})^{T} \phi(x_{i})$],

where e is a matrix of ones and $K(x_i, x_j)$ is a kernel function, and unlike other models, the classes are predicted by the sign of the decision function [Scikit-learn, 1.4.1]. The

value for C is chosen as 10 through available grid search cross validation function. The resultant accuracy score is 86.70% and the AUROC is 77.87%. Its confusion matrix is

summarized in Figure 2.5.

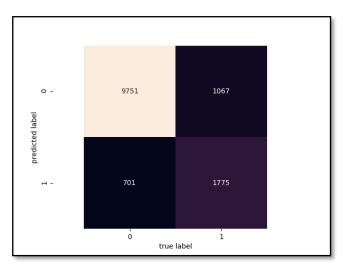


Figure 2.5: Confusion matrix for SVM classifier.

The following model is Neural Networks classifier. It is an interesting model which works as a black box but it also takes some time to compute the weights for each

layer to produce the result. The multilayer perceptron trained with backpropagation is used in this algorithm [Scikit-learn, 1.17.2]. The parameters are tuned manually, choosing the two layer representation with 10 nodes each. Other parameters have lower scores. The accuracy score and AUROC are 87.21% and 78.97%, accordingly. The confusion matrix can be observed in Figure 2.6 below.

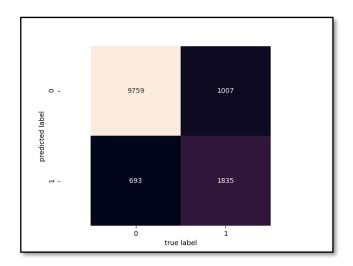


Figure 2.6: Confusion matrix for Neural Networks classifier.

The last model is Extreme Gradient Booster (XGBoost) classifier. It is also an ensemble learning model with boosting and regularization [Chen & Guestrin, 2016]. It has been the winner in many programming contests. The number of estimators is tuned as 13 with the grid search cross validation function. Its accuracy score is 87.73% and AUROC score is 81.07%. Its confusion matrix is given in Figure 2.7.

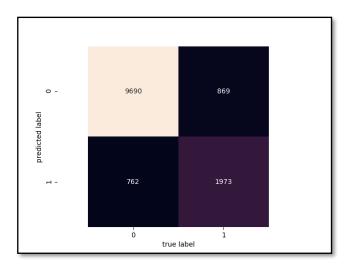


Figure 2.7: Confusion matrix for XGBoost classifier.

To conclude, among all the used models, both the highest accuracy and the highest area under the ROC curve (AUROC) belong to XGBoost classifier, although it is slightly better than its alternatives. To summarize the comparison even better, the ROC curve is demonstrated in Figure 2.8 below.

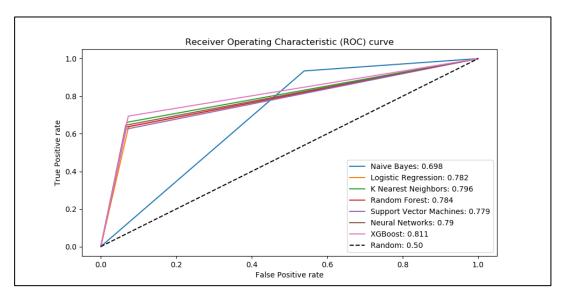


Figure 2.8: ROC curve for binary classification models of churn estimation.

As seen in the graph, there are two peaks where ROC curves concentrate. On the left, the dominant model is XGBoost, followed by KNN, which is followed by the Neural Networks model. The peak on the right represents the Naïve Bayes classifier, which is best in predicting the underrepresented group of exited clients, although its accuracy score in total is the lowest among the models used. All other accuracy scores are greater than 85%, which is far better than guessing all as retained clients (79.1% of accuracy) and this is a significant improvement.

