Batch Normalization

In this notebook, you will implement the batch normalization layers of a neural network to increase its performance. 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 numerica
        l_gradient_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-i
        n-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': 'tr
        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:
          After batch normalization (gamma=1, beta=0)
          mean: [-2.38697950e-16 -2.39808173e-16 -4.60742555e-17]
          std: [1. 1. 1.]
        After batch normalization (nontrivial gamma, beta)
          means: [11. 12. 13.]
                             1.99999999 2.99999999]
          stds: [1.
```

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'}
         qamma = 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: [-0.05967139 \quad 0.00957257 \quad -0.0259153]
          stds: [1.0150653 0.99071727 1.00119239]
```

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: 1.0365309983627649e-08
         dgamma error: 4.235503173805597e-12
         dbeta error: 3.275576553527879e-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.

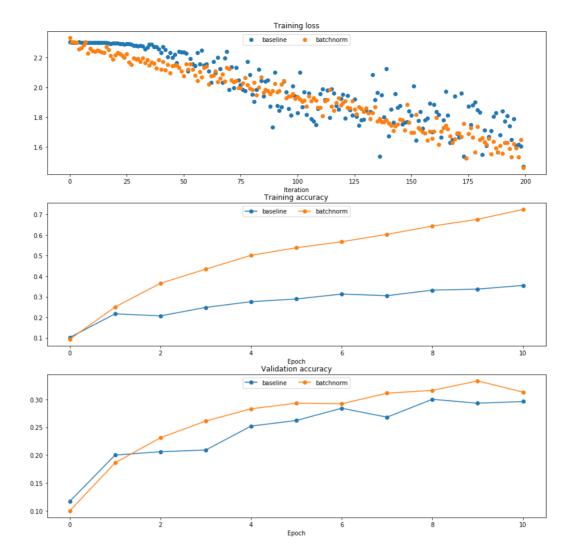
```
In [6]: N, D, H1, H2, C = 2, 15, 20, 30, 10
        X = np.random.randn(N, D)
        y = np.random.randint(C, size=(N,))
        for reg in [0, 3.14]:
            print('Running check with reg = ', reg)
            model = FullyConnectedNet([H1, H2], input_dim=D, num_classes=C,
                                       reg=reg, weight scale=5e-2, dtype=np.float
        64.
                                      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], verbos
        e=False, h=1e-5)
                print('{} relative error: {}'.format(name, rel_error(grad_num, g
        rads[name])))
            if reg == 0: print('\n')
        Running check with reg = 0
        Initial loss: 2.495181649437833
        W1 relative error: 5.7133128371328395e-05
        W2 relative error: 6.152107138686845e-06
        W3 relative error: 3.860542911698204e-10
        b1 relative error: 2.220446049250313e-08
        b2 relative error: 5.551115123125783e-09
        b3 relative error: 1.303595543107751e-10
        beta1 relative error: 4.66520471937565e-08
        beta2 relative error: 1.616541206944714e-09
        gamma1 relative error: 3.0691595985183625e-08
        gamma2 relative error: 2.3542593993916706e-09
        Running check with reg = 3.14
        Initial loss: 6.988149746148828
        W1 relative error: 3.63450880030614e-07
        W2 relative error: 1.4399569445718448e-08
        W3 relative error: 6.072364329088387e-08
        b1 relative error: 1.734723475976807e-10
        b2 relative error: 2.7755575615628914e-09
        b3 relative error: 1.4826365551616593e-10
        betal relative error: 3.12891605875848e-06
        beta2 relative error: 4.003822494515134e-08
        gamma1 relative error: 4.038265200968759e-06
        gamma2 relative error: 3.0722719270750266e-08
```

