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# MLP Coursework 1

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## Abstract

In this report we study the problem of overfitting, which is the training regime where performance increases on the training set but decrease on validation data. **[In overfitting, the model learns too much about the training data and is unable to generalise its learnings for unseen data which leads to poor performance on new data points.]** . We first analyse the given example and discuss the probable causes of the underlying problem. Then we investigate how the depth and width of a neural network can affect overfitting in a feedforward architecture and observe that increasing width and depth **[slightly results in improved model performance overall. If either or both the width and depth are increased substantially, the model starts overfitting the training data and returns low accuracy scores for any unseen data.]** . Next we discuss how two standard methods, Dropout and Weight Penalty, can mitigate overfitting, then describe their implementation and use them in our experiments to reduce the overfitting on the EMNIST dataset. Based on our results, we ultimately find that **[the model which was just using dropout regularisation was performing the best. It was able to properly learn from the training data and return high accuracy scores on the testing data while keeping the error of the model low.]** . Finally, we conclude the report with our observations and related work. Our main findings indicate that **[complex models can easily overfit the data resulting in low accuracy scores on unseen data points. We can use various techniques like Dropout & L1/L2 regularisations to ensure that the model doesn't memorise the data and is instead able to generalise the learnings from it.]** .

## 1. Introduction

In this report we focus on a common and important problem while training machine learning models known as overfitting, or overtraining, which is the training regime where performances increase on the training set but decrease on unseen data. We first start with analysing the given problem in Figure 1, study it in different architectures and then investigate different strategies to mitigate the problem. In particular, Section 2 identifies and discusses the given problem, and investigates the effect of network width and depth

in terms of generalisation gap (see Ch. 5 in Goodfellow et al. 2016) and generalisation performance. Section 3 introduces two regularisation techniques to alleviate overfitting: Dropout (Srivastava et al., 2014) and L1/L2 Weight Penalties (see Section 7.1 in Goodfellow et al. 2016). We first explain them in detail and discuss why they are used for alleviating overfitting. In Section 4, we incorporate each of them and their various combinations to a three hidden layer<sup>1</sup> neural network, train it on the EMNIST dataset, which contains 131,600 images of characters and digits, each of size 28x28, from 47 classes. We evaluate them in terms of generalisation gap and performance, and discuss the results and effectiveness of the tested regularisation strategies. Our results show that **[the model which was just using dropout regularisation was performing the best. It was able to properly learn from the training data and return high accuracy scores on the testing data while keeping the error of the model low.]** .

Finally, we conclude our study in Section 5, noting that **[complex models can easily overfit the data resulting in low accuracy scores on unseen data points. We can use various techniques like Dropout & L1/L2 regularisations to ensure that the model doesn't memorise the data and is instead able to generalise the learnings from it.]** .

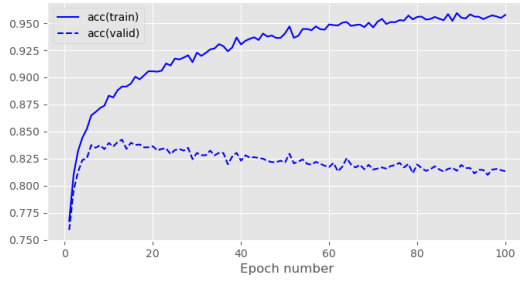
## 2. Problem identification

Overfitting to training data is a very common and important issue that needs to be dealt with when training neural networks or other machine learning models in general (see Ch. 5 in Goodfellow et al. 2016). A model is said to be overfitting when as the training progresses, its performance on the training data keeps improving, while its is degrading on validation data. Effectively, the model stops learning related patterns for the task and instead starts to memorize specificities of each training sample that are irrelevant to new samples.

