Image Classification in Practice: Implementing Multi-layer Neural Network Using Numpy

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Abstract

In this assignment, we implemented a multi-layer neural network, including forward and backward propagation, by using Numpy. We have used this neural network to classify Japanese Hiragana hand written characters in the KMNIST dataset with softmax outputs. We performed **one-fold cross validation** and applied regularization in our model. To visualize the training progress, we have use Matplotlib to perform visualizations on the loss history.

1 Introduction

We continued with the Japanese Hiragana handwriting dataset from Programming Assignment 1. We preprocessed the data by converting all images from 2 dimensional arrays of size (28,28) into one-dimensional arrays of size 784. We applied z-score normalization to the samples and shuffled them before the training process. The data was split into 80% training data and 20% validation data. For saving time purposes, we have only used **one-fold cross validation** according to the write-up. After fine-tuning learning rate, we achieved an accuracy of 0.8546 on the test set with learning rate 0.05. When added L1 and L2 regularization, we are able to optimize our model to have a test accuracy of 0.8725 using L2 Regularization with λ of 1e-3. Then we experimented with different activation functions and network topologies. We obtained the best performing model using ReLU with accuracy 0.8688; By changing number of layers and number of hidden units in hidden layers, we were able to obtain 0.8626 accuracy on our best model on the test set.

2 Back Propagation

We have implemented backward propagation and tested our calculated gradient with both backward propagation and numerical approximation. We have used the cross-entropy as the loss function and the softmax as the output layer activation function.

2.1 Preliminary Math Work

Before implementing, it is essential to understand the math behind the process.

For backward propagation, we need to calculate the derivative of the loss with respect to the weights and bias accordingly.

First, we calculate the gradient with respect to the output of the output layer by using the cross-entropy function $E^n = -\sum_{i=1}^c t_i^n \ln y_i^n$ and the definition of $\delta_i^n = -\frac{\partial E^n}{\partial a_i^n}$ where n represents the n^{th} pattern. We get:

$$\delta_k^n = t_k^n - y_k^n \tag{1}$$

Then, we calculate the gradient with respect to the hidden layer with the output of the same hidden layer. According to the derivation, the δ_i should follow the rules of:

$$\delta_j^n = g'(a_j^n) \sum_k \delta_k^n w_{jk}^n \tag{2}$$

In the above equation, j, k symbolized the current layer and previous layer accordingly. Whereas $g(a_i^n)$ symbolized the activation function of the previous layer.

(Note: we have followed the notation used in the lecture where j is for the hidden layers and k is for the output layer).

2.2 Implementation

With equations 1 and 2, we can derive the following equations for the gradient of loss with respect to the weights and the gradient of loss with respect to the bias with some algebraic manipulations:

$$\frac{\partial E^n}{\partial w_i} = \delta_j a_j^n \tag{3}$$

$$\frac{\partial E^n}{\partial b_j} = \delta_j \tag{4}$$

With the equations 3 and 4 above, we have vectorized the above equations for faster computation in our implementations.

2.3 Testing Introduction

Suggested in the write-up, we have used the following equation to test approximate the gradient of loss with respect to the weights:

$$\frac{d}{dw}E^n(w) \approx \frac{E^n(w+\epsilon) - E^n(w-\epsilon)}{2\epsilon}, \ \epsilon > 0, \epsilon \in \mathbb{R}$$
 (5)

The w here is one weight in the neural network. After setting our $\epsilon=10^{-2}$, we compared the approximated gradient with the gradient calculated by back propagation. We have randomly chosen one output bias weight, one hidden bias weight, and two hidden to output weights and two input to hidden weights to compared the differences. You can check out our implementation in the <code>check_gradient()</code> method in <code>train.py</code>. For each selected weight, we have done the following:

- 1. Add ϵ to one weight of the weights vector, run the forward propagation, and find $E^n(w+\epsilon)$ by using 5.
- 2. Subtract ϵ to one weight of the weights vector, run the forward propagation, and find $E^n(w-\epsilon)$ by using 5.
- Run forward and then backward propagation to the unmodified weights vector and get the backward propagation calculated gradient.
- Compare the values of both approximated and the backward propagation calculated gradient.

