

Unsupervised Learning: Clustering

Some material adapted from slides by Andrew Moore, CMU

14.6



Yann LeCun on Unsupervised Learning

"Most of human and animal learning is *unsupervised learning*. If intelligence was a cake, unsupervised learning would be the cake, *supervised learning* would be the icing on the cake, and *reinforcement learning* would be the cherry on the cake. ... We know how to make the icing and the cherry, but we don't know how to make the cake. We need to solve the

unsupervised learning problem before we can even think of getting to true AI."*



* <u>Yann LeCun</u> (*Head of Facebook AI, NYU CS Prof.*) on AlphaGo's success and AI, 2016

Unsupervised Learning

- Supervised learning used labeled data pairs (x, y) to learn a function f : X→y
- But, what if we don't have labels?
- No labels = unsupervised learning
- Only some points are labeled = semi-supervised
 learning

-Getting labels is expensive, so we only get a few

- Clustering is the unsupervised grouping of data points based on similarity
- It can be used for knowledge discovery

Clustering algorithms

- Many clustering algorithms
- Clustering typically done using a distance measure defined between instances or points
- Distance defined by instance feature space, so it works with numeric features
 - Requires encoding of categorial values; may benefit from normalization
- We'll look at three popular approaches
 - 1. Centroid-based clustering
 - 2. Hierarchical clustering
 - 3. DBSCAN

Clustering Data

Given a collection of points (x,y), group them into one or more clusters based on their distance from one another

How many clusters are there?

How can we find them



(1) K-Means Clustering

- Randomly choose k cluster center locations, aka centroids
- Loop until convergence
 - assign a point to cluster of closest centroid
 - re-estimate cluster centroids
 based on its data assigned
- Convergence: no point is re-assigned to a different cluster

k = 5





- 1. k centerpoints are randomly initialized.
- 2. Observations are assigned to the closest centerpoint.
- 3. Centerpoints are moved to the center of their members.
- 4. Repeat steps 2 and 3 until no observation changes membership in step 2.

Chris Albon

distance, centroids

- Distance between points (X_0, Y_0, Z_0) and (X_1, Y_1, Z_1) is just sqrt $((X_0 - X_1)^2 + (Y_0 - Y_1)^2 + (Z_0 - Z_1)^2)$
- In numpy

>>> import numpy as np
>>> p1 = np.array([0,-2,0,1]) ; p2 = np.array([0,1,2,1]))
>>> np.linalg.norm(p1 - p2)
3.605551275463989

Computing centroid of set of points easy
 >> points = np.array([[1,2,3], [2,1,1], [3,1,0]]) # 3D points
 >> centroid = np.mean(points, axis=0) # get mean across columns
 >> centroid
 array([2.0, 1.33, 1.33])

(1) K-Means Clustering

- Randomly choose k cluster center locations, aka centroids
- Loop until convergence
 - assign a point to cluster of the closest centroid
 - re-estimate cluster centroids
 based on its data assigned
- Convergence: no point is assigned to a different cluster

k = 5



K-Means Clustering

K-Means (k, data)

- Randomly choose k cluster center locations (centroids)
- Loop until convergence
 - Assign each point to the cluster of the closest centroid.
 - Re-estimate the cluster centroids based on the data assigned to each
- Convergence: no point is assigned to a different cluster



veroni diagram: add lines for regions of points closest to each centroid

K-Means Clustering

K-Means (k, data)

- Randomly choose k cluster center locations (centroids)
- Loop until convergence
 - Assign each point to the cluster of the closest centroid
 - Re-estimate the cluster centroids based on the data assigned to each
- Convergence: no point is assigned to a different cluster



K-Means Clustering

K-Means (k, data)

- Randomly choose k cluster center locations (centroids)
- Loop until convergence
 - Assign each point to the cluster of the closest centroid
 - Re-estimate the cluster centroids based on the data assigned to each
- Convergence: no point is assigned to a different cluster



Visualizing k-means: http://bit.ly/471kmean

https://www.naftaliharris.com/blog/visu	alizi 🛣 松 G 🛈 🕟	I 🦎 [□] {=} 🕈 🚍	Ø R 🛱 O 🗹	💼 🤹 🏪 🛛 🔊 🚺
	•			
		• •		
	8.0000	• 600	•	
	~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~	• •	•	
	https://www.naftaliharris.com/blog/visu	https://www.naftaliharris.com/blog/visualizi	https://www.naftaliharris.com/blog/visualizi	https://www.naftaliharris.com/blog/visualizi * ~ G û î î î î î î î î î î î î î î î î î î

### **Clustering the Iris Data**

 Let's try using unsupervised clustering on the Iris Data

	Weka	Explorer				
Preprocess Classify Cluster Associate Select att	ributes Visualize					
Clusterer						
Choose SimpleKMeans -init 0 -max-candidates 3	100 -periodic-pruning 1000	0 -min-density 2.0 -t1	-1.25 -t2 -1.0	– N 3 – A "weka.co	ore.EuclideanDistance	-R first-
Cluster mode	Clusterer output					
<ul> <li>Use training set</li> <li>Supplied test set</li> <li>Set</li> <li>Percentage split</li> <li>Classes to clusters evaluation         <ul> <li>(Nom) class</li> <li>Store clusters for visualization</li> </ul> </li> </ul>	Initial starting poi Cluster 0: 6.1,2.9,4 Cluster 1: 6.2,2.9,4 Cluster 2: 6.9,3.1,5 Missing values globa	nts (random): .7,1.4,Iris-versico .3,1.3,Iris-versico .1,2.3,Iris-virginio lly replaced with mo	lor lor ca ean/mode	// <del>+</del>		
	Final cluster centro	ids:				
Ignore attributes Start Stop	Attribute	Full Data (150.0)	Cluster# 0 (50.0) ==========	1 (50.0)	2 (50.0)	
Result list (right-click for options)	sepallength sepalwidth petallength petalwidth class	5.8433 3.054 3.7587 1.1987 Iris-setosa Iris-	5.936 2.77 4.26 1.326 -versicolor	5.006 3.418 1.464 0.244 Iris-setosa	6.588 2.974 5.552 2.026 Iris-virginica	
	Time taken to build	model (full training	g data) : 0 se	econds		
	=== Model and evalua	tion on training se [.]	t ===			
	Clustered Instances					
	0 50 (33%) 1 50 (33%) 2 50 (33%)	]				
		)				<b>V</b>
Status						