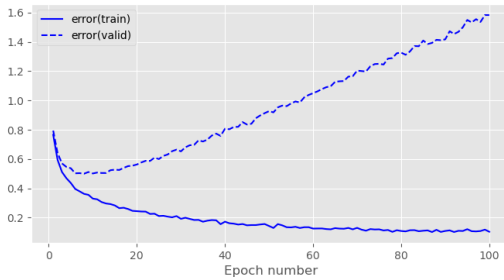
**[Overfitting causes a model to start recognising random deviances and noise present in the training data as a regularity rather than learning the underlying patterns. This poor generalisation leads to the model struggling to capture the trends on unseen data points and hence performing poorly on them. This makes the model less reliable to be used in real world applications. Another direct consequence of overfitting is that it skews the rela-**

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<sup>1</sup>We denote all layers as hidden except the final (output) one. This means that depth of a network is equal to the number of its hidden layers + 1.



(a) accuracy by epoch



(b) error by epoch

Figure 1. Training and validation curves in terms of classification accuracy (a) and cross-entropy error (b) on the EMNIST dataset for the baseline model.

relationship between variables affecting the interpretability of the model and making it difficult to fix the model. Overfitting a model also wastes a lot of resources since it requires a lot of computational power to create a model which will be unable to handle new data points efficiently.] .

Although it eventually happens to all gradient-based training, it is most often caused by models that are too large with respect to the amount and diversity of training data. The more free parameters the model has, the easier it will be to memorise complex data patterns that only apply to a restricted amount of samples. A prominent symptom of overfitting is the generalisation gap, defined as the difference between the validation and training error. A steady increase in this quantity is usually interpreted as the model entering the overfitting regime.

[Network capacity refers to the ability of the model to learn the general characteristics of the data. When a model is more complex, it starts memorising the data rather than learning the trends from it. This usually happens when the model has more room than what is needed for capturing the patterns in the data. If the dataset is small or simple, the basic essence of it can be captured with a small model itself. If in scenarios like this, a complex model is used, the model has the capacity to learn more than the general features of the data and will also start to fit the outliers in the data leading to overfitting on the training set. Thus it is essential to choose an appropriate model for the data provided to us, otherwise choosing a complex model

# Hidden Units	Val. Acc.	Train Error	Val. Error
32	78.9	0.563	0.705
64	80.9	0.345	0.660
128	80.5	0.163	0.924

Table 1. Validation accuracy (%) and training/validation error (in terms of cross-entropy error) for varying network widths on the EMNIST dataset.

with more network capacity will result in overfitting of the data.] .

Figure 1a and 1b show a prototypical example of overfitting. We see in Figure 1a that [the training accuracy of the model increases rapidly in the first few epochs and then gradually slows down as the epochs continue to run. Even though the validation accuracy increases at a good rate in the first few epochs, it starts flattening out after a point and then slowly starts to decline. This is the point where overfitting has started, stating that the model has learned the training data and is unable to generalise it to the validation data. This is further validated by the graph in Figure 1b which plots the error of the model on the training and validation dataset. We can see that the training error reduces rapidly at first and then eventually tends to 0. After a rapid initial decline in the error of the validation set, the error starts increasing at the same time when the accuracy of the validation set starts decreasing. This is due to the fact the model has now overfit and is struggling to process unseen data points.] .

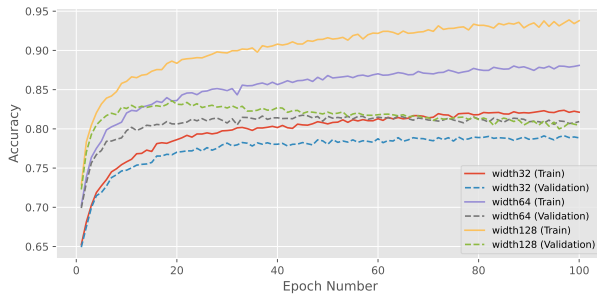
The extent to which our model overfits depends on many factors. For example, the quality and quantity of the training set and the complexity of the model. If we have sufficiently many diverse training samples, or if our model contains few hidden units, it will in general be less prone to overfitting. Any form of regularisation will also limit the extent to which the model overfits.

