Problem Set 4: Neural Networks

This assignment requires a working IPython Notebook installation, which you should already have. If not, please refer to the instructions in the previous problem sets.

In part 2 (programming) of this assignment, you DO NOT need to make any modification code in this IPython Notebook. Instead you will implement your own simple neural network in the mlp.py file. To submit your answers insert mlp.py into this IPython Notebook along with the written answers and submit everything as a single PDF file.

Total: 100 points.

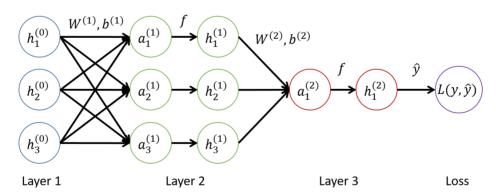
[30pts] Problem 1: Backprop in a simple MLP

This problem asks you to derive all the steps of the backpropagation algorithm for a simple classification network. Consider a fully-connected neural network, also known as a multi-layer perceptron (MLP), with a single hidden layer and a one-node output layer. The hidden and output nodes use an elementwise sigmoid activation function and the loss layer uses cross-entropy loss:

$$f(z) = \frac{1}{1 + exp(-z)}$$

$$L(\hat{y}, y) = -y \ln(\hat{y}) - (1 - y) \ln(1 - \hat{y})$$

The computation graph for an example network is shown below. Note that it has an equal number of nodes in the input and hidden layer (3 each), but, in general, they need not be equal. Also, to make the application of backprop easier, we show the *computation graph* which shows the dot product and activation functions as their own nodes, rather than the usual graph showing a single node for both.



The forward and backward computation for a single example are given below. Note that ⊙ stands for elementwise multiplication.

The forward step is:

```
Require: Network depth, l
Require: \mathbf{W}^{(i)}, i \in \{1, \dots, l\}, the weight matrices of the model
Require: \mathbf{b}^{(i)}, i \in \{1, \dots, l\}, the bias parameters of the model
Require: \mathbf{x}, the input to process
Require: \mathbf{y}, the target output
\mathbf{h}^{(0)} = \mathbf{x}
for k = 1, \dots, l do
\mathbf{a}^{(k)} = \mathbf{b}^{(k)} + \mathbf{W}^{(k)} \mathbf{h}^{(k-1)}
\mathbf{h}^{(k)} = f(\mathbf{a}^{(k)})
end for
\hat{\mathbf{y}} = \mathbf{h}^{(l)}
J = L(\hat{\mathbf{y}}, \mathbf{y}) + \lambda \Omega(\theta)
```

and the backward step is:

```
After the forward computation, compute the gradient on the output layer:  g \leftarrow \nabla_{\hat{y}} J = \nabla_{\hat{y}} L(\hat{y}, y)  for k = l, l - 1, \ldots, 1 do Convert the gradient on the layer's output into a gradient into the prenonlinearity activation (element-wise multiplication if f is element-wise):  g \leftarrow \nabla_{a^{(k)}} J = g \odot f'(a^{(k)})  Compute gradients on weights and biases (including the regularization term, where needed):  \nabla_{b^{(k)}} J = g + \lambda \nabla_{b^{(k)}} \Omega(\theta)  where \partial_{b^{(k)}} J = g h^{(k-1)\top} + \lambda \nabla_{w^{(k)}} \Omega(\theta)  Propagate the gradients w.r.t. the next lower-level hidden layer's activations:  g \leftarrow \nabla_{h^{(k-1)}} J = w^{(k)\top} g  end for
```

Write down each step of the backward pass explicitly for all layers, i.e. compute all intermediate gradients of J above, expressing them as a function of variables x (a vector of inputs), y (a scalar label), $h^{(k)}$ (layer k outputs), $W^{(k)}$ (parameters in layer k), and $b^{(k)}$ (bias parameters of layer k). Notice that $h^{(k)}$ and $b^{(k)}$ can be scalars or vectors depending on layer k, and $W^{(k)}$ can be a matrix or a vector. NOTE: We will assume no regularization, so you can omit the terms involving Ω .

As an example, we will derive the first gradient for you, i.e. the gradient of the cross entropy loss w.r.t. its (scalar) input $\nabla_{\hat{y}} J = \nabla_{\hat{y}} L(\hat{y}, y)$.

$$\nabla_{\hat{y}} L(\hat{y}, y) = \nabla_{\hat{y}} [-y \ln(\hat{y}) - (1 - y) \ln(1 - \hat{y})] = \frac{\hat{y} - y}{(1 - \hat{y})\hat{y}} = \frac{h^{(2)} - y}{(1 - h^{(2)})h^{(2)}}$$

Please derive the remaining gradients, listed below.

Hint: you should substitute the updated values for the gradient *g* in each step and simplify as much as possible.

Hint: Useful information about vectorized chain rule and backpropagation:

If you are struggling with computing the vectorized version of chain rule you may find this example helpful: $\underline{\text{https://web.stanford.edu/class/cs224n/readings/gradient-notes.pdf}}$

(https://web.stanford.edu/class/cs224n/readings/gradient-notes.pdf)

It also contains some helpful shortcuts for computing gradients.

[5pts] Q1.1: Derive $\nabla_{a^{(2)}}J$, where $a^{(2)}$ the (scalar) pre-nonlinearity activation of layer 2.

Hint: to get this scalar value, multiply the partial value g, which should be equal to the $\nabla_{\hat{y}} L(\hat{y}, y)$ that we computed above, with the sigmoid derivative $f'(a^{(2)})$, and note that $f(a^{(2)}) = h^{(2)}$.

