

# Deep Learning (IST, 2024-25)

# **Homework 1**

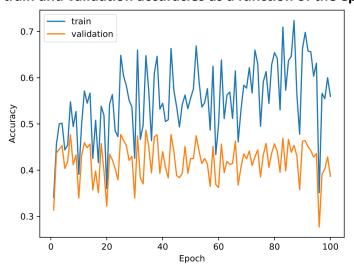
## Question 1

1.

a) After training the perceptron for 100 epochs, we obtained the following results:

train accuracy: 0.5598
validation accuracy: 0.3868
test accuracy: 0.3743

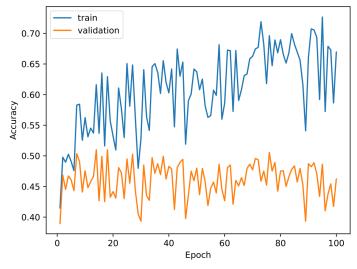
## Perceptron's train and validation accuracies as a function of the epoch number



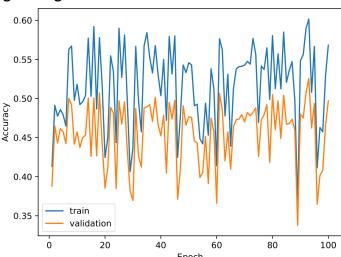
2.

a) After training the logistic regression classifier (without  $\ell_2$  regularization) for 100 epochs, we obtained a final **test accuracy** of 0.4597.

## Non-regularized logistic regression's train and validation accuracies over the epochs



**b)** After training the logistic regression classifier (with  $\ell_2$  regularization) for 100 epochs, we obtained a final **test accuracy** of 0.5053.



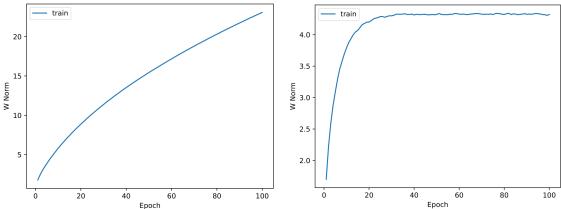
### Regularized logistic regression's train and validation accuracies over the epochs

### Analysis:

Comparing the training and validation accuracies across the stochastic gradient descent epochs, we can see that without regularization the training accuracies tend to be better than with it. Therefore, the differences between training and validation accuracies are greater and tend to increase with the number of epochs, which can make us infer a tendency for overfitting of the model.

So, our analysis enforces that regularization prevents overfitting as expected in theory, reducing and stabilizing the gap between training and testing accuracies. Furthermore, the final test accuracy is better including the regularization.

# $\ell_2$ -norm of the weights of non-regularized (left) and regularized (right) versions of logistic regression as a function of the number of epochs



#### Analysis:

According to the obtained values, it is visible that without regularization the weights tendency is to increase over the epochs and with regularization the weights become more stable and also smaller than non-regularized ones.

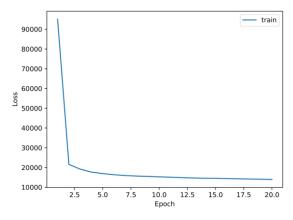
If the weight becomes too large, the regularization term adds a larger penalty, making the loss function increase and leading to the smaller weights seen in the graphics. Therefore, preventing overfitting and reducing the impact of noise in the training data.

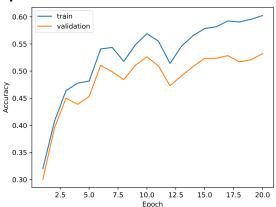
d) As  $\ell_2$  regularization, using  $\ell_1$  regularization the weights are smaller than non-regularized weights, thus preventing overfitting. The main difference between these regularizations is that  $\ell_1$  encourages sparsity. Features with small weights have high probability to be set to zero during training.