## 2.1. Network width

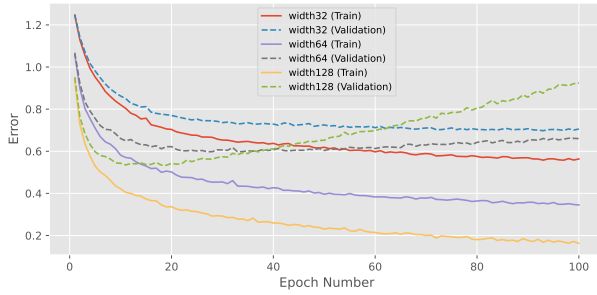
[ Question Table 1 - Fill in Table 1 with the results from your experiments varying the number of hidden units. ] [Question Figure 2 - Replace the images in Figure 2 with figures depicting the accuracy and error, training and validation curves for your experiments varying the number of hidden units. ]

First we investigate the effect of increasing the number of hidden units in a single hidden layer network when training on the EMNIST dataset. The network is trained using the Adam optimiser with a learning rate of  $9 \times 10^{-4}$  and a batch size of 100, for a total of 100 epochs.

The input layer is of size 784, and output layer consists of 47 units. Three different models were trained, with a single hidden layer of 32, 64 and 128 ReLU hidden units respectively. Figure 2 depicts the error and accuracy curves over 100 epochs for the model with varying number of hid-



(a) accuracy by epoch



(b) error by epoch

Figure 2. Training and validation curves in terms of classification accuracy (a) and cross-entropy error (b) on the EMNIST dataset for different network widths.

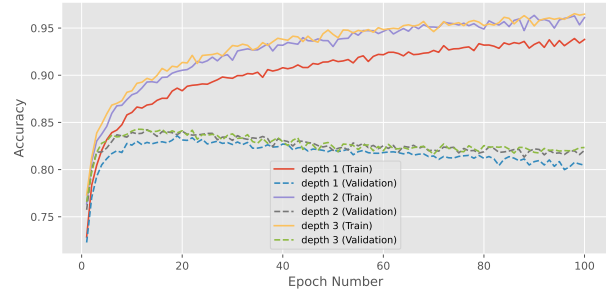
den units. Table 1 reports the final accuracy, training error, and validation error. We observe that [the final validation accuracy of all the models is very close (Figure 2a) indicating that the models perform very similarly. In Figure 2b, if we look at the train and the validation error, we can see the difference in the performance of the models. The training error of the model with 32 hidden units is the highest and the one with 128 hidden units is the lowest. This would ideally mean that the model with 128 neurons would perform the best. But if we look at the validation error, we can see the 128 units model has the highest validation error. This means that model has overfit the training data and is performing relatively worse on data it has not seen before when compared to the 32 units model.

From Table 2, we can see that the model with 64 hidden units has the highest validation accuracy and the lowest validation error signifying that the model has learned from the training data and is able to generalise the best on unseen data out of all the 3 models. This model is relatively less complex and is performing better than a model with more width.] .

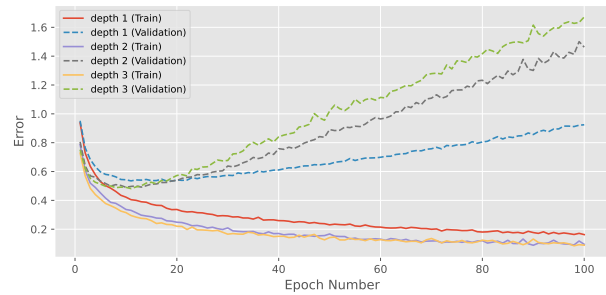
[Changing the width of the model gave us a variety of results. We saw that a model with more width (more complex model) is not necessarily the best model for most scenarios. Even though the model might slightly perform better on the training data, it gives higher error and not a significant increase on the accuracy for new data points. The results we've achieved also prove the point about the network capacity, where a higher

# Hidden Layers	Val. Acc.	Train Error	Val. Error
1	80.5	0.163	0.924
2	82.1	0.095	1.463
3	82.3	0.089	1.672

Table 2. Validation accuracy (%) and training/validation error (in terms of cross-entropy error) for varying network depths on the EMNIST dataset.



