Batch Normalization

In this notebook, you will implement the batch normalization layers of a neural network to increase its performance. If you have any confusion, please review the details of batch normalization from the lecture notes.

CS231n has built a solid API for building these modular frameworks and training them, and we will use their very well implemented framework as opposed to "reinventing the wheel." This includes using their Solver, various utility functions, and their layer structure. This also includes nndl.fc_net, nndl.layers, and nndl.layer_utils. As in prior assignments, we thank Serena Yeung & Justin Johnson for permission to use code written for the CS 231n class (cs231n.stanford.edu).

```
In [1]: ## Import and setups
        import time
        import numpy as np
        import matplotlib.pyplot as plt
        from nndl.fc net import *
        from nndl.layers import *
        from cs231n.data utils import get CIFAR10 data
        from cs231n.gradient_check import eval_numerical_gradient, eval_numerical_gradien
        t array
        from cs231n.solver import Solver
        %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))))
In [2]: # Load the (preprocessed) CIFAR10 data.
        data = get CIFAR10 data()
        for k in data.keys():
          print('{}: {} '.format(k, data[k].shape))
        X train: (49000, 3, 32, 32)
        y train: (49000,)
        X_val: (1000, 3, 32, 32)
        y_val: (1000,)
        X test: (1000, 3, 32, 32)
        y test: (1000,)
```

Batchnorm forward pass

Implement the training time batchnorm forward pass, batchnorm_forward, in nndl/layers.py. After that, test your implementation by running the following cell.

```
In [3]: # Check the training-time forward pass by checking means and variances
        # of features both before and after batch normalization
        # Simulate the forward pass for a two-layer network
        N, D1, D2, D3 = 200, 50, 60, 3
        X = np.random.randn(N, D1)
        W1 = np.random.randn(D1, D2)
        W2 = np.random.randn(D2, D3)
        a = np.maximum(0, X.dot(W1)).dot(W2)
        print('Before batch normalization:')
        print(' means: ', a.mean(axis=0))
        print(' stds: ', a.std(axis=0))
        # Means should be close to zero and stds close to one
        print('After batch normalization (gamma=1, beta=0)')
        a norm, = batchnorm forward(a, np.ones(D3), np.zeros(D3), {'mode': 'train'})
        print(' mean: ', a norm.mean(axis=0))
        print(' std: ', a_norm.std(axis=0))
        # Now means should be close to beta and stds close to gamma
        gamma = np.asarray([1.0, 2.0, 3.0])
        beta = np.asarray([11.0, 12.0, 13.0])
        a_norm, _ = batchnorm_forward(a, gamma, beta, {'mode': 'train'})
        print('After batch normalization (nontrivial gamma, beta)')
        print(' means: ', a norm.mean(axis=0))
        print(' stds: ', a_norm.std(axis=0))
        Before batch normalization:
          means: [-32.01112059 -9.94546852 -41.84035895]
          stds: [35.3568188 32.19471298 34.01701357]
        After batch normalization (gamma=1, beta=0)
          mean: [-4.20774526e-16 8.71264866e-17 -7.19424520e-16]
          std: [1. 1. 1.]
        After batch normalization (nontrivial gamma, beta)
          means: [11. 12. 13.]
          stds: [1.
                             1.99999999 2.99999999]
```

Implement the testing time batchnorm forward pass, batchnorm_forward, in nndl/layers.py. After that, test your implementation by running the following cell.

```
In [4]: # Check the test-time forward pass by running the training-time
        # forward pass many times to warm up the running averages, and then
        # checking the means and variances of activations after a test-time
        # forward pass.
        N, D1, D2, D3 = 200, 50, 60, 3
        W1 = np.random.randn(D1, D2)
        W2 = np.random.randn(D2, D3)
        bn param = {'mode': 'train'}
        gamma = np.ones(D3)
        beta = np.zeros(D3)
        for t in np.arange(50):
          X = np.random.randn(N, D1)
          a = np.maximum(0, X.dot(W1)).dot(W2)
          batchnorm forward(a, gamma, beta, bn param)
        bn param['mode'] = 'test'
        X = np.random.randn(N, D1)
        a = np.maximum(0, X.dot(W1)).dot(W2)
        a_norm, _ = batchnorm_forward(a, gamma, beta, bn_param)
        # Means should be close to zero and stds close to one, but will be
        # noisier than training-time forward passes.
        print('After batch normalization (test-time):')
        print(' means: ', a_norm.mean(axis=0))
        print(' stds: ', a_norm.std(axis=0))
```

```
After batch normalization (test-time):
means: [-0.00385673 0.04523241 -0.05831553]
stds: [1.10601571 1.08595284 1.00428327]
```

