Programming assignment 6: Optimization: Logistic regression

```
In [1]: import numpy as np
   import matplotlib.pyplot as plt
   %matplotlib inline

from sklearn.datasets import load_breast_cancer
   from sklearn.model_selection import train_test_split
   from sklearn.metrics import accuracy_score, fl_score
```

Your task

In this notebook code skeleton for performing logistic regression with gradient descent is given. Your task is to complete the functions where required. You are only allowed to use built-in Python functions, as well as any numpy functions. No other libraries / imports are allowed.

For numerical reasons, we actually minimize the following loss function

$$\mathcal{L}(\mathbf{w}) = rac{1}{N} NLL(\mathbf{w}) + \lambda ||\mathbf{w}||_2^2$$

where $NLL(\mathbf{w})$ is the negative log-likelihood function, as defined in the lecture (Eq. 33)

Load and preprocess the data

In this assignment we will work with the UCI ML Breast Cancer Wisconsin (Diagnostic) dataset https://goo.gl/U2Uwz2).

Features are computed from a digitized image of a fine needle aspirate (FNA) of a breast mass. They describe characteristics of the cell nuclei present in the image. There are 212 malignant examples and 357 benign examples.

```
In [30]: X, y = load_breast_cancer(return_X_y=True)

# Add a vector of ones to the data matrix to absorb the bias term
X = np.hstack([np.ones([X.shape[0], 1]), X])

# Set the random seed so that we have reproducible experiments
np.random.seed(123)

# Split into train and test
test_size = 0.3
X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=test_size)
```

Task 1: Implement the sigmoid function

Task 2: Implement the negative log likelihood

As defined in Eq. 33

Computing the loss function $\mathcal{L}(\mathbf{w})$ (nothing to do here)

```
In [5]: def compute_loss(X, y, w, lmbda):
            Negative Log Likelihood of the Logistic Regression.
            Parameters
            X : array, shape [N, D]
                (Augmented) feature matrix.
            y : array, shape [N]
                Classification targets.
            w : array, shape [D]
                Regression coefficients (w[0] is the bias term).
            lmbda : float
                L2 regularization strength.
            Returns
            loss : float
                Loss of the regularized logistic regression model.
            # The bias term w[0] is not regularized by convention
            return negative_log_likelihood(X, y, w) / len(y) + lmbda * np.linalg.norm(w
         [1:])
```

Task 3: Implement the gradient $abla_{\mathbf{w}} \mathcal{L}(\mathbf{w})$

Make sure that you compute the gradient of the loss function $\mathcal{L}(\mathbf{w})$ (not simply the NLL!)

```
In [39]: def get_gradient(X, y, w, mini_batch_indices, lmbda):
             Calculates the gradient (full or mini-batch) of the negative log likelihood
         w.r.t. w.
             Parameters
             X : array, shape [N, D]
                 (Augmented) feature matrix.
             y : array, shape [N]
                 Classification targets.
             w : array, shape [D]
                 Regression coefficients (w[0] is the bias term).
             mini_batch_indices: array, shape [mini_batch_size]
                 The indices of the data points to be included in the (stochastic) calcu
         lation of the gradient.
                 This includes the full batch gradient as well, if mini batch indices =
         np.arange(n_train).
             lmbda: float
                 Regularization strentgh. lmbda = 0 means having no regularization.
             Returns
             dw : array, shape [D]
                 Gradient w.r.t. w.
             return -(y[mini batch indices] - sigmoid(X[mini batch indices,:].dot(w))).d
         ot(X[mini_batch_indices,:])/len(mini_batch_indices) + 2*lmbda*w
```

