

DATA MINING

CLUSTERING

K-means algorithm

What is clustering?

- **Clustering** is the classification of objects into different groups, or more precisely, the partitioning of a data set into subsets (clusters), so that the data in each subset (ideally) share some common trait - often according to some defined distance measure.

Types of clustering:

1. **Hierarchical algorithms**: these find successive clusters using previously established clusters.
 1. Agglomerative ("bottom-up"): Agglomerative algorithms begin with each element as a separate cluster and merge them into successively larger clusters.
 2. Divisive ("top-down"): Divisive algorithms begin with the whole set and proceed to divide it into successively smaller clusters.
2. **Partitional clustering**: Partitional algorithms determine all clusters at once. They include:
 - **K-means and derivatives**
 - Fuzzy c-means clustering
 - QT (**Quality Threshold**) clustering algorithm

K-MEANS CLUSTERING

- The **k-means algorithm** is an algorithm to cluster n objects based on attributes into k partitions, where $k < n$.
- It assumes that the object attributes form a vector space.

K-MEANS CLUSTERING

- An algorithm for partitioning (or clustering) \mathbf{N} data points into \mathbf{K} disjoint subsets \mathbf{S}_j containing data points so as to minimize the sum-of-squares criterion

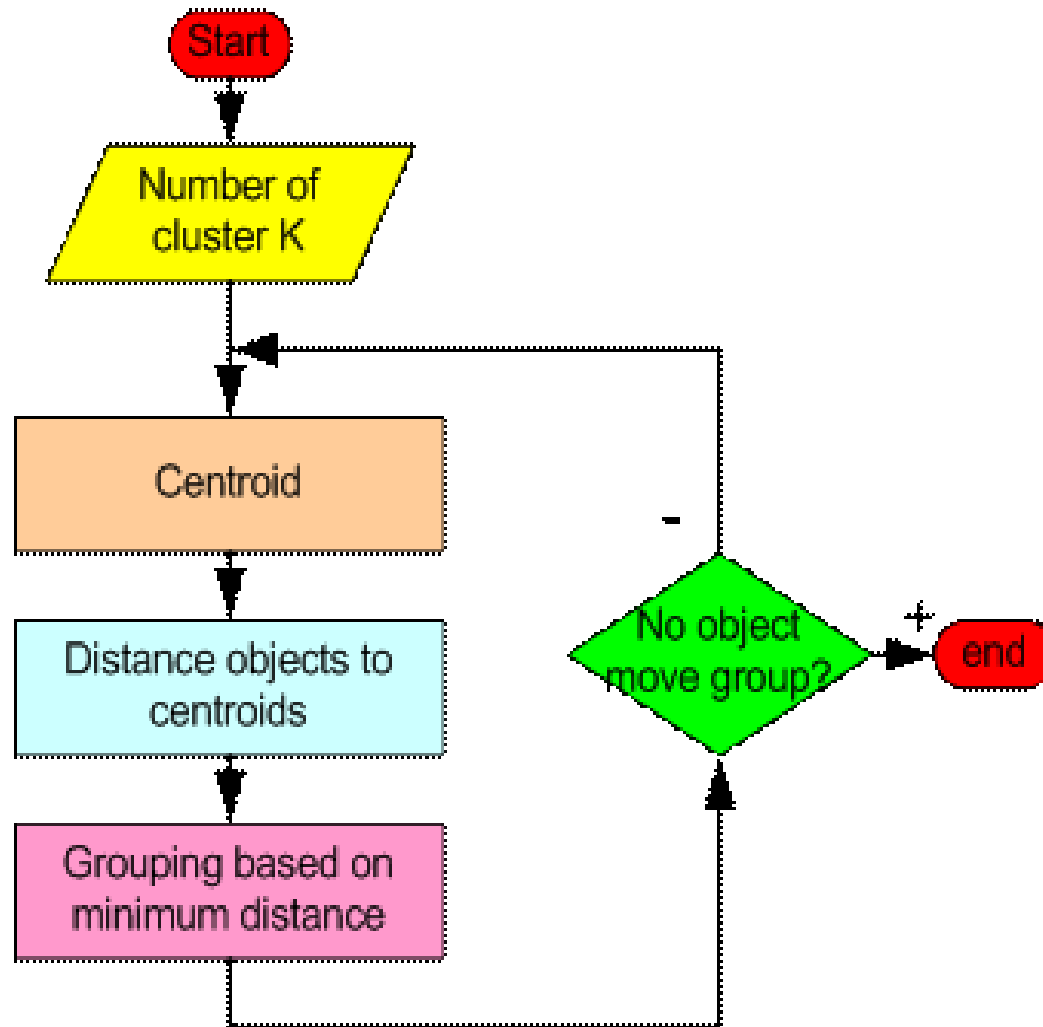
$$J = \sum_{j=1}^K \sum_{n \in \mathbf{S}_j} |x_n - \mu_j|^2,$$

where \mathbf{x}_n is a vector representing the the n^{th} data point and μ_j is the geometric centroid of the data points in \mathbf{S}_j .

K-MEANS CLUSTERING

- Simply speaking k-means clustering is an algorithm to classify or to group the objects based on attributes/features into K number of group.
- K is positive integer number.
- The grouping is done by minimizing the sum of squares of distances between data and the corresponding cluster centroid.

How the K-Mean Clustering algorithm works?



Algorithm Steps

- **Step 1:** Begin with a decision on the value of k = number of clusters .
- **Step 2:** Put any initial partition that classifies the data into k clusters. You may assign the training samples randomly, or systematically as the following:
 1. Take the first k training sample as single- element clusters
 2. Assign each of the remaining $(N-k)$ training sample to the cluster with the nearest centroid. After each assignment, recompute the centroid of the gaining cluster.

Algorithm Steps

- **Step 3:** Take each sample in sequence and compute its [distance](#) from the centroid of each of the clusters. If a sample is not currently in the cluster with the closest centroid, switch this sample to that cluster and update the centroid of the cluster gaining the new sample and the cluster losing the sample.
- **Step 4 .** Repeat step 3 until convergence is achieved, that is until a pass through the training sample causes no new assignments.

A Simple example showing the implementation of k-means algorithm (using K=2)

Individual	Variable 1	Variable 2
1	1.0	1.0
2	1.5	2.0
3	3.0	4.0
4	5.0	7.0
5	3.5	5.0
6	4.5	5.0
7	3.5	4.5

Step 1:

Initialization: Randomly we choose following two centroids (k=2) for two clusters.

In this case the 2 centroid are: $m_1=(1.0,1.0)$ and $m_2=(5.0,7.0)$.