Chapter 3:

HOUSING PRICE PREDICTION

Predicting various prices is another ubiquitous application of statistics and data science in the private sector, since knowing future prices is quite essential for profitability. One of the most famous business cases is to forecast housing prices for potential valuation and investment purposes. Other than in the housing industry itself, real estate investment can be made by miscellaneous companies and institutions as a way to invest their funds. Being considered by managers and investors, real estate is a worthy asset group, which is an ideal hedge against risks and uncertainty [Pinnington, 2020]. In general, housing price prediction constitutes a specific machine learning model of regression with a target variable of the real estate price and independent variables of the related parameters to a housing unit.

The data consist of apartment prices in Moscow, provided by Higher School of Economics and shared on Kaggle in 2018². The dataset has 2040 entries with apartment prices and few related variables. The prices are reported in thousands of US dollars. The independent variables comprise the total space of apartment, the living and kitchen spaces, the distances to the city center and nearest metro station, the categorization of apartments from 1 to 8 depending on observations of data collectors, and several binary variables for the proximity to metro station (within walking distance vs. requiring transportation), the type of building material (brick vs. monolith houses), and the floor indicator (first and last floor vs. the rest). The variables also include the numbering id, which is dropped before modeling since it is useless in predictions.

As a next step before the application of different machine learning algorithms, the data are normalized using the min-max scaler, normalizing by only keeping the relative distances between the data points, since normalization allows some algorithms to approach the solution more efficiently and faster among other benefits. Furthermore, after deciding the best parameters for each model in a grid search of 5-fold cross validation, minimizing the Root Mean Square Error (RMSE), 20% of the data are used for model testing and the other 80% are for model training purposes. The Root Mean Square Error (RMSE) is a popular metric for regression in the literature [Chai & Draxler, 2014]. The

² Can be retrieved from www.kaggle.com/hugoncosta/price-of-flats-in-moscow

bar plots of the RMSE and the correlation coefficient between predicted and actual prices for all models are summarized at the end to compare and decide on the best algorithm.

The first model is Linear Regression, which is a traditional technique for regression problems. It uses the Ordinary Least Squares (OLS) method to minimize the sum of squared errors, i.e. differences between the actual and predicted target variables. The model is as follows:

$$\hat{y}(w,x) = w_0 + w_1 x_1 + w_2 x_2 + \dots + w_N x_N, \tag{3.1}$$

where \hat{y} is the predicted target value, w_0 is the intercept term, and $w = (w_1, w_2, ... w_N)$ is the vector of coefficients [Scikit-learn, 1.1.1]. The RMSE error is 29.41 and the correlation coefficient is 82.29%.

The second model is the Gaussian Process regressor. This method utilizes Gaussian probabilistic processes for regression purposes. In this algorithm, the prior parameters and the noise level, alpha parameter, are the input for the optimization framework. The prior mean is expected to be constant and the prior covariance depends on a kernel to be chosen, such as basic kernel, radial-basis function (RBF) kernel, rational quadratic kernel, and dot-product kernel [Scikit-learn, 1.7.1]. As a result of manual tuning, the combination of rational quadratic kernel and alpha parameter of 1.7 minimizes the error term. This is the rational quadratic kernel:

$$K(x_i, x_j) = (1 + \frac{d(x_i, x_j)^2}{2\alpha l^2})^{-\alpha},$$
 (3.2)

where l>0 and $\alpha>0$ are length-scale and scale mixture parameters, respectively, and d() is a Euclidean distance function [Scikit-learn, 1.7.5]. Consecutively, the calculated RMSE is 27.77 and correlation is 84.45%.

The third model is K Nearest Neighbors regressor. Such nonparametric methods take the average of the closest neighbors by analyzing the distances [Scikit-learn, 1.6.3]. Although usually uniform weights are applied for the average, in this task the inverse distance is used to allow closer neighbors to contribute more to the weights, slightly improving the algorithm performance. The number of neighbors is chosen to be 8, since it minimizes the RMSE score. Although the results are not very promising, the observed RMSE is 29.75 and the correlation coefficient is 82.33%.