As of  $\ell_1$  norm is defined as the summation of absolute values, the gradient changes abruptly and favors a solution where many weights "spike" down to zero (at sparse points) reducing the number of non-zero weights, which leads some features to be completely ignored and to have fewer but more active features. That does not happen in  $\ell_2$  regularization that spreads the penalty evenly across all weights, encouraging them to be small but not zero.

**a)** After training the model (multi-layer perceptron with the parameters given in homework statement) for 20 epochs, we obtained a final **test accuracy** of 0.5417.

# Multi-layer perceptron's train loss (left) and train and validation accuracies (right) over the epochs





## **Question 2**

- **1.** After training the linear model implemented for 100 epochs with 3 different values for learning rate (0.00001, 0.001, 0.1), we obtained the following results:
  - learning rate 0.00001

validation accuracy: 0.2892test accuracy: 0.3087

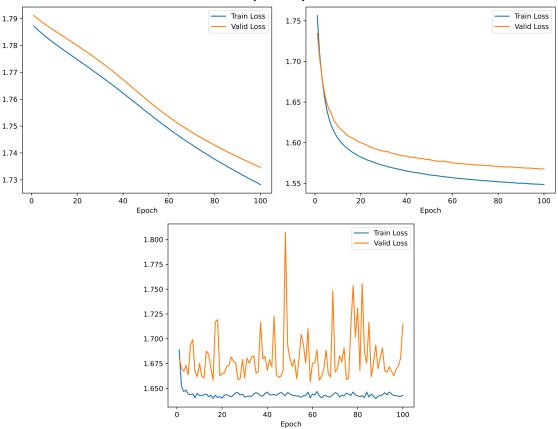
• learning rate 0.001

validation accuracy: 0.4751test accuracy: 0.4827

learning rate 0.00001

validation accuracy: 0.3091test accuracy: 0.3093

Training and validation losses for each configuration – 0.00001 (left); 0.001 (right); 0.1 (bottom)



#### Analysis:

Best validation accuracy was obtained for learning rate 0.001 (final validation accuracy = 0.4751). As it is possible to observe by the validation accuracy results and by the validation loss curves, from learning rates tested the closest to be optimal is 0.001.

It is perceivable that with very low (like 0.00001) learning rates the curve is almost straight, and the weights are being updated steadily but very slowly, which does not make possible to verify any approximation to a convergence point in which the loss curve reaches a minimum.

That is not the case when looking at the curve associated with the learning rate that led to the best result. We can notice the loss decreases smoothly, almost flattening along the epochs which indicates convergence.

In the other hand, a very high learning rate (like 0.1) cause the loss to fluctuate and possible increase, which shows instability, overshooting the minimum and bouncing around instead of settling.

**a)** After training the model for 200 epochs with batch size default (64) and batch size of 512, keeping the other parameters at their default value, we obtained:

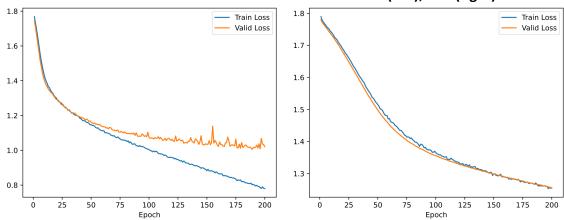
batch size 64 (default)

validation accuracy: 0.6061test accuracy: 0.6073

batch size 512

validation accuracy: 0.5028test accuracy: 0.5190

### Train and validation losses for each model – 64 (left); 512 (right)



### **Analysis:**

The best performance is observed with the default batch size (64). However, the execution time is shorter with the batch size set to 512. This happens because the weights need to be updated less frequently — specifically, dataset\_size / batch\_size times. Having a lower batch size implies a more computationally demanding classification task.

Despite this, a lower batch size allows the model to learn more effectively and capture specific patterns in the data, leading to a better accuracy. Even though, it tends to overfitting (what can be seen in the loss plots, in which the validation loss flattens regardless of the dropping related with the training loss).

