# Programming Homework 1 Matthis Brocheton

## **Problem 1**

### A)

I used the normal equation method to find the parameters of the polynomial  $W_1, W_2, W_3$ , for a polynomial of order m = 2, resulting in a  $3 \times 3$  matrix.

normal equation:

$$W = (X^T X)^{-1} X^T Y$$

To find the inverse  $(A^{-1})$  matrix I used the adjugate (or cofactor) method.

Step 1:

$$A = \begin{bmatrix} a & b & c \\ d & e & f \\ g & h & i \end{bmatrix}$$

To find the determinant

$$\det(A) = a(ei - fh) - b(di - fg) + c(dh - eg)$$

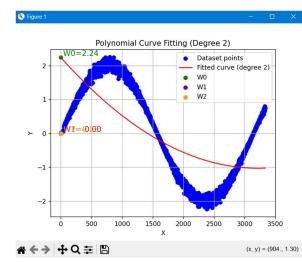
Step 2:

$$C = \begin{bmatrix} e*i - f*h & -(b*i - c*h) & b*f - c*e \\ -(d*i - f*g) & a*i - c*g & -(a*f - c*d) \\ d*h - e*g & -(a*h - b*g) & a*e - b*d \end{bmatrix}$$

Step 3:

$$adj(A) = C^{T}$$

$$A^{-1} = \frac{1}{\det(A)} \times adj(A)$$



W1 = 2.2394943784624255

W2 = -0.0020304497052723164

W3 = 3.1524947624757616e-07

### B)

I used the normal equation to find the parameters of the polynomial.

<u>normal equation:</u>

$$W = (X^T X)^{-1} X^T Y$$

To find the inverse  $(A^{-1})$  matrix I used the gauss jordan inverse.

$$\mathbf{A} = \begin{bmatrix} a_{11} & a_{12} & a_{13} \\ a_{21} & a_{22} & a_{23} \\ a_{31} & a_{32} & a_{33} \end{bmatrix}$$

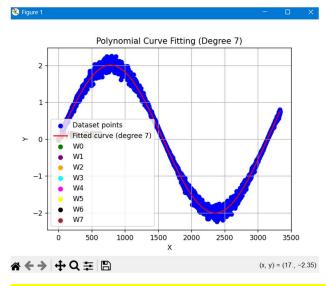
$$[\mathbf{A}|\mathbf{I}] = \begin{bmatrix} a_{11} & a_{12} & a_{13} \\ a_{21} & a_{22} & a_{23} \\ a_{31} & a_{32} & a_{33} \\ \end{bmatrix} \begin{bmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & 1 \end{bmatrix}$$

Transform the first element  $a_{11}$  into 1 by dividing the first row.

Transform the other elements in the first column into 0 by subtracting the first row multiplied by a suitable factor.

Repeat the process for the other columns until you obtain the identity matrix on the left.

Result: 
$$[I | A^{-1}] \to A^{-1}$$



The appropriate order for this polynomial curve is m = 7.

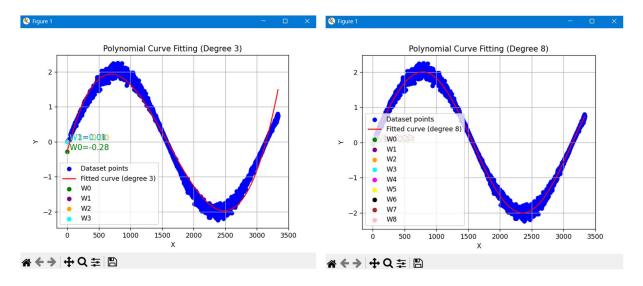
### C)

I use the formula  $\mathit{MSE} = \frac{1}{n} \sum_{i=1}^n (y_1 - \hat{y}_2^{})^2$ 

 $n \rightarrow$  The total number of data points (samples).

 $y_1 \rightarrow$  The actual (true) value of the dependent variable for the i-th data point.