Individual	Variable 1	Variable 2
1	1.0	1.0
2	1.5	2.0
3	3.0	4.0
4	5.0	7.0
5	3.5	5.0
6	4.5	5.0
7	3.5	4.5

	Individual	Mean Vector
Group 1	1	(1.0, 1.0)
Group 2	4	(5.0, 7.0)

Step 2:

- Thus, we obtain two clusters containing:
 $\{1,2,3\}$ and $\{4,5,6,7\}$.
- Their new centroids are:

$$m_1 = \left(\frac{1}{3}(1.0 + 1.5 + 3.0), \frac{1}{3}(1.0 + 2.0 + 4.0) \right) = (1.83, 2.33)$$

$$m_2 = \left(\frac{1}{4}(5.0 + 3.5 + 4.5 + 3.5), \frac{1}{4}(7.0 + 5.0 + 5.0 + 4.5) \right) \\ = (4.12, 5.38)$$

Individual	Centroid 1	Centroid 2
1	0	7.21
2 (1.5, 2.0)	1.12	6.10
3	3.61	3.61
4	7.21	0
5	4.72	2.5
6	5.31	2.06
7	4.30	2.92

$$d(m_1, 2) = \sqrt{|1.0 - 1.5|^2 + |1.0 - 2.0|^2} = 1.12$$

$$d(m_2, 2) = \sqrt{|5.0 - 1.5|^2 + |7.0 - 2.0|^2} = 6.10$$

Step 3:

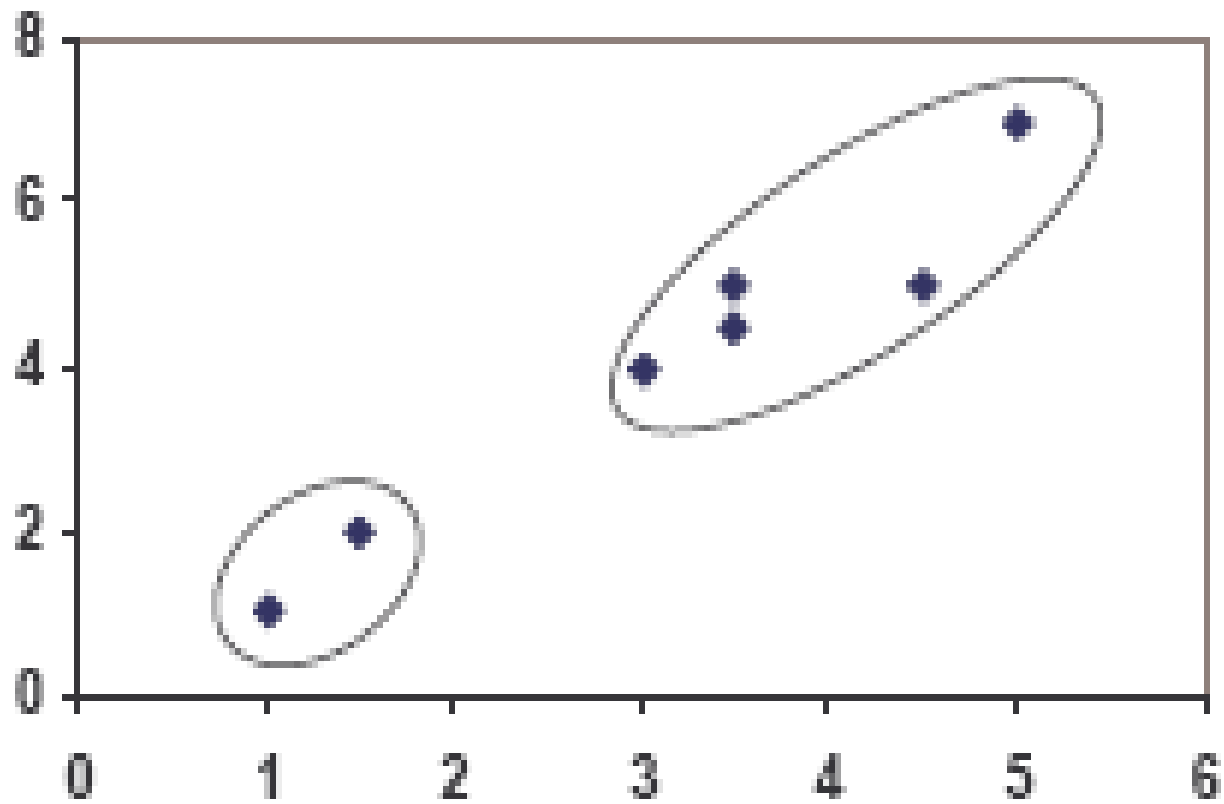
- Now using these centroids we compute the Euclidean distance of each object, as shown in table.
- Therefore, the new clusters are:
 $\{1,2\}$ and $\{3,4,5,6,7\}$
- Next centroids are:
 $m_1=(1.25,1.5)$ and $m_2 = (3.9,5.1)$

Individual	Centroid 1	Centroid 2
1	1.57	5.38
2	0.47	4.28
3	2.04	1.78
4	5.84	1.84
5	3.15	0.73
6	3.78	0.54
7	2.74	1.08

- Step 4 :
The clusters obtained are:
{1,2} and {3,4,5,6,7}
- Therefore, there is no change in the cluster.
- Thus, the algorithm comes to a halt here and final result consist of 2 clusters {1,2} and {3,4,5,6,7}.

Individual	Centroid 1	Centroid 2
1	0.58	5.02
2	0.58	3.92
3	3.05	1.42
4	6.88	2.20
5	4.18	0.41
6	4.78	0.81
7	3.75	0.72

PLOT



(with $K=3$)

Individual	$m_1 = 1$	$m_2 = 2$	$m_3 = 3$	cluster
1	0	1.11	3.61	1
2	1.12	0	2.5	2
3	3.61	2.5	0	3
4	7.21	6.10	3.61	3
5	4.72	3.61	1.12	3
6	5.31	4.24	1.80	3
7	4.30	3.20	0.71	3

C_3

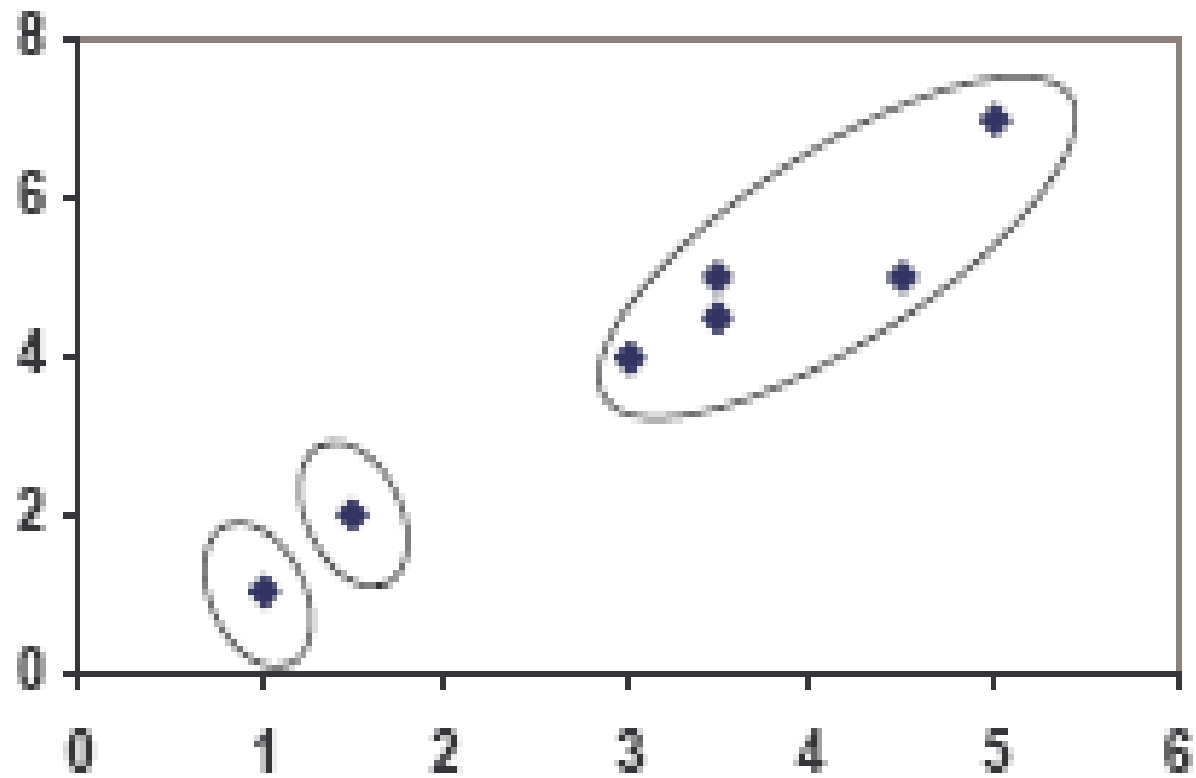
clustering with initial centroids (1, 2, 3)

