Helper_Functions

February 13, 2025

1 Optimization.ipynb Helper Functions

1.1 fc net.py Copied from HW 3

```
[]: import numpy as np
     from .layers import *
     from .layer_utils import *
     class TwoLayerNet(object):
         A two-layer fully-connected neural network with ReLU nonlinearity and
         softmax loss that uses a modular layer design. We assume an input dimension
         of D, a hidden dimension of H, and perform classification over C classes.
         The architecure should be affine - relu - affine - softmax.
         Note that this class does not implement gradient descent; instead, it
         will interact with a separate Solver object that is responsible for running
         optimization.
         The learnable parameters of the model are stored in the dictionary
         self.params that maps parameter names to numpy arrays.
         n n n
         def __init__(self, input_dim=3*32*32, hidden_dims=100, num_classes=10,
                    dropout=0, weight_scale=1e-3, reg=0.0):
             Initialize a new network.
             Inputs:
             - input_dim: An integer giving the size of the input
             - hidden_dims: An integer giving the size of the hidden layer
             - num_classes: An integer giving the number of classes to classify
             - dropout: Scalar between 0 and 1 giving dropout strength.
             - weight_scale: Scalar giving the standard deviation for random
               initialization of the weights.
```

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- req: Scalar giving L2 regularization strength.
      self.params = {}
      self.reg = reg
      # YOUR CODE HERE:
      # Initialize W1, W2, b1, and b2. Store these as self.params['W1'],
      \# self.params['W2'], self.params['b1'] and self.params['b2']. The
      # biases are initialized to zero and the weights are initialized
         so that each parameter has mean O and standard deviation
\neg weight\_scale.
         The dimensions of W1 should be (input dim, hidden dim) and the
      # dimensions of W2 should be (hidden_dims, num_classes)
      # ------ #
      self.params['W1'] = np.random.normal(0, weight_scale, (input_dim,_u
→hidden dims))
      self.params['W2'] = np.random.normal(0, weight_scale, (hidden_dims,_
→num classes))
      self.params['b1'] = np.zeros(hidden_dims)
      self.params['b2'] = np.zeros(num_classes)
      #pass
      # ------ #
      # END YOUR CODE HERE
      # ----- #
  def loss(self, X, y=None):
      Compute loss and gradient for a minibatch of data.
      Inputs:
      - X: Array of input data of shape (N, d_1, \ldots, d_k)
      - y: Array of labels, of shape (N,). y[i] gives the label for X[i].
     Returns:
      If y is None, then run a test-time forward pass of the model and return:
      - scores: Array of shape (N, C) giving classification scores, where
       scores[i, c] is the classification score for X[i] and class c.
     If y is not None, then run a training-time forward and backward pass and
      return a tuple of:
      - loss: Scalar value giving the loss
      - grads: Dictionary with the same keys as self.params, mapping parameter
       names to gradients of the loss with respect to those parameters.
      scores = None
```

```
# ----- #
     # YOUR CODE HERE:
        Implement the forward pass of the two-layer neural network. Store
        the class scores as the variable 'scores'. Be sure to use the
\hookrightarrow layers
     # you prior implemented.
     # ------ #
     hidden, cache_hidden = affine_relu_forward(X, self.params['W1'], self.
→params['b1'])
     scores, cache_scores = affine forward(hidden, self.params['W2'], self.
→params['b2'])
     #pass
     # ------ #
     # END YOUR CODE HERE
     # ============= #
     # If y is None then we are in test mode so just return scores
     if y is None:
       return scores
     loss, grads = 0, \{\}
     # ------ #
     # YOUR CODE HERE:
        Implement the backward pass of the two-layer neural net. Store
        the loss as the variable 'loss' and store the gradients in the
        'grads' dictionary. For the grads dictionary, grads['W1'] holds
        the gradient for W1, grads['b1'] holds the gradient for b1, etc.
        i.e., grads[k] holds the gradient for self.params[k].
     #
     #
       Add L2 regularization, where there is an added cost 0.5*self.reg*W^2
       for each W. Be sure to include the 0.5 multiplying factor to
        match our implementation.
        And be sure to use the layers you prior implemented.
     # ----- #
     loss, dout = softmax_loss(scores,y)
     loss += 0.5 * self.reg * (np.sum(self.params['W1']**2) + np.sum(self.
→params['W2']**2))
     dh, dw2, db2 = affine_backward(dout, cache_scores)
     dx, dw1, db1 = affine_relu_backward(dh, cache_hidden)
     grads['W1'] = dw1 + self.reg * self.params['W1']
     grads['b1'] = db1
     grads['W2'] = dw2 + self.reg * self.params['W2']
     grads['b2'] = db2
```

```
#pass
        # ----- #
        # END YOUR CODE HERE
        # ----- #
       return loss, grads
class FullyConnectedNet(object):
   A fully-connected neural network with an arbitrary number of hidden layers,
   ReLU nonlinearities, and a softmax loss function. This will also implement
    dropout and batch normalization as options. For a network with L layers,
    the architecture will be
    \{affine - [batch norm] - relu - [dropout]\} x (L - 1) - affine - softmax
   where batch normalization and dropout are optional, and the {...} block is
   repeated L - 1 times.
   Similar to the TwoLayerNet above, learnable parameters are stored in the
   self.params dictionary and will be learned using the Solver class.
    11 11 11
   def __init__(self, hidden_dims, input_dim=3*32*32, num_classes=10,
              dropout=0, use batchnorm=False, reg=0.0,
              weight_scale=1e-2, dtype=np.float32, seed=None):
       Initialize a new FullyConnectedNet.
       Inputs:
       - hidden_dims: A list of integers giving the size of each hidden layer.
       - input_dim: An integer giving the size of the input.
        - num_classes: An integer giving the number of classes to classify.
       - dropout: Scalar between 0 and 1 giving dropout strength. If dropout=0\sqcup
 \hookrightarrow then
         the network should not use dropout at all.
       - use\_batchnorm: Whether or not the network should use batch_{\sqcup}
 \neg normalization.
        - reg: Scalar giving L2 regularization strength.
       - weight_scale: Scalar giving the standard deviation for random
         initialization of the weights.
       - dtype: A numpy datatype object; all computations will be performed ⊔
 \hookrightarrow using
         this datatype. float32 is faster but less accurate, so you should use
```