Table 1: Architecture of the baseline model

WEIGHTS	$E(W+\epsilon)$	$E(W-\epsilon)$	NORM OF APPROXIMATED GRADIENT	NORM OF BACKWARD PROPAGATION GRADIENT	ABSOLUTE ERROR
Hidden Weight 1	2.915601	2.915669	-0.003394	-0.003394	4.689468e-10
Hidden Weight 2	1.568530	1.568079	0.02255498	0.02255499	6.128687e-9
Output Weight 1	4.945193	4.944855	0.0168997	0.0168995	2.532141e-7
Output Weight 2	2.789656	2.788244	0.0706445	0.0706437	7.989562e-7
Hidden Bias	2.789018	2.788877	0.0070264	0.0070262	1.803573e-7
Output Bias	1.052991	1.052658	0.016604	0.016603	2.630992e-7
Hidden Weight 2 Output Weight 1 Output Weight 2 Hidden Bias	1.568530 4.945193 2.789656 2.789018	1.568079 4.944855 2.788244 2.788877	-0.003394 0.02255498 0.0168997 0.0706445 0.0070264	-0.003394 0.02255499 0.0168995 0.0706437 0.0070262	6.128687e-9 2.532141e-7 7.989562e-7 1.803573e-7

2.4 Testing result

The Table 1 shows the result of our test. It is apparent that all the absolute error calculated are within big-O of ϵ^2 . Number of decimal points are remained based on whether differences between norm of approximated gradient and norm of backward propagation gradient can be seen.

3 Gradient Descent & Training Procedure

3.1 Minibatch Stochastic Gradient Descent with Momentum

We used minibatch Stochastic Gradient Descent (SGD) throughout all the problems for this programming assignment. SGD in this assignment is similar to programming assignment 1, except that we have also added momentum to our algorithm. A term called velocity is implemented for both weights and bias, denoted as v_w and v_b . The goal of having momentum is to reduce oscillations in the changing gradient, so that loss can be minimized more precisely and efficiently. We set a weight term γ for momentum as 0.9 per assignment instruction. Now the algorithm works as the pseudo code below:

Algorithm 1 Updating rule with momentum

```
1: procedure Updating rule with Momentum
 2:
 3:
           Initialization:
 4:
           w \leftarrow \text{randomized weight}
           d_w \leftarrow \text{weight gradient}
 5:
           b \leftarrow \text{randomized bias}
 6:
 7:
           d_b \leftarrow \text{bias gradient}
 8:
           \alpha \leftarrow learning rate
 9:
           \gamma \leftarrow momentum gamma
10:
           Update velocity:
11:
12:
           v_w \leftarrow \gamma \cdot v_w + (1 - \gamma) \cdot d_w
           v_b \leftarrow \gamma \cdot v_b + (1 - \gamma) \cdot d_b
13:
14:
15:
           Update bias and weight:
16:
           w \leftarrow w - \alpha \cdot v_w
17:
           b \leftarrow b - \alpha \cdot v_b
```

3.2 Training Procedure

The training procedure in this programming assignment follows a very standard deep learning model training method. The training data passes in as minibatches of fixed batch size 128 per assignment instruction, and were then split into training and validation sets. The training set consists of 80% of

the original training data, while the validation set makes up the rest 20%. We were only completing one fold cross validation for this assignment. Batches were forwarded into the network, and their initial loss and accuracy were produced. Then we utilized the loss and accuracy from the input batch to perform backpropagation, thus updating weights and bias. Early stopping is also implemented to prevent overfitting. In general, during training, when validation loss at certain continuous epochs becomes greater than the epochs before, the model must have over-adapted to the training set, thus underfitting the validation set. In such cases, it is justifiable to stop the training process early, rather than receiving a model that fails to recognize general traits in our dataset.