Step 1

Individual	m_1 (1.0, 1.0)	m_2 (1.5, 2.0)	m_3 (3.9, 5.1)	cluster
1	0	1.11	5.02	1
2	1.12	0	3.92	2
3	3.61	2.5	1.42	3
4	7.21	6.10	2.20	3
5	4.72	3.61	0.41	3
6	5.31	4.24	0.61	3
7	4.30	3.20	0.72	3

Step 2

PLOT

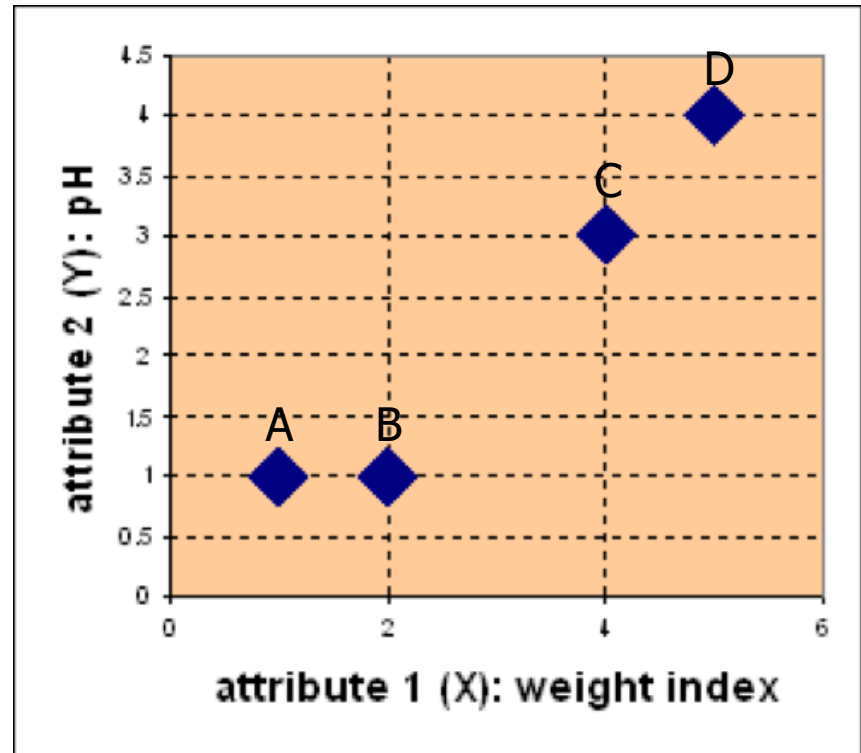


Real-Life Numerical Example

- **Problem**

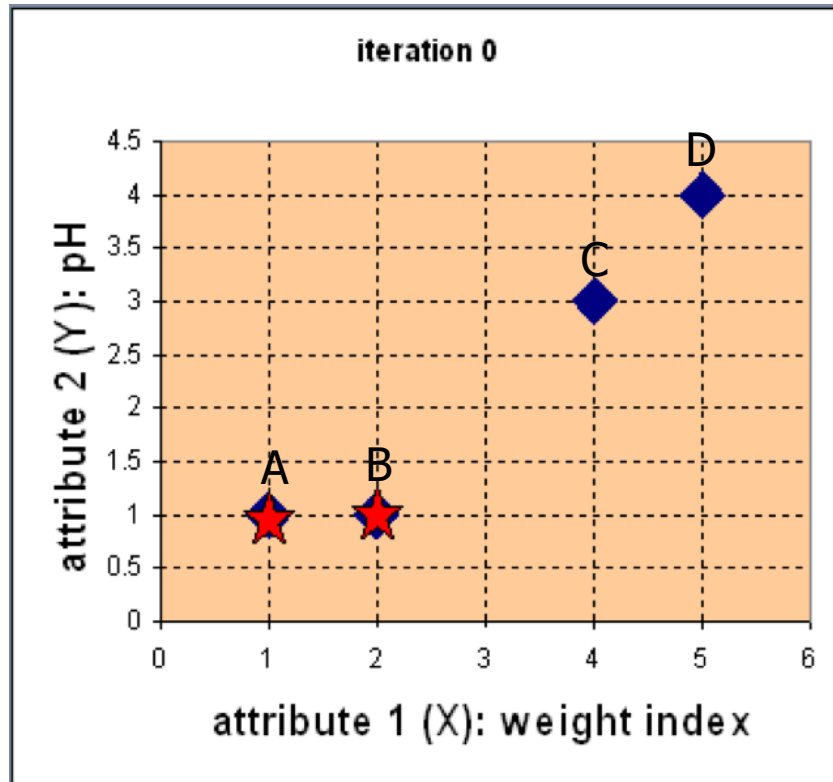
Suppose we have 4 types of medicines and each has two attributes (pH and weight index). Our goal is to group these objects into $K=2$ group of medicine.