Batchnorm backward pass

Implement the backward pass for the batchnorm layer, batchnorm_backward in nndl/layers.py. Check your implementation by running the following cell.

```
In [5]: # Gradient check batchnorm backward pass
        N, D = 4, 5
        x = 5 * np.random.randn(N, D) + 12
        gamma = np.random.randn(D)
        beta = np.random.randn(D)
        dout = np.random.randn(N, D)
        bn param = {'mode': 'train'}
        fx = lambda x: batchnorm forward(x, gamma, beta, bn param)[0]
        fg = lambda a: batchnorm forward(x, gamma, beta, bn param)[0]
        fb = lambda b: batchnorm forward(x, gamma, beta, bn param)[0]
        dx num = eval numerical gradient array(fx, x, dout)
        da num = eval numerical gradient array(fg, gamma, dout)
        db num = eval numerical gradient array(fb, beta, dout)
        , cache = batchnorm forward(x, gamma, beta, bn param)
        dx, dgamma, dbeta = batchnorm backward(dout, cache)
        print('dx error: ', rel_error(dx_num, dx))
        print('dgamma error: ', rel error(da num, dgamma))
        print('dbeta error: ', rel_error(db_num, dbeta))
```

dx error: 1.3992698334533834e-08 dgamma error: 5.746786090518256e-12 dbeta error: 6.2021112657457505e-12

Implement a fully connected neural network with batchnorm layers

Modify the FullyConnectedNet() class in nndl/fc_net.py to incorporate batchnorm layers. You will need to modify the class in the following areas:

- (1) The gammas and betas need to be initialized to 1's and 0's respectively in __init__.
- (2) The batchnorm_forward layer needs to be inserted between each affine and relu layer (except in the output layer) in a forward pass computation in loss. You may find it helpful to write an affine_batchnorm_relu() layer in nndl/layer_utils.py although this is not necessary.
- (3) The batchnorm_backward layer has to be appropriately inserted when calculating gradients.

After you have done the appropriate modifications, check your implementation by running the following cell.

Note, while the relative error for W3 should be small, as we backprop gradients more, you may find the relative error increases. Our relative error for W1 is on the order of 1e-4.

```
In [6]: N, D, H1, H2, C = 2, 15, 20, 30, 10
        X = np.random.randn(N, D)
        y = np.random.randint(C, size=(N,))
        for req in [0, 3.14]:
          print('Running check with reg = ', reg)
          model = FullyConnectedNet([H1, H2], input_dim=D, num_classes=C,
                                    reg=reg, weight scale=5e-2, dtype=np.float64,
                                    use batchnorm=True)
          loss, grads = model.loss(X, y)
          print('Initial loss: ', loss)
          for name in sorted(grads):
            f = lambda : model.loss(X, y)[0]
            grad num = eval numerical gradient(f, model.params[name], verbose=False, h=1e
        -5)
            print('{} relative error: {}'.format(name, rel error(grad num, grads[name])))
          if reg == 0: print('\n')
        Running check with reg = 0
        Initial loss: 2.188787718661933
        W1 relative error: 0.00037497800149038587
        W2 relative error: 2.9802196480990068e-05
        W3 relative error: 4.888275254235284e-10
        b1 relative error: 1.3988810110276972e-06
        b2 relative error: 6.661338147750939e-08
        b3 relative error: 1.4612611236765422e-10
        betal relative error: 5.951068872273817e-08
        beta2 relative error: 8.59207933341679e-09
        gamma1 relative error: 6.559388629007613e-08
        gamma2 relative error: 4.915022499153862e-09
        Running check with reg = 3.14
        Initial loss: 6.387021516091027
        W1 relative error: 2.0159561734290893e-05
        W2 relative error: 7.613736541079917e-06
        W3 relative error: 4.301795856380999e-08
        b1 relative error: 0.004440892098500625
        b2 relative error: 0.004440819934004025
        b3 relative error: 1.887027549448427e-10
        betal relative error: 6.551191124783562e-09
        beta2 relative error: 6.702536344617225e-09
        gamma1 relative error: 6.574369250585113e-09
```

Training a deep fully connected network with batch normalization.

To see if batchnorm helps, let's train a deep neural network with and without batch normalization.