Algorithm 2 Training Procedure with SGD

```
1: procedure Training procedure with SGD
        train\_acc \leftarrow []
 3:
        val\_acc \leftarrow []
 4:
        train_loss \leftarrow []
 5:
        val\_loss \leftarrow []
 6:
        best\_model \leftarrow None
 7:
        best_model_loss \leftarrow float("inf")
 8:
        model \leftarrow NeuralNetwork
 9:
        patience \leftarrow 0
10:
        for epoch \leftarrow range(0,100) do
11:
             Shuffle X_train, y_train, X_val, y_val
12:
             for batch ← batch_generator do
13:
                 model.forward(batch[0], batch[1])
14:
                 model.backward()
15:
             train\_pred, train\_losses \leftarrow model.forward(X\_train, y\_train)
             train_accuracy ← accuracy(train_pred, y_train)
16:
             val\_pred, val\_losses \leftarrow model.forward(X\_val, y\_val)
17:
18:
             val\_accuracy \leftarrow accuracy(val\_pred, y\_val)
             if val_losses < best_model_loss then</pre>
19:
                 best\_model\_loss \leftarrow val\_losses
20:
21:
                 best_mode \leftarrow copy.deepcopy(model)
22:
             if val_losses > best_model_loss then
23:
                 patience+=1
                 if patience > 5 then
24:
25:
                     break
                 else
26:
27:
                     patience \leftarrow 0
28:
             train_loss.append(train_losses)
29:
             train_acc.append(train_accuracy)
30:
             val_loss.append(val_losses)
31:
             val_acc.append(val_accuracy)
32:
        return train_acc, val_acc, train_loss, val_loss, best_model
```

For details of our implementations, please refer to our train.py.

3.3 Training Results

By using the default setup with the following parameters in Table 2:

Table 2: Default parameters of our model

PARAMETER	DEFAULT VALUE		
layer_specs	[784, 128, 10]		
activation	tanh		
learning_rate	0.005		
batch_size	128		
epochs	100		
early_stop	False		
early_stop_epoch	5		
L2_penalty	0.0001		
momentum	True		
momentum_gamma	0.9		

We built our default model. The accuracy and loss on the training and validation set are presented in the line graphs Figure 1 and 2 below.

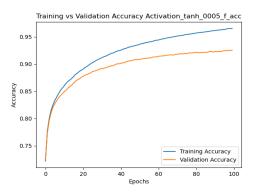


Figure 1: Training and Validation Accuracy For the MLP Using Default Settings

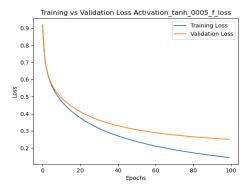


Figure 2: Training and Validation Loss For the MLP Using Default Settings

Afterwards, we tested other learning rates of 0.05, 0.01, and 0.005, and turn on the early stop. Their corresponding losses and accuracy are shown in the line graphs Figure 3 and 4, Figure 5 and 6, and Figure 7 and 8 below:

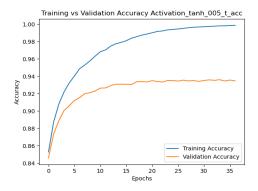


Figure 3: Training and Validation Accuracy For the MLP Using Learning Rate 0.05

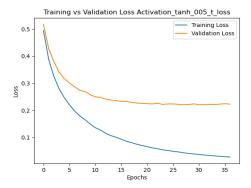


Figure 4: Training and Validation Loss For the MLP Using Learning Rate 0.05

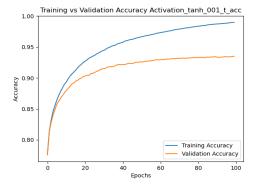


Figure 5: Training and Validation Accuracy For the MLP Using Learning Rate 0.01

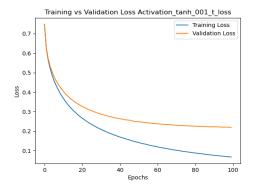


Figure 6: Training and Validation Loss For the MLP Using Learning Rate 0.01

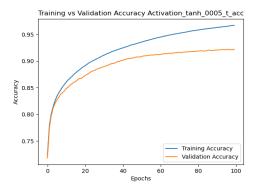


Figure 7: Training and Validation Accuracy For the MLP Using Learning Rate 0.005

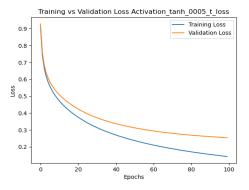


Figure 8: Training and Validation Loss For the MLP Using Learning Rate 0.005

Table 3: Learning rates, accuracy, and loss with default momentum and early stop epoch

LEARNING RATE	TRAIN ACCURACY	TRAIN LOSS	VALIDATION ACCURACY	VALIDATION LOSS
0.005	0.965333	0.143583	0.927833	0.253461
0.05	0.997146	0.036258	0.9345	0.222675
0.01	0.989542	0.067687	0.929833	0.233185

As shown in Table 3 above, we finally chose learning rate 0.05 for testing since its model has the lowest validation loss. Finally, we achieved an accuracy of **0.8546** on the test set.

