CS289A_HW04_Prob4

March 13, 2017

1 CS 289A Homework 4

Start with program overhead: load modules (and reload them as they are modified)

```
In [1]: %load_ext autoreload
In [2]: %autoreload 2
In [3]: import HW04_utils as ut
    import numpy as np
    from scipy import special as spsp
    from matplotlib import pyplot as plt
```

Next, we give a couple paths specifying where to find the data set on the local machine. A user must change this to reflect the path to their data.

Then, load the data using the custom utilities module:

```
In [5]: # Load training data
    descriptions = ut.loaddata(DATA_PATH, BASE_DIR, 'description')
    X = ut.loaddata(DATA_PATH, BASE_DIR, 'X')
    y = ut.loaddata(DATA_PATH, BASE_DIR, 'y')

# Shuffle training data
    data = np.concatenate((X,y),axis=1)
    np.random.shuffle(data)
    X = data[:,:-1]
    y = data[:,-1]

# Normalize training data
    meanX = np.tile(np.mean(X,axis=0),(len(X),1))
    minX = np.tile(np.amin(X,axis=0),(len(X),1))
    maxX = np.tile(np.amax(X,axis=0),(len(X),1))
    X = (X-meanX) / (maxX-minX)
```

```
frac = 1/6
         n = int(len(X) - frac * len(X))
         X_{train} = X[:n]
         X \text{ val} = X[n:]
         y_train = y[:n]
         y_val = y[n:]
         X = X_train
         y = y_train
         # Load test data
         X_test = ut.loaddata(DATA_PATH, BASE_DIR, 'X_test')
         # Normalize test data
         meanXt = np.tile(np.mean(X_test,axis=0),(len(X_test),1))
         minXt = np.tile(np.amin(X_test,axis=0),(len(X_test),1))
         maxXt = np.tile(np.amax(X_test,axis=0),(len(X_test),1))
         X_test = (X_test-meanXt) / (maxXt-minXt)
1.1 Part 1
We use the following procedure to find the optimal w using batch gradient descent:
       w \leftarrow arbitrary starting point
    while J(w) > 0
       w \leftarrow w - \epsilon(2\lambda w - X^T(y - s))
    return w
  Step (1)
In [6]: w = np.zeros(len(descriptions))
  Step (2)
In [7]: def update_w_batch(w, X, y, lam, eps):
             s = spsp.expit(np.dot(X, w))
             w_{prime} = w - eps*(2*lam*w-np.dot(X.T,(y-s)))
             return w_prime
In [8]: def costfnJ(w, X, y, lam):
             s = spsp.expit(np.dot(X,w))
             J = lam*np.linalg.norm(w)**2 - np.sum(y*np.log(s) + (np.ones_like(y)-y)
             return J
In [9]: def whileloop(w, X, y, lam, eps, tol, update_fn):
             iters, Js = [], []
             J = costfnJ(w, X, y, lam)
```

Separate a validation set that is a given fraction of the training data

```
lastJ = J+1 #dummy condition to pass while condition on first run
while J>0 and i<=1e7 and np.absolute(lastJ-J)>tol:
    w_prime = update_fn(w,X,y,lam,eps)
    w = w_prime
    if i%10==0:
        if i%500000==0:
            print(str(i)+":\tJ =",str(J))
        iters.append(i)
        Js.append(J)
    lastJ = J
    J = costfnJ(w,X,y,lam)
    i+=1

return w,iters,Js
```

Step (3)

