Deep Learning — Assignment 2

Second assignment for the 2023 Deep Learning course (NWI-IMC070) of the Radboud University.

Names:			
Group:			

Instructions:

- Fill in your names and the name of your group.
- Answer the questions and complete the code where necessary.
- Keep your answers brief, one or two sentences is usually enough.
- Re-run the whole notebook before you submit your work.
- Save the notebook as a PDF and submit that in Brightspace together with the .ipynb notebook file.
- The easiest way to make a PDF of your notebook is via File > Print Preview and then
 use your browser's print option to print to PDF.

Objectives

In this assignment you will

- 1. Implement a neural network in PyTorch;
- 2. Use automatic differentiation to compute gradients;
- 3. Experiment with SGD and Adam;
- 4. Experiment with hyperparameter optimization;
- 5. Experiment with regularization techniques.

Before we start, if PyTorch or pandas is not installed, install it now with pip install.

```
In []: # !pip install torch
# !pip install torchvision
# !pip install pandas

In []: %config InlineBackend.figure_formats = ['png']
%matplotlib inline
import numpy as np
import sklearn.datasets
import matplotlib.pyplot as plt
import torch
import time
```

```
import torchvision
import tqdm.notebook as tqdm
import collections
import IPython
import pandas as pd

np.set_printoptions(suppress=True, precision=6, linewidth=200)
plt.style.use('ggplot')

# Fix the seed, so outputs are exactly reproducible
torch.manual_seed(12345);
```

2.1 Implementing a model with PyTorch

In the first assignment, you implemented a neural network from scratch in NumPy. In practice, it is more convenient to use a deep learning framework. In this course, we use PyTorch.

In this example, we will use PyTorch to implement and train a simple neural network.

PyTorch tensors

Similar to NumPy, PyTorch works with multi-dimensional tensors. These can be scalars, vectors, matrices, or have an even higher dimension.

Tensors can be created by converting NumPy arrays, or directly in PyTorch:

```
In []: # create a 10x10 matrix filled with zeros
x = np.zeros([10, 10])
x = torch.tensor(x)
print(x)

# create a 3x5 matrix of ones
x = torch.ones([3, 5])
print(x)
```

Note that the tensors have a datatype (dtype) that defines the type of number in the matrix. This can be 32-bit or 64-bit floating point numbers, or various types of integers. (See the PyTorch documentation for a complete list.)

```
In []: print('float32:', torch.tensor([1., 2., 3.], dtype=torch.float32))
    print('float64:', torch.tensor([1., 2., 3.], dtype=torch.float64))
    print('int:', torch.tensor([1, 2, 3], dtype=torch.int))
    print('long:', torch.tensor([1, 2, 3], dtype=torch.long))
```

It can be important to choose the correct datatype for your tensors, because this influences the precision, the computational cost, or the memory requirements. Some functions also specifically require integers.

Moving data to and from the GPU

If you have a GPU, you can use it to increase the speed of PyTorch computations. To do this, you have to move your data to the GPU by calling to ('cuda'). Afterwards, you can move the results back to the CPU by calling to ('cpu') or cpu().

Note: in MacOS you can use the 'mps' backend instead of 'cuda'.

```
In []: # this only works if you have a GPU available
        if torch.cuda.is available():
            # define a variable on the CPU
            x = torch.ones((3,))
            # move the variable to the GPU
            x = x.to('cuda')
            print('x is now on the GPU:', x)
            # move the variable back to the CPU
            x = x.to('cpu')
            print('x is now on the CPU:', x)
        elif torch.backends.mps.is available():
            # define a variable on the CPU
            x = torch.ones((3,))
            # move the variable to the GPU
            x = x.to('mps')
            print('x is now on the GPU:', x)
            # move the variable back to the CPU
            x = x.to('cpu')
            print('x is now on the CPU:', x)
        else:
            print('It looks like you don\'t have a GPU available.')
```

Note: If you want to run a computation on the GPU, all variables of that function should be moved to the GPU first. If some variables are still on the CPU, PyTorch will throw an error.