 $\hat{y}_2\!\rightarrow\!$  The predicted value of the dependent variable for the i-th data point.



This is the average sum of squared errors:

MSE for m = 3: 0.03984

MSE for m = 8: 0.00443

# **Problem 2**

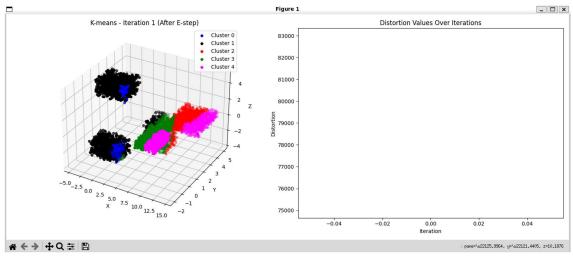
A)

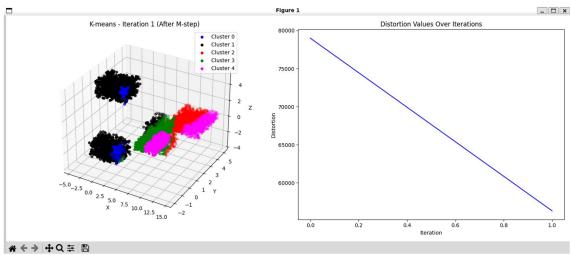
These are the centroids:

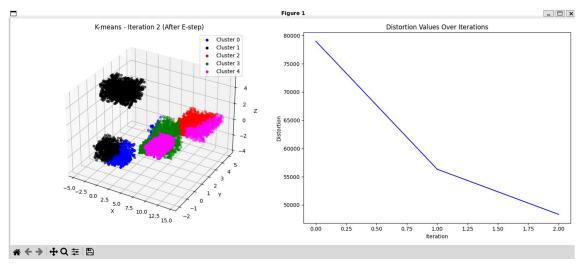
$$y_1 = m_1 = \begin{bmatrix} 4.0 \\ -0.5 \\ 2.0 \end{bmatrix}, \quad y_2 = m_2 = \begin{bmatrix} 2.0 \\ 2.5 \\ 1.0 \end{bmatrix}, \quad y_3 = m_3 = \begin{bmatrix} 10.0 \\ 2.0 \\ -1.0 \end{bmatrix}$$
$$y_4 = m_4 = \begin{bmatrix} 7.0 \\ 0.5 \\ -0.5 \end{bmatrix}, \quad y_5 = m_5 = \begin{bmatrix} 12.0 \\ 1.0 \\ -1.0 \end{bmatrix}$$

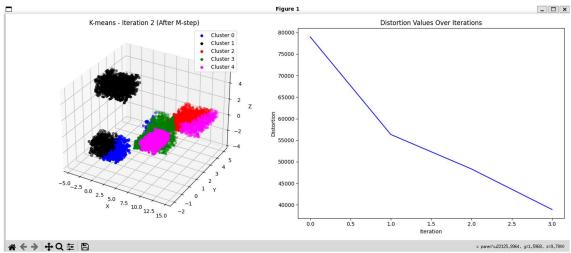
This is the result of the classification of the dataset points using the k-means = 5 algorithm, with the Euclidean distance as the distortion function.

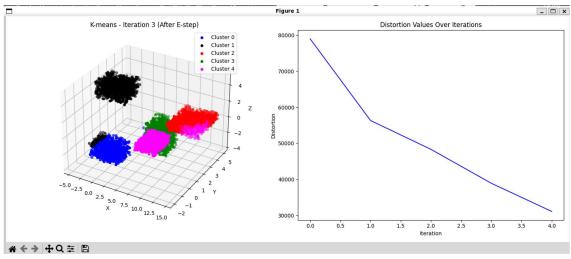
1er iteration:

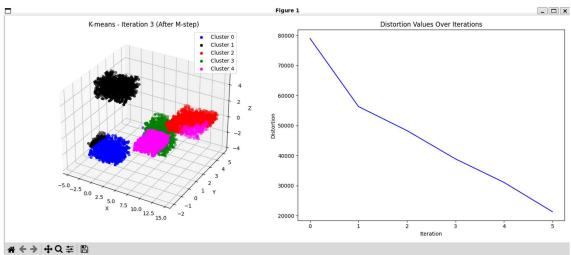




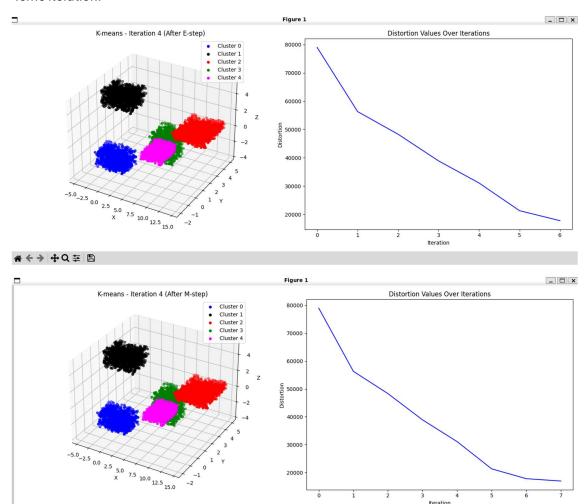


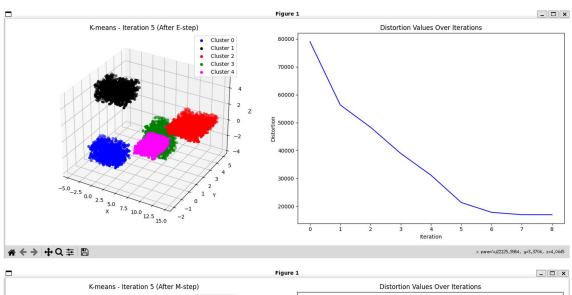


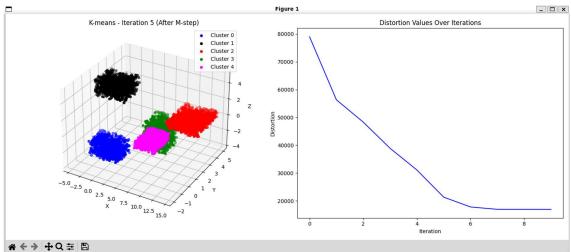




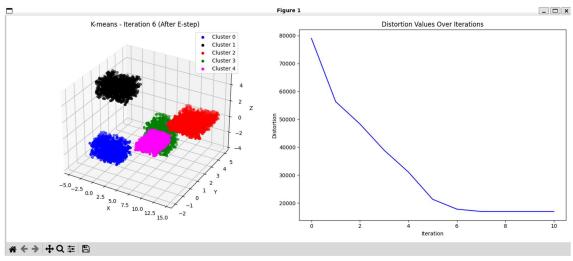
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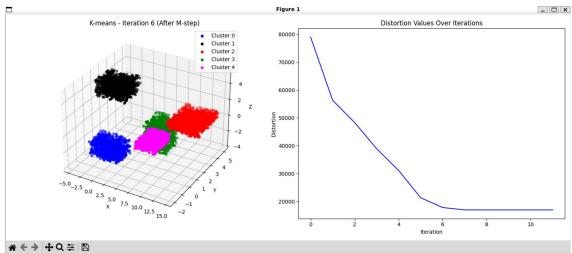