Here we try several values of hyperparameters λ and ϵ to find the optimal values. We also introduce a convergence tolerance that is used in the case that data is not linearly separable.

```
In [10]: lambdas = np.logspace(-3, 1, 5)
         epsilons = np.logspace(-5, -3, 3)
         tol = 1e-6
In [11]: # Collect loss function as f'n of iteration number for each combo
         optima_batch = {}
         LvIs batch = {}
         for lam in lambdas:
             optima batch[lam] = { }
             LvIs_batch[lam] = { }
             for eps in epsilons:
                 print("Lambda:", lam, "\tEpsilon:", eps)
                 w_star,iters,Js = whileloop(w,X,y,lam,eps,tol,update_w_batch)
                 optima_batch[lam][eps] = w_star
                 LvIs_batch[lam][eps] = [iters, Js]
Lambda: 0.001
                      Epsilon: 1e-05
          J = 3465.7359028
0:
500000:
               J = 767.198352822
1000000:
                J = 704.407443175
                J = 662.748898799
1500000:
2000000:
                J = 633.433517029
                J = 612.32966552
2500000:
3000000:
                J = 596.759285111
3500000:
                J = 585.04403722
4000000:
                J = 576.096241789
4500000:
                J = 569.178297095
5000000:
                J = 563.773879883
5500000:
                J = 559.513515784
6000000:
               J = 556.128237005
```

```
6500000: J = 553.419368524
            J = 551.238253803
7000000:
7500000:
            J = 549.472364806
            J = 548.035622922
8000000:
            J = 546.861552511
8500000:
9000000:
            J = 545.898371535
9500000:
             J = 545.105426505
10000000:
             J = 544.450573101
Epsilon: 0.0001
Lambda: 0.001
   J = 3465.7359028
500000:
           J = 563.773821406
            J = 544.450560248
1000000:
             J = 541.715934248
1500000:
Lambda: 0.001
                  Epsilon: 0.001
        J = 3465.7359028
Lambda: 0.01
                 Epsilon: 1e-05
    J = 3465.7359028
500000:
            J = 792.373230383
            J = 744.938598164
1000000:
            J = 718.85368012
1500000:
2000000:
            J = 703.760004371
            J = 694.932644597
2500000:
            J = 689.700802893
3000000:
3500000:
            J = 686.560299793
            J = 684.656004021
4000000:
            J = 683.49248145
4500000:
            J = 682.777510852
5000000:
Lambda: 0.01
                 Epsilon: 0.0001
    J = 3465.7359028
0:
500000:
        J = 682.777498732
Lambda: 0.01
                 Epsilon: 0.001
        J = 3465.7359028
Lambda: 0.1
                 Epsilon: 1e-05
    J = 3465.7359028
500000: J = 967.159795304
            J = 963.570102976
1000000:
Lambda: 0.1
                Epsilon: 0.0001
    J = 3465.7359028
Lambda: 0.1
                Epsilon: 0.001
    J = 3465.7359028
Lambda: 1.0
                Epsilon: 1e-05
      J = 3465.7359028
Lambda: 1.0
                 Epsilon: 0.0001
    J = 3465.7359028
Lambda: 1.0
                Epsilon: 0.001
        J = 3465.7359028
Lambda: 10.0
                  Epsilon: 1e-05
0:
   J = 3465.7359028
```

```
Lambda: 10.0 Epsilon: 0.0001 0: J = 3465.7359028 Lambda: 10.0 Epsilon: 0.001 0: J = 3465.7359028
```

Print out and save a list of the accuracies corresponding to the optimum w* for each combination of λ and ϵ .

```
In [12]: def HyperparameterAccs(lambdas, epsilons, optima, valdata, vallabels):
             Accs = np.zeros((len(lambdas)*len(epsilons),3))
             i=0
             for lam in optima:
                 for eps in optima[lam]:
                     w_star = optima[lam][eps]
                     probs = spsp.expit(np.dot(X_val,w_star))
                     tally = 0
                     total = 0
                      for j in range(len(probs)):
                          if probs[j] >= 0.5:
                              prob = 1
                          if probs[j] < 0.5:
                              prob = 0
                          if prob == y_val[j]:
                              tally += 1
                          total += 1
                     acc = tally/total
                     Accs[i] = [acc, lam, eps]
                     i+=1
                     print('lam = '+str(lam)+'\teps =',eps,'\t\tAccuracy: ',acc)
```