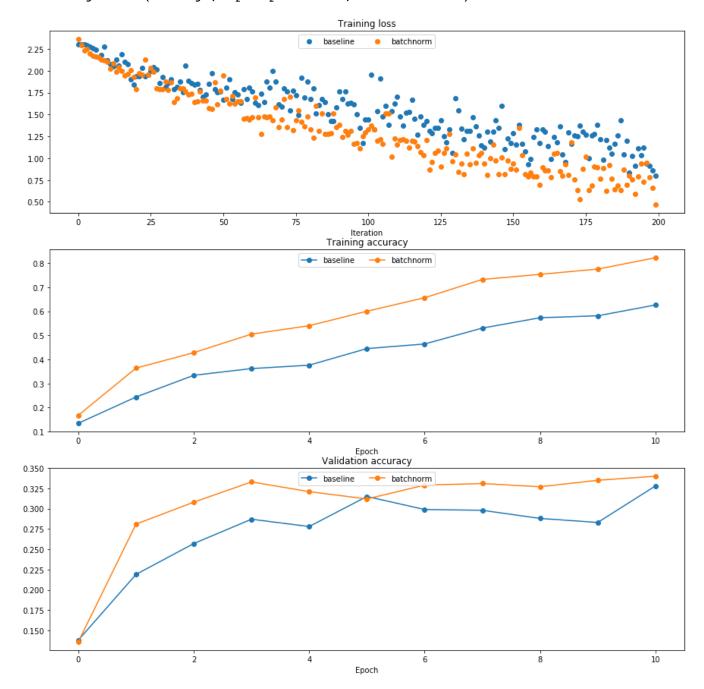
gamma2 relative error: 7.39453183612395e-09

```
In [7]: # Try training a very deep net with batchnorm
        hidden dims = [100, 100, 100, 100, 100]
        num train = 1000
        small data = {
          'X_train': data['X_train'][:num_train],
          'y_train': data['y_train'][:num_train],
          'X_val': data['X_val'],
          'y val': data['y val'],
        weight scale = 2e-2
        bn model = FullyConnectedNet(hidden dims, weight scale=weight scale, use batchnor
        model = FullyConnectedNet(hidden dims, weight scale=weight scale, use batchnorm=F
        alse)
        bn solver = Solver(bn model, small data,
                        num epochs=10, batch size=50,
                        update rule='adam',
                        optim config={
                           'learning rate': 1e-3,
                        },
                        verbose=True, print every=200)
        bn solver.train()
        solver = Solver(model, small data,
                        num epochs=10, batch size=50,
                        update rule='adam',
                        optim config={
                           'learning rate': 1e-3,
                        verbose=True, print every=200)
        solver.train()
        (Iteration 1 / 200) loss: 2.362126
        (Epoch 0 / 10) train acc: 0.168000; val acc: 0.136000
        (Epoch 1 / 10) train acc: 0.364000; val acc: 0.281000
        (Epoch 2 / 10) train acc: 0.428000; val acc: 0.308000
        (Epoch 3 / 10) train acc: 0.505000; val acc: 0.333000
        (Epoch 4 / 10) train acc: 0.540000; val acc: 0.321000
        (Epoch 5 / 10) train acc: 0.600000; val_acc: 0.312000
        (Epoch 6 / 10) train acc: 0.656000; val acc: 0.329000
        (Epoch 7 / 10) train acc: 0.732000; val acc: 0.331000
        (Epoch 8 / 10) train acc: 0.753000; val acc: 0.327000
        (Epoch 9 / 10) train acc: 0.775000; val acc: 0.335000
        (Epoch 10 / 10) train acc: 0.822000; val acc: 0.340000
        (Iteration 1 / 200) loss: 2.302878
        (Epoch 0 / 10) train acc: 0.135000; val acc: 0.138000
        (Epoch 1 / 10) train acc: 0.244000; val acc: 0.219000
        (Epoch 2 / 10) train acc: 0.334000; val acc: 0.257000
        (Epoch 3 / 10) train acc: 0.362000; val acc: 0.287000
        (Epoch 4 / 10) train acc: 0.376000; val acc: 0.278000
        (Epoch 5 / 10) train acc: 0.445000; val_acc: 0.315000
        (Epoch 6 / 10) train acc: 0.464000; val acc: 0.299000
        (Epoch 7 / 10) train acc: 0.530000; val_acc: 0.298000
        (Epoch 8 / 10) train acc: 0.573000; val acc: 0.288000
        (Epoch 9 / 10) train acc: 0.581000; val acc: 0.283000
        (Epoch 10 / 10) train acc: 0.626000; val acc: 0.328000
```

```
In [8]: plt.subplot(3, 1, 1)
        plt.title('Training loss')
        plt.xlabel('Iteration')
        plt.subplot(3, 1, 2)
        plt.title('Training accuracy')
        plt.xlabel('Epoch')
        plt.subplot(3, 1, 3)
        plt.title('Validation accuracy')
        plt.xlabel('Epoch')
        plt.subplot(3, 1, 1)
        plt.plot(solver.loss_history, 'o', label='baseline')
        plt.plot(bn_solver.loss_history, 'o', label='batchnorm')
        plt.subplot(3, 1, 2)
        plt.plot(solver.train_acc_history, '-o', label='baseline')
        plt.plot(bn_solver.train_acc_history, '-o', label='batchnorm')
        plt.subplot(3, 1, 3)
        plt.plot(solver.val_acc_history, '-o', label='baseline')
        plt.plot(bn_solver.val_acc_history, '-o', label='batchnorm')
        for i in [1, 2, 3]:
          plt.subplot(3, 1, i)
          plt.legend(loc='upper center', ncol=4)
        plt.gcf().set size inches(15, 15)
        plt.show()
```