Train the logistic regression model (nothing to do here)

```
In [7]: def logistic_regression(X, y, num_steps, learning_rate, mini_batch_size, lmbda,
        verbose):
            Performs logistic regression with (stochastic) gradient descent.
            Parameters
            X : array, shape [N, D]
                (Augmented) feature matrix.
            y : array, shape [N]
                Classification targets.
            num steps : int
                Number of steps of gradient descent to perform.
            learning rate: float
                The learning rate to use when updating the parameters w.
            mini batch size: int
                The number of examples in each mini-batch.
                If mini_batch_size=n_train we perform full batch gradient descent.
            lmbda: float
                Regularization strentgh. lmbda = 0 means having no regularization.
            verbose : bool
                Whether to print the loss during optimization.
            Returns
            w : array, shape [D]
                Optimal regression coefficients (w[0] is the bias term).
            trace: list
                Trace of the loss function after each step of gradient descent.
            trace = [] # saves the value of loss every 50 iterations to be able to plot
        it later
            n train = X.shape[0] # number of training instances
            w = np.zeros(X.shape[1]) # initialize the parameters to zeros
            # run gradient descent for a given number of steps
            for step in range(num steps):
                permuted idx = np.random.permutation(n train) # shuffle the data
                # go over each mini-batch and update the paramters
                # if mini_batch_size = n_train we perform full batch GD and this loop r
        uns only once
                for idx in range(0, n_train, mini_batch_size):
                    # get the random indices to be included in the mini batch
                    mini_batch_indices = permuted_idx[idx:idx+mini_batch_size]
                    gradient = get_gradient(X, y, w, mini_batch_indices, lmbda)
                    # update the parameters
                    w = w - learning rate * gradient
                # calculate and save the current loss value every 50 iterations
                if step % 50 == 0:
                    loss = compute_loss(X, y, w, lmbda)
                    trace.append(loss)
                    # print loss to monitor the progress
                    if verbose and step % 50 == 0:
                         print('Step \{0\}, loss = \{1:.4f\}'.format(step, loss))
            return w, trace
```

Task 4: Implement the function to obtain the predictions

Full batch gradient descent

```
In [9]: # Change this to True if you want to see loss values over iterations.
         verbose = False
In [36]: n train = X train.shape[0]
         w_full, trace_full = logistic_regression(X_train,
                                                    y train,
                                                    num steps=8000,
                                                    learning rate=1e-5,
                                                   mini batch size=n train,
                                                   lmbda=0.1,
                                                   verbose=verbose)
In [37]: n train = X train.shape[0]
         w minibatch, trace minibatch = logistic regression(X train,
                                                              y train,
                                                              num_steps=8000,
                                                              learning_rate=1e-5,
                                                              mini_batch_size=50,
                                                              lmbda=0.1,
                                                              verbose=verbose)
```

Our reference solution produces, but don't worry if yours is not exactly the same.

```
Full batch: accuracy: 0.9240, f1_score: 0.9384 Mini-batch: accuracy: 0.9415, f1_score: 0.9533
```

0.2

```
In [38]: y_pred_full = predict(X_test, w_full)
          y_pred_minibatch = predict(X_test, w_minibatch)
          print('Full batch: accuracy: {:.4f}, f1_score: {:.4f}'
                .format(accuracy_score(y_test, y_pred_full), f1_score(y_test, y_pred_full)
          )))
          print('Mini-batch: accuracy: {:.4f}, f1_score: {:.4f}'
                .format(accuracy_score(y_test, y_pred_minibatch), f1_score(y_test, y_pred
          minibatch)))
         Full batch: accuracy: 0.9240, f1_score: 0.9384
         Mini-batch: accuracy: 0.9415, f1_score: 0.9528
In [40]: plt.figure(figsize=[15, 10])
          plt.plot(trace full, label='Full batch')
          plt.plot(trace_minibatch, label='Mini-batch')
          plt.xlabel('Iterations * 50')
          plt.ylabel('Loss $\mathcal{L}(\mathbf{w})$')
          plt.legend()
          plt.show()

    Full batch

                                                                                        Mini-batch
            0.9
            0.8
            0.7
          (w)
            0.6
          Loss
            0.5
            0.4
            0.3
```

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60

80

Iterations * 50

100

120

140

160