If you have a GPU, make the following code run without errors.

```
In []: if torch.cuda.is_available():
    x = torch.tensor([1, 2, 3])
    y = torch.tensor([1, 2, 3], device='cuda')
    print('x is on the CPU:', x)
    print('y is on the GPU:', y)

# this will not work
# TODO: make this computation run on the GPU
    z = x * y
    print(z)
```

To use the GPU when it is available, and fall back to the CPU otherwise, a common trick is to define a global device constant. You can then use tensor.to(device) and torch.tensor(device=device).

```
In [ ]: device = torch.device("cuda" if torch.cuda.is_available() else "mps" if torc
```

Converting back to NumPy

Sometimes, it is useful to convert PyTorch tensors back to NumPy arrays, for example, if you want to plot the performance of your network.

Call detach().cpu().numpy() on the tensor variable to convert the variable to NumPy:

```
In [ ]: x = torch.tensor(5.)
x_in_numpy = x.detach().cpu().numpy()
x_in_numpy
```

Note: detach detaches the tensor from the other computation, by among other things, removing gradient information. cpu transfers the tensor from GPU to CPU if needed. Finally numpy converts from a pytorch tensor to a numpy array, these function very similarly, they just come from different libraries.

Computing the gradients of a simple model

You can use the PyTorch tensors to perform computations, such as the function $y=x\cdot w+b$:

```
In [ ]: w = torch.tensor(2.)
b = torch.tensor(1.)
x = torch.tensor(5.)
print('w:', w)
print('b:', b)
print('x:', x)

y = x * w + b
print(y)
```

If we would like to compute the gradient for the parameters w and b, we could derive and compute them manually (as you did last week):

```
In [ ]: y_grad = 1
    w_grad = y_grad * x
    b_grad = y_grad
    print('w_grad:', w_grad)
    print('b_grad:', b_grad)
```

This can be a lot of work, and it is easy to make mistakes. Fortunately, PyTorch (and other deep learning libraries) can compute these gradients automatically using automatic differentiation.

Computing the gradient automatically

You can compute an automatic gradient as follows:

- 1. Tell PyTorch which variables need a gradient. You can do this by setting requires grad=True when you define the variable.
- 2. Perform the computation.
- 3. Use the backward() function on the result to compute the gradients using backpropagation.
- 4. The grad property of your variables will now contain the gradient.

Have a look at this example, and compare the gradients with the gradients we computed manually:

```
In []: w = torch.tensor(2., requires_grad=True)
b = torch.tensor(1., requires_grad=True)
x = torch.tensor(5.)

# compute the function
y = x * w + b

# compute the gradients, given dy = 1
y.backward()

print('w.grad:', w.grad)
print('b.grad:', b.grad)
# x did not have requires_grad, so no gradient was computed
print('x.grad:', x.grad)
```

This also works for much more complicated functions (and even entire neural networks):

```
In []: w = torch.tensor(2., requires_grad=True)
b = torch.tensor(1., requires_grad=True)
x = torch.tensor(5.)

y = torch.exp(torch.sin(x * w) + b)
y.backward()

print('w.grad:', w.grad)
print('b.grad:', b.grad)
```

Inspect the automatic differentiation history

PyTorch can compute these gradients automatically because it keeps track of the operations that generated the result.

While you don't normally need to do this, you can look inside <code>y.grad_fn</code> to see how it works:

```
In []: w = torch.tensor(2., requires_grad=True)
b = torch.tensor(1., requires_grad=True)
x = torch.tensor(5.)

y = x * w + b
```

```
print('y.grad_fn:', y.grad_fn)
print(' \-->', y.grad_fn.next_functions)
print(' \-->', y.grad_fn.next_functions[0][0].next_functions)
```

The grad fn of y contains a tree that reflects how y was computed:

- the last operation was an addition (x * w) plus b: AddBackward0 knows how to compute the gradient of that;
- one of the inputs to the addition was a multiplication x times w: MulBackward0 computes the gradient;
- eventually, the backpropagation reaches the input variables: AccumulateGrad is
 used to store the gradient in the grad property of each variable.

As long as you use operations for which PyTorch knows the gradient, the backward() function can perform automatic backpropagation and the chain rule to compute the gradients. If you want, can read more about this in the PyTorch autograd tutorial.

2.2 PyTorch neural network with torch.nn (1 point)

The torch.nn module of PyTorch contains a large number of building blocks to construct your own neural network architectures. You will need this in this and future assignments. Have a look at the documentation for torch.nn to see what is available. In this assignment, we will use torch.nn.Linear to build networks with linear layers, as well as some activation and loss functions.