```
float64 for numeric gradient checking.
      - seed: If not None, then pass this random seed to the dropout layers. \Box
\hookrightarrow This
       will make the dropout layers deteriminstic so we can gradient check \sqcup
\hookrightarrow the
       model.
      self.use_batchnorm = use_batchnorm
      self.use_dropout = dropout > 0
      self.reg = reg
      self.num_layers = 1 + len(hidden_dims)
      self.dtype = dtype
     self.params = {}
      # YOUR CODE HERE:
        Initialize all parameters of the network in the self.params
\hookrightarrow dictionary.
         The weights and biases of layer 1 are W1 and b1; and in general the
      # weights and biases of layer i are Wi and bi. The
      # biases are initialized to zero and the weights are initialized
      # so that each parameter has mean 0 and standard deviation_
⇔weight scale.
      # ----- #
      for i in np.arange(self.num_layers):
         if(i == 0):
             self.params['W' + str(i+1)] = np.random.normal(0, weight_scale,__
→(input_dim, hidden_dims[i]))
             self.params['b' + str(i+1)] = np.zeros(hidden_dims[i])
         elif(i == self.num_layers - 1):
             self.params['W' + str(i+1)] = np.random.normal(0, weight_scale,__
⇔(hidden_dims[i-1], num_classes))
             self.params['b' + str(i+1)] = np.zeros(num_classes)
         else:
             self.params['W' + str(i+1)] = np.random.normal(0, weight_scale,__
self.params['b' + str(i+1)] = np.zeros(hidden_dims[i])
      #pass
      # ------ #
      # END YOUR CODE HERE
      # ------ #
      # When using dropout we need to pass a dropout_param dictionary to each
      # dropout layer so that the layer knows the dropout probability and the
\rightarrowmode
```

```
\# (train / test). You can pass the same dropout param to each dropout
\hookrightarrow layer.
      self.dropout_param = {}
      if self.use dropout:
        self.dropout_param = {'mode': 'train', 'p': dropout}
        if seed is not None:
           self.dropout_param['seed'] = seed
       # With batch normalization we need to keep track of running means and
       # variances, so we need to pass a special bn param object to each batch
       # normalization layer. You should pass self.bn_params[0] to the forward_
\hookrightarrow pass
       # of the first batch normalization layer, self.bn_params[1] to the
\hookrightarrow forward
       # pass of the second batch normalization layer, etc.
      self.bn params = []
      if self.use_batchnorm:
        self.bn_params = [{'mode': 'train'} for i in np.arange(self.
→num_layers - 1)]
       # Cast all parameters to the correct datatype
      for k, v in self.params.items():
        self.params[k] = v.astype(dtype)
  def loss(self, X, y=None):
       Compute loss and gradient for the fully-connected net.
       Input / output: Same as TwoLayerNet above.
       11 11 11
      X = X.astype(self.dtype)
      mode = 'test' if y is None else 'train'
       # Set train/test mode for batchnorm params and dropout param since they
       # behave differently during training and testing.
      if self.dropout_param is not None:
        self.dropout_param['mode'] = mode
      if self.use_batchnorm:
        for bn_param in self.bn_params:
          bn_param[mode] = mode
      scores = None
       # ----- #
       # YOUR CODE HERE:
          Implement the forward pass of the FC net and store the output
```

```
scores as the variable "scores".
     H = []
     H_{cache} = []
     for i in range(self.num_layers):
        H_app = None
        H_cache_app = None
        if(i==0):
           H_app, H_cache_app = affine_relu_forward(X, self.params['W' +__
⇔str(i+1)], self.params['b' + str(i+1)])
           H.append(H_app)
           H_cache.append(H_cache_app)
        elif(i == self.num_layers - 1):
            scores, H_cache_app = affine_forward(H[i-1], self.params['W' +__
⇔str(i+1)], self.params['b' + str(i+1)])
           H_cache.append(H_cache_app)
        else:
           H_app, H_cache_app = affine_relu_forward(H[i-1], self.
aparams['W' + str(i+1)], self.params['b' + str(i+1)])
           H.append(H_app)
           H_cache.append(H_cache_app)
     #pass
     # ----- #
     # END YOUR CODE HERE
     # ----- #
     # If test mode return early
     if mode == 'test':
       return scores
     loss, grads = 0.0, {}
     # ----- #
     # YOUR CODE HERE:
       Implement the backwards pass of the FC net and store the gradients
     # in the grads dict, so that grads[k] is the gradient of self.
\rightarrow params[k]
       Be sure your L2 regularization includes a 0.5 factor.
     #pass
     loss, dhidden = softmax_loss(scores, y)
     for i in range(self.num_layers,0,-1):
       loss += 0.5*self.reg*np.sum(self.params['W{}'.format(i)]*self.
→params['W{}'.format(i)])
       if i == self.num_layers:
```

```
dH1, dW, db = affine_backward(dhidden,H_cache[i-1])
          grads['W{}'.format(i)] = dW + self.reg*self.params['W{}'.format(i)]
          grads['b{}'.format(i)] = db
          dH1, dW, db = affine_relu_backward(dH1,H_cache[i-1])
          grads['W{}'.format(i)] = dW + self.reg*self.params['W{}'.format(i)]
          grads['b{}'.format(i)] = db
      # loss, dhidden = softmax_loss(scores, y)
      # for i in range(self.num layers,0,-1):
           loss += 0.5 * self.reg*np.sum(self.params['W{}'.format(i)]*self.
\rightarrow params['W{}'.format(i)])
           if i == self.num_layers:
               dFC1, dW, db = affine\_backward(dhidden, FC\_cache[i-1])
               grads['W{}\}'.format(i)] = dW + self.reg*self.params['W{}\}'.
\rightarrow format(i)]
      #
              grads['b\{\}'.format(i)] = db
      #
            else:
             dFC1, dW, db = affine\_relu\_backward(dFC1,FC\_cache[i-1])
               qrads['W{}\}'.format(i)] = dW + self.req*self.params['W{}\}'.
\hookrightarrow format(i)]
               qrads['b{}\}'.format(i)] = db
      # END YOUR CODE HERE
      # ----- #
      return loss, grads
```

1.2 optim.py

```
[]: import numpy as np
     ,, ,, ,,
     This code was originally written for CS 231n at Stanford University
     (cs231n.stanford.edu). It has been modified in various areas for use in the
     ECE 239AS class at UCLA. This includes the descriptions of what code to
     implement as well as some slight potential changes in variable names to be
     consistent with class nomenclature. We thank Justin Johnson & Serena Yeung for
     permission to use this code. To see the original version, please visit
     cs231n.stanford.edu.
     11 11 11
     11 11 11
     This file implements various first-order update rules that are commonly used for
     training neural networks. Each update rule accepts current weights and the
     gradient of the loss with respect to those weights and produces the next set of
     weights. Each update rule has the same interface:
     def update(w, dw, config=None):
```

Inputs:

- w: A numpy array giving the current weights.
- dw: A numpy array of the same shape as w giving the gradient of the loss with respect to w.
- config: A dictionary containing hyperparameter values such as learning rate, momentum, etc. If the update rule requires caching values over many iterations, then config will also hold these cached values.

Returns:

- next_w: The next point after the update.
- config: The config dictionary to be passed to the next iteration of the update rule.

NOTE: For most update rules, the default learning rate will probably not perform well; however the default values of the other hyperparameters should work well for a variety of different problems.

For efficiency, update rules may perform in-place updates, mutating w and setting next_w equal to w.

```
def sgd(w, dw, config=None):
```

11 11 11

Performs vanilla stochastic gradient descent.

```
config format:
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- learning_rate: Scalar learning rate.

