

# ML Final Project

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## 1 Introduction

Kaggle, a data modeling, and data analysis competition platform, holds competition of data science for statisticians' and data mining experts. Participants experimented with different methods, competing against each other for the prize to the best model. [1]

Though Kaggle competitions are incredibly fun and rewarding, they are intimidating for novices like me in data science. [2] Therefore, instead of participating in a Kaggle competition and losing my enthusiasm for machine learning, I joined a playground competition. Tabular Playground Series organizes a playground competition once a month, for my final project, I choose to participate in the one held in March 2021.

The submissions of the competition are evaluated on the area under the ROC curve between the predicted probability and the observed target. There will be 30,000 sets of data for training and 20,000 sets for testing. For each set of data, there are 30 kinds of features, including bool, float, and string.

Since there is no limitation on choosing a model in the competition, I tried models I've learned this semester, including

clustering, linear discrimination, and artificial neural network. The reason why I didn't try nonparametric methods was that after trying clustering, I find it hard to train models efficiently by letting data speak for themselves since the dataset is massive.

## 2 Motivation

Being a student studying biomedical engineering, I considered it is important for me to learn machine learning well on technical level and medical level, that is, I should be able to write a good model while being aware of what kind of clinical problems might happen. Before learning machine learning from the aspect of biomedical engineering, I would like to approve on my technical level. As result, I took part in a playground competition in Kaggle as my final project, hoping I could learn machine learning better by training a synthetic, beginner-friendly dataset (but based on a real dataset and generated using a CTGAN).

## 3 Try and Errors

### 3.1 Preprocessing

In this dataset, the features are composed of different data types, including bool, float, and string (they are categorical).

However, we usually prefer the data type of the features is numeric. Therefore, I tried two methods, respectively to be label encoder and one hot. The former target labels with values between 0 and  $n\_classes - 1$ , transforms non-numerical labels (as long as they are hashable and comparable) to numerical labels. [3] The latter creates a binary column for each category and returns a sparse matrix. [4] After transforming all the non-numerical features into numerical features, I used linear regression to test which method is better. The results show that using one hot could get a lower loss and higher accuracy. Thus, I choose one hot to transform the non-numerical features. After transforming the number of features increased from 30 to 634. Then, I randomly split the dataset into 240,000 training data and 60,000 validation data, and combine every 30,000 data into a loader.

Metrics	Training	Validation
Number of data	240,000	60,000
Loader size	8	2

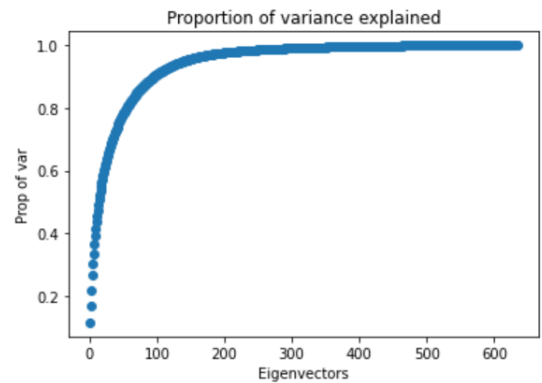
Table 1: Data Distribution

### 3.2 Dimension reduction

To clarify whether using feature selection or feature extraction is better, I plot the histogram of the distribution of each feature and find out that they all look similar. Since they all seem to have a similar distribution, I assumed that each feature attribute is almost equal to the model. Therefore, I choose feature extraction.

At first, I tried PCA and choose to use

100 components because the elbow of the proportion of variance explained approximately happens at 100. After project data on the eigenvectors I choose, I use linear regression to check if I need to select more components or less. However, no matter how I change the numbers of selected components, the result doesn't change much. Later, it occurred to me that maybe I do not necessarily need to do dimension reduction since the ratio of features number and data is  $634/240,000$ , equal to 0.0026, which is lesser than 1%. Therefore, I decided not to do dimension reduction, unless overfitting happens.



Graph 1: Proportion of variance explained

### 3.3 Selecting Model

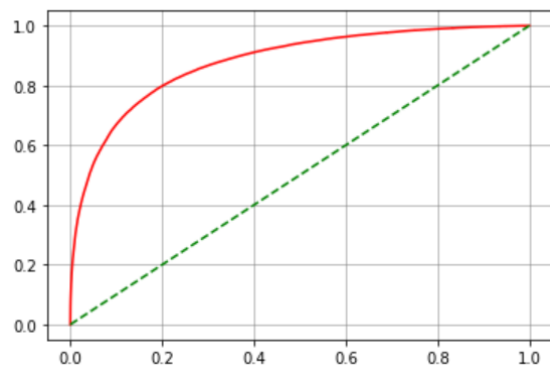
*Linear discrimination:* I tried linear regression and logistic regression.