A network module

As a first example, the two-layer network from last week can be implemented like this:

```
In []:
    class WeeklNet(torch.nn.Module):
        def __init__(self):
            super().__init__()

            self.layer1 = torch.nn.Linear(64, 32)
            self.relu = torch.nn.ReLU()
            self.layer2 = torch.nn.Linear(32, 10)

    def forward(self, x):
            x = self.layer1(x)
            x = self.relu(x)
            x = self.layer2(x)
            return x

net = WeeklNet()
```

Observe the following:

A network in PyTorch is usually implemented as a subclass of torch.nn.Module, as
it is here.

- The init function defines the layers that are used in the network.
- The forward function computes the output of the network given one or more inputs (in this case: x).

Notice that there is no final activation function (such as a sigmoid or softmax) at the end of the network. This is not a mistake: we will do this later, by using a loss function that combines the sigmoid/softmax and the loss in one computation, because this is more numerically stable.

Network parameters

Some of the components of the network have parameters, such as the weight and bias in the Linear layers. We can list them using the parameters or named_parameters function:

```
In [ ]: for name, param in net.named_parameters():
    print(name)
    print(param)
    print()
```

As you can see, these parameters have been initialized to non-zero values.

(a) Why are these weights not zero?

(1 point)

TODO: Your answer here.

Shortcut: use torch.nn.Sequential

Quite often, as in our network above, a network architecture consists of a number of layers that are computed one after the other. PyTorch has a special torch.nn.Sequential function to quickly define these networks, without having to define a new class.

For example, the network we implemented earlier can also be written like this:

2.3 A neural network for Fashion-MNIST (12 points)

In this assignment, we will do experiments with the Fashion-MNIST dataset. First, we download the dataset, and create a random training set with 1000 images and a validation

set with 500 images:

Plot some images

The Fashion-MNIST contains images of 28 by 28 pixels, from 10 different classes. In our experiments we flatten the images to a vector with $28 \times 28 = 784$ features.

```
In []: # plot some of the images
    plt.figure(figsize=(10, 2))
    plt.imshow(np.hstack([fashion_train[i][0].reshape(28, 28) for i in range(9)]
    plt.grid(False)
    plt.tight_layout()
    plt.axis('off')
    plt.title('labels: ' + str([fashion_train[i][1] for i in range(9)]));
```

(a) Can you, as a human, distinguish the different classes? Do you think a neural network should be able to learn to do this as well? (1 point)

TODO: Your answer here.

Use the DataLoader to create batches

We will use the DataLoader from PyTorch (see the documentation) to automatically create random batches of 10 images:

```
In [ ]: data_loader = torch.utils.data.DataLoader(fashion_train, batch_size=10, shuf
```

We will use the data loader to loop over all batches in the dataset.

For each batch, we get x, a tensor containing the images, and y, containing the label for each image:

```
In []: for x, y in data_loader:
    print('x.dtype:', x.dtype, 'x.shape:', x.shape)
    print('y.dtype:', y.dtype, ' y.shape:', y.shape)
    # one batch is enough for now
    break
```

Construct a network

We will construct a network that can classify these images.

(b) Implement a network with the following architecture using torch.nn.Sequential:

(2 points)

- Accept flattened inputs: 28 x 28 images mean an input vector with 784 elements.
- Linear hidden layer 1, ReLU activation, output 128 features.
- Linear hidden layer 2, ReLU activation, output 64 features.
- Linear output layer, to 10 classes.
- No final activation function.

```
In [ ]: def build_net():
    # TODO: construct and return the network
    return None
net = build_net()
print(net)
```

(c) Test your network by creating a data loader and computing the output for one batch: (3 points)

Train a network with PyTorch

To train the network, we need a number of components:

- · A network, like the one you just defined.
- A DataLoader to loop over the training samples in small batches.
- A loss function, such as the cross-entropy loss. See the loss functions in the PyTorch documentation.
- An optimizer, such as SGD or Adam: after we use the backward function to compute
 the gradients, the optimizer computes and applies the updates to the weights of the
 network. See the optimization algorithms in the PyTorch documentation.

As an example, the code below implements all of these components and runs a single update step of the network.

