# Assignment 2, COMP4702

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#### Question 4.2

```
1 % Q2
a = randn(200, 2);
a b = a + 4;
  c = a;
  c(:, 1) = 3 * c(:, 1);
  c = c - 4;
  d = [a; b];
  e = [a; b; c];
12
  % hold on;
  \% plot (a(:, 1), a(:, 2), '+');
  \% plot (b(:, 1), b(:, 2), 'o');
  \% plot (c(:, 1), c(:, 2), '*');
17
  % Use the first dataset
  data = e;
19
20
21
  figure;
22
  subplot (3, 2, 1);
  hold on;
  % ksdensity(data, 'PlotFcn', 'contour');
  plot(a(:, 1), a(:, 2), '+');
  {\tt plot}\,(\,b\,(:\,,\ 1)\,\,,\ b\,(:\,,\ 2)\,\,,\ \lq\circ\,\lq)\,;
27
  plot(c(:, 1), c(:, 2), '*');
29
  title ('Contours Overlay');
31
  % Calculate the grid
  [f, xi] = ksdensity(data); x = linspace(min(xi(:, 1)), max(xi(:, 1)))
     ));
  y = linspace(min(xi(:, 2)), max(xi(:, 2)));
  [xq,yq] = meshgrid(x,y);
  z = griddata(xi(:, 1), xi(:, 2), f, xq, yq);
36
37
  % We now have x, y, z that can be used to get the gradient at any
38
     point
  contour(x, y, z);
```

```
x \lim ([-10, 10]);
  y \lim ([-10, 10]);
  hold off;
43
44
  copy = data;
  new_points = copy;
46
47
   for i = 1:5
48
       new_points = step(new_points, x, y, z);
49
       subplot(3, 2, i + 1);
50
       hold on;
51
       % ksdensity(data, 'PlotFcn', 'contour');
52
       scatter(new_points(:, 1), new_points(:, 2));
53
       title (sprintf ('Step %d', i));
54
       x \lim ([-10, 10]);
55
       y \lim ([-10, 10]);
56
       hold off;
57
58
  end
59
60
61
   function dist = euclid_distance(point1, point2)
62
        dist = \operatorname{sqrt}((\operatorname{point1}(1) - \operatorname{point2}(1))^2 + (\operatorname{point1}(2) - \operatorname{point2}(2))
63
           ) ^2);
  end
64
   function new_points = step(points, x, y, z)
66
       % Calibration factor (lambda)
67
       max_distance = 1.8;
68
69
       new_points = zeros(length(points), 2);
70
       for i = 1:length (points)
71
            point = points(i, :);
72
73
            within_range = zeros(length(points), 1);
74
75
            numerator = 0;
76
            denominator = 0;
77
78
            for j = 1: length (points)
79
                 other_point = points(j, :);
80
                 if euclid_distance(point, other_point) < max_distance
81
                      within_range(j) = 1;
                      \% weight = interp2(x, y, z, other_point(1),
83
                         other_point(2));
```

```
84
                   % Calculate part of sum
85
                   distance = euclid_distance(point, other_point);
86
                   numerator = numerator + (distance * other_point);
87
                    denominator = denominator + distance;
               end
89
           end
91
           mx = numerator / denominator;
92
           new_points(i, :) = mx(1, :);
93
      end
94
  end
95
```