return Accs

In [13]: Accs_batch = HyperparameterAccs(lambdas,epsilons,optima_batch,X_val,y_val)

```
lam = 0.001
                   eps = 1e-05
                                               Accuracy: 0.955
lam = 0.001
                  eps = 0.0001
                                                Accuracy: 0.958
lam = 0.001
                  eps = 0.001
                                              Accuracy: 0.959
lam = 0.01
                 eps = 1e-05
                                              Accuracy: 0.952
lam = 0.01
                 eps = 0.0001
                                              Accuracy: 0.951
lam = 0.01
                 eps = 0.001
                                              Accuracy: 0.952
lam = 0.1
                 eps = 1e-05
                                             Accuracy: 0.935
lam = 0.1
                eps = 0.0001
                                              Accuracy: 0.935
lam = 0.1
                 eps = 0.001
                                             Accuracy: 0.935
lam = 1.0
                eps = 1e-05
                                             Accuracy: 0.922
lam = 1.0
                eps = 0.0001
                                              Accuracy: 0.921
lam = 1.0
                eps = 0.001
                                             Accuracy: 0.921
lam = 10.0
                eps = 1e-05
                                              Accuracy: 0.913
lam = 10.0
                 eps = 0.0001
                                               Accuracy: 0.913
```

```
lam = 10.0 eps = 0.001
                                          Accuracy: 0.913
In [14]: print(optima_batch[0.001][0.001])
        print (Accs_batch)
\begin{bmatrix} -9.76375855 & 13.46225692 & -5.2732488 \end{bmatrix}
                                        -66.13364483
                                                      10.78171502
                                                      2.76490805
  19.51816531 -26.64844479 146.9055984
                                         -5.99243176
  19.14298793
                3.538464251
[[ 9.55000000e-01 1.00000000e-03
                                   1.0000000e-05]
 9.58000000e-01
                   1.00000000e-03 1.0000000e-04]
 [ 9.59000000e-01 1.0000000e-03 1.00000000e-03]
 9.52000000e-01 1.0000000e-02 1.0000000e-05
 9.51000000e-01 1.0000000e-02 1.00000000e-04
 9.52000000e-01 1.0000000e-02 1.0000000e-031
 [ 9.35000000e-01
                   1.00000000e-01 1.0000000e-05]
 9.35000000e-01 1.0000000e-01 1.0000000e-04
 [ 9.35000000e-01 1.00000000e-01 1.00000000e-03]
 9.22000000e-01 1.0000000e+00 1.0000000e-051
 [ 9.21000000e-01 1.00000000e+00 1.00000000e-04]
 [ 9.21000000e-01
                  1.00000000e+00 1.0000000e-03]
 [ 9.13000000e-01 1.00000000e+01 1.00000000e-05]
 9.13000000e-01
                   1.00000000e+01 1.0000000e-04]
 9.13000000e-01 1.00000000e+01 1.00000000e-03]
```

Plot the loss function vs. iteration number for the best combination of λ , ϵ .

In [16]: plot_LFvIt (Accs_batch, LvIs_batch)

