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_numerical_gradie
         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-ipytho
         %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'})
                 mean: ', a_norm.mean(axis=0))
         print('
         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: [ -9.17637808 -2.0580363 -17.78324359]
          stds: [41.16732993 26.50052681 33.07542062]
        After batch normalization (gamma=1, beta=0)
          mean: [ 1.66533454e-16 -7.43849426e-17 1.07136522e-16]
                            0.99999999 1.
          std: [1.
        After batch normalization (nontrivial gamma, beta)
          means: [11. 12. 13.]
```

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

1.99999999 2.999999991

```
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)
```

stds: [1.

```
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: [-0.09378005 0.04614452 -0.02173469]
```

Batchnorm backward pass

stds: [0.97910845 1.04762543 0.98880615]

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: 1.4186096322719448e-08
dgamma error: 8.217669743520836e-12
dbeta error: 3.9185260469964293e-11
```

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.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=1
             print('{} relative error: {}'.format(name, rel error(grad num, grads[name]))
           if reg == 0: print('\n')
        Running check with reg = 0
        Initial loss: 2.207694248543029
        W1 relative error: 0.0005000106339833756
        W2 relative error: 6.579144180337383e-05
        W3 relative error: 3.7053714174736985e-10
        b1 relative error: 1.0495077029659683e-08
        b2 relative error: 2.400857290751901e-07
        b3 relative error: 1.1905639326632003e-10
        betal relative error: 4.773590456956985e-08
        beta2 relative error: 1.0433495321077037e-08
        gamma1 relative error: 3.2019554804305616e-08
        gamma2 relative error: 6.2819204268826725e-09
        Running check with reg = 3.14
        Initial loss: 7.345437888117081
        W1 relative error: 1.549517586548874e-05
        W2 relative error: 7.803342381217168e-06
        W3 relative error: 2.7242808696725256e-08
        b1 relative error: 1.2685165418080402e-08
        b2 relative error: 4.3853809472693683e-07
        b3 relative error: 1.9288315246954152e-10
        beta1 relative error: 2.4476014258843913e-08
        beta2 relative error: 1.2001393219250454e-08
```

```
gamma1 relative error: 2.215518566042456e-08
gamma2 relative error: 2.30108816286757e-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_batchno
         model = FullyConnectedNet(hidden_dims, weight_scale=weight_scale, use_batchnorm=
         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: 2.289108
(Epoch 0 / 10) train acc: 0.120000; val acc: 0.122000
(Epoch 1 / 10) train acc: 0.359000; val acc: 0.267000
(Epoch 2 / 10) train acc: 0.434000; val acc: 0.303000
(Epoch 3 / 10) train acc: 0.513000; val acc: 0.320000
(Epoch 4 / 10) train acc: 0.585000; val acc: 0.324000
(Epoch 5 / 10) train acc: 0.649000; val acc: 0.328000
(Epoch 6 / 10) train acc: 0.717000; val acc: 0.330000
(Epoch 7 / 10) train acc: 0.719000; val acc: 0.311000
(Epoch 8 / 10) train acc: 0.760000; val acc: 0.312000
(Epoch 9 / 10) train acc: 0.789000; val acc: 0.314000
(Epoch 10 / 10) train acc: 0.807000; val acc: 0.328000
(Iteration 1 / 200) loss: 2.302422
(Epoch 0 / 10) train acc: 0.173000; val acc: 0.156000
(Epoch 1 / 10) train acc: 0.224000; val acc: 0.212000
(Epoch 2 / 10) train acc: 0.325000; val acc: 0.287000
(Epoch 3 / 10) train acc: 0.358000; val acc: 0.282000
```

```
(Epoch 4 / 10) train acc: 0.413000; val_acc: 0.310000 (Epoch 5 / 10) train acc: 0.431000; val_acc: 0.297000 (Epoch 6 / 10) train acc: 0.454000; val_acc: 0.308000 (Epoch 7 / 10) train acc: 0.551000; val_acc: 0.306000 (Epoch 8 / 10) train acc: 0.571000; val_acc: 0.340000 (Epoch 9 / 10) train acc: 0.555000; val_acc: 0.288000 (Epoch 10 / 10) train acc: 0.613000; val_acc: 0.319000
```

