## COMP/INDR 421/521 HW03: Multiclass Multilayer Perceptron

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The purpose of this assignment was to implement a multiclass multilayer perceptron algorithm in R

The problem consisted of four main parts:

- 1- Initialization
- 2- Data generation
- 3- <u>Linear Discrimination Algorithm:</u> Where the multiclass multilayer perceptron algorithm is applied on the data set generated in the first part.
- 4- Algorithm performance evaluation

Following, more details will be discussed about each part accompanied by snippets from my source code.

## • Initialization

```
# mean parameters
class_means \leftarrow matrix(c(+2.0, +2.0,
                         -4.0, -4.0,
                         -2.0, +2.0,
                         +4.0, -4.0,
                         -2.0, -2.0.
                         +4.0, +4.0,
                         +2.0, -2.0,
-4.0, +4.0), 2, 8)
# covariance parameters
class_covariances <- array(c(+0.8, -0.6, -0.6, +0.8,
                              +0.4, +0.0, +0.0, +0.4,
                              +0.8, +0.6, +0.6, +0.8,
                              +0.4, +0.0, +0.0, +0.4,
                              +0.8, -0.6, -0.6, +0.8,
                              +0.4, +0.0, +0.0, +0.4,
                              +0.8, +0.6, +0.6, +0.8,
                              +0.4, +0.0, +0.0, +0.4), c(2, 2, 8)
# sample sizes
class_sizes <- c(100, 100, 100, 100)
```

seen in the previous code snippet, actual class parameters are initialized for all four classes.

As can be

$$class\_means[,i] = [\mu_{i1} \quad \mu_{i2}]$$

covariance matrix 
$$\mathbf{\Sigma} \equiv class\_covariances[,,i] = \begin{bmatrix} \sigma_{i1}^2 & \sigma_{i12} \\ \sigma_{i21} & \sigma_{i2}^2 \end{bmatrix}$$

## • Data generation

```
# generate random samples
points1 <- mvrnorm(n = class_sizes[1] / 2, mu = class_means[,1], Sigma = class_covariances[,,1])
points2 <- mvrnorm(n = class_sizes[1] / 2, mu = class_means[,2], Sigma = class_covariances[,,2])
points3 <- mvrnorm(n = class_sizes[2] / 2, mu = class_means[,3], Sigma = class_covariances[,,3])
points4 <- mvrnorm(n = class_sizes[2] / 2, mu = class_means[,4], Sigma = class_covariances[,,4])
points5 <- mvrnorm(n = class_sizes[3] / 2, mu = class_means[,5], Sigma = class_covariances[,,5])
points6 <- mvrnorm(n = class_sizes[3] / 2, mu = class_means[,6], Sigma = class_covariances[,,6])
points7 <- mvrnorm(n = class_sizes[4] / 2, mu = class_means[,7], Sigma = class_covariances[,,7])
points8 <- mvrnorm(n = class_sizes[4] / 2, mu = class_means[,8], Sigma = class_covariances[,,7])
points8 <- mvrnorm(n = class_sizes[4] / 2, mu = class_means[,8], Sigma = class_covariances[,,8])

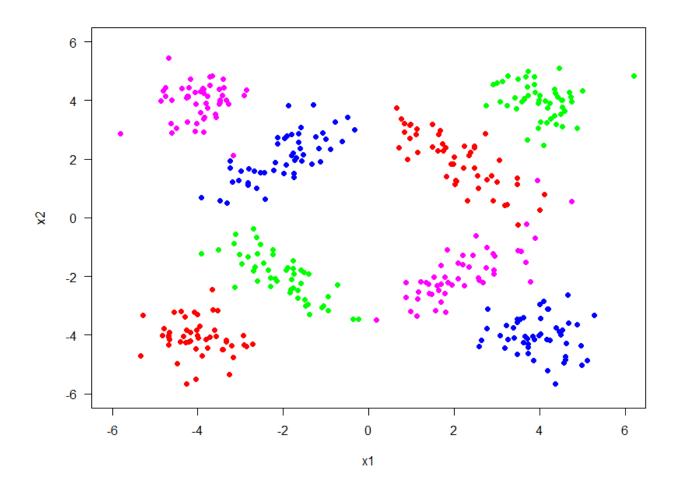
X <- rbind(points1, points2, points3, points4, points5, points6, points7, points8)

colnames(X) <- c("x1", "x2")

# generate corresponding labels
y <- c(rep(1, class_sizes[1]), rep(2, class_sizes[2]), rep(3, class_sizes[3]), rep(4, class_sizes[4]))

# write data to a file
write.csv(x = cbind(X, y), file = "Hw3_data_set.csv", row.names = FALSE)</pre>
```

This part consists of generating the data set (including input/feature values and their corresponding outputs/labels) from bivariate Gaussian densities. Then, the generated data set was plotted as the following graph:



## • Multiclass Multilayer Perceptron Algorithm

This part contains multiple steps as follows:

- Importing data and initializing algorithm parameters

- Defining softmax

$$softmax \equiv y_i = \hat{p}(C_i|x) = \frac{\exp(w_i^T x)}{\sum_{j=1}^K \exp(w_j^T x)}$$
 # define the softmax function softmax <- function(x, w, k, c){ denomSum <- 0 for(i in 1:k){ denomSum <- denomSum +  $\exp(x \% \% w[,i])$ } print( $\exp(x \% \% w[,c])/\text{denomSum})$  return ( $\exp(x \% \% w[,c])/\text{denomSum})$ }

- Classification

Where the class scoring function is the softmax function.