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.

```
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
         batchnorm=True)
        model = FullyConnectedNet(hidden dims, weight scale=weight scale, use ba
        tchnorm=False)
        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: 2.330732
        (Epoch 0 / 10) train acc: 0.093000; val_acc: 0.100000
        (Epoch 1 / 10) train acc: 0.250000; val_acc: 0.186000 (Epoch 2 / 10) train acc: 0.365000; val_acc: 0.231000
        (Epoch 3 / 10) train acc: 0.434000; val_acc: 0.261000
        (Epoch 4 / 10) train acc: 0.501000; val_acc: 0.283000
        (Epoch 5 / 10) train acc: 0.538000; val_acc: 0.293000
        (Epoch 6 / 10) train acc: 0.567000; val_acc: 0.292000
        (Epoch 7 / 10) train acc: 0.603000; val_acc: 0.311000
        (Epoch 8 / 10) train acc: 0.643000; val_acc: 0.316000
        (Epoch 9 / 10) train acc: 0.676000; val_acc: 0.333000
        (Epoch 10 / 10) train acc: 0.724000; val_acc: 0.313000
        (Iteration 1 / 200) loss: 2.303350
        (Epoch 0 / 10) train acc: 0.102000; val_acc: 0.117000
        (Epoch 1 / 10) train acc: 0.217000; val_acc: 0.200000
        (Epoch 2 / 10) train acc: 0.207000; val_acc: 0.206000
        (Epoch 3 / 10) train acc: 0.248000; val_acc: 0.209000
        (Epoch 4 / 10) train acc: 0.276000; val_acc: 0.252000
        (Epoch 5 / 10) train acc: 0.289000; val_acc: 0.262000
        (Epoch 6 / 10) train acc: 0.313000; val_acc: 0.284000
        (Epoch 7 / 10) train acc: 0.305000; val_acc: 0.268000
        (Epoch 8 / 10) train acc: 0.332000; val_acc: 0.300000
        (Epoch 9 / 10) train acc: 0.337000; val_acc: 0.293000
        (Epoch 10 / 10) train acc: 0.355000; val_acc: 0.296000
```

```
In [8]: fig, axes = plt.subplots(3, 1)
         ax = axes[0]
         ax.set title('Training loss')
         ax.set_xlabel('Iteration')
         ax = axes[1]
         ax.set_title('Training accuracy')
ax.set_xlabel('Epoch')
         ax = axes[2]
         ax.set_title('Validation accuracy')
         ax.set_xlabel('Epoch')
         ax = axes[0]
         ax.plot(solver.loss_history, 'o', label='baseline')
ax.plot(bn_solver.loss_history, 'o', label='batchnorm')
         ax = axes[1]
         ax.plot(solver.train_acc_history, '-o', label='baseline')
         ax.plot(bn_solver.train_acc_history, '-o', label='batchnorm')
         ax = axes[2]
         ax.plot(solver.val_acc_history, '-o', label='baseline')
         ax.plot(bn_solver.val_acc_history, '-o', label='batchnorm')
         for i in [1, 2, 3]:
              ax = axes[i - 1]
              ax.legend(loc='upper center', ncol=4)
         plt.gcf().set_size_inches(15, 15)
         plt.show()
```

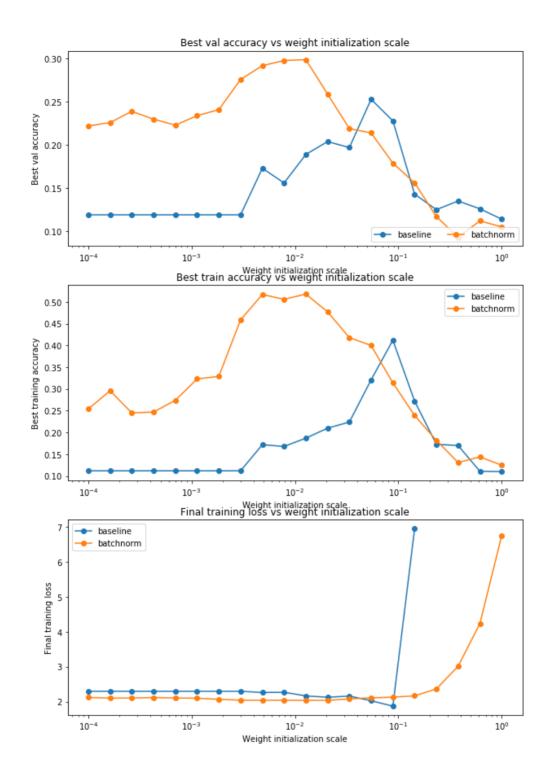


Batchnorm and initialization

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_scale
        s)))
             bn model = FullyConnectedNet(hidden dims, weight scale=weight scale,
        use_batchnorm=True)
            model = FullyConnectedNet(hidden_dims, weight_scale=weight_scale, us
        e 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
        /home/dennis/Documents/PY_PROGRAM/UCLA_C247/HW4-code/HW4-code/nndl/layer
        s.py:446: RuntimeWarning: 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[-10
          0:1))
          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:

On one hand, compared to the case without batch normalization, training with batch normalization allows us to train with a less bigger weight initialization scale, which is 2e-3, while the case without batch normalization is 5e-3. On the other hand, training with batch normalization tolerates greater weight initialization scale, meaning that training loss does not goes to ifinity with 1e-1 weight initialization scale. The main reason is that batch normalization acts as a regularization, which to some extent reduces the possibility for each layer of going to extreme values.