11 11 11

if config is None: config = {}

config.setdefault('learning_rate', 1e-2)

w -= config['learning_rate'] * dw
return w, config

def sgd_momentum(w, dw, config=None):

11 11 11

Performs stochastic gradient descent with momentum.

config format:

- learning_rate: Scalar learning rate.
- momentum: Scalar between 0 and 1 giving the momentum value.

Setting momentum = 0 reduces to sqd.

- velocity: A numpy array of the same shape as w and dw used to store \textbf{a}_{\sqcup} -moving

```
average of the gradients.
   if config is None: config = {}
   config.setdefault('learning_rate', 1e-2)
   config.setdefault('momentum', 0.9) # set momentum to 0.9 if it wasn't there
   v = config.get('velocity', np.zeros_like(w)) # gets velocity, else_u
 ⇔sets it to zero.
   # ----- #
   # YOUR CODE HERE:
   # Implement the momentum update formula. Return the updated weights
   # as next_w, and the updated velocity as v.
   # ============ #
   momentum_update = config['momentum'] * v - config['learning_rate'] * dw
   next_w = w + momentum_update
   v = momentum_update
   # ----- #
   # END YOUR CODE HERE
   config['velocity'] = v
   return next_w, config
def sgd_nesterov_momentum(w, dw, config=None):
   Performs stochastic gradient descent with Nesterov momentum.
   config format:
   - learning_rate: Scalar learning rate.
   - momentum: Scalar between 0 and 1 giving the momentum value.
   Setting momentum = 0 reduces to sqd.
   - velocity: A numpy array of the same shape as w and dw used to store a_{\sqcup}
 \hookrightarrow moving
   average of the gradients.
   if config is None: config = {}
   config.setdefault('learning_rate', 1e-2)
   config.setdefault('momentum', 0.9) # set momentum to 0.9 if it wasn't there
   v = config.get('velocity', np.zeros_like(w)) # gets velocity, else_
 ⇔sets it to zero.
   # =========== #
   # YOUR CODE HERE:
   # Implement the momentum update formula. Return the updated weights
   # as next_w, and the updated velocity as v.
   # ----- #
```

```
v_prev = v
   v = config['momentum']*v - config['learning_rate'] * dw
   next_w = w - config['momentum'] * v_prev + (1 + config['momentum']) * v
   # ----- #
   # END YOUR CODE HERE
   config['velocity'] = v
   return next_w, config
def rmsprop(w, dw, config=None):
   Uses the RMSProp update rule, which uses a moving average of squared \sqcup
 \hookrightarrow gradient
   values to set adaptive per-parameter learning rates.
   config format:
   - learning_rate: Scalar learning rate.
   - decay_rate: Scalar between 0 and 1 giving the decay rate for the squared
   gradient cache.
   - epsilon: Small scalar used for smoothing to avoid dividing by zero.
   - beta: Moving average of second moments of gradients.
   if config is None: config = {}
   config.setdefault('learning_rate', 1e-2)
   config.setdefault('decay_rate', 0.99)
   config.setdefault('epsilon', 1e-8)
   config.setdefault('a', np.zeros_like(w))
   next_w = None
   # ----- #
   # YOUR CODE HERE:
   # Implement RMSProp. Store the next value of w as next_w. You need
   # to also store in config['a'] the moving average of the second
     moment gradients, so they can be used for future gradients. Concretely,
   # config['a'] corresponds to "a" in the lecture notes.
   config['a'] = config['decay_rate'] * config['a'] + (1 -__
 next_w = w - config['learning_rate'] * dw / (np.sqrt(config['a']) +__
 ⇔config['epsilon'])
   # END YOUR CODE HERE
   # ------ #
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return next_w, config
def adam(w, dw, config=None):
   Uses the Adam update rule, which incorporates moving averages of both the
   gradient and its square and a bias correction term.
   config format:
   - learning_rate: Scalar learning rate.
   - beta1: Decay rate for moving average of first moment of gradient.
   - beta2: Decay rate for moving average of second moment of gradient.
   - epsilon: Small scalar used for smoothing to avoid dividing by zero.
   - m: Moving average of gradient.
   - v: Moving average of squared gradient.
   - t: Iteration number.
   if config is None: config = {}
   config.setdefault('learning_rate', 1e-3)
   config.setdefault('beta1', 0.9)
   config.setdefault('beta2', 0.999)
   config.setdefault('epsilon', 1e-8)
   config.setdefault('v', np.zeros_like(w))
   config.setdefault('a', np.zeros_like(w))
   config.setdefault('t', 0)
   next_w = None
   # ----- #
   # YOUR CODE HERE:
   # Implement Adam. Store the next value of w as next w. You need
   # to also store in config['a'] the moving average of the second
   # moment gradients, and in config['v'] the moving average of the
      first moments. Finally, store in config['t'] the increasing time.
   # ----- #
   config['t'] += 1
   config['v'] = config['beta1'] * config['v'] + (1 - config['beta1']) * dw
   config['a'] = config['beta2'] * config['a'] + (1 - config['beta2']) * dw**2
   v_corrected = config['v'] / (1 - config['beta1']**config['t'])
   a_corrected = config['a'] / (1 - config['beta2']**config['t'])
   next_w = w - config['learning_rate'] * v_corrected / (np.sqrt(a_corrected)⊔
 →+ config['epsilon'])
   # ========= #
   # END YOUR CODE HERE
   # ------ #
   return next_w, config
```