(d) Have a look at the code to understand how it works. Then make the following changes: (4 points)

- · Set the batch size to 16
- Use Adam as the optimizer and set the learning rate to 0.01
- · For each minibatch, compute the output of the network
- Compute and optimize the cross-entropy loss

```
In [ ]: # initialize a new instance of the network
        net = build net()
        # construct a data loader for the training set
        data loader = torch.utils.data.DataLoader(fashion train, shuffle=True)
        # initialize the SGD optimizer
        # we pass the list of parameters of the network
        optimizer = torch.optim.SGD(net.parameters(), lr=0)
        # TODO: Initialize cross-entropy loss function
        loss function = ...
        # repeat for multiple epochs
        for epoch in range(10):
            # compute the mean loss and accuracy for this epoch
            loss sum = 0.0
            accuracy sum = 0.0
            steps = 0
            # loop over all minibatches in the training set
            for x, y in data loader:
                # compute the prediction given the input x
                # TODO: compute the output
                output = ...
                # compute the loss by comparing with the target output y
                # TODO: use loss function to compute the loss
                loss = \dots
                # for a one-hot encoding, the output is a score for each class
                # we assign each sample to the class with the highest score
                pred class = torch.argmax(output, dim=1)
                # compute the mean accuracy
                accuracy = torch.mean((pred class == y).to(float))
                # reset all gradients to zero before backpropagation
                optimizer.zero grad()
                # compute the gradient
                loss.backward()
                # use the optimizer to update the parameters
                optimizer.step()
```

```
accuracy_sum += accuracy.detach().cpu().numpy()
    loss_sum += loss.detach().cpu().numpy()
    steps += 1

# print('y:', y)
# print('pred_class:', pred_class)
# print('accuracy:', accuracy)
print('epoch:', epoch,
    'loss:', loss_sum / steps,
    'accuracy:', accuracy_sum / steps)
```

(e) Run the optimization for a few epochs. Does the loss go down? Has the training converged? (1 point)

TODO: Your answer here.

(f) Looking back at the network, we did not include a SoftMax activation function after the last linear layer. But typically you need to use a softmax activation when using cross-entropy loss. Was there a mistake? (1 point)

Hint: Look at the documentation of the cross-entropy loss function. Is the formula there the same as in the slides?

TODO: Your answer here.

(optional) Why do you think the developers of PyTorch did it this way?

OPTIONAL TODO: Your answer here.

2.4 Training code for the rest of this assignment

For the rest of this assignment, we will use a slightly more advanced training function. It runs the training loop for multiple epochs, and at the end of each epoch evaluates the network on the validation set.