* Solution: *

Given that the partial value g is:

$$\nabla_{\hat{y}} L(\hat{y}, y) = \frac{\partial}{\partial \hat{y}} [-y \ln(\hat{y}) - (1 - y) \ln(1 - \hat{y})] = \frac{\hat{y} - y}{(1 - \hat{y})\hat{y}}$$

$$\nabla_{a^{(2)}} f(a^{(2)}) = \frac{\partial}{\partial a^{(2)}} (\frac{1}{1 + exp(-a^{(2)})}) = \frac{exp(-a^{(2)})}{[1 + exp(-a^{(2)})]^2} = \hat{y}(1 - \hat{y})$$

$$\nabla_{a^{(2)}} J = \nabla_{\hat{y}} L(\hat{y}, y) \nabla_{a^{(2)}} f(a^2) = \frac{\hat{y} - y}{(1 - \hat{y})\hat{y}} \hat{y}(1 - \hat{y}) = \hat{y} - y = h^{(2)} - y$$

[5pts] Q1.2: $\nabla_{b}(2)J$

* Solution: *

$$\nabla_b(2J = \nabla_a(2J * \nabla_b(2)(b^{(2)} + W^{(2)}h^{(1)}) = \nabla_a(2J * 1 = h^{(2)} - y$$

[5pts] Q1.3: $\nabla_{W}(2)J$

Hint: this should be a vector, since $W^{(2)}$ is a vector.

* Solution: *

$$\nabla_{W}(2)J = \nabla_{a}(2)J * \nabla_{W}(2)(b^{(2)} + W^{(2)}h^{(1)}) = (h^{(2)} - y)h^{(1)}$$

[5pts] Q1.4: $\nabla_{b}(1)J$

* Solution: *

$$\nabla_{W}(2)J = \nabla_{a}(2)J * \nabla_{h}(1)(b^{(2)} + W^{(2)}h^{(1)}) = (h^{(2)} - y)W^{(2)}$$

[5pts] Q1.5: ∇_b (1)J, ∇_W (1)J

* **Solution:** * Similar to the $\nabla_a(2)J$, we can derive $\nabla_a(1)J$ as:

$$\nabla_{a^{(1)}} J = \nabla_{h^{(1)}} J * \frac{\partial}{\partial a^{(1)}} (\frac{1}{1 + exp(-a^{(1)})}) = (h^{(2)} - y) W^{(2)} h^{(1)} (1 - h^{(1)})$$

Therefore, we have:

$$\nabla_{b}^{(1)}J = \nabla_{a}^{(1)}J * \nabla_{b}^{(1)}(b^{(1)} + W^{(1)}h^{(0)}) = \nabla_{a}^{(1)}J * 1 = (h^{(2)} - y)W^{(2)}h^{(1)}(1 - h^{(1)})$$

$$\nabla_{W}^{(1)}J = \nabla_{a}^{(1)}J * \nabla_{W}^{(1)}(b^{(1)} + W^{(1)}h^{(0)}) = \nabla_{a}^{(1)}J * h^{(0)} = (h^{(2)} - y)W^{(2)}h^{(1)}(1 - h^{(1)})h^{(0)}$$

[5pts] Q1.6 Briefly, explain how the computational speed of backpropagation would be affected if it did not include a forward pass

* Solution: *

If the forward pass was not performed, the local gradient of each layer will not be calculated, which makes it impossible to derive the gradient of every step directly by backpropagation. As a result, we have to calculate the initial local gradient of each layer from the input layer to the output layer, and then update the gradient by backpropagation. Consequently, the computational speed will decrease significantly as the number of layers increase.

[50pts] Problem 2 (Programming): Implementing a simple MLP

In this problem we will develop a neural network with fully-connected layers, or Multi-Layer Perceptron (MLP). We will use it in classification tasks. This question is adapted from Stanford CS231n (http://cs231n.stanford.edu/).

In the current directory, you can find a file mlp.py, which contains the definition for class TwoLayerMLP. As the name suggests, it implements a 2-layer MLP, or MLP with 1 *hidden* layer. You will implement your code in the same file, and call the member functions in this notebook. Below is some initialization. The autoreload command makes sure that mlp.py is periodically reloaded.

In [1]:

```
# setup
import numpy as np
import matplotlib.pyplot as plt
from mlp import TwoLayerMLP

%matplotlib inline
plt.rcParams['figure.figsize'] = (10.0, 8.0) # set default size of plots
plt.rcParams['image.interpolation'] = 'nearest'
plt.rcParams['image.cmap'] = 'gray'

# for auto-reloading external modules
# see http://stackoverflow.com/questions/1907993/autoreload-of-modules-in-ipython
%load_ext autoreload
%autoreload 2

def rel_error(x, y):
    """ returns relative error """
    return np.max(np.abs(x - y) / (np.maximum(1e-8, np.abs(x) + np.abs(y))))
```

Next we initialize a toy model and some toy data, the task is to classify five 4-d vectors.

In [2]:

```
# Create a small net and some toy data to check your implementations.
# Note that we set the random seed for repeatable experiments.
input size = 4
hidden size = 10
num classes = 3
num inputs = 5
def init toy model(actv, std=1e-1):
    np.random.seed(0)
    return TwoLayerMLP(input size, hidden size, num classes, std=std, activation=act
def init toy data():
    np.random.seed(1)
    X = 10 * np.random.randn(num inputs, input size)
    y = np.array([0, 1, 2, 2, 1])
    return X, y
X, y = init toy data()
print('X = ', X)
print()
print('y = ', y)
X = \begin{bmatrix} 16.24345364 & -6.11756414 & -5.28171752 & -10.72968622 \end{bmatrix}
 [ 8.65407629 -23.01538697 17.44811764 -7.61206901]
```

```
 \begin{bmatrix} 8.65407629 & -23.01538697 & 17.44811764 & -7.61206901 \\ [ 3.19039096 & -2.49370375 & 14.62107937 & -20.60140709 ] \\ [ -3.22417204 & -3.84054355 & 11.33769442 & -10.99891267 ] \\ [ -1.72428208 & -8.77858418 & 0.42213747 & 5.82815214 ] ] 
 y = [ 0 1 2 2 1 ]
```

[5pts] Q2.1 Forward pass: Sigmoid

Our 2-layer MLP uses a softmax output layer (**note**: this means that you don't need to apply a sigmoid on the output) and the multiclass cross-entropy loss to perform classification.