### Question 4.3

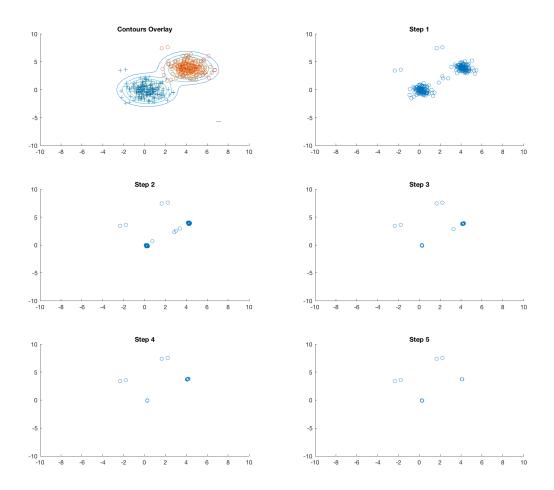


Figure 1: 2 Classes, Lambda = 1.8

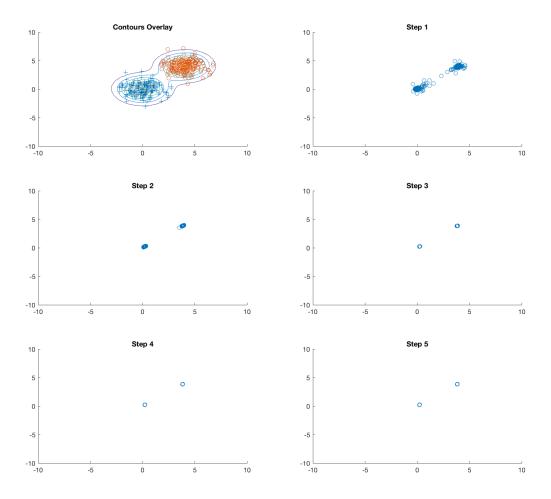


Figure 2: 2 Classes, Lambda = 3

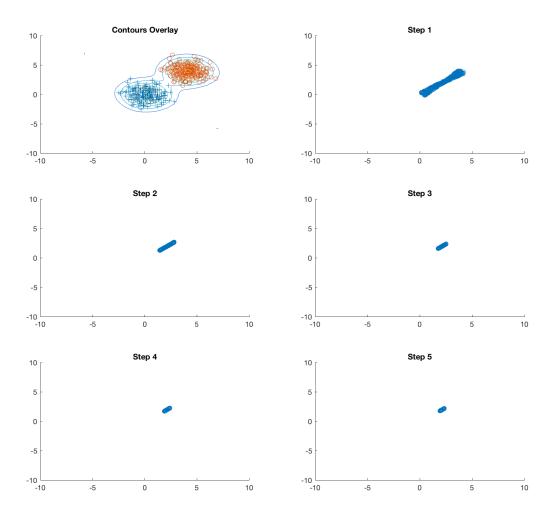


Figure 3: 2 Classes, Lambda = 5

For the 2 class problem, a lambda value set to 3 provided the best result. A lambda of 1.8 cause a few outliers not to shift towards the mean value. Whereas a lambda value of 5 caused all the points to shift towards one of the means, which is not desired. A lambda value of 3 shifted all the points to the two mean values without leaving outliers, thus is the best lambda value out of the three.

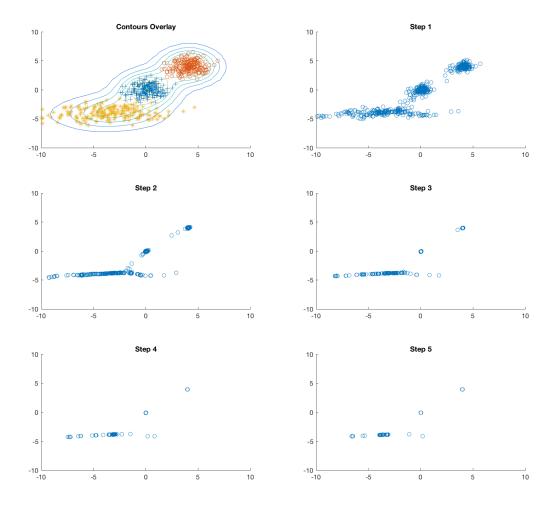


Figure 4: 3 Classes, Lambda = 1.8

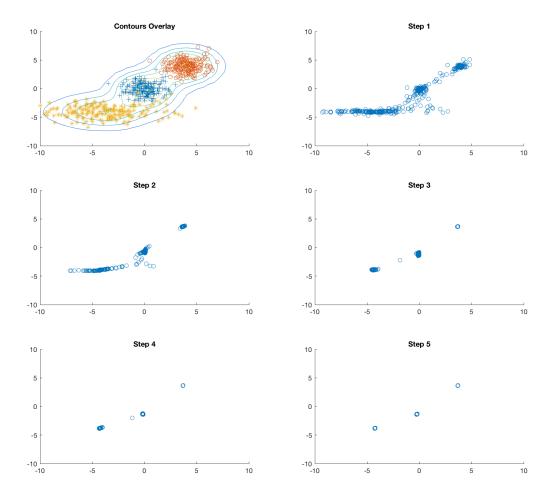


Figure 5: 3 Classes, Lambda = 3

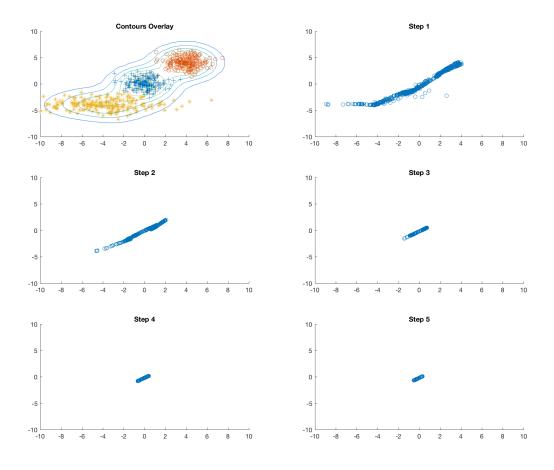


Figure 6: 3 Classes, Lambda = 5

The 3 class problem showed similar results with the same lambda values 1.8, 3 and 5. Again the 1.8 lambda value failed to shift some outliers towards the mean and the 5 lambda value shifted all of them towards a single point. The lambda value 3 correctly shifted all the points to their respective means, thus the lambda value of 3 was the best choice out of the three numbers.