Feel free to look inside, but keep in mind that some of this code is only needed to generate the plots in this assignment.

```
# move the network parameters to the gpu, if necessary
net = net.to(device)
# initialize the loss and accuracy history
history = collections.defaultdict(list)
epoch stats, phase = None, None
# initialize the data loaders
data loader = {
    'train':
                torch.utils.data.DataLoader(train, batch size=batch si
    'validation': torch.utils.data.DataLoader(validation, batch size=bat
}
# measure the length of the experiment
start time = time.time()
# some advanced PyTorch to look inside the network and log the outputs
# you don't normally need this, but we use it here for our analysis
def register measure hook(idx, module):
    def hook(module, input, output):
        with torch.no grad():
            # store the mean output values
            epoch stats['%s %d: %s output mean' % (phase, idx, type(modu
                output.mean().detach().cpu().numpy()
            # store the mean absolute output values
            epoch stats['%s %d: %s output abs mean' % (phase, idx, type(
                output.abs().mean().detach().cpu().numpy()
            # store the std of the output values
            epoch_stats['%s %d: %s output std' % (phase, idx, type(modul
                output.std().detach().cpu().numpy()
    module.register forward hook(hook)
# store the output for all layers in the network
for layer idx, layer in enumerate(net):
    register measure hook(layer idx, layer)
# end of the advanced PyTorch code
for epoch in tqdm.tqdm(range(epochs), desc='Epoch', leave=False):
    # initialize the loss and accuracy for this epoch
    epoch stats = collections.defaultdict(float)
    epoch stats['train steps'] = 0
    epoch stats['validation steps'] = 0
    epoch_outputs = {'train': [], 'validation': []}
    # first train on training data, then evaluate on the validation data
    for phase in ('train', 'validation'):
        # switch between train and validation settings
        net.train(phase == 'train')
        epoch steps = 0
        epoch loss = 0
        epoch_accuracy = 0
        # loop over all minibatches
        for x, y in data loader[phase]:
```

```
# move data to gpu, if necessary
            x = x.to(device)
            y = y.to(device)
           # compute the forward pass through the network
            pred y = net(x)
            # compute the current loss and accuracy
            loss = torch.nn.functional.cross entropy(pred y, y)
            pred class = torch.argmax(pred y, dim=1)
            accuracy = torch.mean((pred class == y).to(float))
            # add to epoch loss and accuracy
            epoch stats['%s loss' % phase] += loss.detach().cpu().numpy(
            epoch stats['%s accuracy' % phase] += accuracy.detach().cpu(
            # store outputs for later analysis
            epoch outputs[phase].append(pred y.detach().cpu().numpy())
            # only update the network in the training phase
            if phase == 'train':
                # set gradients to zero
                optimizer.zero grad()
                # backpropagate the gradient through the network
                loss.backward()
                # track the gradient and weight of the first layer
                # (not standard; we only need this for the assignment)
                epoch stats['train mean abs grad'] += \
                    torch.mean(torch.abs(net[0].weight.grad)).detach().c
                epoch stats['train mean abs weight'] += \
                    torch.mean(torch.abs(net[0].weight)).detach().cpu().
                # update the weights
                optimizer.step()
            epoch stats['%s steps' % phase] += 1
        # compute the mean loss and accuracy over all minibatches
        for key in epoch stats:
            if phase in key and not 'steps' in key:
                epoch stats[key] = epoch stats[key] / epoch stats['%s st
                history[key].append(epoch stats[key])
        # count the number of update steps
        history['%s steps' % phase].append((epoch + 1) * epoch stats['%s
        # store the outputs
        history['%s outputs' % phase].append(np.concatenate(epoch output
    history['epochs'].append(epoch)
    history['time'].append(time.time() - start time)
return history
```

```
In [ ]: # helper code to plot our results
        class HistoryPlotter:
            def init (self, plots, table, rows, cols):
                self.plots = plots
                self.table = table
                self.rows = rows
                self.cols = cols
                self.histories = {}
                self.results = []
                self.fig, self.axs = plt.subplots(ncols=cols * len(plots), nrows=row
                                                   sharex='col', sharey='none',
                                                   figsize=(3.5 * cols * len(plots),
                plt.tight layout()
                IPython.display.display(self.fig)
                IPython.display.clear output(wait=True)
            # add the results of an experiment to the plot
            def add(self, title, history, row, col):
                self.histories[title] = history
                self.results.append((title, {key: history[key][-1] for key in self.t
                for plot idx, plot xy in enumerate(self.plots):
                    ax = self.axs[row, col * len(self.plots) + plot idx]
                    for key in plot xy['y']:
                        ax.plot(history[plot_xy['x']], history[key], label=key)
                    if 'accuracy' in plot xy['y'][0]:
                        ax.set ylim([0, 1.01])
                    ax.legend()
                    ax.set xlabel(plot xy['x'])
                    ax.set title(title)
                plt.tight layout()
                IPython.display.clear output(wait=True)
                IPython.display.display(self.fig)
            # print a table of the results for all experiments
            def print table(self):
                df = pd.DataFrame([
                    { 'experiment': title, **{key: row[key] for key in self.table} }
                    for title, row in self.results
                IPython.display.display(df)
            def done(self):
                plt.close()
                self.print table()
```

2.5 Optimization and hyperparameters (10 points)

An important part of training a neural network is hyperparameter optimization: finding good learning rates, minibatch sizes, and other parameters to train an efficient and effective network.

In this part, we will explore some of the most common hyperparameters.

Learning rate with SGD and Adam

First, we will investigate optimizers and learning rates:

- The choice of optimizer: SGD and especially Adam are common choices.
- The learning rate determines the size of the updates by the optimizer.

Optimizing hyperparameters is often a matter of trial-and-error.

We will run an experiment to train our network with the following settings:

- Optimizer: SGD or Adam
- Learning rate: 0.1, 0.01, 0.001, 0.0001
- Minibatch size: 32
- 150 epochs

For each setting, we will plot:

- The train and validation accuracy
- The train and validation loss

We will also print a table with the results of the final epoch.