Softmax function:

For class j:

$$P(y_{(j)}|x) = \frac{\exp(z_j)}{\sum_{c=1}^{C} \exp(z_c)}$$

Where C is the number of classes and z is class-wise output of the network.

Multiclass cross-entropy loss function:

$$J = \frac{1}{m} \sum_{i=1}^{m} \sum_{c=1}^{C} \left[-y_{(c)} log(P(y_{(c)}|x^{(i)})) \right]$$

 $y_{(c)} = 1$ for the ground truth class and 0 otherwise.

m is the number of inputs in a batch and C is the number of classes.

Please take a look at method TwoLayerMLP.loss in the file mlp.py. This function takes in the data and weight parameters, and computes the class scores (aka logits), the loss L, and the gradients on the parameters.

• Complete the implementation of forward pass (up to the computation of scores) for the sigmoid activation: $\sigma(x) = \frac{1}{1 + exp(-x)}$.

Note 1: Softmax cross entropy loss involves the <u>log-sum-exp operation</u> (https://en.wikipedia.org/wiki/LogSumExp). This can result in numerical underflow/overflow. Read about the solution in the link, and try to understand the calculation of loss in the code.

Note 2: You're strongly encouraged to implement in a vectorized way and avoid using slower for loops. Note that most numpy functions support vector inputs.

Check the correctness of your forward pass below. The difference should be very small (<1e-6).

In [3]:

```
net = init_toy_model('sigmoid')
loss, _ = net.loss(X, y, reg=0.1)
correct_loss = 1.182248
print(loss)
print('Difference between your loss and correct loss:')
print(np.sum(np.abs(loss - correct_loss)))
```

```
1.1822479803941373
Difference between your loss and correct loss:
1.9605862711102873e-08
```

[10pts] Q2.2 Backward pass: Sigmoid

• For sigmoid activation, complete the computation of grads, which stores the gradient of the loss with respect to the variables W1, b1, W2, and b2.

Now debug your backward pass using a numeric gradient check. Again, the differences should be very small.

In [4]:

```
# Use numeric gradient checking to check your implementation of the backward pass.
# If your implementation is correct, the difference between the numeric and
# analytic gradients should be less than 1e-8 for each of W1, W2, b1, and b2.
from utils import eval_numerical_gradient

loss, grads = net.loss(X, y, reg=0.1)

# these should all be very small
for param_name in grads:
    f = lambda W: net.loss(X, y, reg=0.1)[0]
    param_grad_num = eval_numerical_gradient(f, net.params[param_name], verbose=Fals
    print('%s max relative error: %e'%(param_name, rel_error(param_grad_num, grads[r])

W2 max relative error: 1.527006e-09
b2 max relative error: 3.092960e-11
W1 max relative error: 1.126756e-08
```

[5pts] Q2.3 Train the Sigmoid network

b1 max relative error: 2.035406e-06

To train the network we will use stochastic gradient descent (SGD), implemented in TwoLayerNet.train. Then we train a two-layer network on toy data.

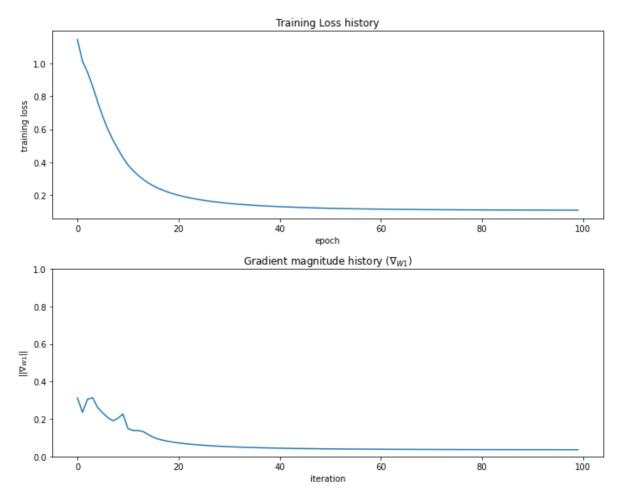
• Implement the prediction function TwoLayerNet.predict, which is called during training to keep track of training and validation accuracy.

You should get the final training loss around 0.1, which is good, but not too great for such a toy problem. One problem is that the gradient magnitude for W1 (the first layer weights) stays small all the time, and the neural net doesn't get much "learning signals". This has to do with the saturation problem of the sigmoid activation function.

In [5]:

```
net = init_toy_model('sigmoid', std=1e-1)
stats = net.train(X, y, X, y,
                  learning_rate=0.5, reg=1e-5,
                  num_epochs=100, verbose=False)
print('Final training loss: ', stats['loss history'][-1])
# plot the loss history and gradient magnitudes
fig, (ax1, ax2) = plt.subplots(2, 1)
ax1.plot(stats['loss history'])
ax1.set xlabel('epoch')
ax1.set ylabel('training loss')
ax1.set title('Training Loss history')
ax2.plot(stats['grad magnitude history'])
ax2.set xlabel('iteration')
ax2.set_ylabel(r'$||\nabla_{W1}||$')
ax2.set title('Gradient magnitude history ' + r'($\nabla {W1}$)')
ax2.set ylim(0,1)
fig.tight layout()
plt.show()
```

Final training loss: 0.10926794610680683



[5pts] Q2.4 Using ReLU activation

The Rectified Linear Unit (ReLU) activation is also widely used: ReLU(x) = max(0, x).