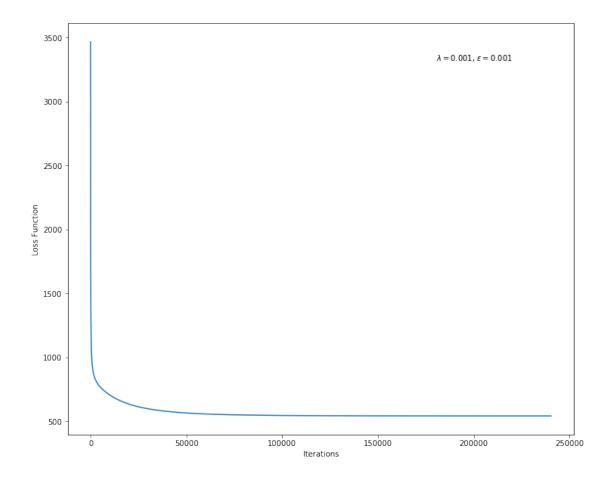
```
In [15]: def plot_LFvIt(Accs,LvIs):
    fig = plt.figure(figsize=(12,10))
    plt.clf()

# Find the most accurate lam, eps combo

imax = np.argmax(Accs[:,0])

lam = Accs[imax,1]
    eps = Accs[imax,2]

Iters = LvIs[lam][eps][0]
    LossFn = LvIs[lam][eps][1]
    plt.plot(Iters,LossFn)
    plt.xlabel('Iterations')
    plt.ylabel('Loss Function')
    plt.text(0.75*(np.amax(Iters)-np.amin(Iters))+np.amin(Iters),0.95*(np.amin))
```



1.2 Part 2

Instead of batch descent, we can use the following procedure to find the optimal w using stochastic gradient descent:

- (1) $w \leftarrow \text{arbitrary starting point}$
- (2) while J(w) > 0 $w \leftarrow w - \epsilon(2\lambda w - X_i^T(y_i - s_i))$
- (3) return w

Step (1) - Same as in batch gradient descent

```
In [17]: w = np.zeros(len(descriptions))
```

Step (2) - we can reuse the functions for calculating the cost function and the looping process defined in part 1, step 2; we define a new function for the stochastic update rule

```
w_prime = w - eps*(2*lam*w-X_i*(y_i-s_i))
return w_prime
```

Step (3)