The following model is Random Forest regressor, which is an ensemble of decision trees, predicting the target by taking the average of each leaf node. A combination of numerous decision trees using the boosting aggregation technique converts a single decision tree, a weak nonparametric model, into a powerful regressor. In a sample with replacement, the algorithm takes the average of target predictions made by each decision tree to make the final prediction [Scikit-learn, 1.11.2.1]. The number of estimators is chosen to be 115, since it minimizes the error term. The RMSE score and correlation are 26.18 and 86.42%, accordingly.

The fifth model is Support Vector Machines regression. Basing on framework of computed optimizations, the model involves primal and dual optimization steps with such specifications:

$$\min_{w,b,\zeta,\zeta^*} \frac{1}{2} w^T w + C \sum_{i=1}^n (\zeta_i + \zeta_i^*)$$
subject to $y_i - w^T \phi(x_i) - b \le \varepsilon + \zeta_i$,
$$w^T \phi(x_i) + b - y_i \le \varepsilon + \zeta_i^*,$$

$$\zeta_i, \zeta_i^* \ge 0, \qquad i = 1, ..., n$$
(3.3)

* Dual:

$$\min_{\alpha,\alpha^*} \frac{1}{2} (\alpha - \alpha^*)^T Q(\alpha - \alpha^*) + \varepsilon e^T (\alpha + \alpha^*) - y^T (\alpha - \alpha^*)$$

$$subject \ to \quad e^T (\alpha - \alpha^*) = 0,$$

$$0 \le \alpha_i, \alpha_i^* \le C, \qquad i = 1, ..., n$$

$$[Q_{ij} = K(x_i, x_j) = \phi(x_i)^T \phi(x_j)],$$

where e is a matrix of ones and $K(x_i, x_j)$ is a kernel function [Scikit-learn, 1.4.2]. The C coefficient is chosen to be 10000 within the grid search cross validation function; greater values of C do not improve the performance significantly. The resultant RMSE score is 25.29 and the correlation coefficient is 87.25%.

The next model is Neural Networks regressor. Being a black box model, it can achieve great results without detailed interpretations. Here, the multilayer perceptron is trained using backpropagation with no activation function at the output level to have the regression format in predictions [Scikit-learn, 1.17.3]. The parameters are tuned manually, choosing the neural network of five layers with 150 nodes each. Other parameters have higher error terms. The RMSE error and correlation are 25.74 and 86.88%, respectively.

The last model is Extreme Gradient Booster (XGBoost) regressor, a popular technique. It is a powerful ensemble algorithm with regularized boosting of regression trees [Chen & Guestrin, 2016]. The number of estimators is tuned as 21 with the help of grid search cross validation function. Its RMSE score is 26.09 and correlation coefficient is 86.30%. The comparison of models can be clearly seen within the bar plots summarized in Figures 3.1-3.2 below.

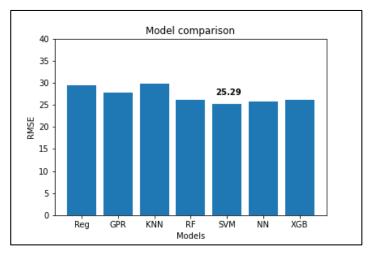


Figure 3.1: RMSE scores across regression models of housing price prediction.

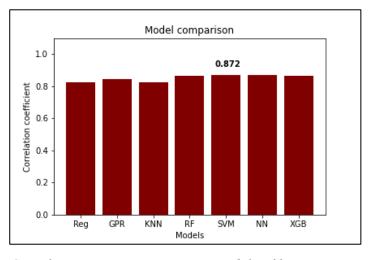


Figure 3.2: Correlations across regression models of housing price prediction.

In sum, among all the applied algorithmic models, both the lowest error term (RMSE score) and the best linear relationship between predictions and actual prices (correlation coefficient) belong to SVM regressor, although it is slightly better than its alternatives.