### Question 5.1

```
1 %
2 % Principle Component Analysis
3 %
4
5 function result = pca(data)
6 m = mean(data);
```

```
S = cov(data - m);
       [\operatorname{evec}, \operatorname{eval}] = \operatorname{eigs}(S);
       \% Sort the eigenvalues
10
       [y, i] = sort(diag(eval), 'descend');
       \% Sort the eigenvectors columns by the eigenvalue indexes
12
       evec = evec(:, i);
13
14
       % PCA only works if you subtract the mean
15
       result = evec' * (data - m)';
16
  end
^{17}
```

### Question 5.2

 $\mathbf{a}$ 

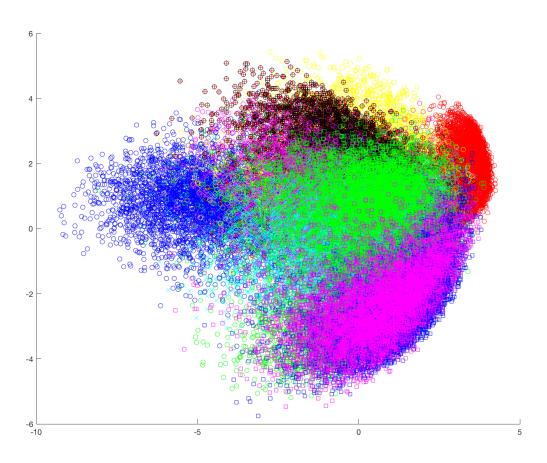


Figure 7: PCA on MNIST dataset

#### $\mathbf{b}$

The first principle component accounts for 5.116% of the data, whereas the second principle component accounts for 3.7414% of the data.

 $\mathbf{c}$ 

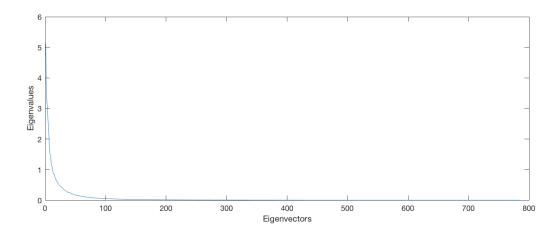


Figure 8: Scree Graph

## Question 5.6

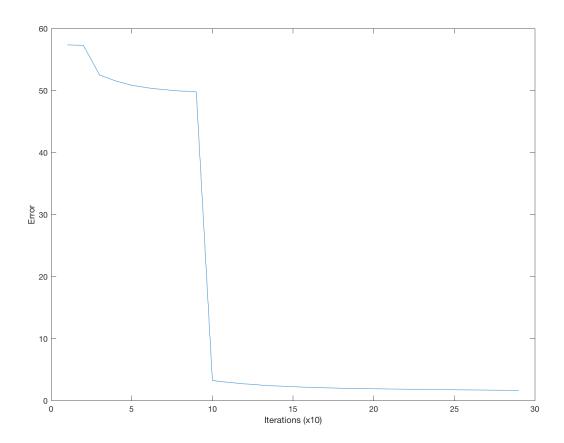


Figure 9: Error vs Iteration

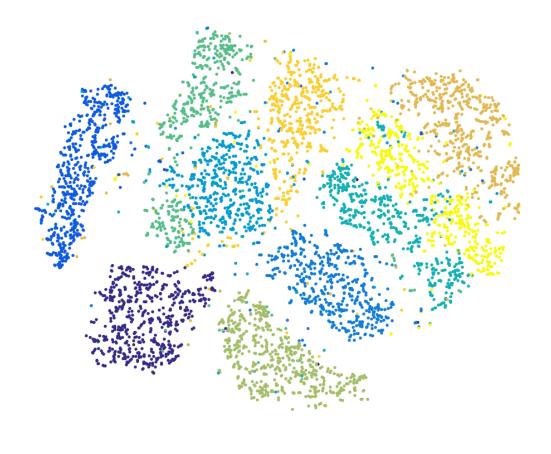


Figure 10: Iteration 300

## Question 5.8