1.3 layers.py

```
[]: import numpy as np
    import pdb
    11 11 11
    This code was originally written for CS 231n at Stanford University
    (cs231n.stanford.edu). It has been modified in various areas for use in the
    ECE 239AS class at UCLA. This includes the descriptions of what code to
    implement as well as some slight potential changes in variable names to be
    consistent with class nomenclature. We thank Justin Johnson & Serena Yeung for
    permission to use this code. To see the original version, please visit
    cs231n.stanford.edu.
    def affine_forward(x, w, b):
       Computes the forward pass for an affine (fully-connected) layer.
       The input x has shape (N, d_1, \ldots, d_k) and contains a minibatch of N
       examples, where each example x[i] has shape (d_1, \ldots, d_k). We will
       reshape each input into a vector of dimension D = d_1 * ... * d_k, and
       then transform it to an output vector of dimension M.
       Inputs:
       -x: A numpy array containing input data, of shape (N, d_1, \ldots, d_k)
       - w: A numpy array of weights, of shape (D, M)
       - b: A numpy array of biases, of shape (M,)
       Returns a tuple of:
       - out: output, of shape (N, M)
       - cache: (x, w, b)
        11 11 11
       # ----- #
       # YOUR CODE HERE:
       # Calculate the output of the forward pass. Notice the dimensions
          of w are D x M, which is the transpose of what we did in earlier
         assignments.
       X = x.reshape((x.shape[0], -1))
       out = np.dot(X,w) + b
       # ----- #
       # END YOUR CODE HERE
       # ------ #
```

```
cache = (x, w, b)
   return out, cache
def affine_backward(dout, cache):
   Computes the backward pass for an affine layer.
   Inputs:
   - dout: Upstream derivative, of shape (N, M)
   - cache: Tuple of:
   - x: A numpy array containing input data, of shape (N, d_1, \ldots, d_k)
   - w: A numpy array of weights, of shape (D, M)
   - b: A numpy array of biases, of shape (M,)
   Returns a tuple of:
   - dx: Gradient with respect to x, of shape (N, d1, ..., d_k)
   - dw: Gradient with respect to w, of shape (D, M)
   - db: Gradient with respect to b, of shape (M,)
   x, w, b = cache
   dx, dw, db = None, None, None
   # ----- #
   # YOUR CODE HERE:
   # Calculate the gradients for the backward pass.
   # Notice:
   # dout is N x M
     dx should be N x d1 x ... x dk; it relates to dout through_
 \rightarrowmultiplication with w, which is D x M
     dw should be D x M; it relates to dout through multiplication with x_{,\cup}
 \rightarrowwhich is N x D after reshaping
   # db should be M; it is just the sum over dout examples
   X = x.reshape((x.shape[0], -1))
   db = np.sum(dout, axis=0)
   dw = np.dot(X.T, dout)
   dx = np.dot(dout, w.T).reshape(x.shape)
   # ----- #
   # END YOUR CODE HERE
   # ----- #
   return dx, dw, db
def relu_forward(x):
   11 11 11
```

```
Computes the forward pass for a layer of rectified linear units (ReLUs).
  Input:
  - x: Inputs, of any shape
  Returns a tuple of:
  - out: Output, of the same shape as x
  - cache: x
  # ----- #
  # YOUR CODE HERE:
  # Implement the ReLU forward pass.
  # ----- #
  out = np.maximum(0,x)
  # ============ #
  # END YOUR CODE HERE
  # ======== #
  cache = x
  return out, cache
def relu_backward(dout, cache):
  Computes the backward pass for a layer of rectified linear units (ReLUs).
  Input:
  - dout: Upstream derivatives, of any shape
  - cache: Input x, of same shape as dout
  Returns:
  - dx: Gradient with respect to x
  x = cache
  # YOUR CODE HERE:
  # Implement the ReLU backward pass
  # ----- #
  dx = dout * (x>0)
  # ----- #
  # END YOUR CODE HERE
  return dx
```

```
def batchnorm_forward(x, gamma, beta, bn_param):
   Forward pass for batch normalization.
   During training the sample mean and (uncorrected) sample variance are
    computed from minibatch statistics and used to normalize the incoming data.
   During training we also keep an exponentially decaying running mean of the \sqcup
 ⊶mean
   and variance of each feature, and these averages are used to normalize data
   at test-time.
   At each timestep we update the running averages for mean and variance using
    an exponential decay based on the momentum parameter:
   running mean = momentum * running mean + (1 - momentum) * sample mean
   running var = momentum * running var + (1 - momentum) * sample var
   Note that the batch normalization paper suggests a different test-time
    behavior: they compute sample mean and variance for each feature using a
    large number of training images rather than using a running average. For
    this implementation we have chosen to use running averages instead since
    they do not require an additional estimation step; the torch7 implementation
    of batch normalization also uses running averages.
   Input:
    - x: Data of shape (N, D)
    - gamma: Scale parameter of shape (D,)
    - beta: Shift paremeter of shape (D,)
    - bn_param: Dictionary with the following keys:
    - mode: 'train' or 'test'; required
    - eps: Constant for numeric stability
    - momentum: Constant for running mean / variance.
    - running_mean: Array of shape (D,) giving running mean of features
    - running_var Array of shape (D,) giving running variance of features
   Returns a tuple of:
    - out: of shape (N, D)
    - cache: A tuple of values needed in the backward pass
    n n n
   mode = bn_param['mode']
   eps = bn_param.get('eps', 1e-5)
   momentum = bn_param.get('momentum', 0.9)
   N, D = x.shape
   running_mean = bn_param.get('running_mean', np.zeros(D, dtype=x.dtype))
   running_var = bn_param.get('running_var', np.zeros(D, dtype=x.dtype))
```

```
out, cache = None, None
 if mode == 'train':
  # ----- #
  # YOUR CODE HERE:
    A few steps here:
      (1) Calculate the running mean and variance of the minibatch.
      (2) Normalize the activations with the sample mean and variance.
      (3) Scale and shift the normalized activations. Store this
         as the variable 'out'
     (4) Store any variables you may need for the backward pass in
        the 'cache' variable.
  sample_mean = np.mean(x, axis = 0)
    sample_variance = np.var(x, axis = 0)
    x_hat = (x - sample_mean) / np.sqrt(sample_variance + eps)