Again, we try several values of hyperparameters λ and ϵ to find the optimal values and we introduce a convergence tolerance that is used in the case that data is not linearly separable.

```
In [19]: tol = 1e-9
In [20]: # Collect loss function as f'n of iteration number for each combo
        optima_stoch = {}
        LvIs stoch = {}
        for lam in lambdas:
            optima_stoch[lam] = {}
            LvIs_stoch[lam] = { }
            for eps in epsilons:
                print("Lambda:", lam, "\tEpsilon:", eps)
                w_star,iters,Js = whileloop(w,X,y,lam,eps,tol,update_w_stoch)
                optima_stoch[lam][eps] = w_star
                LvIs_stoch[lam][eps] = [iters, Js]
                print(iters[len(iters)-1])
Lambda: 0.001
                     Epsilon: 1e-05
     J = 3465.7359028
500000: J = 3364.07844802
751460
Lambda: 0.001
                     Epsilon: 0.0001
        J = 3465.7359028
0:
407340
Lambda: 0.001
                    Epsilon: 0.001
    J = 3465.7359028
500000:
             J = 1780.02392783
957690
Lambda: 0.01
                    Epsilon: 1e-05
    J = 3465.7359028
            J = 3368.30796372
500000:
672060
Lambda: 0.01
                    Epsilon: 0.0001
         J = 3465.7359028
0:
359080
Lambda: 0.01
                    Epsilon: 0.001
         J = 3465.7359028
0:
500000:
             J = 2785.21595913
1000000:
              J = 2787.32786271
1500000:
              J = 2785.51122138
2000000:
              J = 2785.11129489
             J = 2787.0705809
2500000:
3000000:
              J = 2785.28963853
```

```
J = 2785.803959
3500000:
             J = 2787.22372734
4000000:
4500000:
             J = 2788.00847534
5000000:
             J = 2786.29573236
            J = 2785.48439103
5500000:
6000000:
             J = 2788.66480945
6005030
Lambda: 0.1 Epsilon: 1e-05
   J = 3465.7359028
214990
Lambda: 0.1 Epsilon: 0.0001
   J = 3465.7359028
           J = 3365.64013098
500000:
            J = 3365.10773544
1000000:
             J = 3365.41298886
1500000:
             J = 3365.44933867
2000000:
            J = 3365.16683129
2500000:
            J = 3365.2569922
3000000:
3238330
Lambda: 0.1
                Epsilon: 0.001
J = 3465.7359028
            J = 3365.02642054
500000:
1000000:
             J = 3364.96104161
             J = 3365.04799404
1500000:
2000000:
             J = 3365.24070245
            J = 3364.91826271
2500000:
             J = 3365.17746645
3000000:
             J = 3365.1079893
3500000:
             J = 3364.48461212
4000000:
4500000:
             J = 3366.13618097
             J = 3364.66628881
5000000:
             J = 3365.56950205
5500000:
6000000:
             J = 3365.13812904
             J = 3366.38725617
6500000:
7000000:
             J = 3366.13648221
             J = 3365.20638599
7500000:
             J = 3365.1508241
8000000:
8500000:
             J = 3364.84247723
9000000:
             J = 3365.38729735
             J = 3365.47776594
9500000:
10000000:
              J = 3364.7427037
10000000
Lambda: 1.0 Epsilon: 1e-05
  J = 3465.7359028
         J = 3455.19947615
500000:
892190
Lambda: 1.0 Epsilon: 0.0001
J = 3465.7359028
```

```
500000: J = 3455.22516758
806130
Lambda: 1.0
                   Epsilon: 0.001
         J = 3465.7359028
               J = 3455.45095814
500000:
1000000:
                J = 3455.32709589
1500000:
               J = 3454.99078137
2000000:
                J = 3455.46400094
2500000:
               J = 3454.41694654
3000000:
                J = 3454.82719222
                J = 3455.04570563
3500000:
4000000:
                J = 3455.11178217
4500000:
                J = 3455.63697838
                J = 3455.37833326
5000000:
                J = 3455.22887265
5500000:
6000000:
                J = 3454.9626366
6500000:
                J = 3455.19824855
7000000:
                J = 3455.31077341
7500000:
                J = 3455.36412691
8000000:
                J = 3455.8766627
8500000:
                J = 3455.37583182
                J = 3455.11135474
9000000:
9500000:
               J = 3455.25644384
10000000:
                J = 3455.08109678
10000000
Lambda: 10.0
                     Epsilon: 1e-05
         J = 3465.7359028
355830
Lambda: 10.0
                     Epsilon: 0.0001
         J = 3465.7359028
500000:
             J = 3464.65317259
974090
Lambda: 10.0
                     Epsilon: 0.001
         J = 3465.7359028
500000:
               J = 3464.63245865
                J = 3464.71713815
1000000:
1500000:
               J = 3464.75740153
2000000:
                J = 3464.74406301
2500000:
                J = 3464.59253951
3000000:
                J = 3464.63204545
3500000:
                J = 3464.68100638
4000000:
                J = 3464.7078865
4500000:
                J = 3464.61943668
5000000:
                J = 3464.63747874
5500000:
                J = 3464.67095198
6000000:
               J = 3464.73626162
6500000:
               J = 3464.63546903
7000000:
               J = 3464.63242981
```

```
7500000: J = 3464.47215421

8000000: J = 3464.76204115

8500000: J = 3464.5792482

9000000: J = 3464.60974749

9500000: J = 3464.77891622

10000000: J = 3464.72769749

10000000
```

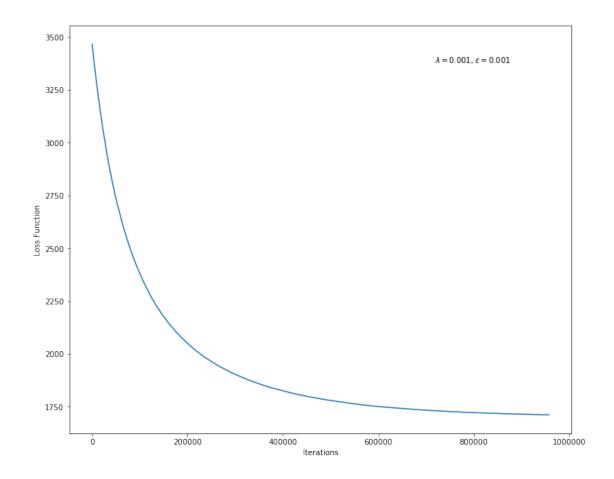
Print out a list of the optimum w* for each combination of λ and ϵ . Save the accuracies in a list.