(a) Run the experiment and have a look at the results.

```
In [ ]: plotter = HistoryPlotter(plots=[{'x': 'epochs', 'y': ['train loss', 'validat']
                                        {'x': 'epochs', 'y': ['train accuracy', 'val
                                 table=['train accuracy', 'validation accuracy'],
                                 rows=4, cols=2)
        epochs = 150
        batch size = 32
        for row, lr in enumerate((0.1, 0.01, 0.001, 0.0001)):
            net = build net()
            optimizer = torch.optim.SGD(net.parameters(), lr=lr)
            history = fit(net, fashion_train, fashion_validation, optimizer=optimize
            plotter.add('SGD lr=%s' % str(lr), history, row=row, col=0)
            net = build net()
            optimizer = torch.optim.Adam(net.parameters(), lr=lr)
            history = fit(net, fashion train, fashion validation, optimizer=optimize
            plotter.add('Adam lr=%s' % str(lr), history, row=row, col=1)
        plotter.done()
```

As you can see, not every combination of hyperparameters works equally well.

(b) Was 150 epochs long enough to train the network with all settings? List the experiments that have/have not yet converged. (2 points)

Experiments that have converged to a good result: TODO

Experiments that need more training: TODO

Other experiments: TODO

(c) How does the learning rate affect the speed of convergence? (1 point)

TODO Your answer here.

(d) A larger learning rate does not always lead to better or faster training. What happened to Adam with a learning rate of 0.1? (1 point)

TODO Your answer here.

(e) It seems that Adam works reasonably well with learning rates 0.01, 0.001, and 0.0001. Can you explain the difference between the three learning curves, in terms of performance, stability, and speed? (2 points)

Final performance: TODO

Stability: TODO

Speed: TODO

Same accuracy, increasing loss

You may have noticed something interesting in the curves for "Adam Ir=0.001": after 10 to 20 epochs, the loss on the validation set starts increasing again, while the accuracy remains the same. How is this possible?

We can find a clue by looking at the output of the network. We will plot the final outputs: the prediction just before the softmax activation function. These values are also called 'logits'.

(f) Run the code below to generate the plots.

```
axs[0, 1].plot(plotter.histories['Adam lr=0.001']['epochs'],
                   plotter.histories['Adam lr=0.001']['train loss'],
                   label='train loss')
    axs[0, 1].plot(plotter.histories['Adam lr=0.001']['epochs'],
                   plotter.histories['Adam lr=0.001']['validation loss'],
                   label='validation loss')
   axs[0, 1].set xlabel('epochs')
   axs[0, 1].set ylabel('loss')
   axs[0, 1].set title('Adam lr=0.001')
   axs[0, 1].legend()
   # plot curve of mean absolute output values
   axs[1, 0].plot(plotter.histories['Adam lr=0.001']['epochs'],
                   plotter.histories['Adam lr=0.001']['train 4: Linear outpu
                   label='output before softmax (training set)')
   axs[1, 0].set xlabel('epochs')
   axs[1, 0].set ylabel('mean absolute output')
   axs[1, 0].set title('Output before softmax (Adam lr=0.001)')
   axs[1, 0].legend()
   # plot distributions of output values
   for epoch in (149, 80, 25, 5, 1):
        axs[1, 1].hist(plotter.histories['Adam lr=0.001']['train outputs'][e
                       label='epoch %d' % epoch)
   axs[1, 1].set xlabel('output before softmax (training set)')
   axs[1, 1].set_ylabel('number of values')
   axs[1, 1].set title('Output before softmax (Adam lr=0.001)')
   axs[1, 1].legend()
   plt.tight layout()
plot output stats()
```

Bottom left: the mean of the final outputs for all training images at different epochs.

Bottom right: histograms showing the distribution of the output values at different epochs.

You should now be able to answer this question:

(g) Why does the accuracy remain stable while the loss keeps increasing? (1 points)

Perhaps you can combine these plots with your knowledge of the softmax activation and cross-entropy loss function to explain this curious behaviour.

TODO: Your answer here.

Minibatch size

Another important hyperparameter is the minibatch size. Sometimes your minibatch size is limited by the available memory in your GPU, but you can often choose different values.

We will run an experiment to train our network with different minibatch sizes:

Minibatch size: 4, 16, 32, 64

We will fix the other hyperparameters to values that worked well in the previous experiment:

- Optimizer: Adam
- Learning rate: 0.0001
- 150 epochs

For each setting, we will plot:

- The train and validation accuracy vs number of epochs
- The train and validation loss vs number of epochs
- The train and validation accuracy vs the number of gradient descent update steps
- The train and validation accuracy vs the training time

We will also print a table with the results of the final epoch.

(h) Run the experiment and have a look at the results.