- Complete the implementation for the ReLU activation (forward and backward) in mlp.py.
- · Train the network with ReLU, and report your final training loss.

Make sure you first pass the numerical gradient check on toy data.

In [6]:

```
net = init_toy_model('relu', std=le-1)
loss, grads = net.loss(X, y, reg=0.1)
print('loss = ', loss) # correct_loss = 1.320973

# The differences should all be very small
print('checking gradients')
for param_name in grads:
    f = lambda W: net.loss(X, y, reg=0.1)[0]
    param_grad_num = eval_numerical_gradient(f, net.params[param_name], verbose=Fals
    print('%s max relative error: %e'%(param_name, rel_error(param_grad_num, grads[param_name]))
```

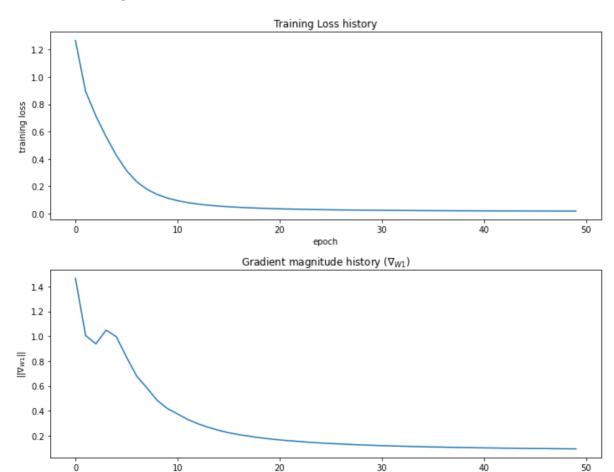
```
loss = 1.3037878913298206
checking gradients
W2 max relative error: 3.440708e-09
b2 max relative error: 4.447646e-11
W1 max relative error: 3.561318e-09
b1 max relative error: 2.738421e-09
```

Now that it's working, let's train the network. Does the net get stronger learning signals (i.e. gradients) this time? Report your final training loss.

In [7]:

```
net = init toy model('relu', std=1e-1)
stats = net.train(X, y, X, y,
                  learning rate=0.1, reg=1e-5,
                  num epochs=50, verbose=False)
print('Final training loss: ', stats['loss_history'][-1])
# plot the loss history
fig, (ax1, ax2) = plt.subplots(2, 1)
ax1.plot(stats['loss history'])
ax1.set xlabel('epoch')
ax1.set ylabel('training loss')
ax1.set_title('Training Loss history')
ax2.plot(stats['grad magnitude history'])
ax2.set xlabel('iteration')
ax2.set ylabel(r'$||\nabla {W1}||$')
ax2.set_title('Gradient magnitude history ' + r'($\nabla_{W1}$)')
fig.tight layout()
plt.show()
```

Final training loss: 0.0178562204869839



iteration

Application to a real Problem

Load MNIST data

Now that you have implemented a two-layer network that works on toy data, let's try some real data. The MNIST dataset is a standard machine learning benchmark. It consists of 70,000 grayscale handwritten digit images, which we split into 50,000 training, 10,000 validation and 10,000 testing. The images are of size 28x28, which are flattened into 784-d vectors.

Note 1: the function <code>get_MNIST_data</code> requires the <code>scikit-learn</code> package. If you previously did anaconda installation to set up your Python environment, you should already have it. Otherwise, you can install it following the instructions here: http://scikit-learn.org/stable/install.html (http://scikit-learn.org/stable/install.html)

Note 2: If you encounter a HTTP 500 error, that is likely temporary, just try again.

Note 3: Ensure that the downloaded MNIST file is 55.4MB (smaller file-sizes could indicate an incomplete download - which is possible)

In [8]:

```
# load MNIST
from utils import get MNIST data
X_train, y_train, X_val, y_val, X_test, y_test = get_MNIST_data()
print('Train data shape: ', X_train.shape)
print('Train labels shape: ', y_train.shape)
print('Validation data shape: '
                               , X val.shape)
print('Validation labels shape: ', y_val.shape)
print('Test data shape: ', X test.shape)
print('Test labels shape: '
                           ', y_test.shape)
                   (50000, 784)
Train data shape:
Train labels shape: (50000,)
                       (10000, 784)
Validation data shape:
Validation labels shape: (10000,)
```

Train a network on MNIST

Test data shape: (10000, 784)
Test labels shape: (10000,)

We will now train a network on MNIST with 64 hidden units in the hidden layer. We train it using SGD, and decrease the learning rate with an exponential rate over time; this is achieved by multiplying the learning rate with a constant factor <code>learning_rate_decay</code> (which is less than 1) after each epoch. In effect, we are using a high learning rate initially, which is good for exploring the solution space, and using lower learning rates later to encourage convergence to a local minimum (or saddle.point (http://www.offconvex.org/2016/03/22/saddle.points/), which may happen more often).

Train your MNIST network with 2 different activation functions: sigmoid and ReLU.

We first define some variables and utility functions. The plot_stats function plots the histories of gradient magnitude, training loss, and accuracies on the training and validation sets. The show_net_weights function visualizes the weights learned in the first layer of the network. In most neural networks trained on visual data, the first layer weights typically show some visible structure when visualized. Both functions help you to diagnose the training process.