(a) accuracy by epoch



(b) error by epoch

Figure 3. Training and validation curves in terms of classification accuracy (a) and cross-entropy error (b) on the EMNIST dataset for different network depths.

width model has the more than the required capacity to learn the features of the data and thus it starts fitting the outliers to the data leading to overfitting. This won't always be the case as this test was conducted on a relatively simple dataset. If the data to be processed had more features, we would require a model with more width to learn the underlying patterns in the data. A complex model in this scenario won't overfit because its capacity is being completely utilised in learning the features of the data.] .

## 2.2. Network depth

[ Question Table 2 - Fill in Table 2 with the results from your experiments varying the number of hidden layers. ] [Question Figure 3 - Replace these images with figures depicting the accuracy and error, training and validation curves for your experiments varying the number of hidden layers. ]

Next we investigate the effect of varying the number of hidden layers in the network. Table 2 and Figure 3 depict results from training three models with one, two and three

hidden layers respectively, each with 128 ReLU hidden units. As with previous experiments, they are trained with the Adam optimiser with a learning rate of  $9 \times 10^{-4}$  and a batch size of 100.

We observe that **[models of different depth are all capable of classifying classes of EMNIST dataset. From Figure 3a, providing the models with more depth has enabled them to extract features from the data efficiently and give excellent accuracy scores. Irrespective of their depth, in case of validation data, all of them are performing similarly and giving accuracy of around 80% (Table 2). The rise of accuracy is very steep in the first 10 epochs, and then it is slow gradual improvement in training accuracy and barely any change in the validation accuracy.]**

**The error behaves very differently and gives a good understanding of the models. In Figure 3b, we can see that the error of all the models on the training set approaches 0, but on the validation set starts increasing as the epochs increase. This is the point of overfitting; and the magnitude of the error indicates how much the model has learned the training data. The model with the highest depth has the highest capacity and hence has been able to learn the train set the best resulting in it having the highest error on the validation set (from Table 2).]**

**[Increasing the depth of the model can be helpful in processing large or complex datasets. But in other situations it can turn out to be counter productive and as the model will overfit the data. This is because certain capability is required for the model to process the data, but if the model has more room to capture features, it'll start fitting the noise to the dataset. As the depth of the model increases, there are more neurons available which allow the model to extract more features from the dataset allowing it to overfit the training data and result in poor accuracy scores for unseen data. It also becomes computationally expensive to train the model and the resources will be wasted if the model overfits on the training data. We can tackle the problem of overfitting by using methods like Dropout and Regularisation which penalise the model for generalising too much to the training set preventing overfitting.]**

### 3. Dropout and Weight Penalty

In this section, we investigate three regularisation methods to alleviate the overfitting problem, specifically dropout layers and the L1 and L2 weight penalties.

#### 3.1. Dropout

Dropout (Srivastava et al., 2014) is a stochastic method that randomly inactivates neurons in a neural network according to an hyperparameter, the inclusion rate (*i.e.* the rate that an unit is included). Dropout is commonly represented by an additional layer inserted between the linear layer

and activation function. Its forward pass during training is defined as follows:

$$\text{mask} \sim \text{bernoulli}(p) \quad (1)$$

$$\mathbf{y}' = \text{mask} \odot \mathbf{y} \quad (2)$$

where  $\mathbf{y}, \mathbf{y}' \in \mathbb{R}^d$  are the output of the linear layer before and after applying dropout, respectively.  $\text{mask} \in \mathbb{R}^d$  is a mask vector randomly sampled from the Bernoulli distribution with inclusion probability  $p$ , and  $\odot$  denotes the element-wise multiplication.

At inference time, stochasticity is not desired, so no neurons are dropped. To account for the change in expectations of the output values, we scale them down by the inclusion probability  $p$ :

$$\mathbf{y}' = \mathbf{y} * p \quad (3)$$

As there is no nonlinear calculation involved, the backward propagation is just the element-wise product of the gradients with respect to the layer outputs and mask created in the forward calculation. The backward propagation for dropout is therefore formulated as follows:

$$\frac{\partial \mathbf{y}'}{\partial \mathbf{y}} = \text{mask} \quad (4)$$

Dropout is an easy to implement and highly scalable method. It can be implemented as a layer-based calculation unit, and be placed on any layer of the neural network at will. Dropout can reduce the dependence of hidden units between layers so that the neurons of the next layer will not rely on only few features from of the previous layer. Instead, it forces the network to extract diverse features and evenly distribute information among all features. By randomly dropping some neurons in training, dropout makes use of a subset of the whole architecture, so it can also be viewed as bagging different sub networks and averaging their outputs.