Medicine	Weight	pH-Index
A	1	1
B	2	1
C	4	3
D	5	4



Example

- Step 1: Use initial seed points for partitioning



$$c_1 = A, c_2 = B$$

$D^0 =$	$\begin{bmatrix} 0 & 1 & 3.61 & 5 \\ 1 & 0 & 2.83 & 4.24 \end{bmatrix}$	$c_1 = (1,1)$ group - 1
		$c_2 = (2,1)$ group - 2
	$\begin{matrix} A & B & C & D \\ \begin{bmatrix} 1 & 2 & 4 & 5 \\ 1 & 1 & 3 & 4 \end{bmatrix} \end{matrix}$	Euclidean distance
		X
		Y

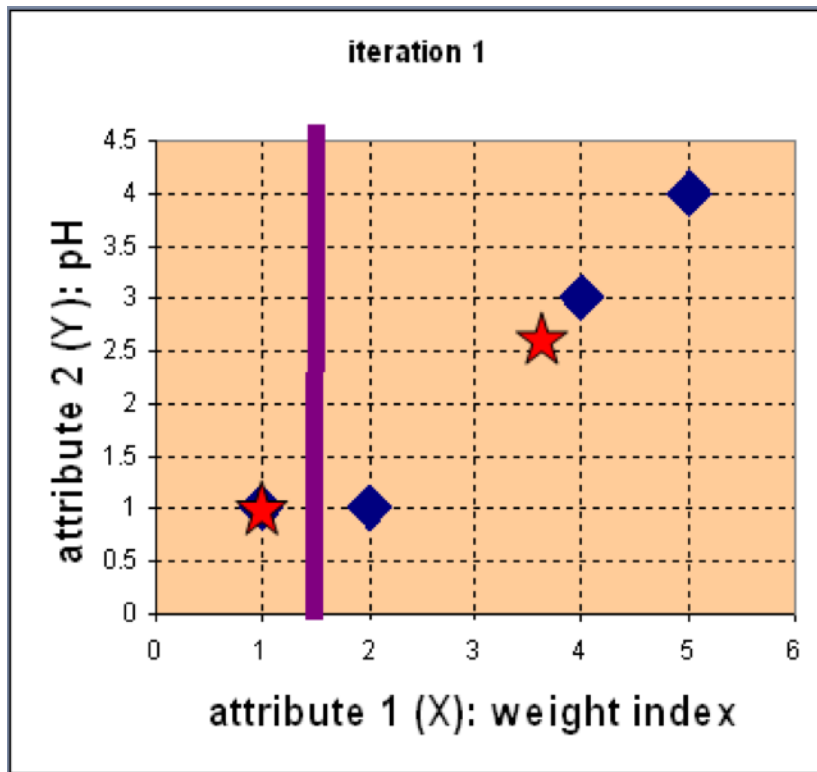
$$d(D, c_1) = \sqrt{(5-1)^2 + (4-1)^2} = 5$$

$$d(D, c_2) = \sqrt{(5-2)^2 + (4-1)^2} = 4.24$$

Assign each object to the cluster with the nearest seed point

Example

- Step 2: Compute new centroids of the current partition



Knowing the members of each cluster, now we compute the new centroid of each group based on these new memberships.

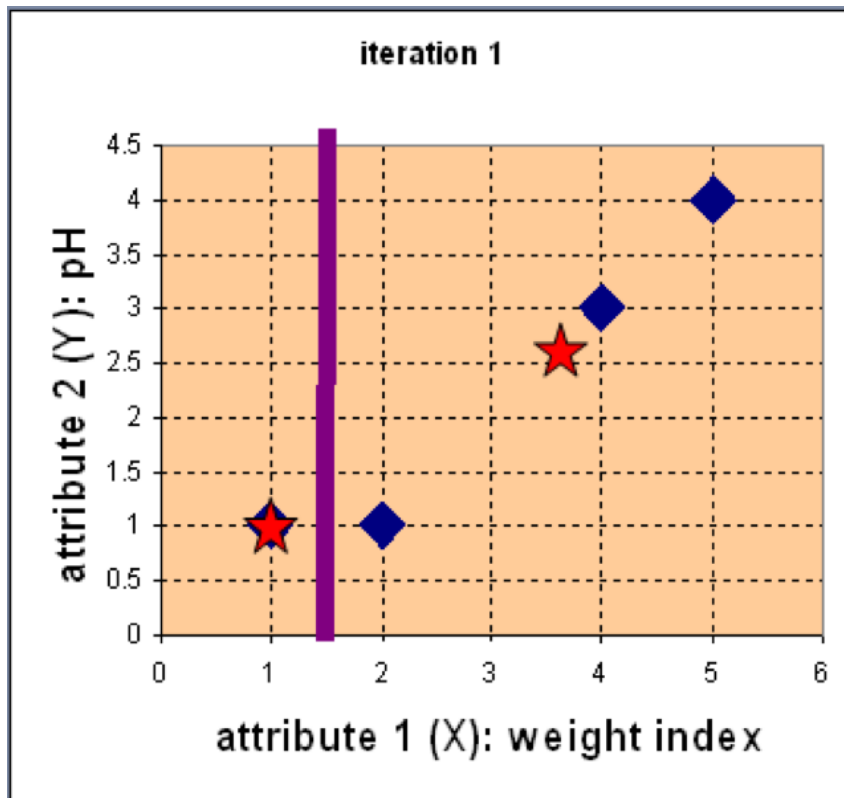
$$c_1 = (1, 1)$$

$$c_2 = \left(\frac{2 + 4 + 5}{3}, \frac{1 + 3 + 4}{3} \right)$$

$$= \left(\frac{11}{3}, \frac{8}{3} \right)$$

Example

- Step 2: Renew membership based on new centroids



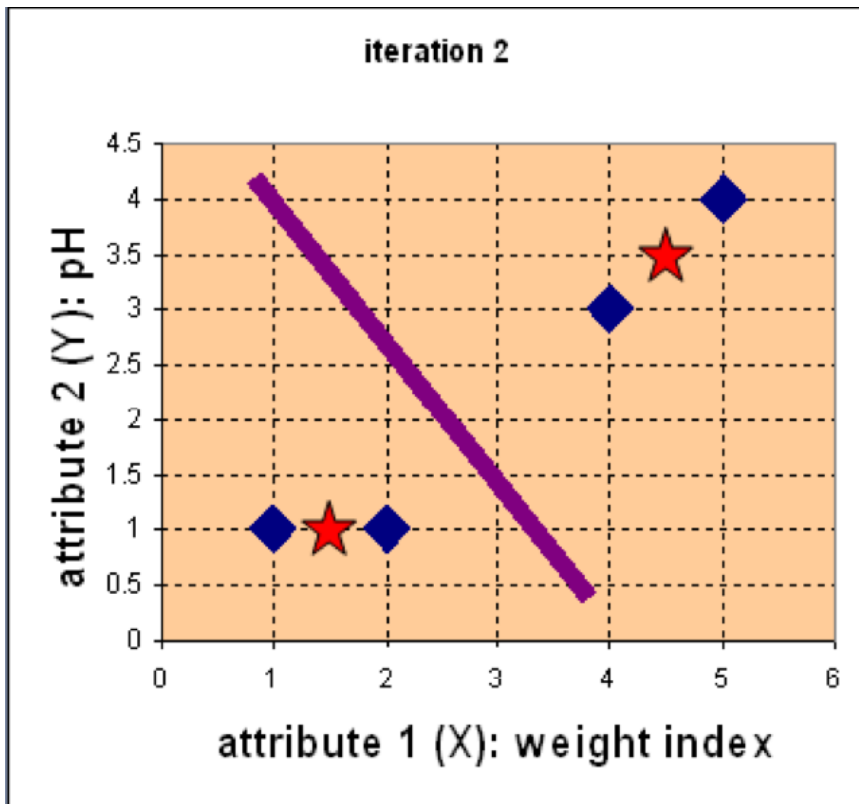
Compute the distance of all objects to the new centroids

$$D^1 = \begin{bmatrix} 0 & 1 & 3.61 & 5 \\ 3.14 & 2.36 & 0.47 & 1.89 \\ A & B & C & D \\ \begin{bmatrix} 1 & 2 & 4 & 5 \\ 1 & 1 & 3 & 4 \end{bmatrix} & X & Y \end{bmatrix} \quad \begin{array}{l} \mathbf{c}_1 = (1, 1) \text{ group-1} \\ \mathbf{c}_2 = (\frac{11}{3}, \frac{8}{3}) \text{ group-2} \end{array}$$