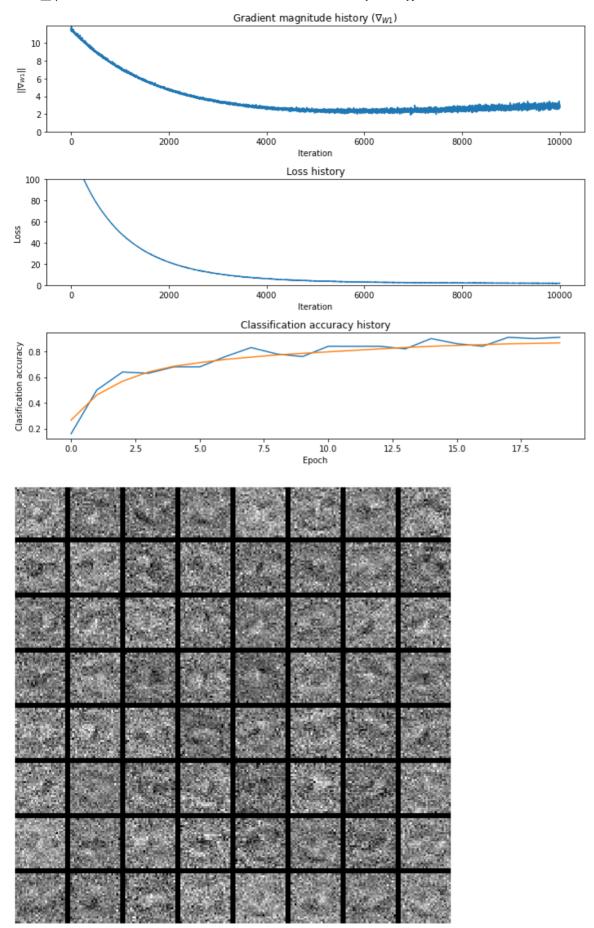
In [9]:

```
input size = 28 * 28
hidden size = 64
num classes = 10
# Plot the loss function and train / validation accuracies
def plot stats(stats):
    fig, (ax1, ax2, ax3) = plt.subplots(3, 1)
    ax1.plot(stats['grad_magnitude_history'])
    ax1.set title('Gradient magnitude history ' + r'$(\nabla {W1})$')
    ax1.set xlabel('Iteration')
    ax1.set ylabel(r'$|\nabla {W1}||$')
    ax1.set ylim(0, np.minimum(100,np.max(stats['grad magnitude history'])))
    ax2.plot(stats['loss history'])
    ax2.set title('Loss history')
    ax2.set xlabel('Iteration')
    ax2.set ylabel('Loss')
    ax2.set ylim(0, 100)
    ax3.plot(stats['train acc history'], label='train')
    ax3.plot(stats['val acc history'], label='val')
    ax3.set title('Classification accuracy history')
    ax3.set xlabel('Epoch')
    ax3.set ylabel('Clasification accuracy')
    fig.tight_layout()
    plt.show()
# Visualize the weights of the network
from utils import visualize grid
def show_net_weights(net):
    W1 = net.params['W1']
    W1 = W1.T.reshape(-1, 28, 28)
    plt.imshow(visualize grid(W1, padding=3).astype('uint8'))
    plt.gca().axis('off')
    plt.show()
```

[10pts] Q2.5 Train a Sigmoid network

In [10]:

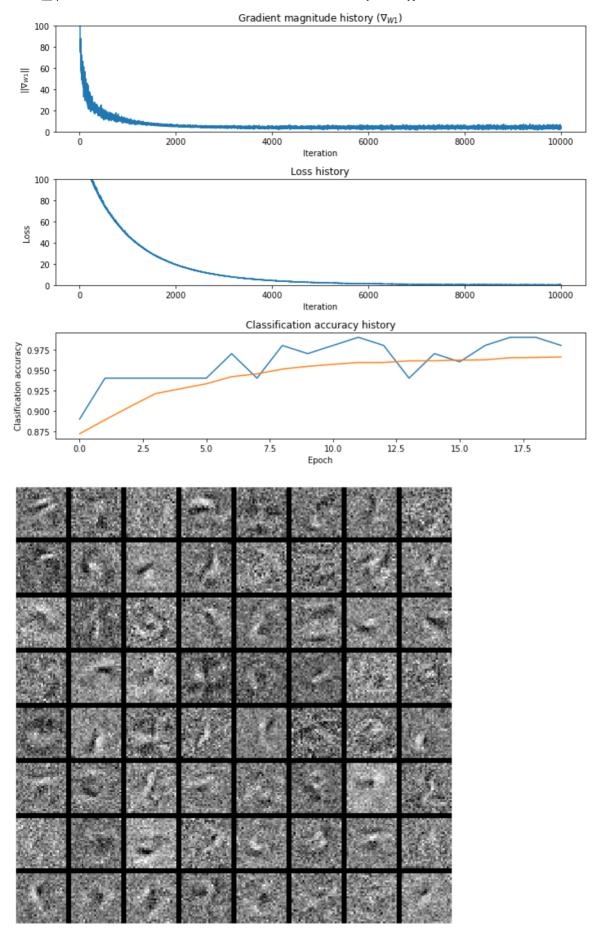
```
sigmoid net = TwoLayerMLP(input size, hidden size, num classes, activation='sigmoid
# Train the network
sigmoid stats = sigmoid net.train(X train, y train, X val, y val,
                                  num epochs=20, batch size=100,
                                  learning rate=1e-3, learning rate decay=0.95,
                                  req=0.5, verbose=True)
# Predict on the training set
train acc = (sigmoid net.predict(X_train) == y_train).mean()
print('Sigmoid final training accuracy: ', train acc)
# Predict on the validation set
val acc = (sigmoid net.predict(X val) == y val).mean()
print('Sigmoid final validation accuracy: ', val acc)
# Predict on the test set
test acc = (sigmoid net.predict(X test) == y test).mean()
print('Sigmoid test accuracy: ', test acc)
# show stats and visualizations
plot stats(sigmoid stats)
show net weights(sigmoid net)
/Users/simple/Desktop/pset4/mlp.py:88: RuntimeWarning: overflow encoun
tered in exp
  hidden = 1 / (1 + np.exp(-z1))
Epoch 1: loss 79.040004, train acc 0.160000, val acc 0.266300
Epoch 2: loss 49.814996, train acc 0.500000, val acc 0.461100
Epoch 3: loss 32.419904, train acc 0.640000, val acc 0.568300
Epoch 4: loss 21.756599, train_acc 0.630000, val_acc 0.639700
Epoch 5: loss 15.148895, train acc 0.680000, val acc 0.685000
Epoch 6: loss 10.909900, train acc 0.680000, val acc 0.712600
Epoch 7: loss 8.078106, train_acc 0.760000, val_acc 0.737900
Epoch 8: loss 6.166522, train_acc 0.830000, val_acc 0.755600
Epoch 9: loss 4.948016, train acc 0.780000, val acc 0.772900
Epoch 10: loss 4.113118, train acc 0.760000, val acc 0.785000
Epoch 11: loss 3.455138, train_acc 0.840000, val_acc 0.797000
Epoch 12: loss 3.026239, train acc 0.840000, val acc 0.808100
Epoch 13: loss 2.702231, train acc 0.840000, val acc 0.819600
Epoch 14: loss 2.438965, train acc 0.820000, val acc 0.830900
Epoch 15: loss 2.258613, train_acc 0.900000, val_acc 0.839900
Epoch 16: loss 2.166625, train acc 0.860000, val acc 0.846800
Epoch 17: loss 2.098843, train acc 0.840000, val acc 0.852800
Epoch 18: loss 1.975990, train acc 0.910000, val acc 0.859300
Epoch 19: loss 1.898398, train acc 0.900000, val acc 0.862300
Epoch 20: loss 1.876564, train acc 0.910000, val acc 0.866400
Sigmoid final training accuracy: 0.8721
Sigmoid final validation accuracy: 0.8664
Sigmoid test accuracy: 0.8639
```