(i) Why is it useful to look at the number of training steps and the training time? How is this visible in the plots? (1 point)

TODO Your answer here.

(j) What are the effects of making the minibatches very small? (1 point)

TODO Your answer here.

(k) What are the effects of making the minibatches very large? (1 point)

TODO Your answer here.

2.6 Regularization (13 points)

Besides choosing the hyperparameters, we can include other components to improve the training of the model.

In this section, we will experiment with batch normalization, weight decay, and data augmentation.

Batch normalization

Batch normalization can be implemented with the batch normalization modules from torch.nn (documentation).

For a network with 1D feature vectors, you can use torch.nn.BatchNorm1d (documentation).

(a) Construct a network with batch normalization:

(1 point)

Use the same structure as before, but include batchnorm after the hidden linear layers. So we have:

- A linear layer from 784 to 128 features, followed by batchnorm and a ReLU activation.
- A linear layer from 128 to 64 features, followd by batchnorm and ReLU activation.
- A final linear layer from 64 features to 10 outputs, no activation.

```
In [ ]: def build_net_bn():
    # TODO implement the network
net = build_net_bn()
print(net)
```

We will run an experiment to compare a network without batch normalization with a network with batch normalization.

We will fix the other hyperparameters to values that worked well in the previous experiment:

Optimizer: Adam

• Learning rate: 0.0001

Minibatch size: 32

150 epochs

(b) Run the experiment and have a look at the results.

(c) Does batch normalization improve the performance of the network? (1 point)

TODO Your answer here.

(d) Does batch normalization affect the training or convergence speed? (1 point)

TODO Your answer here.

Let us look a bit closer at how batch normalization changes the network.

We will plot some statistics about the values inside the network.

```
In [ ]: def plot layer stats(layers, history):
            fig, axs = plt.subplots(ncols=layers, nrows=2,
                                    figsize=(3.5 * layers, 3 * 2)
            for layer in range(layers):
                i = 0
                for stat in ('mean', 'std'):
                    for phase in ('train', 'validation'):
                        keys = [key for key in history.keys()
                                 if '%s %d:' % (phase, layer) in key and 'output %s'
                        if len(keys) == 1:
                            key = keys[0]
                            ax = axs[i, layer]
                            ax.plot(history['epochs'], history[key], label='%s outpu
                            ax.set xlabel('epochs')
                            ax.legend()
                            ax.set title(key.replace(' output %s' % stat, '').replace
                    i += 1
            plt.tight layout()
```

First, we will plot the statistics of the network without batch normalization.

For each layer, the plots show the mean and standard deviation of the output values of that layer:

```
In [ ]: plot_layer_stats(5, histories['without batchnorm'])
```

We make similar plots for the network with batch normalization: (Note that the number of layers is slightly larger.)

```
In [ ]: plot_layer_stats(7, histories['with batchnorm'])
```

(e) Compare the mean training values in the batch normalization network with those in the network without batch normalization. Can you explain this with how batch normalization works? (1 point)

TODO Your answer here.

(f) Compare the train and validation curves for the batch normalization layers. The training curves are smooth, but the validation curve is noisy. Why does this happen?

(1 point)

TODO Your answer here.

(g) Batch normalization is supposed to normalize the values to $\mu=0$ and $\sigma=1$, but in layer 4, the mean and standard deviation are steadily increasing over time. Why and how does this happen? (1 point)

TODO Your answer here.

Weight decay

The training can also be regularized using weight decay. This option is built-in in many of the PyTorch optimizers (documentation).

We will set up an experiment to investigate how this affects the training of the model.

We use the good settings from before:

Optimizer: Adam

Learning rate: 0.0001

• Minibatch size: 32

150 epochs

and apply L2 weight decay with a factor 0, 0.0001, 0.001, 0.01, or 0.1.

(h) Complete the code below and run the experiment.

(1 point)

```
batch_size = 32
weight_decays = [0, 0.0001, 0.001, 0.01, 0.1]

for row, weight_decay in enumerate(weight_decays):
    net = build_net()
    # TODO: Set up optimizer with the given weight_decay
    history = fit(net, fashion_train, fashion_validation, optimizer=optimize
    plotter.add('weight_decay=%s' % str(weight_decay), history, row=row, col

plotter.done()
```

(i) How can you observe the amount of overfitting in the plots?

(1 point)

TODO Your answer here.