#### 3.2. Weight penalty

L1 and L2 regularisation (Ng, 2004) are simple but effective methods to mitigate overfitting to training data. The application of L1 and L2 regularisation strategies could be formulated as adding penalty terms with L1 and L2 norm square of weights in the cost function without changing other formulations. The idea behind this regularisation method is to penalise the weights by adding a term to the cost function, and explicitly constrain the magnitude of the weights with either the L1 and L2 norms. The optimisation problem takes a different form:

$$\text{L1: } \min_{\mathbf{w}} E_{\text{data}}(\mathbf{X}, \mathbf{y}, \mathbf{w}) + \lambda \|\mathbf{w}\|_1 \quad (5)$$

$$\text{L2: } \min_{\mathbf{w}} E_{\text{data}}(\mathbf{X}, \mathbf{y}, \mathbf{w}) + \lambda \|\mathbf{w}\|_2^2 \quad (6)$$

where  $E_{\text{data}}$  denotes the cross entropy error function, and  $\{\mathbf{X}, \mathbf{y}\}$  denotes the input and target training pairs.  $\lambda$  controls the strength of regularisation.



Weight penalty works by constraining the scale of parameters and preventing them to grow too much, avoiding overly sensitive behaviour on unseen data. While L1 and L2 regularisation are similar to each other in calculation, they have different effects. Gradient magnitude in L1 regularisation does not depend on the weight value and tends to bring small weights to 0, which can be used as a form of feature selection, whereas L2 regularisation tends to shrink the weights to a smaller scale uniformly.

[The L1 penalty also known as Lasso regularisation and L2 penalty known as Ridge regularisation can be used together to tackle the overfitting that occurs during the training of the model. One of the approaches to combining the L1 and L2 regularisations is called the Elastic Net. The equation for the Elastic Net is given by:

$$E = E_{\text{data}} + \lambda_1 \sum_{i=1}^N |w_i| + \lambda_2 \sum_{i=1}^N w_i^2 \quad (7)$$

where  $\lambda_1$  and  $\lambda_2$  are the hyperparameters for the L1 and L2 regularisation respectively.

The L1 regularisation is used to reduce the number of features by shrinking the coefficients of the less important features to 0. The L2 regularisation penalises the model for having large coefficients and shrinks them to a smaller scale. By adjusting the values of  $\lambda_1$  and  $\lambda_2$  appropriately, we can find a right balance of L1 and L2 penalties that will help the model generalise over new or unseen data points reducing the overall variance and preventing overfitting.

Sometimes while using L1 regularisation, the model can get stuck at a local minima due to some of the parameters being set to 0. This problem can be mitigated by using L2 norms along with L1 norms. This is because the L2 norm is a smooth function and hence the model can easily escape the local minima.] .

## 4. Balanced EMNIST Experiments

[ Question Table 3 - Fill in Table 3 with the results from your experiments for the missing hyperparameter values for each of L1 regularisation, L2 regularisation, and Dropout (use the values shown on the table). ]

[Question Figure 4 - Replace these images with figures depicting the Validation Accuracy and Generalisation Gap (difference between validation and training error) for each of the experiment results varying the Dropout inclusion rate, and L1/L2 weight penalty depicted in Table 3 (including any results you have filled in). ]

Here we evaluate the effectiveness of the given regularisation methods for reducing the overfitting on the EMNIST dataset. We build a baseline architecture with three hidden layers, each with 128 neurons, which suffers from overfitting as shown in section 2.

Here we train the network with a lower learning rate of  $10^{-4}$ , as the previous runs were overfitting after only a handful of

epochs. Results for the new baseline (c.f. Table 3) confirm that lower learning rate helps, so all further experiments are run using it.