**b)** After training the model for 200 epochs with 3 different dropout values (0.01, 0.25, 0.5), keeping the other parameters at their default value, we obtained:

dropout 0.01

validation accuracy: 0.5762test accuracy: 0.5803

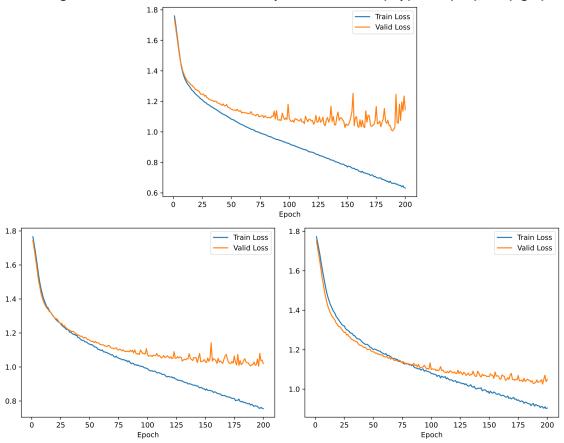
dropout 0.25

validation accuracy: 0.6083test accuracy: 0.6057

dropout 0.5

validation accuracy: 0.5990test accuracy: 0.5960

Training and validation losses for each dropout value - 0.01 (top); 0.25 (left); 0.5 (right)



#### **Analysis:**

Using a low dropout value (e.g. 0.01) only a small fraction of neurons is inactive during training, which translates in a minimal impact in terms of preventing overfitting (this is evident in the loss graph, where for this dropout rate it shows the most pronounced signs of overfitting).

When dropout value is 0.5, half neurons are inactive during each training epoch, besides improving in generalization, it may start to excessively limit the network's capacity to learn patterns, particularly in some cases where the model struggles to converge effectively within each epoch, what is translated in a decreasing of performance. The drop in accuracy from 0.25 enforces this fact.

The best results are from training with dropout 0.25, which may indicate a better trade-off between overfitting and underfitting risks.

c) After training the model for 200 epochs with batch size of 1024 and setting momentum to 0.0 and 0.9, while keeping the other parameters at their default value, we obtained:

• momentum 0.0

validation accuracy: 0.4701test accuracy: 0.4883

momentum 0.9

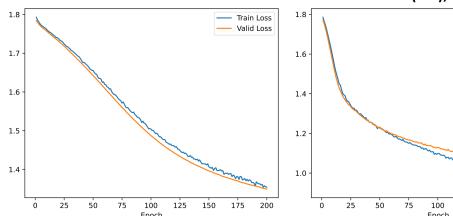
validation accuracy: 0.5990test accuracy: 0.6010

## Train and validation losses for each momentum – 0.0 (left); 0.9 (right)

Train Loss

Valid Loss

150



#### **Analysis:**

It is possible to verify that with a higher momentum it is achieved a better performance.

From the plots it can be seen that the convergence is faster when a momentum is applied. Hence, less epochs are needed for assisting to smaller loss values.

Moreover, the account for the weight of past gradients in the updates helps the algorithm to avoid being stuck in local minimum or saddle points and to get closer to the global minimum, attaining higher accuracy values.

## **Question 3**

1.

2.

Even Question 3 We have: 
$$\hat{y} = n^{-1}k + no$$
and from the previous claim: 
$$k = A_0 \, \hat{p}(x)$$
replacing  $k$  in the  $\hat{y}$  expression: 
$$\hat{y} = n^{-1}(A_0 \, \hat{p}(x)) + no$$

$$= (n^{-1}A_0) \, \hat{p}(x) + [n_0 \, 0 \cdots \, 0] \, \hat{p}(x)$$

$$= (n^{-1}A_0 + [n_0 \, 0 \cdots \, 0]) \, \hat{p}(x)$$
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## **TODO**

IMPORTANT: Please write 1 paragraph indicating clearly what was the contribution of each member of the group in this project. A penalization of 10 points will be applied if this information is missing.