- These steps are repeated iteratively with randomly initialized w and v values, and an update rule is equal to gradient function multiplied by the learning rate  $(\eta)$ 

```
while (1) {
 print(pasteO("running iteration#", iteration))
for (i in sample(N)) { # sample(N) gives a random ordering of numbers 1 to N, we want to access data in a random order
      calculate hidden nodes
    Z[i,] \leftarrow sigmoid(c(1, X[i,]) %*% W)
    y_predicted[i,] \leftarrow sapply(1:K, function(c){softmax(matrix(c(1, Z[i,]), 1, H + 1), v, K, c)})
      if(y_truth[i]
      v[,e] <- v[,e] + eta * (1 - y_predicted[i, y_truth[i]]) * c(1, z[i,])
} else [</pre>
        v[,e] "<- v[,e] + eta * -1 * y_predicted[i, e] * c(1, Z[i,])</pre>
      }
    for (h in 1:H) {
      sum_error <- 0
      for(class in 1:K){
        if(y_truth[i] == class){
          sum_error <- sum_error + (1 - y_predicted[i, class]) * v[h, class]
          sum_error <- sum_error - y_predicted[i, class] * v[h, class]</pre>
      W[,h] \leftarrow W[,h] + eta * sum_error * Z[i, h] * (1 - Z[i, h]) * c(1, X[i,])
    }
  error
  for(i in 1:N){
    error <- error + log(y_predicted[i, y_truth[i]] + 1e-100)
  objective_values <- c(objective_values, -error)
     (iteration != 1){}
    if ((abs(objective_values[iteration] - objective_values[iteration - 1]) < epsilon || iteration >= max_iteration)) {
  iteration <- iteration + 1
```

Then, class classification is determined as same as the best scoring class.

```
y_predicted_class <- c()
for(i in 1:N){
   y_predicted_class <- rbind(y_predicted_class, which.max(y_predicted[i,]))
}</pre>
```

The values of w parameters after the iterations were as follows:

```
> print(W)
[1,] [,2] [,3] [,4] [,5] [,6] [,7] [,8] [1,] 6.6421114 16.183535 13.568445 9.606574 10.882318 9.457999 10.258945 4.246377
                                                                                                  0.9366378 7.515359 7.851477
[2,] -0.3973043 -3.247718 -2.738522 -1.675305 1.872582 2.172038 2.069357 -3.938547 -6.1934437 1.340761 -1.434061 [3,] -1.3745071 -1.888773 -2.414250 -1.643067 2.734930 1.660503 -1.532246 5.863815 0.2611013 1.965851 1.609761
                                                                                             [,19]
                                 [,14]
                                                         [,16]
                                                                     [,17]
                                                                                   [,18]
          [,12]
                       [,13]
                                              [,15]
                                                                                                       [,20]
[1,] 7.549591 6.08786684 8.358916 9.0584142
                                                    9.649272
                                                                6.542305 3.73750780 8.169214 6.101761
[2,] -1.556821 0.33129887 2.149746 -0.4373296 -2.395928 -3.045919 -0.09817704 2.596878 1.877990 [3,] 1.552035 0.03377539 1.107313 2.8371040 1.684598 2.114441 7.78099965 4.469213 3.155037
> print(v)
                [.1]
                            [,2]
                                         Γ.31
        1.22578898 -3.2752719 -0.8138384
                                                 2.85491376
 [1.]
         6.45466756 -0.5562685 -6.0946715
                                                 0.21017778
 [2.]
         6.48803439 0.2287617 -5.1419805
 [3,]
                                                -1.57672859
         4.42965542
                      2.4213326 -6.8581354
                                                 0.01189458
 [5,] -10.46428810 9.3609146 -1.9226784
                                                 3.02511974
 [6,]
       -6.29895900 1.9403376 2.3865458
                                                 1.98564019
       -3.88727056 1.7979948 3.2354870
                                                -1.14258910
       3.62658762 3.9069454 3.4525445 -10.98362312
 [8,]
 [9.]
      -2.29755614 1.9565008
                                  2.0482115
                                                -1.70898779
                                  2.7969166
[10,]
      -6.45489239
                     3.2503260
                                                 0.38510835
[11,]
         0.05263743 -4.2880708
                                  2.0518933
                                                 2.18187222
[12.]
        1.22027299 -5.4567859 2.4293527
                                                 1.81034826
[13,]
       -0.11906993 -4.5325438 2.4379738
                                                 2.21387609
       -3.72530229 -1.6997933 2.2524160
[14,]
                                                 3.18287697
       -3.43953756 -2.1092468 2.6597211
[15,]
                                                 2.87995519
         2.86544434 -8.5651483 2.6571417
                                                 3.05699048
[16,]
         3.35122658 -4.6881655
                                  2.2584355
Γ17. ]
                                                -0.92444250
[18,]
         0.64322297
                      2.7412882 2.9941344
                                                -6.39329884
         3.02454686 5.3580628 -3.2904141
                                                -5.07990564
[19,]
[20,]
         0.34256288 3.7710632 -5.5250503
                                                 1.41472120
         2.66228065 2.3268913 -6.3755187
[21,]
                                                1.37370073
```

- The last step is to evaluate the performance of the algorithm by finding the confusion matrix, plotting the objective values to check convergence and visualizing the classification by plotting the points based on their predicted labels, with the decision boundaries.