Then, we apply the L1 or L2 regularisation with dropout to our baseline and search for good hyperparameters on the validation set. We summarise all the experimental results in Table 3. For each method, we plot the relationship between generalisation gap and validation accuracy in Figure 4.

First we analyze three methods separately, train each over a set of hyperparameters and compare their best performing results.

[We introduce a new hyperparameter for controlling the dropout in the model which is the inclusion probability. This hyperparameter specifies the percentage of the neurons that should be considered and the rest are converted to 0. This helps the model become less reliant on particular neurons and prevent overfitting. We tested with different inclusion probabilities ranging from 0.6 to almost 1 and found that the validation accuracy didn't increase much even though more neurons were available for the model to extract the patterns from the data. The validation accuracy increased at a slow pace with availability of more neurons, but the increase in generalisation gap was relatively high, hinting at overfitting of the model (Figure 4a).

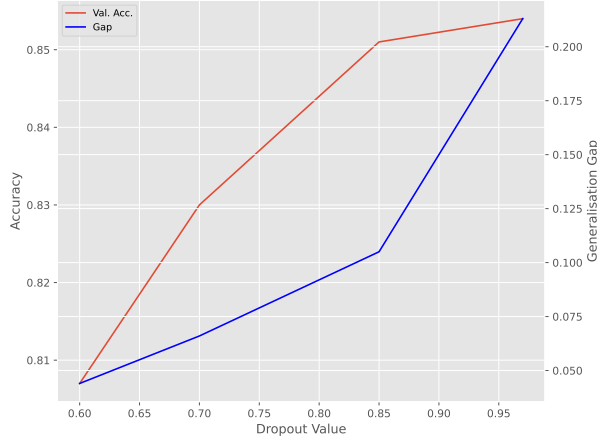
Using another hyperparameter, we controlled the amount of weight penalties that were applied to the model using the L1 & L2 norms. In case of L1 penalties, as the coefficients of the penalty increased, the error increased while simultaneously decreasing the accuracy of the model. Since L1 penalties add the absolute values of the weights to the error, this told us that some features are relatively more important than others for prediction of the model. L2 penalties had a similar but not as extreme effect when compared to L1 penalties. Even though Figure 4b shows that the generalisation gap is decreasing, it is because the training error is also increasing signifying that the model is unable to fit the training data well.

From these testings, we concluded that the dropout model with inclusion probability of 0.85 was performing the best. Even though the model with inclusion probability of 0.97 has a slightly higher validation accuracy, it comes at the cost of increased generalisation gap. This model was then used to classify the test set of the EMNIST data. The accuracy of the model on the test data is 84.0% and the error is 0.471. This is because the model is able to generalise well over unseen data and has not overfit on the training set. We also noted that the model with 0.97 inclusion probability had a slightly higher accuracy on the test set of 84.55%. Although the accuracy is higher, this cannot be a factor in choosing a model because we can't check the accuracy scores on the test set in real life scenarios.] .

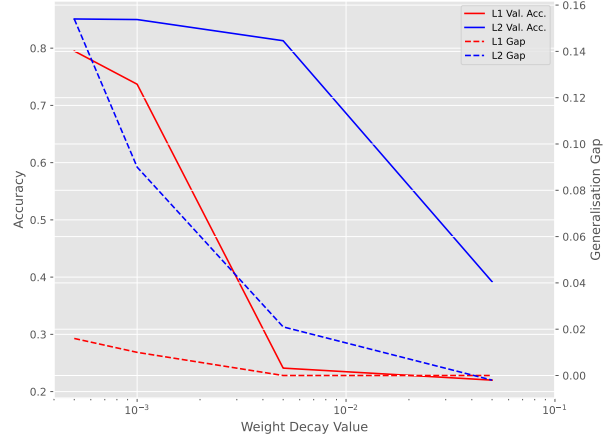
[We can extend this experiment by testing a plethora