For linear regression, I simply solve the linear equation without using loss function or optimizer to adjust the model. The training area under the ROC curve (simply AUC) is 0.8814, while the validation AUC is 0.8799. With this result, I was having confidence that the AUC would at least reach 0.9 if loss

function and gradient descent is applied.

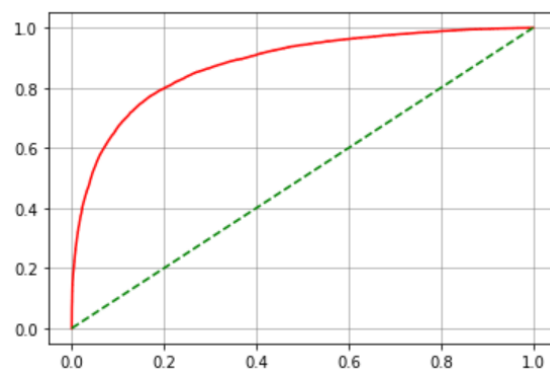
For logistic regression, I used the module in sklearn to build a model. The result was fine, too. The training AUC is 0.8791 and the validation AUC is 0.8794.

```
Accuracy 0.842
Sensitivity= 0.93
Specificity= 0.599
Precision= 0.865
Negative Predictive Value= 0.754
AUC: 0.8791
```



*Graph 2: performance of training data using logistic regression*

```
Accuracy 0.843
Sensitivity= 0.93
Specificity= 0.601
Precision= 0.867
Negative Predictive Value= 0.753
AUC: 0.8794
```

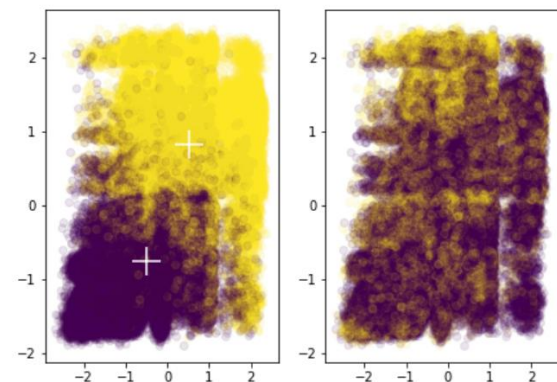


*Graph 3: performance of validation data using logistic regression*

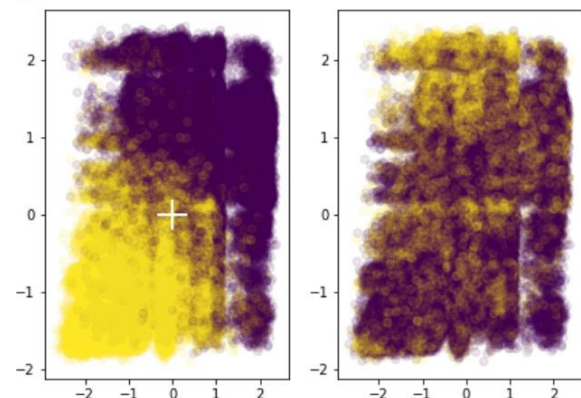
**Clustering:** I tried K-means, fuzzy C-Means, and Spectral clustering.

For K-means, I used the module in sklearn to build a model. The accuracy for training data ranges from 0.5 to 0.8. I assumed that it is because the first two centers are chosen randomly, thus the accuracy flows.

For the fuzzy C-mean, I used the module in fcmeans to build the model. The accuracy for training data ranges from 0 to 0.5. To figure out what happened, I plot the scatter plot to see the distribution of the data.



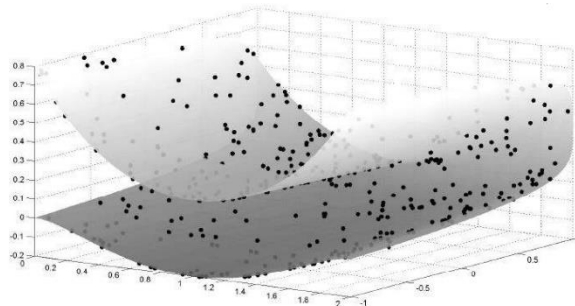
*Graph 4, 5: K-means centers and data scatter plot with prediction of k-means (4: left) and data scatter plot with ground truth (5: right)*



*Graph 6, 7: fuzzy C-means data scatter plot with prediction of fuzzy C-means (6: left) and data scatter plot with ground truth (7: right)*

Graph 4 and graph 6 change in every run. In graph 5 (same as 7), the data with different

labels mix together, therefore, I conclude that this dataset is not suitable for clustering. However, the plot reminded me of the graph down below:



*Picture 1: the 3-D from slide clustering\_v2*

I thought, maybe the data seems close in 2-D but actually far in 3-D. Therefore, I wanted to try spectral clustering. I used a module in sklearn to build a model. However, it occupied too many RAM resources on Colab and was forced to shut down every time I tried to run the code. I tried to consume the number of features and data, but it still wouldn't work. Thus, I gave up on this option.

