# $\mathbf{PRML}$

Assignement - 3 Report

by

P Yashwanth Sai(cs22b002)

Dr Arun Rajkumar



Computer Science and Engineering Indian Institute of Technology Madras  ${\bf Jan\ -\ May\ 2024}$ 

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## 1 Datasets

### 1.1 emails.csv

The following dataset is used from kaggle {emails.csv}. The emails were pre-processed by removing the numbers and punctuation marks. The final text in the email contains only words/alphabets in lowercase and whitespaces.

## 1.2 Preprocessing

The resulting data-frame was pre-processed using sklearn.feature\_extraction.text

### 1.2.1 CountVectorizer(binary = True, stop\_words = 'english')

- The dataset has been converted into a DataFrame where each column header represents a unique word/feature present in the dataset.
- Each row represents a data point or an email. The values are binary, either 0 or 1, indicating whether the corresponding word is present in the email (1) or not (0).
- stop\_words are the common words in english that have no impact on our training methods an hence are removed to decrease the number of features/dimensions.
- This DataFrame will be represented as binary\_data\_array or something similar further in the report.

### 1.2.2 CountVectorizer(binary = False, stop\_words = 'english')

- The dataset has been converted into a DataFrame where each column header represents a unique word/feature present in the dataset
- Each row represents a data point/email the values are the frequencies of how many times that word has occurred in the mail.
- This DataFrame will be represented as frequency\_data\_array or something similar further in the report.

#### 1.2.3 TfidfVectorizer(stop\_words = 'english')

- The dataset has been converted into a DataFrame where each column header represents a unique word/feature present in the dataset
- Each value in the row represents the value as calculated by the tfidf transformer over the dataset.
- This DataFrame will be represented as tfidf\_data\_array or something similar further in the report.

#### 1.3 Notations

- X is the data matrix of dimensions  $n \times d$ . It can obtained from rawdata by using any of the pre-processing methods 1.2.1 , 1.2.2 , 1.2.3 whichever is applicable in the usage context.
- Y is the labels matrix of the data. It is of dimensions  $n \times 1$

## 2 Naive Bayes Classifier

## 2.1 Training

- The maximum likelihood estimators for the parameters of naive bayes algo yields the following equations
- binary\_data\_array is used for Naive Bayes, Bayesian distribution is used.

$$\hat{p} = \frac{1}{n} \sum_{i=1}^{n} y_i$$

$$\hat{p}_j^y = \frac{\sum_{i=1}^{n} \mathbb{1}(f_j^i = 1, y_i = y)}{\sum_{i=1}^{n} (y_i = y)}$$

The first equation is the fraction of spam emails in the data.

The second equation is the fraction of y-labelled emails that contain the  $j^{th}$  word. This is a 2D matrix of dimensions  $2 \times d$ .

#### 2.2 Prediction

The label of a test data point is obtained from the application of Baye's rule

$$P(y^{test} = 1 | x^{test}) \propto \left( \prod_{k=1}^{d} (\hat{p}_k^1)^{f_k} (1 - \hat{p}_k^1)^{1 - f_k} \right) \cdot \hat{p}$$

$$P(y^{test} = 0 | x^{test}) \propto \left( \prod_{k=1}^{d} (\hat{p}_k^0)^{f_k} (1 - \hat{p}_k^0)^{1 - f_k} \right) \cdot (1 - \hat{p})$$

if 
$$P(y^{test} = 1 | x^{test}) > P(y^{test} = 0 | x^{test})$$
 predict  $y^{test} = 1$  else  $y_{test} = 0$ 

As the values of probabilities are very low and the number of features are  $\approx 30000$  the product will be too small and might get ignored even in *float 64* data types. Hence logarithm is applied to the above inequality and on simplification the following equation is obtained.

$$\sum_{i=1}^{d} f_i \log \left( \frac{\hat{p}_i^1 (1 - \hat{p}_i^0)}{\hat{p}_i^0 (1 - \hat{p}_i^1)} \right) + \left( \sum_{i=1}^{d} \log \left( \frac{1 - \hat{p}_i^1}{\hat{p}_i^0} \right) \right) + \log \left( \frac{\hat{p}}{1 - \hat{p}} \right) \ge 0$$

- The  $2^{nd}$  and  $3^{rd}$  terms are constant for all points so let that be denoted by b.
- In the first term, put the log terms in the summation into a  $1 \times d$  matrix. Let that matrix be W.
- Also  $X_{test} = [f_1, f_2 \dots f_d]$ . Hence we can write  $y_{pred} = sign(X_{test}W^T + b)$
- If  $X_{test}$  has n data points even then the same method is applied and we'll get an  $n \times 1$  prediction matrix.

#### 2.3 Evaluation

- Rawdata is processed and split into train\_data and test\_data with a ratio of 4:1.
- The model is trained on the train split and W, b are calculated.
- On the train split, 98.6% and on the test split 96.6% accuracy was obtained

## 3 Logistic Regression Classifier

## 3.1 Training

- tfidf\_data\_array is used for this classifier. The reasons will be discussed after tuning the hyper-parameters and cross validation which will be discussed in 3.3
- The likelihood function is given by the following equation. Here g is the sigmoid function.

$$\mathcal{L}(w; Data) = \prod_{i=1}^{n} (g(w^{T}x_{i}))^{y_{i}} (1 - g(w^{T}x_{i}))^{1-y_{i}}$$

• Apply log on both sides and further simplification, we obtain the log-likelihood function as follows

$$\log \mathcal{L}(w; Data) = \sum_{i=1}^{n} \left( (1 - y_i)(-w^T x_i) - \log(1 + \exp(-w^T x_i)) \right)$$