Model	Hyperparameter value(s)	Validation accuracy	Train Error	Validation Error
Baseline	-	0.837	0.241	0.533
Dropout	0.6	80.7	0.549	0.593
	0.7	82.7	0.447	0.509
	<b>0.85</b>	<b>85.1</b>	<b>0.329</b>	<b>0.434</b>
	0.97	85.4	0.244	0.457
L1 penalty	5e-4	79.5	0.642	0.658
	1e-3	75.0	0.850	0.856
	5e-3	2.41	3.850	3.850
	5e-2	2.20	3.850	3.850
L2 penalty	5e-4	85.1	0.306	0.460
	1e-3	84.8	0.362	0.460
	5e-3	81.3	0.586	0.607
	5e-2	39.2	2.258	2.256

Table 3. Results of all hyperparameter search experiments. *italics* indicate the best results per series (Dropout, L1 Regularisation, L2 Regularisation) and **bold** indicates the best overall.



(a) Accuracy and error by inclusion probability.



(b) Accuracy and error by weight penalty.

Figure 4. Accuracy and error by regularisation strength of each method (Dropout and L1/L2 Regularisation).

of combinations of the dropout regularisation and L1 & L2 penalties to fine tune the model further. We have already seen that each regularisation method is used for a specific task and has its own advantages and disadvantages. By combining these methods, we can try to mitigate the disadvantages of one method by using the other. Some of the combinations for the models that we can use are:

- $\alpha = 1e-4$ ; Dropout = 0.75; L1 Penalty = 5e-4
- $\alpha = 1e-4$ ; Dropout = 0.97; L1 Penalty = 1e-3
- $\alpha = 1e-4$ ; Dropout = 0.75; L2 Penalty = 5e-4
- $\alpha = 5e-4$ ; Dropout = 0.97; L2 Penalty = 1e-3
- $\alpha = 1e-4$ ; L1 Penalty = 5e-4; L2 Penalty = 5e-4
- $\alpha = 1e-4$ ; L1 Penalty = 3e-4; L2 Penalty = 9e-4

- $\alpha = 1e-4$ ; L1 Penalty = 5e-4; L2 Penalty = 1e-3; Dropout = 0.9

- $\alpha = 1e-4$ ; L1 Penalty = 1e-4; L2 Penalty = 9e-4; Dropout = 0.9

These experiments would give us an idea about which regularisation methods work well together. The dropouts control the number of neurons the model is using to train. Controlling this directly controls the capacity of the network and hence it is a crucial tool in preventing overfitting.

The L1 norms penalise the model for having too large gradients and bring them down to 0. A combination of L1 norms and Dropouts would promote a simpler model with small architecture which would help with generalisation and faster convergence. L2 norms also penalise the model for having large weights, but instead of reducing the weights to 0, it scales it down in a uniform

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scale. Dropouts with L2 will lead to a smoother model which will converge in fewer instances while still helping the performance of the model. The values of hyperparameters above will check well performing regularisation methods in combination with another relatively low performing regulariser to find the ideal balance of the parameters. We found that the model with learning rate  $1e-4$ , L1 penalty  $5e-4$ , L2 penalty  $1e-3$  and dropout inclusion probability of 0.9 performed the best giving an validation accuracy of 83.5%].

## 5. Conclusion

[In this report, we have discussed the problem of overfitting in machine learning models. We studied how increasing the width and depth of the model increased the model's ability to learn from the data and found that a model with more than required capacity is guaranteed to overfit the training data eventually. Consequentially, we looked at methods to tackle this problem of overfitting. Experiments with dropout regularisation were carried out which help in removing the dependency of the model from certain specific neurons. Considering too many neurons caused the model to overfit whereas considering too few neurons made the model underfit the data. We then looked at L1/L2 norms which penalise the weights of the model helping in preventing overfitting. Various experiments were formulated to combine the power of dropout and L1/L2 norms to get the best of all the regularisation methods and help the model performance. These formulations can help us get a good idea about combinations of regularisation methods that will work well together. Further studies should focus on using combined regularisation methods like Elastic Net to prevent overfitting. Early stopping is another approach that can be explored to combat overfitting in machine learning models.]

## References

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