[10pts] Q2.6 Train a ReLU network

In [11]:

```
relu_net = TwoLayerMLP(input_size, hidden_size, num_classes, activation='relu', std=
# Train the network
relu stats = relu net.train(X train, y train, X val, y val,
                            num epochs=20, batch size=100,
                            learning rate=1e-3, learning rate decay=0.95,
                            reg=0.5, verbose=True)
# Predict on the training set
train acc = (relu net.predict(X train) == y train).mean()
print('ReLU final training accuracy: ', train acc)
# Predict on the validation set
val_acc = (relu_net.predict(X_val) == y_val).mean()
print('ReLU final validation accuracy: ', val acc)
# Predict on the test set
test acc = (relu net.predict(X test) == y test).mean()
print('ReLU test accuracy: ', test acc)
# show stats and visualizations
plot stats(relu stats)
show net weights(relu net)
/Users/simple/Desktop/pset4/mlp.py:268: RuntimeWarning: overflow encou
ntered in true divide
  y pred = np.argmax(np.exp(loss1) / np.exp(np.sum(loss1, axis=1)).res
hape(-1,1),axis=1)
Epoch 1: loss 77.112349, train acc 0.890000, val acc 0.872100
Epoch 2: loss 47.480978, train acc 0.940000, val acc 0.889000
Epoch 3: loss 29.866142, train_acc 0.940000, val_acc 0.905400
Epoch 4: loss 19.525541, train acc 0.940000, val acc 0.921200
Epoch 5: loss 13.036480, train acc 0.940000, val acc 0.927200
Epoch 6: loss 8.991456, train_acc 0.940000, val_acc 0.933200
Epoch 7: loss 6.142489, train acc 0.970000, val acc 0.941800
Epoch 8: loss 4.524229, train acc 0.940000, val acc 0.945300
Epoch 9: loss 3.237475, train acc 0.980000, val acc 0.951100
Epoch 10: loss 2.411581, train acc 0.970000, val acc 0.954500
Epoch 11: loss 1.850353, train acc 0.980000, val acc 0.956900
Epoch 12: loss 1.393945, train acc 0.990000, val acc 0.959200
Epoch 13: loss 1.163271, train_acc 0.980000, val_acc 0.959200
Epoch 14: loss 0.994304, train_acc 0.940000, val acc 0.961200
Epoch 15: loss 0.816019, train acc 0.970000, val acc 0.961300
Epoch 16: loss 0.682771, train acc 0.960000, val acc 0.962100
Epoch 17: loss 0.626259, train_acc 0.980000, val_acc 0.962400
Epoch 18: loss 0.488502, train acc 0.990000, val acc 0.965000
Epoch 19: loss 0.450650, train acc 0.990000, val acc 0.965300
Epoch 20: loss 0.405613, train acc 0.980000, val acc 0.965900
ReLU final training accuracy: 0.973
ReLU final validation accuracy: 0.9659
ReLU test accuracy: 0.9655
```



[5pts] Q2.7

Which activation function would you choose in practice? Why?

* Solution: *

I will choose the "Relu" activation function.

- (1) For one thing, the "Relu" is more computational efficient since it does not contain complex math operations. On the contrary, the exponential operation in "sigmoid" costs a lot more time, especially when the scale of data is large.
- (2) Using the "Relu function" can make the converge rate more stable so that we can avoid the Vanishing Gradient Problem. However, when the number of network layers grow, the "sigmoid" function in the saturable region is too flat that the gradient will be approaching zero, which makes it hard to converge.

