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

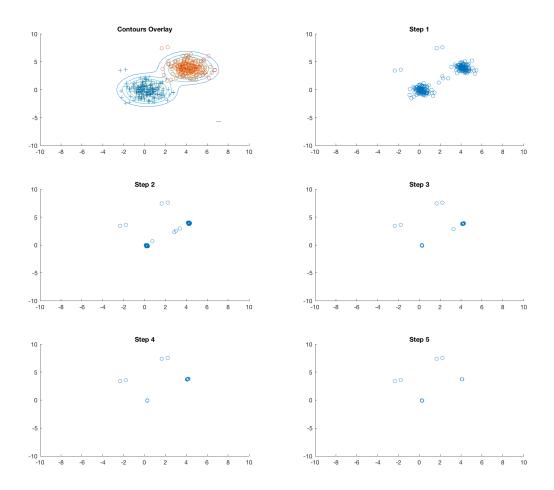


Figure 1: 2 Classes, Lambda = 1.8

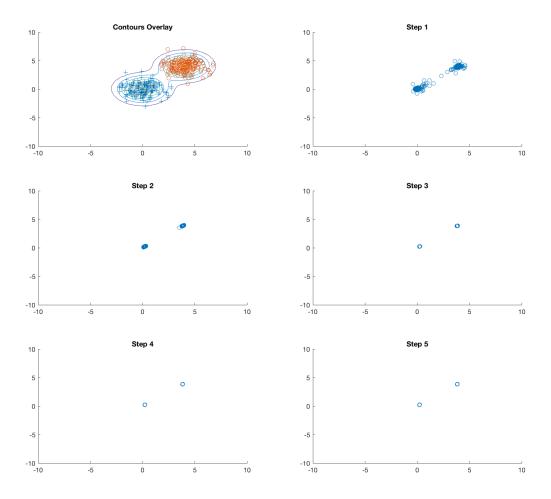


Figure 2: 2 Classes, Lambda = 3

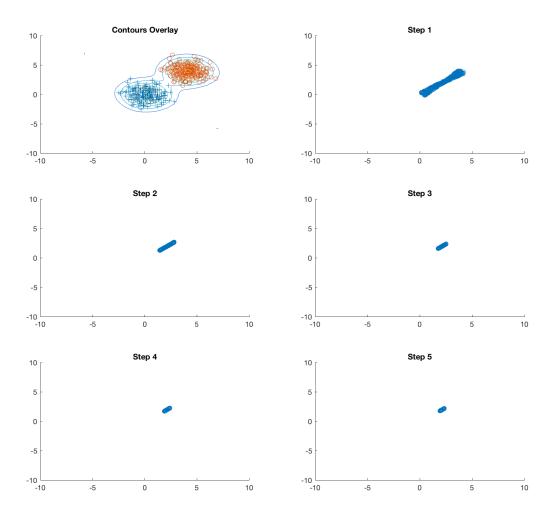


Figure 3: 2 Classes, Lambda = 5

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.