Sentiment Analysis 14

Chapter 4:

SENTIMENT ANALYSIS

Another notable application of machine learning methodology in business is sentiment analysis technique, also known as opinion mining. It is a type of natural language processing (NLP) algorithms, commonly constituting text analysis. Given a piece of writing, usually voice of the customer material such as a comment or a review, the algorithm aims to understand its overall tone and classify it according to some acceptable measure. For instance, an opinion could be classified as positive, negative, or neutral. Hence, it is usually a classification task with various number of targeted classes and preprocessed text as an input, obtained through the feature extraction process. In practice, the sentiment analysis is deployed in a range of aspects in business, such as customer support and service, market research, public relations, and human resources management in numerous industries [Puschmann & Powell, 2018].

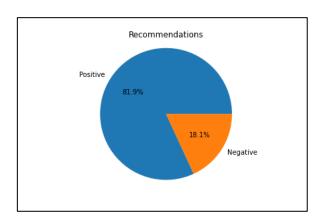


Figure 4.1: Distribution of customer recommendations.

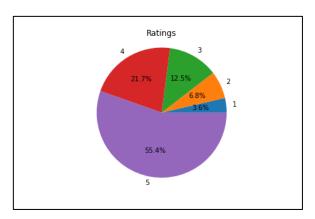


Figure 4.2: Distribution of customer ratings.

The dataset is the collection of women's e-commerce clothing reviews from a real but anonymous source, shared on Kaggle in 2018³. The data initially consisted of 23486 non-null entries depicting the customer reviews with related information, including product and customer descriptions. Other columns than text reviews, customer recommendations and rankings were dropped to concentrate on a sole sentiment analysis task. The text reviews column is the main variable for the learning model. The two others are possible classification targets with the customer recommendations, being either yes or no in a binary representation, and the rankings on a scale from 1 to 5. Since text reviews were omitted in some entries, they were dropped, resulting in a total of 22641 rows. The distribution for recommendations and rankings are represented as pie charts in Figures 4.1-4.2 above. Only 18.1% of all reviews have negative recommendations. Similarly, the majority of reviews have maximum ranking with 55.4% of proportion. In a consistent manner, each of next ranking level has lesser proportion than the previous. Since there are more classes, a multiclass classification task with a target of rankings is expected to be less successful in predictions than a binary classification task with a target of recommendations, in a general algorithmic setup.

First of all, the text reviews have to be preprocessed. Thus, the text reviews are filtered from numbers and special characters, since most of them are irrelevant and do not possess any important information. The cases of remaining alphabetical characters are lowered to enforce uniformity. Taking out the stopwords from the current data, the tag clouds are produced for both positive and negative recommendation labels. Both clouds are demonstrated in Figures 4.3-4.4 below.



Figure 4.3: Tag cloud of text reviews with positive recommendation.

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³ Can be retrieved from www.kaggle.com/nicapotato/womens-ecommerce-clothing-reviews



Figure 4.4: Tag cloud of text reviews with negative recommendation.

The stopwords, such as prepositions and conjunctions, are omitted because they are usually meaningless in such an analysis. In both tag clouds, although there are many common words related to the e-commerce of women's clothing in general, some words related to either positive or negative reaction can also be observed. As a next step, the text reviews with omitted stopwords are tokenized into bag-of-words (BoW) representation, which is a simplified matrix with the word counts. Then, the BoW representation is converted to term frequency – inverse document frequency (TF-IDF) measure, which is quite popular for text analysis. This measure reflects the relevance of words to a document in a collection of documents, i.e. corpus [Qaiser & Ali, 2018]. The unigram version of BoW and TF-IDF measures are compared, where unigram is a single word representation from the original text. Alternatively, their bigram version, which is a representation as pairs of consecutive words, being also considered, is not included in the analysis due to low effectiveness in comparison.