/Users/quentintruong/Desktop/UCLA-ECE239AS-W19/hw4/env/lib/python3.6/site-packag es/matplotlib/cbook/deprecation.py:106: MatplotlibDeprecationWarning: Adding an axes using the same arguments as a previous axes currently reuses the earlier in stance. In a future version, a new instance will always be created and returne d. Meanwhile, this warning can be suppressed, and the future behavior ensured, by passing a unique label to each axes instance.

warnings.warn(message, mplDeprecation, stacklevel=1)



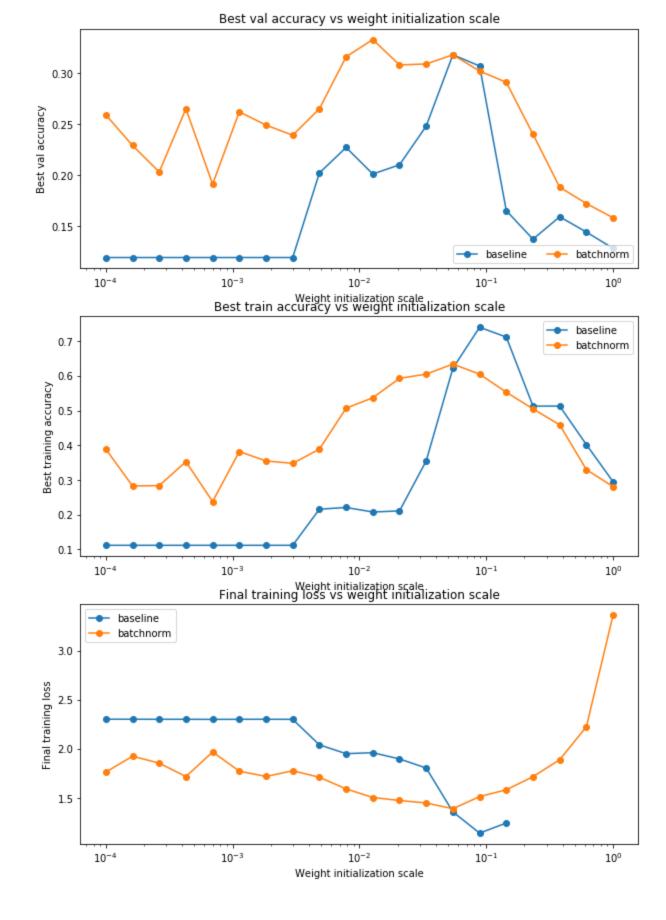
Batchnorm and initialization

The following cells run an experiment where for a deep network, the initialization is varied. We do training for when batchnorm layers are and are not included.

```
In [9]: # Try training a very deep net with batchnorm
        hidden dims = [50, 50, 50, 50, 50, 50, 50]
        num train = 1000
        small data = {
          'X_train': data['X_train'][:num_train],
          'y_train': data['y_train'][:num_train],
          'X_val': data['X_val'],
          'y_val': data['y_val'],
        bn solvers = {}
        solvers = {}
        weight scales = np.logspace(-4, 0, num=20)
        for i, weight scale in enumerate(weight scales):
          print('Running weight scale {} / {}'.format(i + 1, len(weight scales)))
          bn model = FullyConnectedNet(hidden dims, weight scale=weight scale, use batchn
        orm=True)
          model = FullyConnectedNet(hidden dims, weight scale=weight scale, use batchnorm
        =False)
          bn solver = Solver(bn model, small data,
                          num_epochs=10, batch_size=50,
                          update rule='adam',
                          optim config={
                             'learning rate': 1e-3,
                           },
                          verbose=False, print every=200)
          bn solver.train()
          bn solvers[weight scale] = bn solver
          solver = Solver(model, small data,
                          num epochs=10, batch size=50,
                          update rule='adam',
                          optim config={
                             'learning rate': 1e-3,
                          verbose=False, print every=200)
          solver.train()
          solvers[weight scale] = solver
```

```
Running weight scale 1 / 20
Running weight scale 2 / 20
Running weight scale 3 / 20
Running weight scale 4 / 20
Running weight scale 5 / 20
Running weight scale 6 / 20
Running weight scale 7 / 20
Running weight scale 8 / 20
Running weight scale 9 / 20
Running weight scale 10 / 20
Running weight scale 11 / 20
Running weight scale 12 / 20
Running weight scale 13 / 20
Running weight scale 14 / 20
Running weight scale 15 / 20
Running weight scale 16 / 20
/Users/quentintruong/Desktop/UCLA-ECE239AS-W19/hw4/code/nndl/layers.py:409: Runt
imeWarning: divide by zero encountered in log
  loss = -np.sum(np.log(probs[np.arange(N), y])) / N
Running weight scale 17 / 20
Running weight scale 18 / 20
Running weight scale 19 / 20
Running weight scale 20 / 20
```

```
In [10]: # Plot results of weight scale experiment