4 Regularization

In this section, we have experimented with adding different weight decay value to the update rule by using both L1 and L2 regularization starting from the best model we found in part c. After using the equations 6 and 7 for both regularizations, we have aimed to find the optimal amount of regularization λ given the number of epochs increased to 110 and the rest of the hyperparemeters for the best model in part (c) stays unchanged.

4.1 L2 Regularization

L2 Regularization =
$$\sum_{i=1}^{N} l(y_i, \hat{y}_i) + \lambda \sum_{j=1}^{N} w_j^2$$
 (6)

Table 4: Accuracy, and loss with 110 epochs and other parameters in part (c) at different L2 λ

PENALTY λ	TRAIN ACCURACY	TRAIN LOSS	VALIDATION ACCURACY	VALIDATION LOSS	TEST ACCURACY
1e-3	0.9938	0.0690	0.9489	0.1781	0.8725
1e-6	0.9999	0.0051	0.9333	0.2591	0.8469

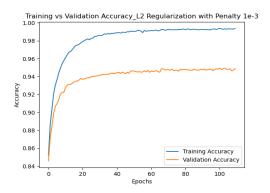


Figure 9: Training and Validation Accuracy For the MLP Using L2 Regularization 1e-3

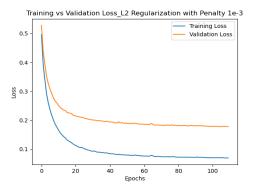


Figure 10: Training and Validation Loss For the MLP Using L2 Regularization 1e-3

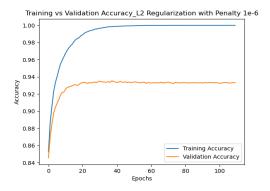


Figure 11: Training and Validation Accuracy For the MLP Using L2 Regularization 1e-6

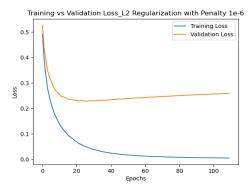


Figure 12: Training and Validation Loss For the MLP Using L2 Regularization 1e-6

4.2 L1 Regularization

L1 Regularization =
$$\sum_{i=1}^{N} l(y_i, \hat{y}_i) + \lambda \sum_{j=1}^{N} |w_j|$$
 (7)

Table 5: Accuracy, and loss with 110 epochs and other parameters in part (c) at different L1 λ

PENALTY λ	TRAIN	TRAIN	VALIDATION	VALIDATION	TEST
	ACCURACY	LOSS	ACCURACY	LOSS	ACCURACY
1e-3	0.9302	0.2515	0.9091	0.3094	0.8154
1e-6	0.9999	0.0052	0.9337	0.2574	0.8495

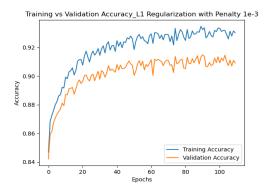


Figure 13: Training and Validation Accuracy For the MLP Using L1 Regularization 1e-3

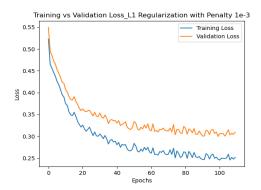


Figure 14: Training and Validation Loss For the MLP Using L1 Regularization 1e-3

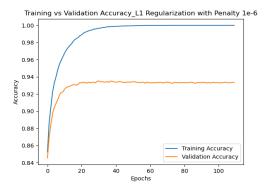


Figure 15: Training and Validation Accuracy For the MLP Using L1 Regularization 1e-6

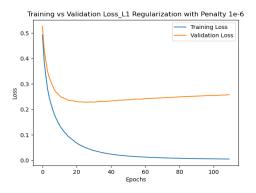


Figure 16: Training and Validation Loss For the MLP Using L1 Regularization 1e-6

4.3 Differences and Conclusion on Regularization

From the Table 4, 5 and graphs above, we have optimized L2 regularization with penalty $\lambda = 1e-3$, it produces a minimized validation loss of 0.1781, and a final test accuracy of 0.8725. For L1 regularization, it is optimized at penalty $\lambda = 1e-6$ with a validation loss of 0.2574 and a test accuracy of 0.8495. According to the graphs, L2 regularization has very steady loss decreasing and accuracy increasing graphs without too much oscillation. However, much oscillation is shown for L1 regularization with a larger penalty. This might be due to the fact that L1 regularization has a sparse solution, thus would be more likely to miss the minimum with a larger penalty. When penalty is reduced, we can observe significant improvement in model's performance.