Moreover, after seeing graph 5, I assume that the non-parametric methods such as K-NN wouldn't work well as well. In addition, there are 300,000 data, which would take lots of RAM resources to run a non-parametric model since every data needs to speak for itself. Therefore, I decided not to try non-parametric method.

After I get the conclusion above, it occurred to me that the predictions of the clustering are not probabilities. Therefore, AUC couldn't be evaluated.

*Artificial Neural Network:* I tried DNN.

I took the code from homework 3 and

tried 2 hidden layers, 3 hidden layers, different nodes number, and different learning rates. However, the accuracy couldn't be higher than 60% no matter how hard I tried.

*Comparison:* Among all kinds of models, logistic regression seems to have the best performance. To tune the parameters, I decided to build a model myself instead of using the module.

## 4 Experiments

### 4.1 Building model

Before building a logistic regression model, I added an all-ones column as a feature, which acts as the constant in the linear problem. In formulas down below,  $y$  is ground truth,  $x$  is the dataset, and  $w$  is weight in the linear model.

*Prediction:* The function for prediction is the transpose matrix of  $w$  dot transpose matrix of  $x$  plus the constant  $w^0$ , and then put the result into sigmoid function.

$$\text{prediction } y^t = \text{sigmoid}(w^T x^T + w^0)$$

*Loss function:* I tried the Lasso function, MSE function, and cross-entropy. After testing three kinds of function with validation data, cross-entropy had the best performance. Therefore, I choose cross-entropy as my loss function.

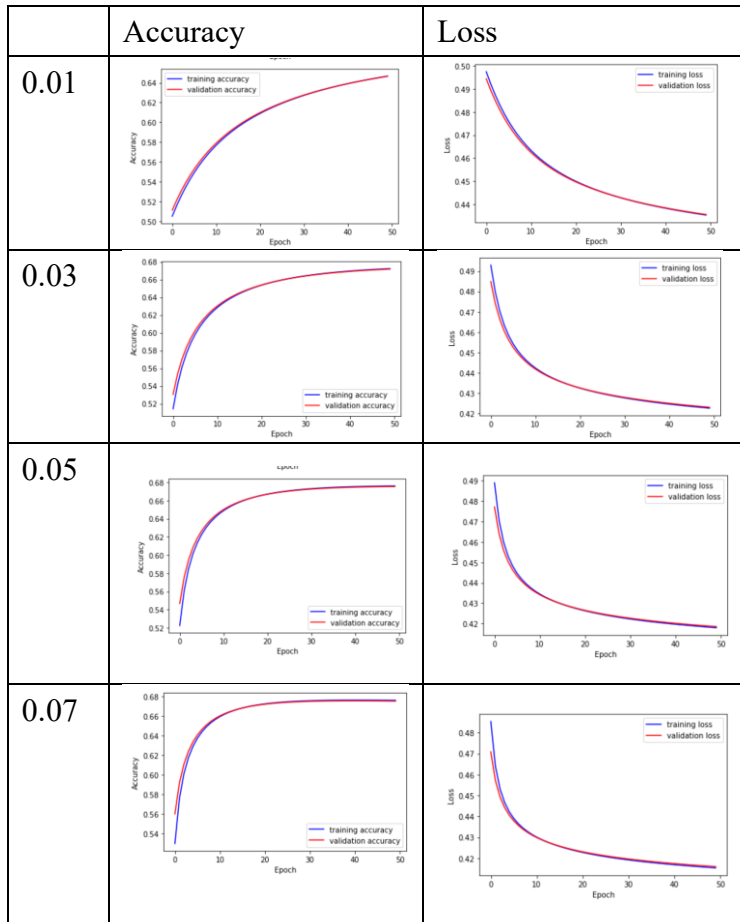
$$\text{loss} = \frac{1}{N} \sum_{n=1}^N \ln(1 + \exp(-y_n w^T x_n))$$

*Gradient:* I do the gradient decent by trying to minimize the gradient written down below. The weight matrix will then minus learning rate  $\times$  gradient to carry out gradient descent.

$$gradient = \frac{1}{N} \sum_{n=1}^N \partial(-y_n w^T x_n)(-y_n x_n)$$

## 4.2 Experiment 1

I tested different learning rates with 0.01, 0.03, 0.05, 0.07, for each learning rate, I set the epoch to be 50, and the result is shown in Table 2. As Table 2 shows, the lower the learning rate is, the more gently the curves are, and eventually converge to a horizontal



line.