In [21]: Accs_stoch = HyperparameterAccs(lambdas,epsilons,optima_stoch,X_val,y_val)

lam =	0 001	eps = 1e-05	Accuracy: 0.902
-		i	<u> -</u>
lam =	0.001	eps = 0.0001	Accuracy: 0.903
lam =	0.001	eps = 0.001	Accuracy: 0.913
lam =	0.01	eps = 1e-05	Accuracy: 0.902
lam =	0.01	eps = 0.0001	Accuracy: 0.903
lam =	0.01	eps = 0.001	Accuracy: 0.904
lam =	0.1	eps = 1e-05	Accuracy: 0.901
lam =	0.1	eps = 0.0001	Accuracy: 0.901
lam =	0.1	eps = 0.001	Accuracy: 0.9
lam =	1.0	eps = 1e-05	Accuracy: 0.903
lam =	1.0	eps = 0.0001	Accuracy: 0.902
lam =	1.0	eps = 0.001	Accuracy: 0.899
lam =	10.0	eps = 1e-05	Accuracy: 0.904
lam =	10.0	eps = 0.0001	Accuracy: 0.889
lam =	10.0	eps = 0.001	Accuracy: 0.893

Plot the loss function vs. iteration number for the best combination of λ , ϵ .

```
In [22]: plot_LFvIt(Accs_stoch, LvIs_stoch)
```



1.3 Part 3

Now we wish to repeat part 2 (stochastic gradient descent) but using a variable ϵ . We can accomplish this by redefining our while loop to decrease ϵ such that $\epsilon \propto 1/t$.

return w, iters, Js

Then, we call that function using the same procedure used before.

```
In [24]: # Collect loss function as f'n of iteration number for each combo
        optima_stoch_deceps = {}
        LvIs_stoch_deceps = {}
        for lam in lambdas:
            optima_stoch_deceps[lam] = { }
            LvIs_stoch_deceps[lam] = {}
            for eps in epsilons:
               print("Lambda:", lam, "\tEpsilon:", eps)
               w_star,iters,Js = whileloop_deceps(w,X,y,lam,eps,tol,update_w_stoc
               optima_stoch_deceps[lam][eps] = w_star
               LvIs_stoch_deceps[lam][eps] = [iters, Js]
               print(iters[len(iters)-1])
Lambda: 0.001
                    Epsilon: 1e-05
J = 3465.7359028
Lambda: 0.001
                   Epsilon: 0.0001
    J = 3465.7359028
0:
1400
Lambda: 0.001 Epsilon: 0.001
0:
        J = 3465.7359028
3500
Lambda: 0.01
                  Epsilon: 1e-05
      J = 3465.7359028
0:
Lambda: 0.01
                  Epsilon: 0.0001
       J = 3465.7359028
0:
1430
Lambda: 0.01 Epsilon: 0.001
      J = 3465.7359028
0:
Lambda: 0.1
                 Epsilon: 1e-05
     J = 3465.7359028
0:
840
Lambda: 0.1 Epsilon: 0.0001
0:
     J = 3465.7359028
990
Lambda: 0.1
                 Epsilon: 0.001
      J = 3465.7359028
0:
23150
Lambda: 1.0
                 Epsilon: 1e-05
```