The same list of classification models, described in chapter 1, is utilized for binary classifications of sentiment analysis. The most successful ones are also used for the multiclass classification task. Similarly, AUROC measure and accuracy score are employed for the binary classification. For the multiclass classification, however, another measure has to be used in addition to accuracy and confusion matrix, since there is no well-designed ROC curve for this case. One common measure is Macro F1 score, a macro averaged F1 measure of each class, which combines both popular precision and recall measures at once [Grandini, Bagli, & Visani, 2020]. The precision measures how good the model is in predicting a particular class and the recall measures how well the particular class is predicted overall [Grandini, Bagli, & Visani, 2020]. The F1 score, which is a

tradeoff of both, geometrically averaged for each class, is a good substitute for the AUROC measure. The macro average, rather than the weighted one, is used to deem each class equally; otherwise, guessing the most probable class at random would increase the score, as it does for accuracy measure. Having decided the best parameters for each model, 80% of the data are employed for model training and the remaining 20% are for model testing purposes. The grid search of 5-fold cross validation is used to decide for the best model parameters, where it is required.

All seven models are applied for binary classification task of sentiment analysis. Firstly, the Naïve Bayes classifier [Scikit-learn, 1.9.1] and Logistic regression [Scikit-learn, 1.1.11] models do not have any decision variables, as before. The number of neighbors is chosen to be 50 for K Nearest Neighbors classifier [Scikit-learn, 1.6.2]. Moreover, the number of estimators for Random Forest classifier [Scikit-learn, 1.11.2.1] is chosen as 100; more estimators do not improve the results significantly. Then, the Support Vector Machines classifier [Scikit-learn, 1.4.1] maximizes the AUROC score at the C value of 100 for the BoW version and C of 10000 for the TF-IDF one. Being tuned manually, the Neural Networks classifier [Scikit-learn, 1.17.2] is decided to have six layer representation with 5 nodes each for BoW and four layers with 7 nodes per each for TF-IDF. Lastly, the Extreme Gradient Booster classifier [Chen & Guestrin, 2016] is chosen to have 1000 estimators for both cases. The outputs of the binary classification modeling are summarized in Table 4.1 below.

Table 4.1: Comparison of binary classification models of sentiment analysis.

Model	BoW		TF-IDF	
Wiodel	accuracy	AUROC	accuracy	AUROC
Naïve Bayes	45.9%	55.15%	45.93%	55.03%
Logistic Regression	88.54%	78.29%	88.23%	74.38%
K Nearest Neighbors	80.95%	51.49%	82.69%	56.13%
Random Forest	82.65%	55.66%	82.73%	55.89%
Support Vector Machines	87.99%	74.88%	88.7%	77.87%
Neural Networks	87.46%	82.48%	85.45%	80.8%
XGBoost	88.87%	77.72%	87.88%	77.33%

As seen in the table, the outputs for BoW and TF-IDF input versions are more or less similar, especially for the accuracy scores. In the case of Logistic Regression and Neural Networks, the AUROC score is significantly better for BoW. On the other hand, KNN and SVM models perform better for TF-IDF. The ROC curves for both BoW and TF-IDF versions are also demonstrated in Figures 4.5-4.6.

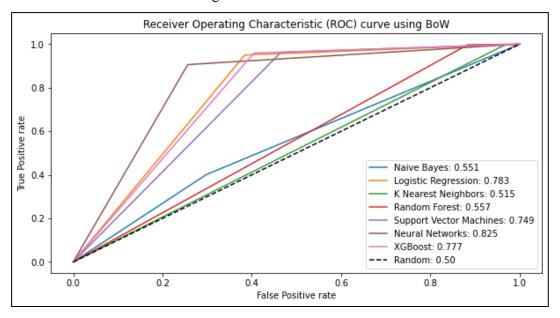


Figure 4.5: ROC curve for binary classification models using BoW.

As a result, despite the indication of higher accuracy scores for some models, which is less important, according to AUROC measures, the best performing model for both BoW and TF-IDF inputs is Neural Network classifier.