    # eps is the epsilon value needed to make sure it is a nonzero value
    out = gamma * x_hat + beta
    cache = (x, x_hat, sample_mean, sample_variance, gamma, beta, eps)
    running_mean = momentum * running_mean + (1 - momentum) * sample_mean
    running_variance = momentum * running_variance + (1 - momentum) *__
→sample_variance
  # ----- #
  # END YOUR CODE HERE
  # ============ #
 elif mode == 'test':
  # ------ #
  # YOUR CODE HERE:
    Calculate the testing time normalized activation. Normalize using
    the running mean and variance, and then scale and shift appropriately.
    Store the output as 'out'.
  # ----- #
    x_hat = (x - running_mean)/np.sqrt(running_variance + eps)
    out = gamma * x_hat + beta
  # ============ #
  # END YOUR CODE HERE
  else:
    raise ValueError('Invalid forward batchnorm mode "%s"' % mode)
 # Store the updated running means back into bn_param
```

```
bn_param['running_mean'] = running_mean
   bn_param['running_var'] = running_var
   return out, cache
def batchnorm_backward(dout, cache):
   Backward pass for batch normalization.
   For this implementation, you should write out a computation graph for
   batch normalization on paper and propagate gradients backward through
   intermediate nodes.
   Inputs:
   - dout: Upstream derivatives, of shape (N, D)
   - cache: Variable of intermediates from batchnorm_forward.
   Returns a tuple of:
   - dx: Gradient with respect to inputs x, of shape (N, D)
   - dgamma: Gradient with respect to scale parameter gamma, of shape (D,)
   - dbeta: Gradient with respect to shift parameter beta, of shape (D,)
   11 11 11
   dx, dgamma, dbeta = None, None, None
   # ------ #
   # YOUR CODE HERE:
   # Implement the batchnorm backward pass, calculating dx, dgamma, and_{\square}
 \hookrightarrowdbeta.
   # ============ #
   x, x_hat, sample_mean, sample_variance, gamma, beta, eps = cache
   N, D = x.shape
   dbeta = np.sum(dout, axis = 0)
   dgamma = np.sum(dout * x_hat, axis = 0)
   dx hat = dout * gamma
   dsample_variance = np.sum(dx_hat * (x - sample_mean) * (-0.5) *__
 \Rightarrow (sample_variance + epsilon)**(-1.5), axis = 0)
   dsample_mean = np.sum(dx_hat * (-1)/np.sqrt(sample_variance + eps) + _U

dsample_variance *
                       np.mean(-2*(x-sample_mean), axis = 0))
   dx = dx_hat / np.sqrt(sample_variance + eps) + dsample_variance * 2 * (x - <math>\Box
 ⇒sample_mean)/N + dsample_mean/N
   # ----- #
   # END YOUR CODE HERE
   # ----- #
   return dx, dgamma, dbeta
```

```
def dropout_forward(x, dropout_param):
   Performs the forward pass for (inverted) dropout.
   Inputs:
   - x: Input data, of any shape
   - dropout_param: A dictionary with the following keys:
   - p: Dropout parameter. We keep each neuron output with probability p.
   - mode: 'test' or 'train'. If the mode is train, then perform dropout;
     if the mode is test, then just return the input.
   - seed: Seed for the random number generator. Passing seed makes this
    function deterministic, which is needed for gradient checking but not in
    real networks.
   Outputs:
   - out: Array of the same shape as x.
   - cache: A tuple (dropout_param, mask). In training mode, mask is the \sqcup
 \hookrightarrow dropout
   mask that was used to multiply the input; in test mode, mask is None.
   p, mode = dropout_param['p'], dropout_param['mode']
   if 'seed' in dropout_param:
      np.random.seed(dropout_param['seed'])
   mask = None
   out = None
   if mode == 'train':
   # ======== #
   # YOUR CODE HERE:
      Implement the inverted dropout forward pass during training time.
     Store the masked and scaled activations in out, and store the
      dropout mask as the variable mask.
   mask = (np.random.rand(*x.shape) < p)/p</pre>
      out = x*mask
   # ------ #
   # END YOUR CODE HERE
   # ------ #
   elif mode == 'test':
   # ------ #
   # YOUR CODE HERE:
      Implement the inverted dropout forward pass during test time.
   # ------ #
```

```
out = x
  # ------ #
  # END YOUR CODE HERE
  # ----- #
  cache = (dropout_param, mask)
  out = out.astype(x.dtype, copy=False)
  return out, cache
def dropout_backward(dout, cache):
  Perform the backward pass for (inverted) dropout.
  Inputs:
  - dout: Upstream derivatives, of any shape
  - cache: (dropout_param, mask) from dropout_forward.
  dropout_param, mask = cache
  mode = dropout_param['mode']
  dx = None
  if mode == 'train':
  # ----- #
  # YOUR CODE HERE:
   Implement the inverted dropout backward pass during training time.
  # ------ #
    dx = dout * mask
  # END YOUR CODE HERE
  # ============ #
  elif mode == 'test':
  # ======== #
    Implement the inverted dropout backward pass during test time.
  dx = dout
  # END YOUR CODE HERE
  # ------ #
  return dx
def svm_loss(x, y):
  Computes the loss and gradient using for multiclass SVM classification.
  Inputs:
```

```
- x: Input data, of shape (N, C) where x[i, j] is the score for the jth_{\sqcup}
 \hookrightarrow class
    for the ith input.
    - y: Vector of labels, of shape (N,) where y[i] is the label for x[i] and
    0 <= y[i] < C
    Returns a tuple of:
    - loss: Scalar giving the loss
    - dx: Gradient of the loss with respect to x
    N = x.shape[0]
    correct_class_scores = x[np.arange(N), y]
    margins = np.maximum(0, x - correct_class_scores[:, np.newaxis] + 1.0)
    margins[np.arange(N), y] = 0
    loss = np.sum(margins) / N
    num_pos = np.sum(margins > 0, axis=1)
    dx = np.zeros_like(x)
    dx[margins > 0] = 1
    dx[np.arange(N), y] -= num_pos
    dx /= N
    return loss, dx
def softmax_loss(x, y):
    11 11 11
    Computes the loss and gradient for softmax classification.
    Inputs:
    - x: Input data, of shape (N, C) where x[i, j] is the score for the jth_{\sqcup}
 \hookrightarrow class
    for the ith input.
    - y: Vector of labels, of shape (N,) where y[i] is the label for x[i] and
    0 <= y[i] < C
    Returns a tuple of:
    - loss: Scalar giving the loss
    - dx: Gradient of the loss with respect to x
    probs = np.exp(x - np.max(x, axis=1, keepdims=True))
    probs /= np.sum(probs, axis=1, keepdims=True)
    N = x.shape[0]
    loss = -np.sum(np.log(probs[np.arange(N), y])) / N
    dx = probs.copy()
    dx[np.arange(N), y] = 1
    dx /= N
    return loss, dx
```