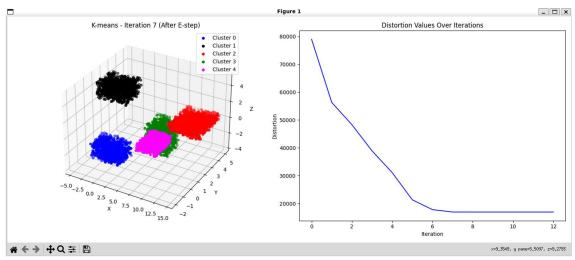


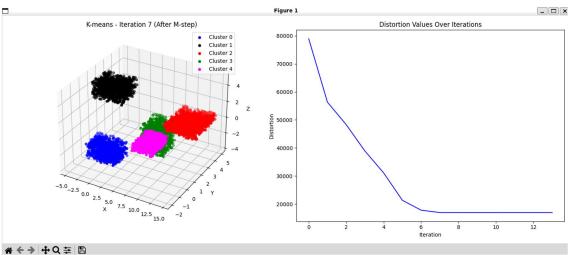


### 6er iteration:









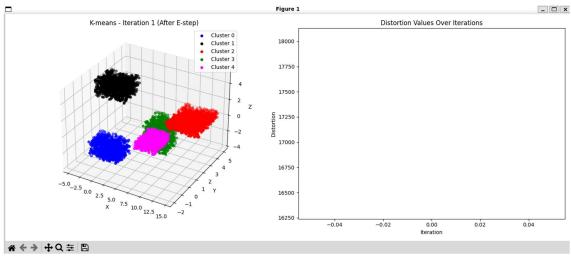
These are the centroids:

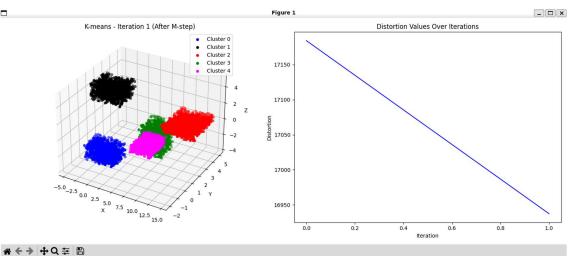
$$y_1 = m_1 = \begin{bmatrix} 0.0 \\ -0.3 \\ -2.0 \end{bmatrix}, \quad y_2 = m_2 = \begin{bmatrix} -1.3 \\ 1.5 \\ 4.0 \end{bmatrix}, \quad y_3 = m_3 = \begin{bmatrix} 11.3 \\ 3.0 \\ 0.2 \end{bmatrix}$$

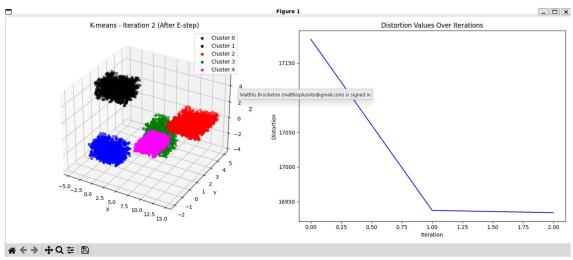
$$y_4 = m_4 = \begin{bmatrix} 5.7 \\ 3.0 \\ -2.0 \end{bmatrix}, \quad y_5 = m_5 = \begin{bmatrix} 10.0 \\ -1.0 \\ 1.2 \end{bmatrix}$$

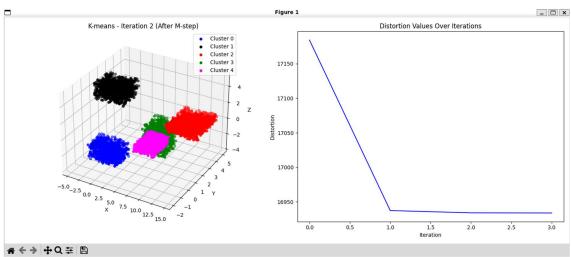
This is the behavior of the K-means (K = 5):

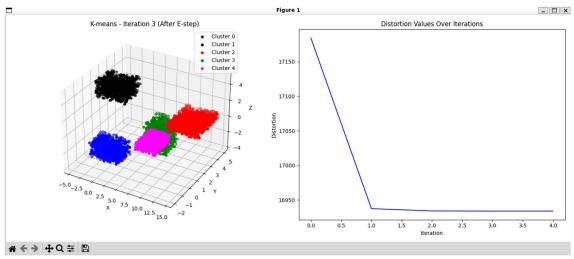
1er iteration:

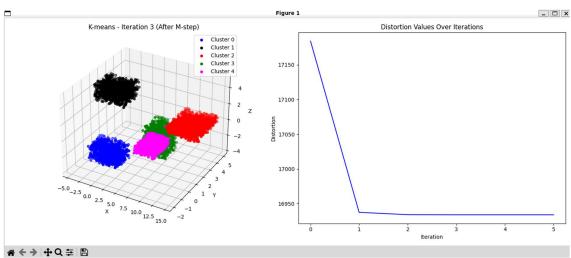








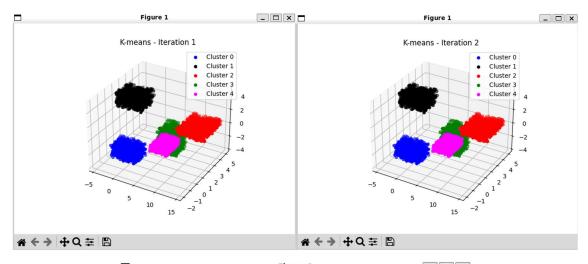


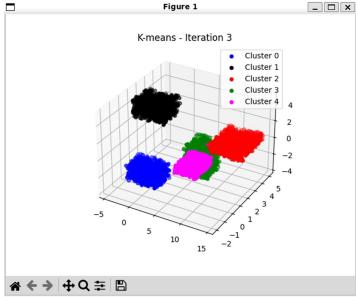


These are the centroids:

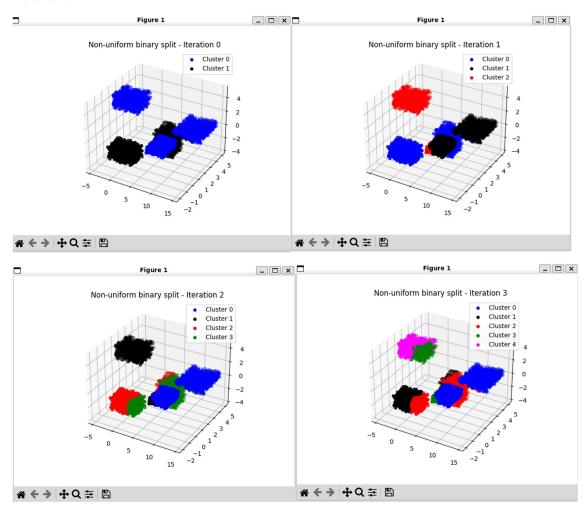
$$y_1 = m_1 = \begin{bmatrix} 0.0 \\ -0.3 \\ -2.0 \end{bmatrix}, \quad y_2 = m_2 = \begin{bmatrix} -1.3 \\ 1.5 \\ 4.0 \end{bmatrix}, \quad y_3 = m_3 = \begin{bmatrix} 11.3 \\ 3.0 \\ 0.2 \end{bmatrix}$$
$$y_4 = m_4 = \begin{bmatrix} 5.7 \\ 3.0 \\ -2.0 \end{bmatrix}, \quad y_5 = m_5 = \begin{bmatrix} 10.0 \\ -1.0 \\ 1.2 \end{bmatrix}$$

Whereas the **k-means algorithm** was able, from the very first iteration, to classify the dataset into clusters with a high degree of accuracy, convergence was reached quickly in only a few steps.





In contrast, the **non-uniform binary split algorithm** performs clustering in a **hierarchical and incremental** manner. As shown in the sequence of figures, the algorithm begins by splitting the dataset into two clusters (iteration 0), and then continues subdividing them in the subsequent iterations.



In conclusion, while **k-means** is fast and effective for well-separated and roughly spherical clusters, the **non-uniform binary split** method offers more flexibility at the cost of longer convergence times.