Assign the membership to objects

Example

- Step 3: Repeat the first two steps until its convergence



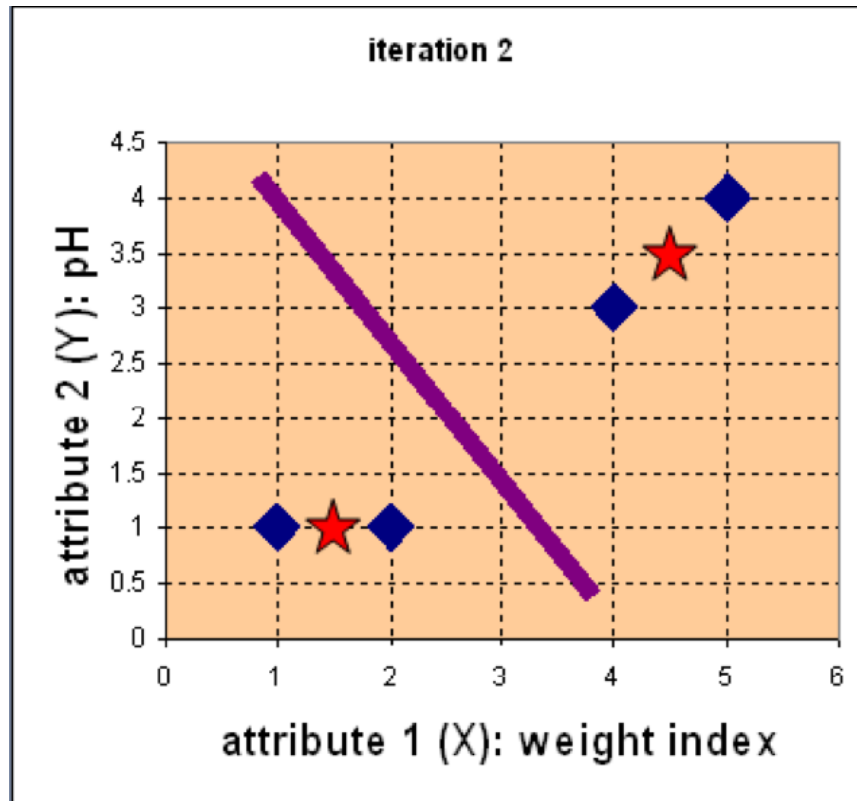
Knowing the members of each cluster, now we compute the new centroid of each group based on these new memberships.

$$c_1 = \left(\frac{1+2}{2}, \frac{1+1}{2} \right) = \left(1\frac{1}{2}, 1 \right)$$

$$c_2 = \left(\frac{4+5}{2}, \frac{3+4}{2} \right) = \left(4\frac{1}{2}, 3\frac{1}{2} \right)$$

Example

- Step 3: Repeat the first two steps until its convergence



Compute the distance of all objects to the new centroids

$$D^2 = \begin{bmatrix} 0.5 & 0.5 & 3.20 & 4.61 \\ 4.30 & 3.54 & 0.71 & 0.71 \end{bmatrix} \quad \begin{array}{l} \mathbf{c}_1 = (1\frac{1}{2}, 1) \text{ group-1} \\ \mathbf{c}_2 = (4\frac{1}{2}, 3\frac{1}{2}) \text{ group-2} \end{array}$$

	A	B	C	D	
	1	2	4	5	X
	1	1	3	4	Y

Stop due to no new assignment
Membership in each cluster no longer change

We get the final grouping as the results as:

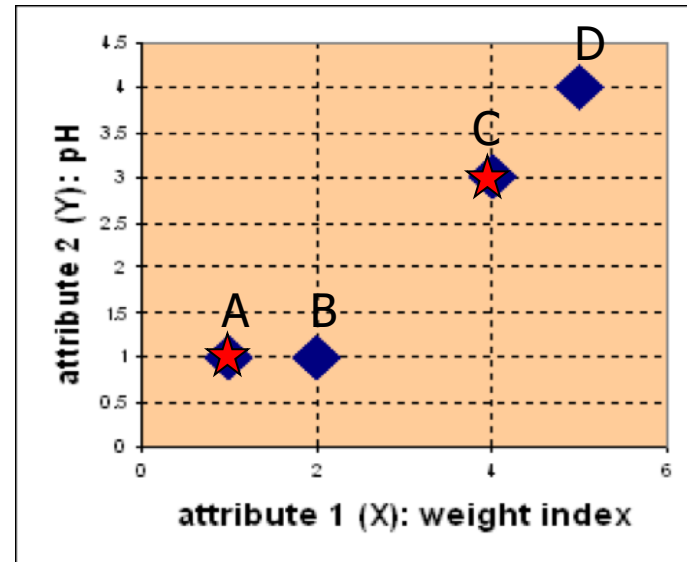
<u>Object</u>	<u>Feature1(X): weight index</u>	<u>Feature2 (Y): pH</u>	<u>Group (result)</u>
Medicine A	1	1	1
Medicine B	2	1	1
Medicine C	4	3	2
Medicine D	5	4	2

Exercise

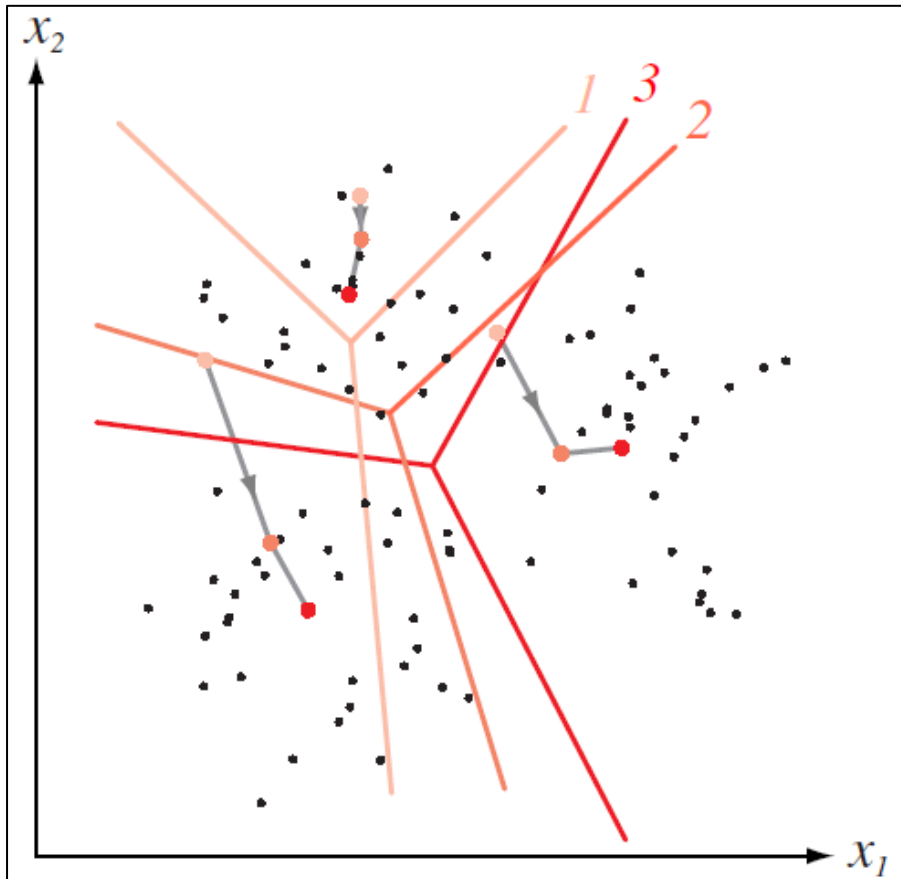
For the medicine data set, use K-means with the **Manhattan** distance metric for clustering analysis by setting $K=2$ and initialising seeds as $C_1 = A$ and $C_2 = C$. Answer three questions as follows:

1. How many steps are required for convergence?
2. What are memberships of two clusters after convergence?
3. What are centroids of two clusters after convergence?

Medicine	Weight	pH-Index
A	1	1
B	2	1
C	4	3
D	5	4



How K-means partitions?



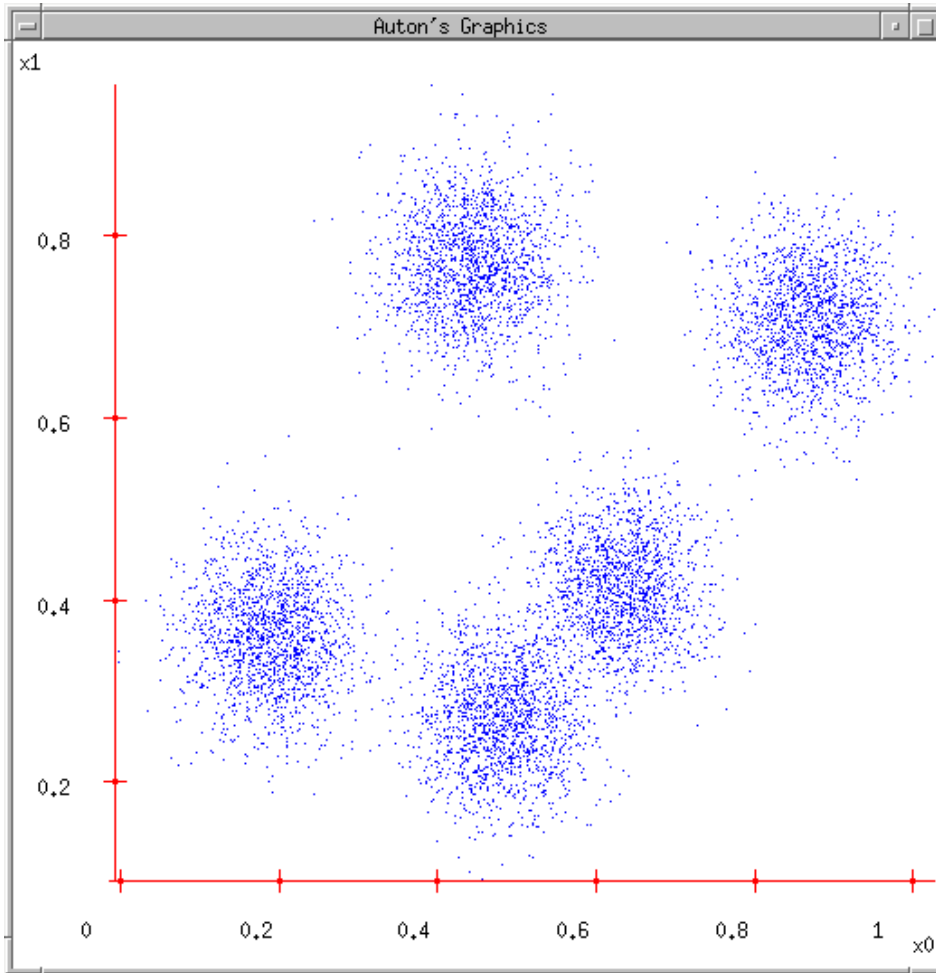
When K centroids are set/fixed, they partition the whole data space into K mutually exclusive subspaces to form a partition.

A partition amounts to a

[Voronoi Diagram](#)

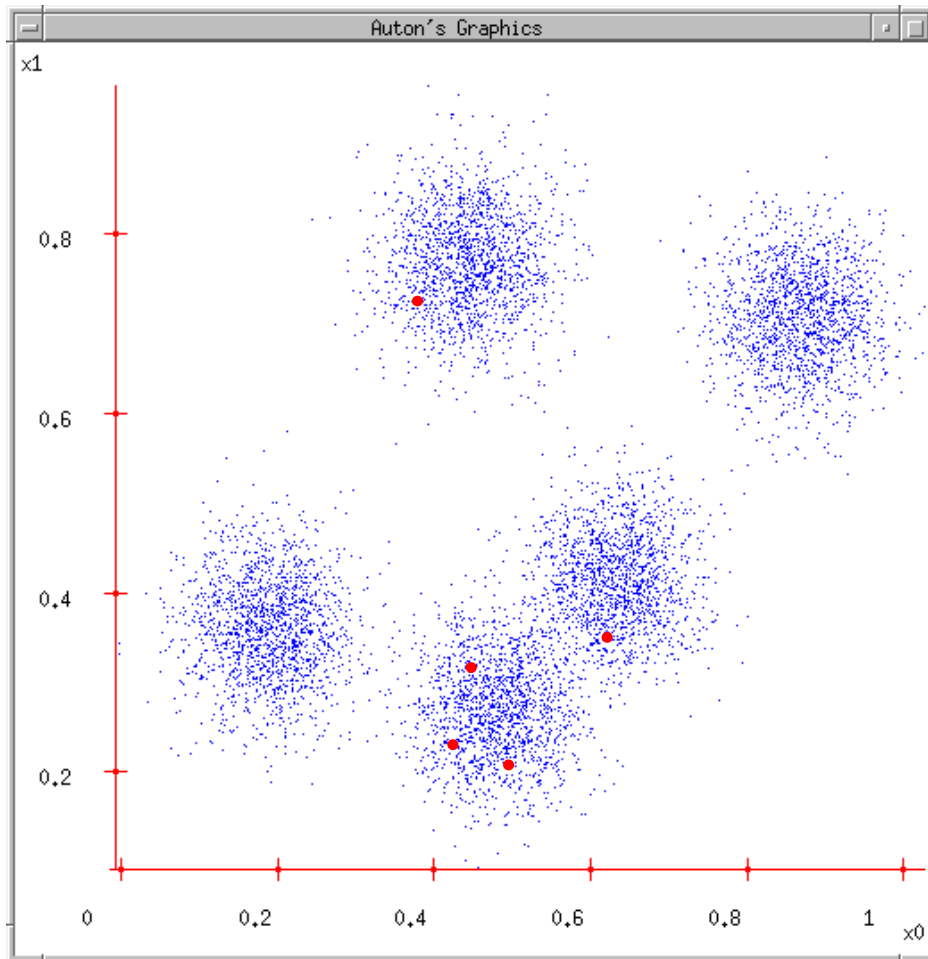
Changing positions of centroids leads to a new partitioning.