1.4 data utils.py

I kept the CIFAR data in the HW3 folder

```
[]: from __future__ import print_function
    from six.moves import cPickle as pickle
    import numpy as np
    import os
    from imageio import imread
    import platform
    def load_pickle(f):
        version = platform.python_version_tuple()
        if version[0] == '2':
             return pickle.load(f)
        elif version[0] == '3':
             return pickle.load(f, encoding='latin1')
        raise ValueError("invalid python version: {}".format(version))
    def load CIFAR batch(filename):
       """ load single batch of cifar """
      with open(filename, 'rb') as f:
        datadict = load_pickle(f)
        X = datadict['data']
        Y = datadict['labels']
        X = X.reshape(10000, 3, 32, 32).transpose(0,2,3,1).astype("float")
        Y = np.array(Y)
        return X, Y
    # def load_CIFAR10(ROOT):
       """ load all of cifar """
       xs = []
     #
     #
       ys = []
        for b in range (1,6):
     #
         f = os.path.join(ROOT, 'data_batch_%d' % (b, ))
     #
         X, Y = load_CIFAR_batch(f)
     #
         xs.append(X)
     #
         ys.append(Y)
     # Xtr = np.concatenate(xs)
     #
       Ytr = np.concatenate(ys)
     \# del X, Y
     # Xte, Yte = load CIFAR batch(os.path.join(ROOT, 'test batch'))
       return Xtr, Ytr, Xte, Yte
    def load_CIFAR10(ROOT):
         """Load all of CIFAR-10 using absolute paths."""
```

```
xs = []
    ys = []
    """ NOTE FOR THE GRADERS: I had something going on with my join ROOT_{\sqcup}
 \hookrightarrow function
        so I decided to simply manually join the file directory with a similar ...
 ⇔for loop
        because I kept getting the same error despite having the correct_{\sqcup}
 \hookrightarrow directory
        You can see I tested to see if the directory is present in the normal \sqcup
 ⇔code"""
    for b in range(1, 6):
        f = f"/Users/ctang/Desktop/ECE_C147/HW3/cifar-10-batches-py/

data_batch_{b}"

        if not os.path.exists(f):
            raise FileNotFoundError(f"File not found: {f}")
        X, Y = load_CIFAR_batch(f)
        xs.append(X)
        ys.append(Y)
    Xtr = np.concatenate(xs)
    Ytr = np.concatenate(ys)
    del X, Y
    test_file = "/Users/ctang/Desktop/ECE_C147/HW3/cifar-10-batches-py/
 ⇔test_batch"
    if not os.path.exists(test_file):
        raise FileNotFoundError(f"File not found: {test_file}")
    Xte, Yte = load_CIFAR_batch(test_file)
    return Xtr, Ytr, Xte, Yte
def get_CIFAR10_data(num_training=49000, num_validation=1000, num_test=1000,
                      subtract_mean=True):
    Load the CIFAR-10 dataset from disk and perform preprocessing to prepare
    it for classifiers. These are the same steps as we used for the SVM, but
    condensed to a single function.
    # Load the raw CIFAR-10 data
    cifar10_dir = '/Users/ctang/Desktop/ECE_C147/HW3/cifar-10-batches-py/
 \hookrightarrowtest_batch'
    X_train, y_train, X_test, y_test = load_CIFAR10(cifar10_dir)
    # Subsample the data
    mask = list(range(num_training, num_training + num_validation))
```

```
X_val = X_train[mask]
    y_val = y_train[mask]
    mask = list(range(num_training))
    X_train = X_train[mask]
    y_train = y_train[mask]
    mask = list(range(num_test))
    X_test = X_test[mask]
    y_test = y_test[mask]
    # Normalize the data: subtract the mean image
    if subtract mean:
     mean_image = np.mean(X_train, axis=0)
     X_train -= mean_image
      X_val -= mean_image
     X_test -= mean_image
    # Transpose so that channels come first
    X_train = X_train.transpose(0, 3, 1, 2).copy()
    X_{val} = X_{val.transpose}(0, 3, 1, 2).copy()
    X_test = X_test.transpose(0, 3, 1, 2).copy()
    # Package data into a dictionary
    return {
      'X train': X train, 'y train': y train,
      'X_val': X_val, 'y_val': y_val,
      'X_test': X_test, 'y_test': y_test,
    }
def load tiny imagenet(path, dtype=np.float32, subtract_mean=True):
 Load TinyImageNet. Each of TinyImageNet-100-A, TinyImageNet-100-B, and
  TinyImageNet-200 have the same directory structure, so this can be used
  to load any of them.
  Inputs:
  - path: String giving path to the directory to load.
  - dtype: numpy datatype used to load the data.
  - subtract_mean: Whether to subtract the mean training image.
 Returns: A dictionary with the following entries:
  - class_names: A list where class_names[i] is a list of strings giving the
    WordNet names for class i in the loaded dataset.
  - X_train: (N_tr, 3, 64, 64) array of training images
  - y_train: (N_tr,) array of training labels
  - X_val: (N_val, 3, 64, 64) array of validation images
  - y_val: (N_val,) array of validation labels
```

```
- X_test: (N_test, 3, 64, 64) array of testing images.
- y test: (N test,) array of test labels; if test labels are not available
  (such as in student code) then y test will be None.
- mean_image: (3, 64, 64) array giving mean training image
# First load wnids
with open(os.path.join(path, 'wnids.txt'), 'r') as f:
  wnids = [x.strip() for x in f]
# Map wnids to integer labels
wnid to label = {wnid: i for i, wnid in enumerate(wnids)}
# Use words.txt to get names for each class
with open(os.path.join(path, 'words.txt'), 'r') as f:
  wnid_to_words = dict(line.split('\t') for line in f)
  for wnid, words in wnid_to_words.iteritems():
    wnid_to_words[wnid] = [w.strip() for w in words.split(',')]
class_names = [wnid_to_words[wnid] for wnid in wnids]
# Next load training data.
X_train = []
y train = []
for i, wnid in enumerate(wnids):
  if (i + 1) % 20 == 0:
    print('loading training data for synset %d / %d' % (i + 1, len(wnids)))
  # To figure out the filenames we need to open the boxes file
  boxes_file = os.path.join(path, 'train', wnid, '%s_boxes.txt' % wnid)
  with open(boxes_file, 'r') as f:
    filenames = [x.split('\t')[0] for x in f]
  num_images = len(filenames)
  X_train_block = np.zeros((num_images, 3, 64, 64), dtype=dtype)
  y_train_block = wnid_to_label[wnid] * np.ones(num_images, dtype=np.int64)
  for j, img_file in enumerate(filenames):
    img_file = os.path.join(path, 'train', wnid, 'images', img_file)
    img = imread(img_file)
    if img.ndim == 2:
      ## grayscale file
      img.shape = (64, 64, 1)
    X_train_block[j] = img.transpose(2, 0, 1)
  X train.append(X train block)
  y_train.append(y_train_block)
# We need to concatenate all training data
X_train = np.concatenate(X_train, axis=0)
y_train = np.concatenate(y_train, axis=0)
```

```
# Next load validation data
with open(os.path.join(path, 'val', 'val_annotations.txt'), 'r') as f:
  img files = []
  val_wnids = []
  for line in f:
    img_file, wnid = line.split('\t')[:2]
    img_files.append(img_file)
    val_wnids.append(wnid)
  num val = len(img files)
  y_val = np.array([wnid_to_label[wnid] for wnid in val_wnids])
  X_val = np.zeros((num_val, 3, 64, 64), dtype=dtype)
  for i, img_file in enumerate(img_files):
    img_file = os.path.join(path, 'val', 'images', img_file)
    img = imread(img_file)
    if img.ndim == 2:
      img.shape = (64, 64, 1)
    X_{val}[i] = img.transpose(2, 0, 1)
# Next load test images
# Students won't have test labels, so we need to iterate over files in the
# images directory.
img_files = os.listdir(os.path.join(path, 'test', 'images'))
X_test = np.zeros((len(img_files), 3, 64, 64), dtype=dtype)
for i, img file in enumerate(img files):
  img_file = os.path.join(path, 'test', 'images', img_file)
  img = imread(img file)
  if img.ndim == 2:
    img.shape = (64, 64, 1)
  X_test[i] = img.transpose(2, 0, 1)
y_test = None
y_test_file = os.path.join(path, 'test', 'test_annotations.txt')
if os.path.isfile(y_test_file):
  with open(y_test_file, 'r') as f:
    img_file_to_wnid = {}
    for line in f:
      line = line.split('\t')
      img_file_to_wnid[line[0]] = line[1]
  y_test = [wnid_to_label[img_file_to_wnid[img_file]] for img_file in_
→img_files]
  y_test = np.array(y_test)
mean_image = X_train.mean(axis=0)
if subtract_mean:
  X_train -= mean_image[None]
  X_val -= mean_image[None]
  X_test -= mean_image[None]
```

```
return {
    'class_names': class_names,
    'X_train': X_train,
    'y_train': y_train,
    'X_val': X_val,
    'y_val': y_val,
    'X_test': X_test,
    'y_test': y_test,
    'class_names': class_names,
    'mean_image': mean_image,
 }
def load_models(models_dir):
  HHHH
 Load saved models from disk. This will attempt to unpickle all files in a
  directory; any files that give errors on unpickling (such as README.txt) will
  be skipped.
  Inputs:
  - models_dir: String giving the path to a directory containing model files.
    Each model file is a pickled dictionary with a 'model' field.
  Returns:
  A dictionary mapping model file names to models.
 models = \{\}
  for model_file in os.listdir(models_dir):
    with open(os.path.join(models_dir, model_file), 'rb') as f:
      try:
        models[model_file] = load_pickle(f)['model']
      except pickle.UnpicklingError:
        continue
  return models
```