Table 2: performances for different learning rates

The accuracy for different learning rate with 0.01, 0.03, 0.05, 0.07 is 65%, 67%, 68%, 68% (for both training and validation data) respectively. I would choose 0.05 as the final learning rate since the slope of the curve looks small enough and won't reach the highest too fast. However, the slope of the loss curve for neither 0.05 nor 0.07 isn't small enough. Therefore, I would like to try running a bigger epoch.

## 4.3 Experiment 2

After experiment 1, the learning rate was fixed, and the next step is to see if the loss could be lower, and the result is shown in Table 3.

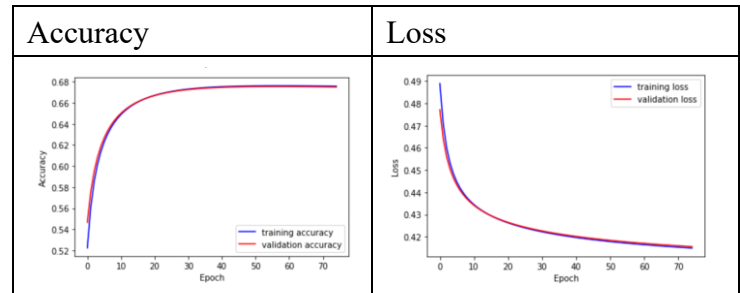


Table 3: performances when there are 75 epochs

As the result shown in graph 3, the slope of the loss curve approaches zero, and the accuracy remains the same.

## 5 Result

### 5.1 Dataset

Since the features in the dataset need to be numerical, I used one hot method to transform non-numerical features into numerical. However, after transforming, the feature number for the training dataset is not equal to the one for the testing dataset. The

testing dataset is four feature lesser, thus, I added four all-ones columns as constant in logistic regression to make sure the size of the training dataset equals to the testing dataset.

	Train	Test
Number of data	300,000	200,000
Number of features	634	634

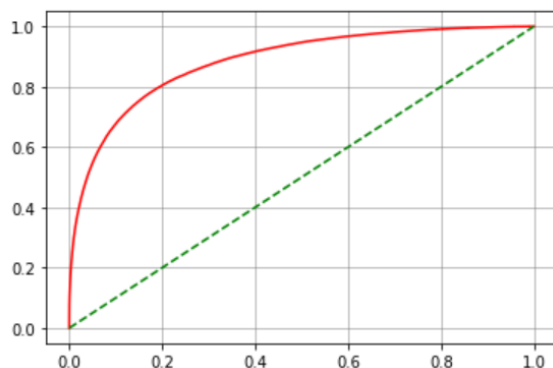
## 5.2 Training Model

In the experiment, it is obvious to see that the logistic regression I built myself doesn't work better than the one in sklearn no matter in result or time complexity, so I use the module to get the result for the competition. The penalty used in the model is l2 [5], and since the module's default is to add constant automatically, I didn't add an extra column.

## 5.3 Performance

The AUC of the testing dataset is 0.8841, and the performance is shown down below.

Accuracy 0.845  
Sensitivity= 0.931  
Specificity= 0.605  
Precision= 0.868  
Negative Predictive Value= 0.761  
AUC: 0.8841



The score of the testing data is 0.73815. Score  
0.73815

## 6 Conclusion and Discussion

There is still some difference between my model and sklearn, which I believe is the reason why the area under the ROC curve my model evaluates is lower.

*Loss function:* The module only uses cross-entropy as the loss function in multiclassification while I use it in a 2 classes problem.

*Regularization:* The module uses the l2 regularization to make sure the model won't be too complicated while I didn't. However, since there is no sign of overfitting, I believe this is not the reason for bad performance.

There are still other parameters in the module that I'm not familiar with. There is still a big process to make and knowledge to learn.

## 8 Reference

- [1] Kaggle wiki <https://en.wikipedia.org/wiki/Kaggle>
- [2] Tabular Playground Series - Mar 2021  
<https://www.kaggle.com/c/tabular-playground-series-mar-2021/overview>
- [3] Sklearn web for encoder <https://scikit-learn.org/stable/modules/generated/sklearn.preprocessing.LabelEncoder.html>
- [4] Sklearn web for one\_hot <https://scikit-learn.org/stable/modules/generated/sklearn.preprocessing.OneHotEncoder.html>
- [5] Sklearn web for logistic regression  
[https://scikit-learn.org/stable/modules/generated/sklearn.linear\\_model.LogisticRegression.html](https://scikit-learn.org/stable/modules/generated/sklearn.linear_model.LogisticRegression.html)