- We cant obtain a closed form solution for w from the above equation hence we need to apply gradient ascent to maximize the log-likelihood function.
- On calculation, we get the graident as

$$\nabla \log \mathcal{L}(w) = \sum_{i=1}^{n} x_i \left( y_i - \frac{1}{1 + \exp(-w^T x_i)} \right)$$

• Gradient Ascent step:

$$w_{t+1} = w_t + \eta \cdot \nabla \log \mathcal{L}(w_t)$$

$$w_{t+1} = w_t + \eta \cdot \sum_{i=1}^{n} x_i \left( y_i - \frac{1}{1 + \exp(-w^T x_i)} \right)$$

- $\bullet$  The above equation can be computationally expensive due to the calculation of gradient at each step which is loop of length n
- We can use the some properties of the np arrays in python and do the following. Let  $\sigma$  denote the sigmoid function.
- $\sigma(w^T \cdot X)$  yields an matrix of size  $n \times 1$ , where each element represents the corresponding sigmoid function value.
- The Gradient Ascent step can be modified into:

$$w_{t+1} = w_t + \eta(X^T \cdot (Y - \sigma(w^T \cdot X)))$$

#### 3.2 Prediction

• The predication for a test data point is given by the sign of  $w^T \cdot x_i$ 

$$y_{pred} = \begin{cases} 1 & \text{if } w^T \cdot x_{test} \ge 0\\ 0 & \text{if } w^T \cdot x_{test} < 0 \end{cases}$$

## 3.3 Evaluation and Inferences

- The following data was obtained for 100 iterations of gradient ascent.
- '-' Indicates that an overflow has occured.

Data	$\eta$ (step size)	Time	Accuracy	
Data	ij (step size)		Train split	Test split
tfidf_data	1	10s	100%	98.8%
	$10^{-1}$	10s	100%	99.4%
tiidi_data	$10^{-2}$	10s	99.8%	99.1%
	$10^{-3}$	10s	94.9%	91.2%
	$10^{-3}$	_	_	_
frequency_data	$10^{-4}$	90s	99.54%	99.56%
nequency_data	$10^{-5}$	90s	98.4%	97.6%
	$10^{-6}$	90s	89.6%	86.1%

- From the above table, we can see that Logistic regression runs **faster** with tfidf\_data than frequency\_data. It is also more accurate.
- Hence the features are extracted from the text using tfidf\_vectorizer (1.2.3).
- $\eta = 10^{-1}$  produces a better result over both test and train splits.
- Train accuracy decreases on decreasing  $\eta$  and test accuracy increases and then decreases, the maximum occurs at  $\eta=10^{-1}$

## 4 SVM Classifier

## 4.1 Training

• SVM is implemented using the sklearn module in python.

```
from sklearn.svm import LinearSVC
from sklearn.metrics import accuracy_score
svm = LinearSVC(dual=True, max_iter=5000, C = 1)
svm.fit(tfidf_data_array, data['label'])
```

- The above code initializes an SVM object with a linear kernel and trains the model on the input feature matrix and corresponding labels.
- tfidf\_data(1.2.3) and C = 1 is chosen which is explained in 4.2

#### 4.2 Evaluation and Inferences

• Labels are predicted from the test data and the accuracy is calculated as follows.

```
train_predictions_svm = svm.predict(tfidf_train_array)
train_accuracy_svm = accuracy_score(train_data['label'], train_predictions_svm)

test_predictions_svm = svm.predict(tfidf_test_array)
test_accuracy_svm = accuracy_score(test_data['label'], test_predictions_svm)
```

Data	C	Time	Accuracy		
			Train split	Test split	
${ m tfidf\_data}$	0.1	1.1s	99.7%	98%	
	1	1.1s	100%	99.1%	
	5	1.1s	100%	99.1%	
	10	1.1s	100%	99.1%	
	100	1.1s	100%	99.2%	
frequency_data	0.1	2s	100%	98.7%	
	1	2s	100%	98.4%	
nequency_data	5	2s	100%	98.3%	
	10	2s	100%	98.3%	
	100	2s	100%	98.3%	

- The input features are processed using tfidf\_data (1.2.3) rather than frequency\_data (1.2.2) because the former runs relatively faster than latter as evident from the above table. Also test accuracies are better with tfidf\_data.
- Hence tfidf\_data (1.2.3) is used.
- We can see that there isn't much difference with the change of the hyper parameter C hence the default value of C = 1.0 is chosen. A high value gave a slight increase in test accuracy but that may lead to hard margin classifier. Hence they arent chosen.

## 5 Final Predictor

- Entire data, 1.1 is used for training all the models i.e, Naive Bayes, Logistic Regression and SVM. No train and test splits are created.
- The models are initialized with the hyper-parameters and data sets as discussed in the previous sections.
- For an email{n}.txt which is present in the test folder under the same directory, the output is predicted by the mode of predictions of Naive Bayes, Logistic Regression and SVM.

- That is we give the most occurring prediction in Naive Bayes, Logistic Regression and SVM as the final output.
- The outputs are written out through the print command into the standard output stream with each prediction (1 or 0) in a seperate line. Also the individual predictions of each algorithm is written into a 'predictions.txt' file

```
for i in range(len(final_predictions)):
    print(final_predictions[i])

with open('predictions.txt', 'w') as output_file:
    for i, prediction in enumerate(final_predictions):
        output_file.write(f'Email {i + 1} : NB: {bnb_predictions[i]},
        SVM : {svm_predictions[i]}, LR: {lr_predictions[i]},
        Final prediction: {prediction}\n')
```