K-means Demo



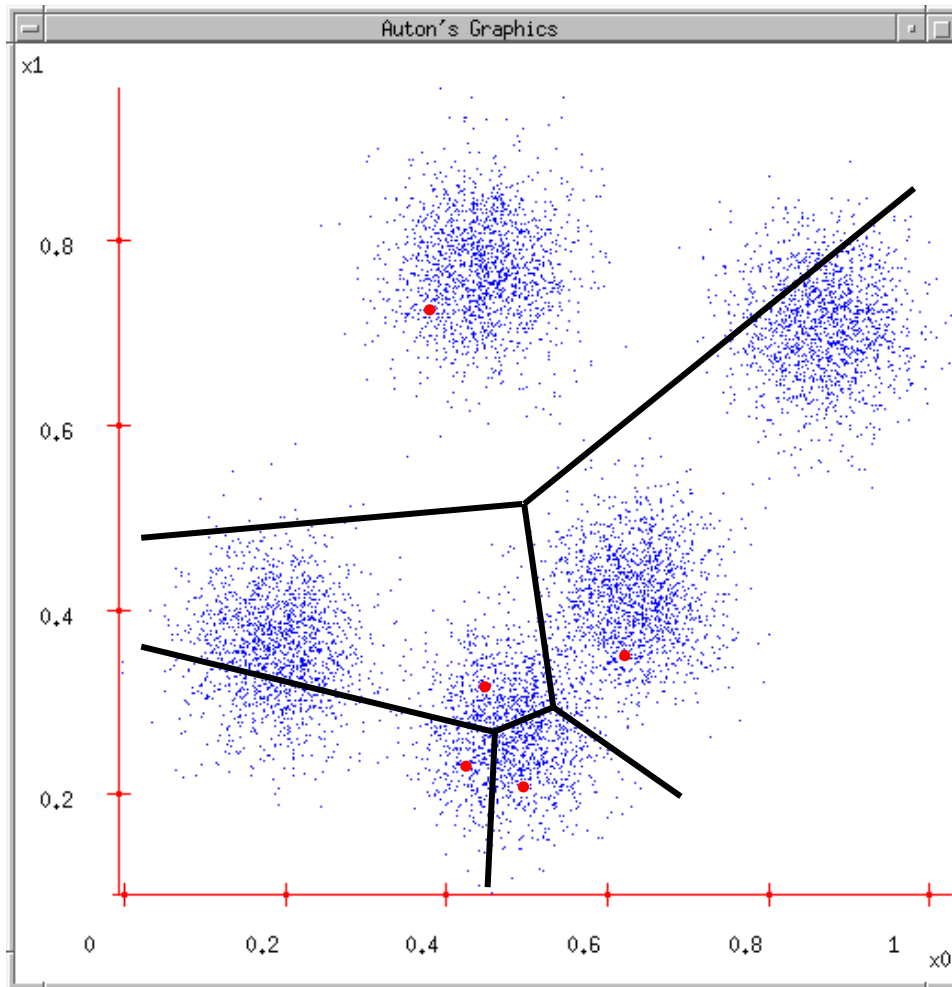
1. User set up the number of clusters they'd like. (*e.g.* $k=5$)

K-means Demo



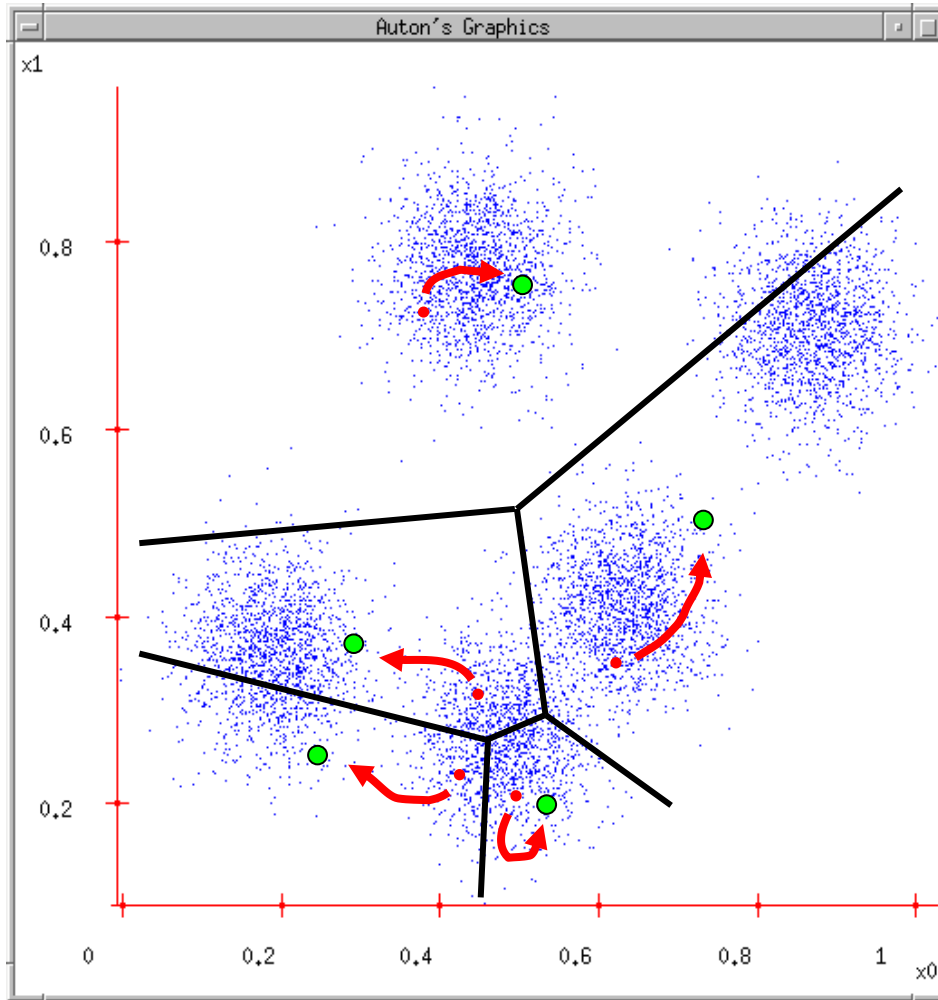
1. User set up the number of clusters they'd like. (*e.g.* $K=5$)
2. Randomly guess K cluster Center locations