J = 3465.7359028

```
290
Lambda: 1.0 Epsilon: 0.0001
     J = 3465.7359028
2940
Lambda: 1.0 Epsilon: 0.001
     J = 3465.7359028
3020
Lambda: 10.0
                Epsilon: 1e-05
    J = 3465.7359028
1180
Lambda: 10.0
                 Epsilon: 0.0001
0:
    J = 3465.7359028
3930
                Epsilon: 0.001
Lambda: 10.0
0:
        J = 3465.7359028
3880
```

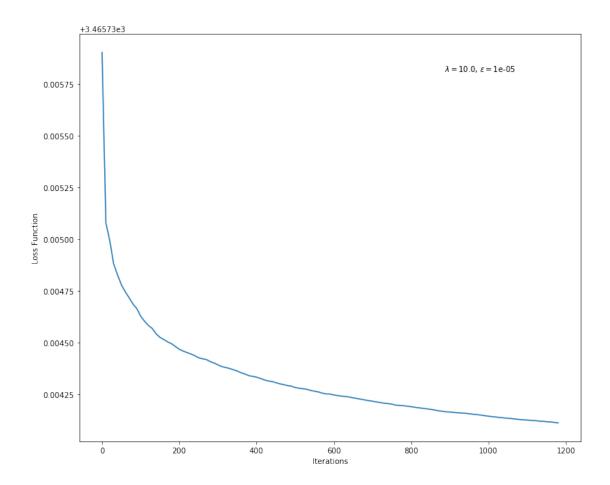
Print out a list of the optimum w* for each combination of λ and ϵ . Save the accuracies in a list.

In [25]: Accs_stoch_deceps = HyperparameterAccs(lambdas,epsilons,optima_stoch_deceps

lam =	0.001	eps = 1e-05	Accuracy: 0.88
lam =	0.001	eps = 0.0001	Accuracy: 0.884
lam =	0.001	eps = 0.001	Accuracy: 0.875
lam =	0.01	eps = 1e-05	Accuracy: 0.884
lam =	0.01	eps = 0.0001	Accuracy: 0.769
lam =	0.01	eps = 0.001	Accuracy: 0.854
lam =	0.1	eps = 1e-05	Accuracy: 0.877
lam =	0.1	eps = 0.0001	Accuracy: 0.871
lam =	0.1	eps = 0.001	Accuracy: 0.883
lam =	1.0	eps = 1e-05	Accuracy: 0.761
lam =	1.0	eps = 0.0001	Accuracy: 0.874
lam =	1.0	eps = 0.001	Accuracy: 0.869
lam =	10.0	eps = 1e-05	Accuracy: 0.892
lam =	10.0	eps = 0.0001	Accuracy: 0.864
lam =	10.0	eps = 0.001	Accuracy: 0.887

Plot the loss function vs. iteration number for the best combination of λ, ϵ .

```
In [26]: plot_LFvIt(Accs_stoch_deceps, LvIs_stoch_deceps)
```



Finally, we use our most successful training algorithm (in this case, batch gradient descent for $\lambda = 0.001$ and $\epsilon = 0.001$) to predict on the test data.

```
In [51]: lam,eps = 0.001,0.001
    w_star = optima_batch[lam][eps]
    preds = spsp.expit(np.dot(X_test,w_star))
    predictions = np.rint(preds)
```

We save these predictions to the csv file for Kaggle submission.

Noting that this submission on Kaggle produced a test error of almost 12% (significantly more than the training error of 3% for the same hyperparameters), I chose a new set of hyperparameters with λ greater than the first submission. The intention of this was to reduce overfitting, as I had presumably been overfitting, giving an excellent training error but mediocre test error.

```
In [50]: lam,eps = 0.1,0.001
    w_star2 = optima_batch[lam][eps]
```

```
preds2 = spsp.expit(np.dot(X_test,w_star2))
predictions2 = np.rint(preds2)
```

Again, we save these predictions to the csv file for Kaggle submission.

The hypothesis seems accurate. This submission (username mnegus) gave a score of 95.565%.