```
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()
```

<ipython-input-8-8e49aa315b6d>:13: MatplotlibDeprecationWarning: Adding an axes
using the same arguments as a previous axes currently reuses the earlier instanc
e. In a future version, a new instance will always be created and returned. Me
anwhile, this warning can be suppressed, and the future behavior ensured, by pas
sing a unique label to each axes instance.

plt.subplot(3, 1, 1)

<ipython-input-8-8e49aa315b6d>:17: MatplotlibDeprecationWarning: Adding an axes
using the same arguments as a previous axes currently reuses the earlier instanc
e. In a future version, a new instance will always be created and returned. Me
anwhile, this warning can be suppressed, and the future behavior ensured, by pas
sing a unique label to each axes instance.

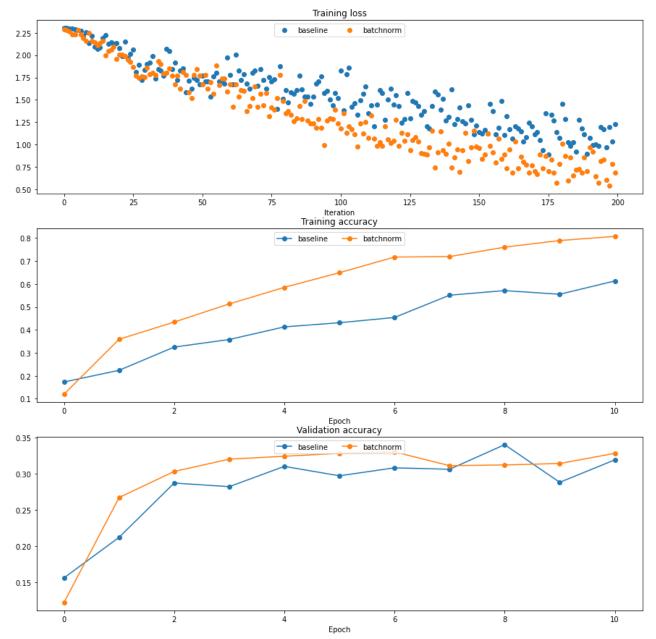
plt.subplot(3, 1, 2)

<ipython-input-8-8e49aa315b6d>:21: MatplotlibDeprecationWarning: Adding an axes
using the same arguments as a previous axes currently reuses the earlier instanc
e. In a future version, a new instance will always be created and returned. Me
anwhile, this warning can be suppressed, and the future behavior ensured, by pas
sing a unique label to each axes instance.

plt.subplot(3, 1, 3)

<ipython-input-8-8e49aa315b6d>:26: MatplotlibDeprecationWarning: Adding an axes
using the same arguments as a previous axes currently reuses the earlier instanc
e. In a future version, a new instance will always be created and returned. Me
anwhile, this warning can be suppressed, and the future behavior ensured, by pas
sing a unique label to each axes instance.

plt.subplot(3, 1, i)



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.

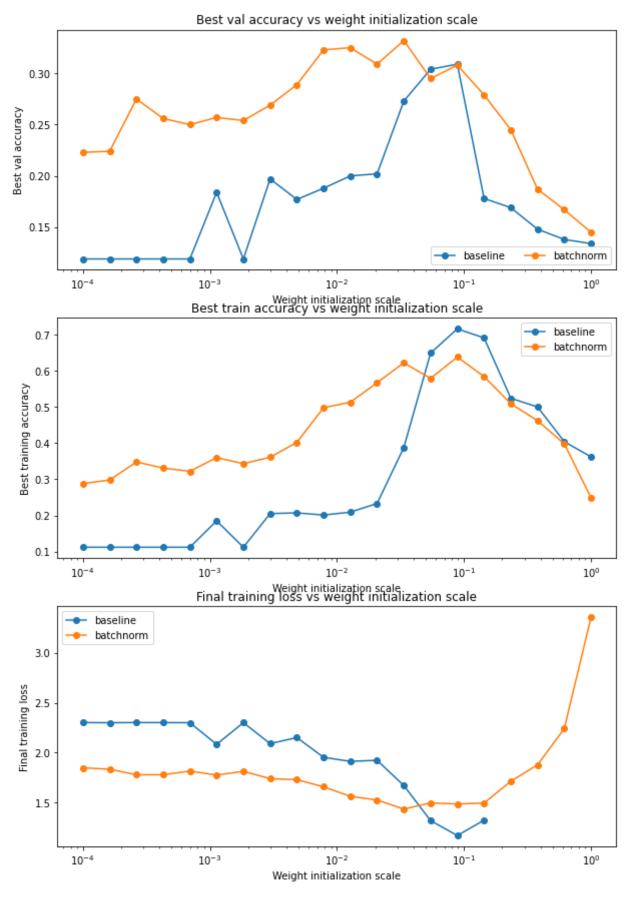
```
In [9]: # Try training a very deep net with batchnorm
hidden_dims = [50, 50, 50, 50, 50, 50]

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_batch
            model = FullyConnectedNet(hidden_dims, weight_scale=weight_scale, use_batchnor
            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/Jonathanchang/Downloads/HW4-code/nndl/layers.py:422: RuntimeWarning: divi
         de 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:

We can find out weight initialization will effect a lot on finding best training and validation accuracies. But we can found out that in some batch normalization performs the loss is worse then the original training. Because batch normalization modulated signal in each layers, and some information will lost in the process. But it have improvements at certain situations and we can find the best condition for such training.

In []:		