5 Activations

In this section, we have experimented different actication functions and compared the differences in terms of the validation accuracy and loss. We applied the best performing model's parameter to test on different activation functions. However, due to different activation functions, some of the models may not converge in 100 epochs, so we adjusted the epochs to 200.

5.1 Sigmoid

The Sigmoid function is given by equation 8:

$$f(z) = \frac{1}{1 + e^{-z}} \tag{8}$$

The key characteristic is that the Sigmoid function has a range of 0 to 1. i.e. $f(z) \in (0,1)$. As a result, it has a great vanishing gradient problem when the input is too large, which will affect the effectiveness of updating the weights.

5.2 Tanh

The Tanh function is given by equation 9:

$$f(z) = \frac{e^z - e^{-z}}{e^z + e^{-z}} \tag{9}$$

The key characteristic is that the Tanh function has a range of -1 to 1. i.e. $f(z) \in (-1,1)$. As a result, it also has a great vanishing gradient problem when the input is too large, which will affect the effectiveness of updating the weights. However, the difference is that this activation function is centered at 0.

5.3 ReLU

The ReLU function is given by equation 10:

$$f(z) = \max(0, z) \tag{10}$$

The key characteristic is that the ReLU function has a range of 0 to infinity. i.e. $f(z) \in (0, \infty)$. As a result, it will be able to prevent the vanishing gradient problem in comparison to the previous two activation functions.

5.4 Different Activation Results

Table 6: Accuracy, and loss with different activation functions

ACTIVATION	TRAIN	TRAIN	VALIDATION	VALIDATION	TEST
FUNCTION	ACCURACY	LOSS	ACCURACY	LOSS	ACCURACY
Sigmod	0.993438	0.05249	0.94058	0.20047	0.8587
Tanh	0.998396	0.02781	0.93225	0.22849	0.8516
ReLU	0.997729	0.01941	0.94483	0.20109	0.8688

Table 6 contains the accuracy and loss of test set trained with different activation functions. It is clear that the ReLU activation function has the best performance because of its ability to resolve the vanishing gradient problem.

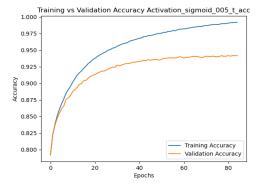


Figure 17: Training and Validation Accuracy For the MLP Using Sigmoid as Activation Function

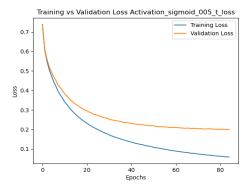


Figure 18: Training and Validation Loss For the MLP Using Sigmoid as Activation Function

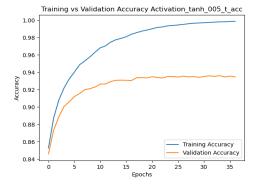


Figure 19: Training and Validation Accuracy For the MLP Using Tanh as Activation Function

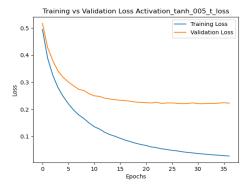


Figure 20: Training and Validation Loss For the MLP Using Tanh as Activation Function

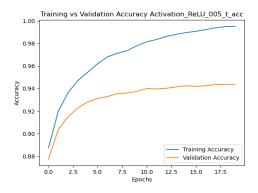


Figure 21: Training and Validation Accuracy For the MLP Using ReLU as Activation Function

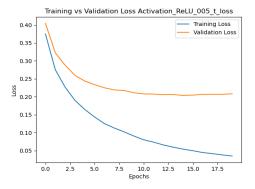


Figure 22: Training and Validation Loss For the MLP Using ReLU as Activation Function

5.5 Conclusion

By using the best performing parameter obtained from assignment part (c), We achieved best test accuracy of **0.8688** with ReLU activation function. According to the graphs, both tanh and ReLu activation functions take approximately 20 to 30 epochs to converge, while Sigmoid activation function requires more than 80 epochs to converge.

6 Network Topology

In this section, we have experimented with different topology of the neural network starting from the best model from part c. Training and validation loss and accuracy, and each topology's accuracy are shown in the graphs below. The topologies include halving the number of units in the hidden layer, doubling the number of hidden units in the hidden layer, and adding a second hidden layer. To make sure that using two hidden layers of equal size and has approximately the same number of parameters for the best model choice in the previous experiment, we set the number of hidden units in both hidden layers to 203.