[20pts] Problem 3: Simple Regularization Methods

You may have noticed the <code>reg</code> parameter in <code>TwoLayerMLP.loss</code>, controlling "regularization strength". In learning neural networks, aside from minimizing a loss function $\mathcal{L}(\theta)$ with respect to the network parameters θ , we usually explicitly or implicitly add some regularization term to reduce overfitting. A simple and popular regularization strategy is to penalize some *norm* of θ .

[10pts] Q3.1: L2 regularization

We can penalize the L2 norm of θ : we modify our objective function to be $\mathcal{L}(\theta) + \lambda \|\theta\|^2$ where λ is the weight of regularization. We will minimize this objective using gradient descent with step size η . Derive the update rule: at time t+1, express the new parameters θ_{t+1} in terms of the old parameters θ_t , the gradient $g_t = \frac{\partial \mathcal{L}}{\partial \theta_t}$, η , and λ .

* Solution: *

Given the objective function $\mathcal{L}(\theta) + \lambda \|\theta\|^2$:

$$\theta_{t+1} = \theta_t - \eta (\frac{\partial}{\partial \theta_t} [\mathcal{L}(\theta) + \lambda \|\theta\|^2])$$

$$= \theta_t - \eta (\frac{\partial \mathcal{L}}{\partial \theta_t}) - \eta \frac{\partial}{\partial \theta_t} \lambda \|\theta\|^2$$

$$= \theta_t - \eta g_t - 2\eta \lambda \theta_t$$

[10pts] Q3.2: L1 regularization

Now let's consider L1 regularization: our objective in this case is $\mathcal{L}(\theta) + \lambda \|\theta\|_1$. Derive the update rule.

(Technically this becomes *Sub-Gradient* Descent since the L1 norm is not differentiable at 0. But practically it is usually not an issue.)

* Solution: *

Similar to the L2 regularization, given the objective function $\mathcal{L}(\theta) + \lambda \|\theta\|_1$:

$$\theta_{t+1} = \theta_t - \eta(\frac{\partial}{\partial \theta_t} [\mathcal{L}(\theta) + \lambda \|\theta\|_1])$$

$$= \theta_t - \eta(\frac{\partial \mathcal{L}}{\partial \theta_t}) - \eta \frac{\partial}{\partial \theta_t} \lambda \|\theta\|_1$$

$$= \theta_t - \eta g_t - \eta \lambda sign(\theta_t)$$

The mlp.py is shown below.