K-means Demo



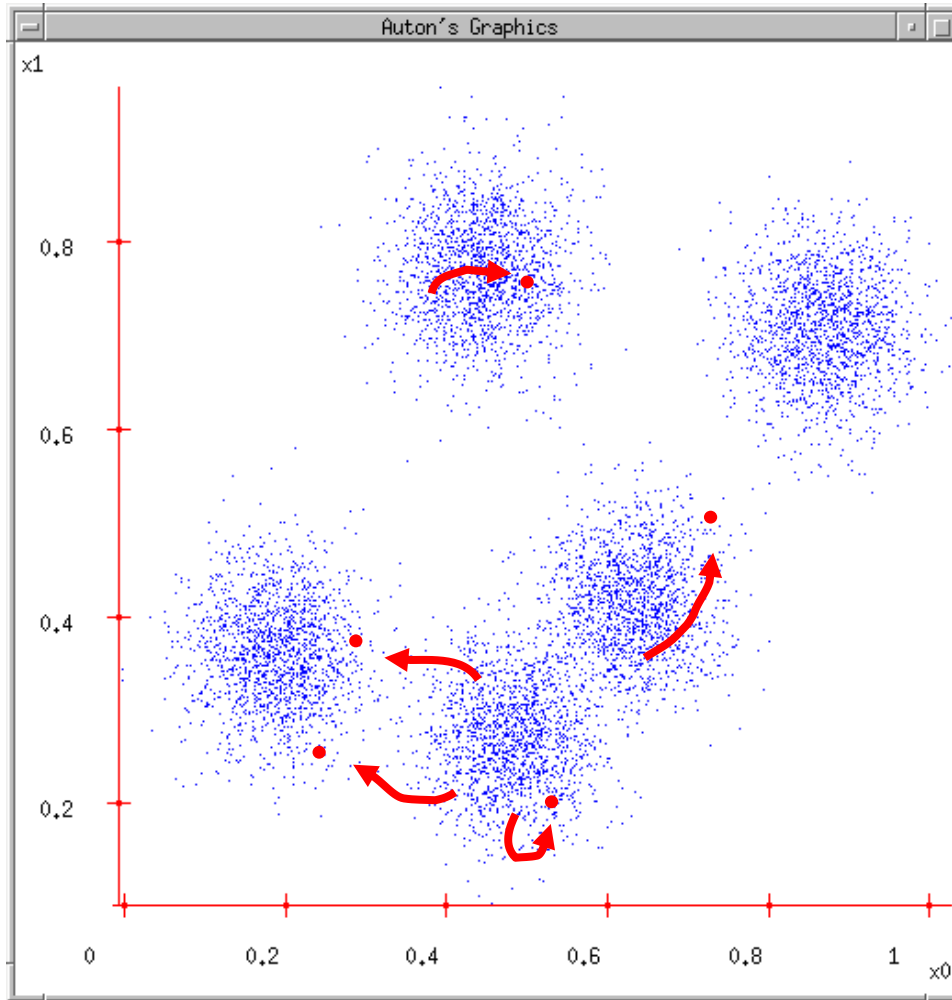
1. User set up the number of clusters they'd like. (*e.g.* $K=5$)
2. Randomly guess K cluster Center locations
3. Each data point finds out which Center it's closest to. (Thus each Center "owns" a set of data points)

K-means Demo



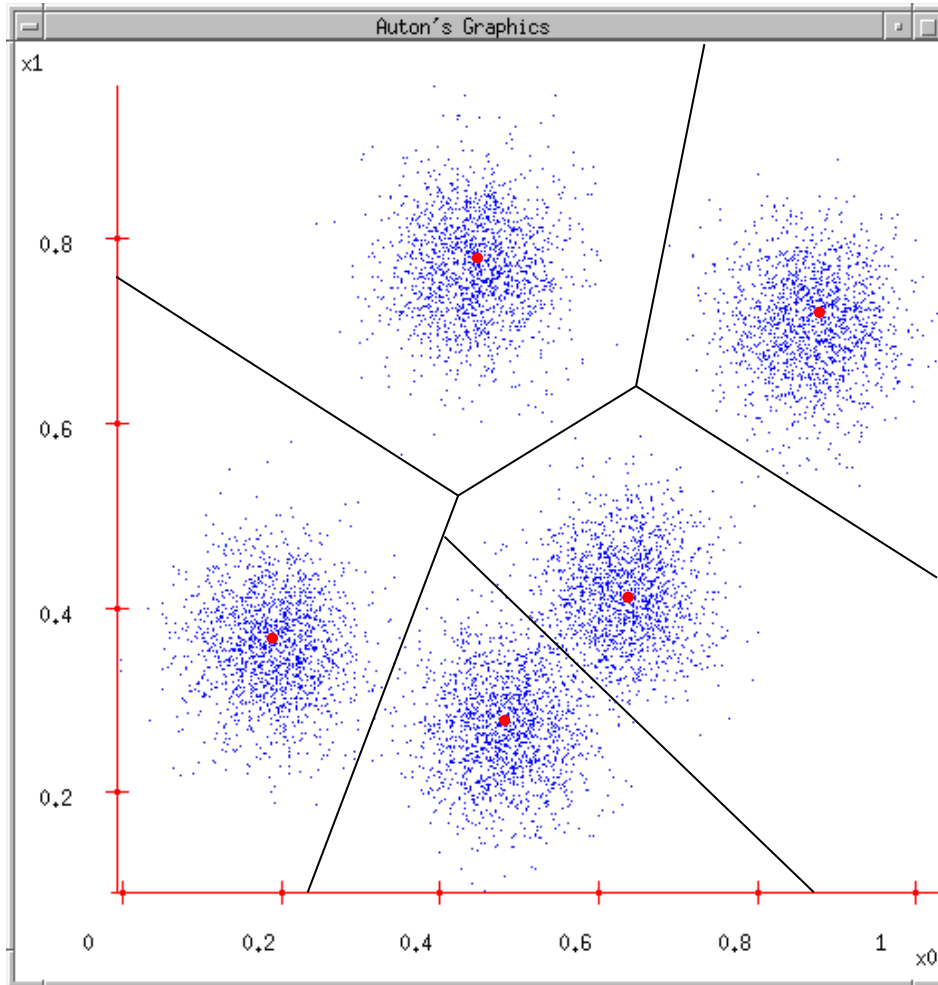
1. User set up the number of clusters they'd like. (e.g. $K=5$)
2. Randomly guess K cluster centre locations
3. Each data point finds out which centre it's closest to. (Thus each Center "owns" a set of data points)
4. Each centre finds the centroid of the points it owns

K-means Demo



1. User set up the number of clusters they'd like. (e.g. $K=5$)
2. Randomly guess K cluster centre locations
3. Each data point finds out which centre it's closest to. (Thus each centre "owns" a set of data points)
4. Each centre finds the centroid of the points it owns
5. ...and jumps there

K-means Demo



1. User set up the number of clusters they'd like. (e.g. $K=5$)
2. Randomly guess K cluster centre locations
3. Each data point finds out which centre it's closest to. (Thus each centre "owns" a set of data points)
4. Each centre finds the centroid of the points it owns
5. ...and jumps there
6. ...Repeat until terminated!