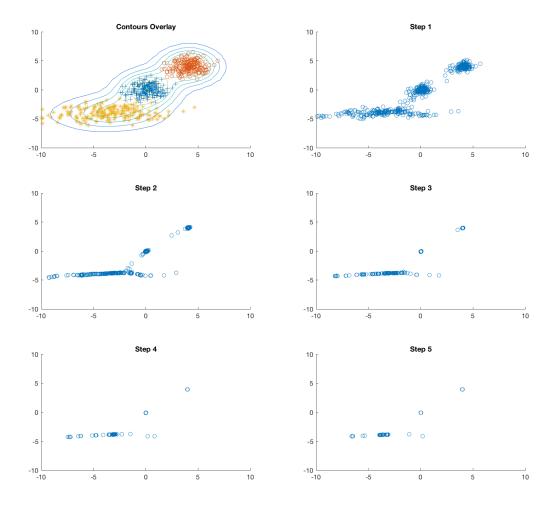


Figure 4: 3 Classes, Lambda = 1.8

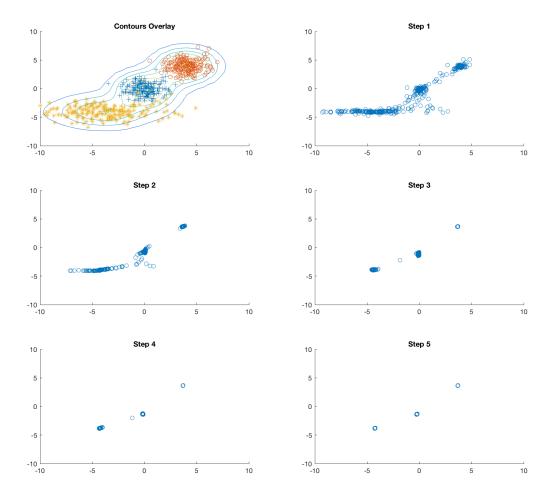


Figure 5: 3 Classes, Lambda = 3

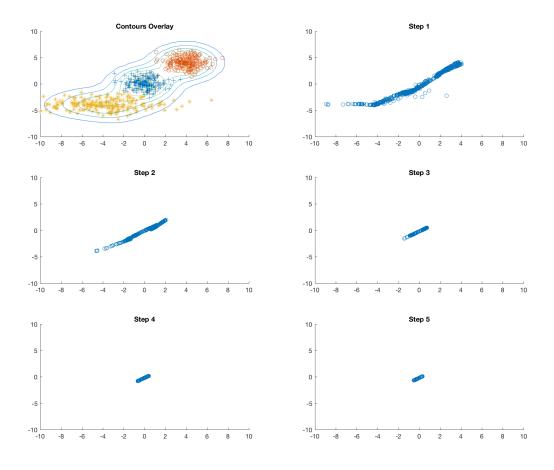


Figure 6: 3 Classes, Lambda = 5

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}

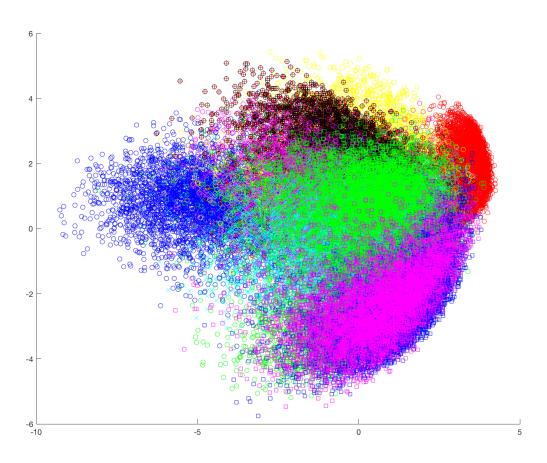


Figure 7: PCA on MNIST dataset

\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}

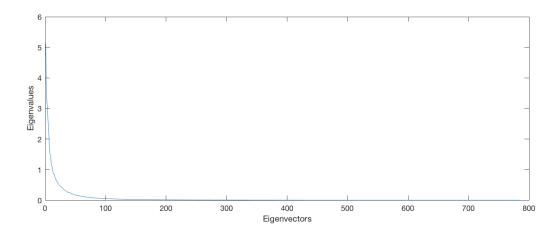


Figure 8: Scree Graph

Question 5.6

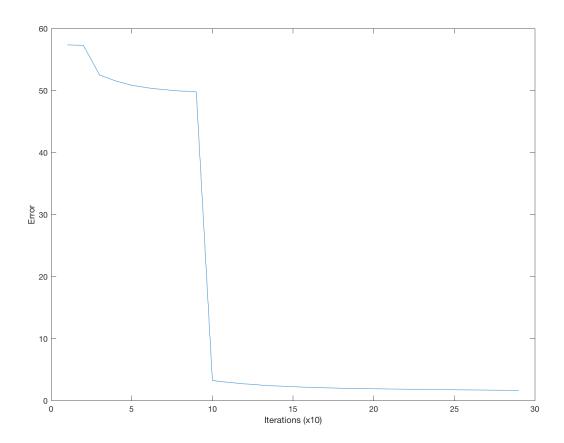


Figure 9: Error vs Iteration

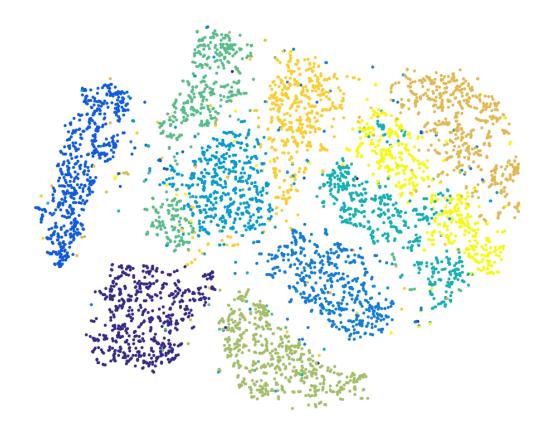


Figure 10: Iteration 300

Question 5.8

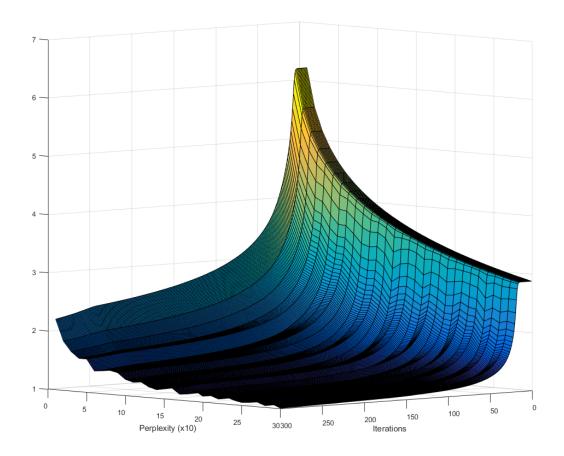


Figure 11: Perplexity and Iterations on Cost

The perplexity determines how to balance the attention between local and global aspects of the dataset. The graph shows the error change on iterations, thus we should choose a perplexity the reduces the error as quickly as possible, such that the algorithm requires less iterations to find a good solution. A simple heuristic to choose a good perplexity would be to look for the greatest rate of change in the error, by inspecting the chart we can see that the lower perplexities have far greater rate of change. Further inspection showed that a perplexity value of 2 provided the best rate of change.

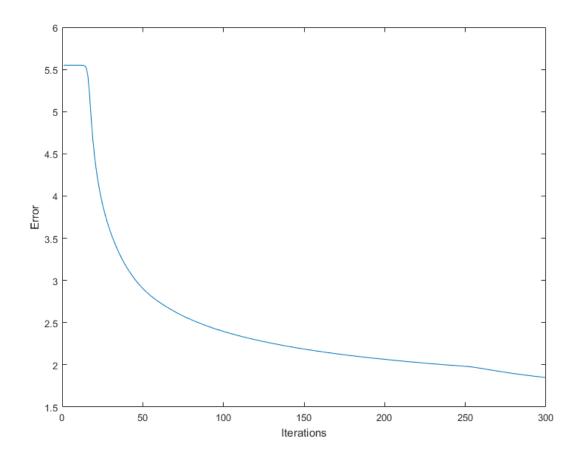


Figure 12: Error on Perplexity