In []:

```
import numpy as np
import matplotlib.pyplot as plt
class TwoLayerMLP(object):
 A two-layer fully-connected neural network. The net has an input dimension of
 N, a hidden layer dimension of H, and performs classification over C classes.
 We train the network with a softmax loss function and L2 regularization on the
 weight matrices. The network uses a ReLU nonlinearity after the first fully
 connected layer.
 In other words, the network has the following architecture:
  input - fully connected layer - ReLU - fully connected layer - softmax
  The outputs of the second fully-connected layer are the scores for each class.
 def __init__(self, input_size, hidden_size, output_size, std=1e-4, activation='rel
    Initialize the model. Weights are initialized to small random values and
    biases are initialized to zero. Weights and biases are stored in the
    variable self.params, which is a dictionary with the following keys:
   W1: First layer weights; has shape (D, H)
    b1: First layer biases; has shape (H,)
    W2: Second layer weights; has shape (H, C)
   b2: Second layer biases; has shape (C,)
    Inputs:
    - input size: The dimension D of the input data.
    - hidden size: The number of neurons H in the hidden layer.
    - output size: The number of classes C.
    self.params = {}
    self.params['W1'] = std * np.random.randn(input size, hidden size)
    self.params['b1'] = np.zeros(hidden size)
    self.params['W2'] = std * np.random.randn(hidden_size, output_size)
    self.params['b2'] = np.zeros(output_size)
    self.activation = activation
  def loss(self, X, y=None, reg=0.0):
    Compute the loss and gradients for a two layer fully connected neural
    network.
    Inputs:
    - X: Input data of shape (N, D). Each X[i] is a training sample.
    - y: Vector of training labels. y[i] is the label for X[i], and each y[i] is
     an integer in the range 0 \le y[i] \le C. This parameter is optional; if it
      is not passed then we only return scores, and if it is passed then we
      instead return the loss and gradients.
    - reg: Regularization strength.
   Returns:
    If y is None, return a matrix scores of shape (N, C) where scores[i, c] is
    the score for class c on input X[i].
```

```
If y is not None, instead return a tuple of:
- loss: Loss (data loss and regularization loss) for this batch of training
 samples.
- grads: Dictionary mapping parameter names to gradients of those parameters
 with respect to the loss function; has the same keys as self.params.
# Unpack variables from the params dictionary
W1, b1 = self.params['W1'], self.params['b1']
W2, b2 = self.params['W2'], self.params['b2']
_{\text{, C}} = W2.shape
N, D = X.shape
# Compute the forward pass
# write your own code where you see [PLEASE IMPLEMENT]
# Perform the forward pass, computing the class scores for the input.
# Store the result in the scores variable, which should be an array of
# shape (N, C).
z1 = np.dot(X, W1) + b1 # 1st layer activation, N*H
# 1st layer nonlinearity, N*H
if self.activation is 'relu':
   # [PLEASE IMPLEMENT]
   hidden = np.maximum(0, z1)
elif self.activation is 'sigmoid':
   # [PLEASE IMPLEMENT]
   hidden = 1 / (1 + np.exp(-z1))
else:
   raise ValueError('Unknown activation type')
# [PLEASE IMPLEMENT] 2nd layer activation, N*C
# hint: involves W2, b2
scores = (hidden @ W2) + b2
END OF YOUR CODE
# If the targets are not given then jump out, we're done
if y is None:
 return scores
# cross-entropy loss with log-sum-exp
A = np.max(scores, axis=1) \# N*1
F = np.exp(scores - A.reshape(N, 1)) # N*C
P = F / np.sum(F, axis=1).reshape(N, 1) # N*C
loss = np.mean(-np.choose(y, scores.T) + np.log(np.sum(F, axis=1)) + A)
# add regularization terms
loss += 0.5 * reg * np.sum(W1 * W1)
loss += 0.5 * reg * np.sum(W2 * W2)
# Backward pass: compute gradients
grads = {}
# write your own code where you see [PLEASE IMPLEMENT]
# Compute the backward pass, computing the derivatives of the weights
# and biases. Store the results in the grads dictionary. For example,
# grads['W1'] should store the gradient on W1, and be a matrix of same size.
```

```
# You should define the hidden variable in the part where you implement the
 # different activation functions. "hidden" is the output after you apply
 # the activation function for the hidden layer.
 # Hint: you should apply different activation functions on "z1" and get "hidden"
  # output layer
 dscore = P # [PLEASE IMPLEMENT] partial derivative of loss wrt. the logits (dL)
  for i in range(N):
     dscore[i,y[i]] == 1
 dW2 = np.dot(hidden.T, dscore)/N # partial derivative of loss wrt. W2
 db2 = np.mean(dscore, axis=0) # partial derivation of loss wrt. b2
  # hidden layer
 dhidden = dscore @ (W2.T)
  if self.activation is 'relu':
     # [PLEASE IMPLEMENT]
     dz1 = dhidden
     for i in range(dz1.shape[0]):
         for j in range(dz1.shape[1]):
            if hidden[i,j] <= 0:</pre>
                dz1[i][j] = 0
     for i in range(dz1.shape[0]):
         for j in range(dz1.shape[1]):
            if z1[i,j] <= 0:</pre>
                dz1[i][j] = 0
 elif self.activation is 'sigmoid':
     # [PLEASE IMPLEMENT]
     dz1 = dhidden * hidden * (1-hidden)
     raise ValueError('Unknown activation type')
  # first layer
 dW1 = (X.T) @ dz1 / N # [PLEASE IMPLEMENT]
 db1 = np.mean(dz1, axis=0) # [PLEASE IMPLEMENT]
  END OF YOUR CODE
  grads['W2'] = dW2 + reg*W2
  grads['b2'] = db2
 grads['W1'] = dW1 + reg*W1
 grads['b1'] = db1
 return loss, grads
def train(self, X, y, X_val, y_val,
         learning rate=1e-3, learning rate decay=0.95,
         reg=1e-5, num epochs=10,
         batch size=200, verbose=False):
 Train this neural network using stochastic gradient descent.
 Inputs:
 - X: A numpy array of shape (N, D) giving training data.
 - y: A numpy array f shape (N,) giving training labels; y[i] = c means that
   X[i] has label c, where 0 <= c < C.
  - X val: A numpy array of shape (N val, D) giving validation data.
```

```
- y val: A numpy array of shape (N val,) giving validation labels.
- learning rate: Scalar giving learning rate for optimization.
- learning rate decay: Scalar giving factor used to decay the learning rate
  after each epoch.
- reg: Scalar giving regularization strength.
- num iters: Number of steps to take when optimizing.
- batch size: Number of training examples to use per step.
- verbose: boolean; if true print progress during optimization.
num train = X.shape[0]
iterations per epoch = max(int(num train / batch size), 1)
epoch num = 0
# Use SGD to optimize the parameters in self.model
loss history = []
grad magnitude history = []
train acc history = []
val acc history = []
np.random.seed(1)
for epoch in range(num epochs):
    # fixed permutation (within this epoch) of training data
    perm = np.random.permutation(num train)
    # go through minibatches
    for it in range(iterations per epoch):
        X batch = None
        y_batch = None
        # Create a random minibatch
        idx = perm[it*batch size:(it+1)*batch size]
        X \text{ batch} = X[idx, :]
        y_batch = y[idx]
        # Compute loss and gradients using the current minibatch
        loss, grads = self.loss(X batch, y=y batch, reg=reg)
        loss history.append(loss)
        # do gradient descent
        for param in self.params:
            self.params[param] -= grads[param] * learning rate
        # record gradient magnitude (Frobenius) for W1
        grad magnitude history.append(np.linalg.norm(grads['W1']))
    # Every epoch, check train and val accuracy and decay learning rate.
    # Check accuracy
    train acc = (self.predict(X batch) == y batch).mean()
    val acc = (self.predict(X val) == y val).mean()
    train_acc_history.append(train_acc)
    val_acc_history.append(val_acc)
    if verbose:
        print('Epoch %d: loss %f, train acc %f, val acc %f'%(
            epoch+1, loss, train acc, val acc))
    # Decay learning rate
    learning_rate *= learning_rate_decay
return {
  'loss history': loss history,
  'grad_magnitude_history': grad_magnitude_history,
```

```
'train acc history': train acc history,
   'val acc history': val acc history,
def predict(self, X):
 Use the trained weights of this two-layer network to predict labels for
 data points. For each data point we predict scores for each of the C
 classes, and assign each data point to the class with the highest score.
 Inputs:
 - X: A numpy array of shape (N, D) giving N D-dimensional data points to
   classify.
 Returns:
 - y_pred: A numpy array of shape (N,) giving predicted labels for each of
   the elements of X. For all i, y_pred[i] = c means that X[i] is predicted
   to have class c, where 0 <= c < C.
 0.00
 # [PLEASE IMPLEMENT]
 # hint: it should be very easy
 loss1 = self.loss(X)
 y pred = np.argmax(np.exp(loss1) / np.exp(np.sum(loss1, axis=1)).reshape(-1,1), a
 return y pred
```