2 Batch Normalization

2.1 fc_net.py

```
[]: import numpy as np
import pdb

from .layers import *
from .layer_utils import *
```

This code was originally written for CS 231n at Stanford University (cs231n.stanford.edu). It has been modified in various areas for use in the ECE 239AS class at UCLA. This includes the descriptions of what code to implement as well as some slight potential changes in variable names to be consistent with class nomenclature. We thank Justin Johnson & Serena Yeung for permission to use this code. To see the original version, please visit cs231n.stanford.edu. 11 11 11 class TwoLayerNet(object): A two-layer fully-connected neural network with ReLU nonlinearity and softmax loss that uses a modular layer design. We assume an input dimension of D, a hidden dimension of H, and perform classification over C classes. The architecure should be affine - relu - affine - softmax. Note that this class does not implement gradient descent; instead, it will interact with a separate Solver object that is responsible for running optimization. The learnable parameters of the model are stored in the dictionary self.params that maps parameter names to numpy arrays. def __init__(self, input_dim=3*32*32, hidden_dims=100, num_classes=10, dropout=1, weight_scale=1e-3, reg=0.0): Initialize a new network. Inputs: - input_dim: An integer giving the size of the input - hidden dims: An integer giving the size of the hidden layer - num_classes: An integer giving the number of classes to classify - dropout: Scalar between 0 and 1 giving dropout strength. - weight_scale: Scalar giving the standard deviation for random initialization of the weights. - reg: Scalar giving L2 regularization strength. self.params = {} self.reg = reg # YOUR CODE HERE: Initialize W1, W2, b1, and b2. Store these as self.params['W1'], # self.params['W2'], self.params['b1'] and self.params['b2']. The