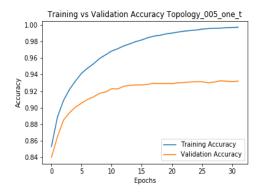


Figure 23: Training and Validation Loss For the MLP Using 128 Hidden Units

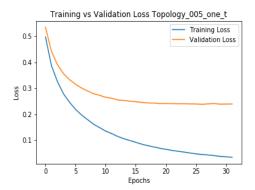


Figure 24: Training and Validation Loss For the MLP Using 128 Hidden Units

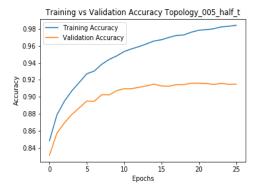


Figure 25: Training and Validation Accuracy For the MLP Using 64 Hidden Units

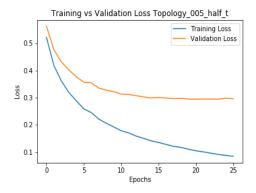


Figure 26: Training and Validation Loss For the MLP Using 64 Hidden Units

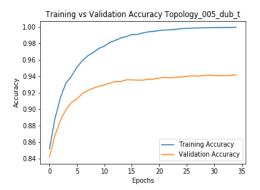


Figure 27: Training and Validation Accuracy For the MLP Using 256 Hidden Units

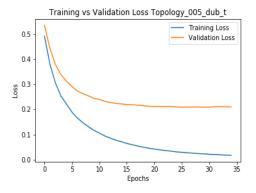


Figure 28: Training and Validation Loss For the MLP Using 256 Hidden Units

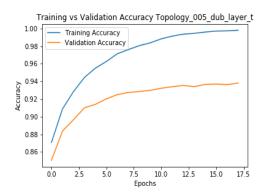


Figure 29: Training and Validation Accuracy For the MLP Using 2 Hidden Layers

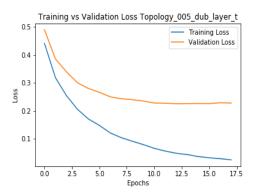


Figure 30: Training and Validation Loss For the MLP Using 2 Hidden Layers

Table 7: Accuracy, and loss with different topologies

TOPOLOGY	TRAIN	TRAIN	VALIDATION	VALIDATION	TEST
	ACCURACY	LOSS	ACCURACY	LOSS	ACCURACY
64 Units	0.98425	0.085203	0.915	0.296058	0.826
128 Units	0.99725	0.034221	0.932167	0.239340	0.8514
256 Units	0.999542	0.017683	0.941833	0.210170	0.8626
203x203 Units	0.998083	0.025471	0.938083	0.228035	0.8583

6.1 Conclusion

As shown in Table 7, we had minimum validation loss on the model with 1 hidden layer of 256 hidden units. Per assignment instruction that we should only "Report final accuracy", we obtained final accuracy on the test set of **0.8626**. According to the graphs, all of them converged after 20 to 30 epochs. The accuracy increases, and the loss decreases as the number of hidden units in the hidden layer increases. However, adding extra hidden layer, even though improved model's performance, but started to overfit much earlier much faster than other topologies, which is reflected in the graph as it stopped at less than 20 epochs. Thus we can reach the conclusion that increasing number of hidden units in a hidden layer and increasing number of hidden layers help improve model's performance, while increasing number of hidden layers may tend to have overfitting issue.

7 Team Contributions

- Linghang Kong: In this project, I worked on implementation of neuralnet.py together with Yi Li and Weiyue Li. Then I initialized train(), train_mlp(), activation_experiment() and topology_experiment(), as well as the plotting functions affiliated with them. We did equal amount of work on the assignment report.
- Weiyue Li: In this project, I worked on implementing the neuralnet.py and data.py, and debugging the train.py with my teammates. I was also in charge of the main.py, README.md, visualization methods of our results, as well as the documentations in our.py files. We did equal amount of work on this report.
- Yi Li: In this project, I was in charge of the neuralnet.py, train.py, data.py, especially focusing on the parts of Regularizations and Momentum. I was also in charge of the part 2 table, part3 plots, part4 table and plots, and algorithm descriptions on final report.