(j) How does weight decay affect the performance of the model in the above experiments? Give an explanation in terms of the amount of overfitting. (1 point)

TODO Your answer here.

Data augmentation

Finally, we will look at data augmentation.

We will run experiments with three types of data augmentation:

- Adding random noise to the pixels, taken from a normal distribution $\mathcal{N}(0,\sigma)$ with σ = 0, 0.01, 0.1, or 0.2.
- Flipping the image horizontally.
- Shifting the image up, down, left, or right by one pixel.

We create a new dataset class that generates noisy images and use this instead of our normal training set.

```
# shift the image by one pixel in the horizontal or vertical directi
if self.shift:
    x = x.reshape(28, 28)
    # shift max one pixel
    shifts = [*torch.randint(-1, 2, (2,)).numpy()]
    x = torch.roll(x, shifts=shifts, dims=(0, 1))
    x = x.flatten()
return x, y
```

We set up an experiment to see if data augmentation improves our results. We use combinations of the three augmentations: noise, flipping, and shifting.

We will train for 250 epochs.

We keep the other settings as before:

Optimizer: AdamLearning rate: 0.0001Minibatch size: 32

(k) Run the experiment and have a look at the results.

```
In [ ]: plotter = HistoryPlotter(plots=[{'x': 'epochs', 'y': ['train loss', 'validat
                                        {'x': 'epochs', 'y': ['train accuracy', 'val
                                 table=['train accuracy', 'validation accuracy', 'ti
                                  rows=8, cols=2)
        epochs = 250
        lr = 0.0001
        batch size = 32
        for row, noise sigma in enumerate((0, 0.01, 0.1, 0.2)):
            for row2, shift in enumerate([False,True]):
                for col, flip horizontal in enumerate([False,True]):
                    noisy fashion train = NoisyDataset(fashion train, noise sigma, f
                    net = build net()
                    optimizer = torch.optim.Adam(net.parameters(), lr=lr)
                    history = fit(net, noisy_fashion_train, fashion_validation, opti
                    label = ('noise=%s' % str(noise sigma)) + \
                            (', shift' if shift else '') + \
                            (', flip horizontal' if flip horizontal else '')
                    plotter.add(label, history, row=row*2+row2, col=col)
        plotter.done()
```

(I) How does data augmentation affect overfitting in the above experiment? Discuss each of the augmentation types. (3 points)

Adding noise: TODO Your answer here.

Horizontal flips: TODO Your answer here.

Shifting: TODO Your answer here.

(m) Why do we have to train the networks with data augmentation a bit longer than networks without data augmentation? (1 points)

TODO: Your answer here.

2.7 Network architecture (5 points)

An often overlooked hyperparameter is the architecture of the neural network itself. Here you can think about the width (the size of each hidden layer) or the depth (the number of layers).

(a) Copy the build_net function from 2.3b and change it to take a parameter for the width of the first hidden layer. (1 point)

The second hidden layer should have half that width, so the network we have been using so far has width=128.

Hint: a // b is the Python notation for integer division rounding down.

```
In [ ]: def build_net(width = 128):
    # TODO: construct and return the network
    return None
```

(b) Set up an experiment to see how the size of the network affects our results.

(1 point)

We keep the other settings as before:

- Optimizer: Adam
- Epochs: 150
- Learning rate: 0.0001
- Minibatch size: 32
- Widths: 16, 32, 64, 128, 256, 512

TODO: Your answer here.

(d) Do you see overfitting for the largest networks? How can you see this from the plots? (1 point)

TODO: Your answer here.

(e) How many parameters are there in a network with width 128? How does that compare to the number of training samples? (1 point)

You don't have to give an exact value, as long as you are in the right order of magnitude it is okay.

(Feel free to write some python code to do computations.)

TODO: Your answer here.

2.8 Discussion (3 points)

(a) Several of the experiments have included a baseline with exactly the same hyperparameters (batch_size=32, weight_decay=0, network_size=128). Are the results exactly the same? What does this tell you about comparing results for picking the best hyperparameters? (2 points)

TODO: Your answer here.

(b) Throughout this assignment we have used a validation set of 500 samples for selecting hyperparameters. Do you think that you will see the same results on an independent test set? Would the best results be obtained with the hyperparameters that are optimal on the validation set? (1 point)

TODO: Your answer here.

The end

Well done! Please double check the instructions at the top before you submit your results.

This assignment has 44 points.

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