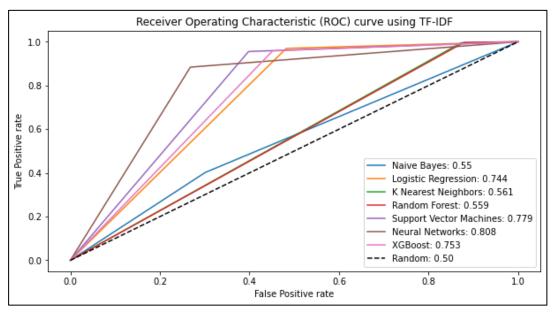


Figure 4.6: ROC curve for binary classification models using TF-IDF.

The classifiers such as Naïve Bayes, Random Forest, and K Nearest Neighbors, show poor performance for this type of problem with AUROC score close to random guessing and, therefore, are omitted in the multiclass classification case.

Then, four best models are applied for multiclass classification task of sentiment analysis. The Logistic regression [Scikit-learn, 1.1.11] classifier is used without any decision variable. Secondly, the SVM classifier [Scikit-learn, 1.4.1] maximizes the Macro F1 score at C value of 1000 for the BoW case and C of 10000 for TF-IDF. With a manual tuning, the Neural Networks classifier [Scikit-learn, 1.17.2] is chosen to have one layer representation with 100 nodes for both cases. Lastly, the XGBoost classifier [Chen & Guestrin, 2016] is decided to have 150 estimators for both BoW and TF-IDF input versions. Table 4.2 summarizes the resultant accuracy and macro averaged F1 scores.

Model	BoW		TF-IDF	
Wiodei	accuracy	Macro F1 score	accuracy	Macro F1 score
Logistic Regression	61.54%	0.42	63.66%	0.39
Support Vector Machines	62.4%	0.42	63.52%	0.42
Neural Networks	58.8%	0.40	57.52%	0.39
XGBoost	62.35%	0.40	61.87%	0.38

Table 4.2: Comparison of multiclass classification models of sentiment analysis.

As expected, the resultant scores are much lower than in the binary classification case, since there are five different classes. With a priority given to Macro F1 score, the best performing model here is Support Vector Machines classifier. Its confusion matrices for two versions are depicted in Figures 4.7-4.8 below.

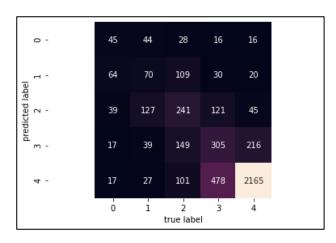


Figure 4.7: Confusion matrix for SVM classifier using BoW.

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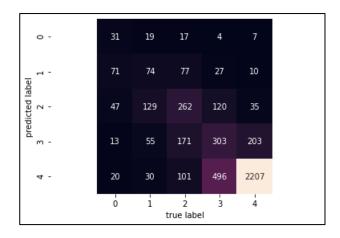


Figure 4.8: Confusion matrix for SVM classifier using TF-IDF.

In general, the BoW input representation has slightly better performance for multiclass classification, although in the case of SVM classifier, both BoW and TF-IDF have equal macro averaged F1 scores. Thus, the sentiment analysis problem with classifications of different forms has multiple aspects, showing a range of machine learning tools and methods.

Conclusion 21

Chapter 5:

CONCLUSION

To conclude, the machine learning methodology of three distinct business applications is analyzed and compared in the paper. The first chapter applies the binary classification models for the churn estimation task. The second chapter deploys the regression analysis techniques to predict the housing prices. The third chapter incorporates the sentiment analysis tools along with models of binary and multiclass classification for e-commerce reviews and ratings. Hence, the diverse range of machine learning and data science methodology is very beneficial to be applied in business.

Even though the lack of a solid theoretical basis and a certain dependence of results on the specific datasets employed are possible drawbacks of this research, it presents a practical application of common machine learning algorithms, demonstrating their relative strengths and weaknesses in particular types of problem. Optimizing the feature selection method and adding more complex techniques would potentially improve the results. Future works in this direction are very promising and inspiring.

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