biases are initialized to zero and the weights are initialized

```
so that each parameter has mean O and standard deviation weight scale.
    The dimensions of W1 should be (input_dim, hidden_dim) and the
     dimensions of W2 should be (hidden dims, num classes)
  # ------ #
  self.params['W1'] = np.random.normal(0, weight_scale, (input_dim,_
→hidden_dims))
  self.params['W2'] = np.random.normal(0, weight_scale, (hidden_dims,_

¬num_classes))
  self.params['b1'] = np.zeros(hidden_dims)
  self.params['b2'] = np.zeros(num_classes)
  # END YOUR CODE HERE
  def loss(self, X, y=None):
      Compute loss and gradient for a minibatch of data.
     Inputs:
      - X: Array of input data of shape (N, d_1, \ldots, d_k)
      - y: Array of labels, of shape (N,). y[i] gives the label for X[i].
     Returns:
     If y is None, then run a test-time forward pass of the model and return:
      - scores: Array of shape (N, C) giving classification scores, where
       scores[i, c] is the classification score for X[i] and class c.
     If y is not None, then run a training-time forward and backward pass and
     return a tuple of:
      - loss: Scalar value giving the loss
      - grads: Dictionary with the same keys as self.params, mapping parameter
       names to gradients of the loss with respect to those parameters.
     scores = None
      # ----- #
      # YOUR CODE HERE:
      # Implement the forward pass of the two-layer neural network. Store
      \# the class scores as the variable 'scores'. Be sure to use the
\hookrightarrow layers
      # you prior implemented.
      # ----- #
     hidden, cache_hidden = affine_relu_forward(X, self.params['W1'], self.
→params['b1'])
      scores, cache_scores = affine forward(hidden, self.params['W2'], self.
→params['b2'])
```

```
#pass
      # ----- #
      # END YOUR CODE HERE
      # ----- #
      # If y is None then we are in test mode so just return scores
      if y is None:
         return scores
      loss, grads = 0, \{\}
                    ----- #
      # YOUR CODE HERE:
         Implement the backward pass of the two-layer neural net. Store
         the loss as the variable 'loss' and store the gradients in the
         'grads' dictionary. For the grads dictionary, grads['W1'] holds
        the gradient for W1, grads['b1'] holds the gradient for b1, etc.
         i.e., grads[k] holds the gradient for self.params[k].
        Add L2 regularization, where there is an added cost 0.5*self.reg*W^2
         for each W. Be sure to include the 0.5 multiplying factor to
      #
        match our implementation.
      # And be sure to use the layers you prior implemented.
      # ------ #
      loss, dout = softmax_loss(scores,y)
      loss += 0.5 * self.reg * (np.sum(self.params['W1']**2) + np.sum(self.
 →params['W2']**2))
      dh, dw2, db2 = affine_backward(dout, cache_scores)
      dx, dw1, db1 = affine_relu_backward(dh, cache_hidden)
      grads['W1'] = dw1 + self.reg * self.params['W1']
      grads['b1'] = db1
      grads['W2'] = dw2 + self.reg * self.params['W2']
      grads['b2'] = db2
      # ----- #
      # END YOUR CODE HERE
      # ----- #
      return loss, grads
class FullyConnectedNet(object):
   A fully-connected neural network with an arbitrary number of hidden layers,
   ReLU nonlinearities, and a softmax loss function. This will also implement
   dropout and batch normalization as options. For a network with L layers,
```

```
the architecture will be
  \{affine - [batch norm] - relu - [dropout]\}\ x\ (L - 1) - affine - softmax
  where batch normalization and dropout are optional, and the {...} block is
  repeated L - 1 times.
  Similar to the TwoLayerNet above, learnable parameters are stored in the
  self.params dictionary and will be learned using the Solver class.
  def __init__(self, hidden_dims, input_dim=3*32*32, num_classes=10,
             dropout=1, use batchnorm=False, reg=0.0,
             weight_scale=1e-2, dtype=np.float32, seed=None):
  11 11 11
  Initialize a new FullyConnectedNet.
  Inputs:
  - hidden_dims: A list of integers giving the size of each hidden layer.
  - input_dim: An integer giving the size of the input.
  - num_classes: An integer giving the number of classes to classify.
  - dropout: Scalar between 0 and 1 giving dropout strength. If dropout=1 then
    the network should not use dropout at all.
  - use batchnorm: Whether or not the network should use batch normalization.
  - reg: Scalar giving L2 regularization strength.
  - weight scale: Scalar giving the standard deviation for random
    initialization of the weights.
  - dtype: A numpy datatype object; all computations will be performed using
    this datatype. float32 is faster but less accurate, so you should use
    float64 for numeric gradient checking.
  - seed: If not None, then pass this random seed to the dropout layers. This
    will make the dropout layers deteriminstic so we can gradient check the
    model.
  11 11 11
      self.use_batchnorm = use_batchnorm
      self.use_dropout = dropout < 1</pre>
      self.reg = reg
      self.num_layers = 1 + len(hidden_dims)
      self.dtype = dtype
      self.params = {}
      # ----- #
      # YOUR CODE HERE:
      # Initialize all parameters of the network in the self.params_
\hookrightarrow dictionary.
         The weights and biases of layer 1 are W1 and b1; and in general the
      # weights and biases of layer i are Wi and bi. The
```

```
biases are initialized to zero and the weights are initialized
          so that each parameter has mean 0 and standard deviation
\rightarrow weight_scale.
      #
        BATCHNORM: Initialize the gammas of each layer to 1 and the beta
      # parameters to zero. The gamma and beta parameters for layer 111
\hookrightarrowshould
        be self.params['qamma1'] and self.params['beta1']. For layer 2,
\hookrightarrow they
          should be gamma2 and beta2, etc. Only use batchnorm if self.
\neg use\_batchnorm
          is true and DO NOT do batch normalize the output scores.
      # ------ #
      for i in range(self.num_layers):
          if i == 0:
             self.params['W1'] = weight_scale * np.random.randn(input_dim,__
→hidden_dims[i])
              self.params['b1'] = np.zeros(hidden_dims[i])
              if self.use_batchnorm:
                 self.params['gamma1'] = np.ones(hidden_dims[i])
                 self.params['beta1'] = np.zeros(hidden_dims[i])
          elif i == self.num_layers - 1:
              self.params['W' + str(i + 1)] = weight_scale * np.random.
→randn(hidden_dims[i - 1], num_classes)
             self.params['b' + str(i + 1)] = np.zeros(num_classes)
          else:
              self.params['W' + str(i + 1)] = weight_scale * np.random.
→randn(hidden_dims[i - 1], hidden_dims[i])
              self.params['b' + str(i + 1)] = np.zeros(hidden_dims[i])
              if self.use_batchnorm:
                 self.params['gamma' + str(i + 1)] = np.ones(hidden_dims[i])
                 self.params['beta' + str(i + 1)] = np.zeros(hidden dims[i])
      # END YOUR CODE HERE
      # ------ #
      # When using dropout we need to pass a dropout_param dictionary to each
      # dropout layer so that the layer knows the dropout probability and the
⊶mode
      # (train / test). You can pass the same dropout param to each dropout
\hookrightarrow layer.
      self.dropout_param = {}
      if self.use_dropout:
        self.dropout_param = {'mode': 'train', 'p': dropout}
        if seed is not None:
          self.dropout_param['seed'] = seed
```

```
# With batch normalization we need to keep track of running means and
       # variances, so we need to pass a special bn param object to each batch
       # normalization layer. You should pass self.bn_params[0] to the forward_{\sqcup}
→pass
      # of the first batch normalization layer, self.bn params[1] to the
\hookrightarrow forward
      # pass of the second batch normalization layer, etc.
      self.bn_params = []
      if self.use_batchnorm:
        self.bn_params = [{'mode': 'train'} for i in np.arange(self.
→num layers - 1)]
      # Cast all parameters to the correct datatype
      for k, v in self.params.items():
        self.params[k] = v.astype(dtype)
  def loss(self, X, y=None):
       11 11 11
      Compute loss and gradient for the fully-connected net.
      Input / output: Same as TwoLayerNet above.
      X = X.astype(self.dtype)
      mode = 'test' if y is None else 'train'
      # Set train/test mode for batchnorm params and dropout param since they
      # behave differently during training and testing.
      if self.dropout_param is not None:
           self.dropout_param['mode'] = mode
      if self.use_batchnorm:
          for bn_param in self.bn_params:
               bn_param['mode'] = mode
      scores = None
       # YOUR CODE HERE:
          Implement the forward pass of the FC net and store the output
       #
          scores as the variable "scores".
       #
         BATCHNORM: If self.use_batchnorm is true, insert a bathnorm layer
          between the affine_forward and relu_forward layers. You may
       #
          also write an affine_batchnorm_relu() function in layer_utils.py.
       #
       #
          DROPOUT: If dropout is non-zero, insert a dropout layer after
```

```
every ReLU layer.
      # ----- #
      caches = []
      for i in range(self.num_layers):
         if i == 0:
             a, cache = affine_forward(X, self.params['W1'], self.
→params['b1'])
             caches.append(cache)
             if self.use_batchnorm:
                 a, cache = batchnorm_forward(a, self.params['gamma1'],
                                           self.params['beta1'], self.
→bn_params[0])
                 caches.append(cache)
             h, cache = relu_forward(a)
             caches.append(cache)
             if self.use_dropout:
                 h, cache = dropout_forward(h, self.dropout_param)
                 caches.append(cache)
         elif i == self.num_layers - 1:
             scores, cache = affine_forward(h, self.params['W' + str(i + 1)],
                                         self.params['b' + str(i + 1)])
             caches.append(cache)
         else:
             a, cache = affine_forward(h, self.params['W' + str(i + 1)],
                                     self.params['b' + str(i + 1)])
             caches.append(cache)
             if self.use batchnorm:
                 a, cache = batchnorm_forward(a, self.params['gamma' + str(i_
+ 1)],
                                           self.params['beta' + str(i + u
caches.append(cache)
             h, cache = relu_forward(a)
             caches.append(cache)
             if self.use dropout:
                 h, cache = dropout_forward(h, self.dropout_param)
                 caches.append(cache)
      # END YOUR CODE HERE
      # ------ #
      # If test mode return early
      if mode == 'test':
        return scores
      loss, grads = 0.0, {}
```

```
# ----- #
      # YOUR CODE HERE:
         Implement the backwards pass of the FC net and store the gradients
         in the grads dict, so that grads[k] is the gradient of self.
\rightarrow params [k]
      #
         Be sure your L2 regularization includes a 0.5 factor.
        BATCHNORM: Incorporate the backward pass of the batchnorm.
      #
      #
       DROPOUT: Incorporate the backward pass of dropout.
      loss, dscores = softmax_loss(scores, y)
     for i in range(self.num_layers):
         loss += 0.5 * self.reg * np.sum(self.params['W' + str(i + 1)] ** 2)
     dh, grads['W' + str(self.num_layers)], grads['b' + str(self.
onum_layers)] = affine_backward(dscores,
                 caches.pop())
     grads['W' + str(self.num_layers)] += self.reg * self.params['W' +

str(self.num_layers)]
     for i in range(self.num_layers - 1, 0, -1):
         if self.use dropout:
            dh = dropout_backward(dh, caches.pop())
         da = relu_backward(dh, caches.pop())
         if self.use_batchnorm:
            da, dgamma, dbeta = batchnorm_backward(da, caches.pop())
            grads['gamma' + str(i)] = dgamma
            grads['beta' + str(i)] = dbeta
         dh, grads['W' + str(i)], grads['b' + str(i)] = affine_backward(da,__
⇔caches.pop())
         grads['W' + str(i)] += self.reg * self.params['W' + str(i)]
                     ------ #
      # END YOUR CODE HERE
      # ----- #
     return loss, grads
```