         best train accs, bn best train accs = [], []
         best val accs, bn best val accs = [], []
         final train loss, bn final train loss = [], []
         for ws in weight scales:
           best_train_accs.append(max(solvers[ws].train_acc_history))
           bn best train accs.append(max(bn solvers[ws].train acc history))
           best val accs.append(max(solvers[ws].val acc history))
           bn best val accs.append(max(bn solvers[ws].val acc history))
           final train loss.append(np.mean(solvers[ws].loss history[-100:]))
           bn final train loss.append(np.mean(bn solvers[ws].loss history[-100:]))
         plt.subplot(3, 1, 1)
         plt.title('Best val accuracy vs weight initialization scale')
         plt.xlabel('Weight initialization scale')
         plt.ylabel('Best val accuracy')
         plt.semilogx(weight_scales, best_val_accs, '-o', label='baseline')
         plt.semilogx(weight scales, bn best val accs, '-o', label='batchnorm')
         plt.legend(ncol=2, loc='lower right')
         plt.subplot(3, 1, 2)
         plt.title('Best train accuracy vs weight initialization scale')
         plt.xlabel('Weight initialization scale')
         plt.ylabel('Best training accuracy')
         plt.semilogx(weight scales, best train accs, '-o', label='baseline')
         plt.semilogx(weight_scales, bn_best_train_accs, '-o', label='batchnorm')
         plt.legend()
         plt.subplot(3, 1, 3)
         plt.title('Final training loss vs weight initialization scale')
         plt.xlabel('Weight initialization scale')
         plt.ylabel('Final training loss')
         plt.semilogx(weight scales, final train loss, '-o', label='baseline')
         plt.semilogx(weight scales, bn final train loss, '-o', label='batchnorm')
         plt.legend()
         plt.gcf().set size inches(10, 15)
         plt.show()
```



Question:

In the cell below, summarize the findings of this experiment, and WHY these results make sense.

Answer:

Batchnorm is more capable of achieving higher validation accuracy and training accuracy across a wider range of weight initializations.

First, this makes sense because it prevents the network's weights from going to zero (vanishing); we can see that the network without batchnorm probably died when using weights less than 10^-3. Moreover, it prevents the networks weights from going to infinity (exploding); we can see that the network without batchnorm probably exploded when using weights greater than 10^-1. Batchnorm helps with weight initializations because it normalizes the data by encouraging zero mean and unit variance, which helps prevent the weights from needing to shrink or grow too much.

Second, batchnorm's higher performance (compared to the network without batchnorm for weights that didn't vanish or explode) is related to the fact that the data is encouraged to be zero mean and unit variance. This unit statistics makes it easier for the network to learn because during gradient descent, the previous layer's data has similar statistics across updates, making the updates more relevant.