ОК

Log 🔨 x 0





Weka Explorer

Preprocess Classify Cluster Associate Select attributes Visualize

#### Clusterer

Choose SimpleKMeans -init 0 -max-candidates 100 -periodic-pruning 10000 -min-density 2.0 -t1 -1.25 -t2 -1.0 -N 3 -A "weka.core.EuclideanDistance -R first-

Cluster mode	Clusterer output
<ul> <li>Use training set</li> <li>Supplied test set</li> <li>Percentage split % 66</li> <li>Classes to clusters evaluation</li> </ul>	sepallength         5.8433         5.8885         5.006         6.8462           sepalwidth         3.054         2.7377         3.418         3.0821           petallength         3.7587         4.3967         1.464         5.7026           petalwidth         1.1987         1.418         0.244         2.0795
Store clusters for visualization	Time taken to build model (full training data) : 0 seconds
Ignore attributes       Start     Stop       Result list (right-click for options)       11:17:51 - SimpleKMeans	=== Model and evaluation on training set === Clustered Instances 0 61 ( 41%) 1 50 ( 33%) 2 39 ( 26%)
11:21:09 - SimpleKMeans	Class attribute: class Classes to Clusters: 0 1 2 < assigned to cluster 0 50 0   Iris-setosa 47 0 3   Iris-versicolor 14 0 36   Iris-virginica Cluster 0 < Iris-versicolor Cluster 1 < Iris-setosa Cluster 2 < Iris-virginica Incorrectly clustered instances : 17.0 11.3333 %

👞 х О

Log

Status



Please **cite us** if you use the software.

#### 2.3. Clustering

- 2.3.1. Overview of clustering methods
- 2.3.2. K-means
- 2.3.2.1. Mini Batch K-Means
- 2.3.3. Affinity Propagation
- 2.3.4. Mean Shift
- 2.3.5. Spectral clustering
- 2.3.5.1. Different label assignment strategies
- 2.3.5.2. Spectral Clustering Graphs

2.3.6. Hierarchical clustering

- 2.3.6.1. Different linkage type: Ward, complete, average, and single linkage
- 2.3.6.2. Adding connectivity constraints
- 2.3.6.3. Varying the metric
- 2.3.7. DBSCAN
- 2.3.8. Birch

Clustering

Each clustering algorithm comes in two variants: a class, that implements the fit method to learn the clusters on train data, and a function, that, given train data, returns an array of integer labels corresponding to the different clusters. For the class, the labels over the training data can be found

in the labels attribute.

#### Input data

One important thing to note is that the algorithms implemented in this module can take different kinds of matrix as input. All the methods accept standard data matrices of shape [n_samples, n_features]. These can be obtained from the classes in the **sklearn.feature_extraction** module. For **AffinityPropagation**, **SpectralClustering** and **DBSCAN** one can also input similarity matrices of shape [n_samples, n_samples]. These can be obtained from the functions in the **sklearn.metrics.pairwise** module.

#### 2.3.1. Overview of clustering methods

MiniBatchKMeansAffinityPropagation MeanShift SpectralClustering

Ward AgglomerativeClustering DBSCAN

Birch Gaussian

GaussianMixture





#### 





### **Problems with K-Means**

- Only works for numeric data (typically reals)
- Very sensitive to the initial points
  - -fix: Do many runs, each with different initial centroids
  - fix: Seed centroids with non-random method, e.g., farthest-first sampling
- Sensitive to outliers
  - -E.g.: find three
  - -fix: identify and remove outliers
- Must manually choose k
  - -Learn optimal k using some performance measure

#### **Problems with K-Means**

• How do you tell it which clustering you want?



Constrained clustering technique



## (2) Hierarchical clustering

- Agglomerative
  - Bottom-up approach: elements start as individual clusters & clusters are merged as one moves up the hierarchy
- Divisive
  - –Top-down approach: elements start as a single cluster & clusters are split as one moves down the hierarchy

#### **Hierarchical Clustering**

Recursive partitioning/merging of a data set



#### **Dendogram**

- Tree structure representing all data partitionings
- Constructed as clustering proceeds



Nine items

#### **Dendogram**

- Tree structure representing all data partitionings
- Constructed as clustering proceeds
- Get a K-clustering by looking at connected components at any given level
- Often binary dendograms, but n-ary ones easy to obtain with minor algorithm changes



Four clusters

#### **Hierarchical clustering advantages**

- Need not specify number of clusters
- Good for data visualization
  - See how data points interact at many levels
  - Can view data at multiple granularity levels
  - Understand how all points interact
- Specifies all of the K clusterings/partitions

## **Divisive hierarchical clustering**

- Top-down technique to find best partitioning of data, generally exponential in time
- Common approach:
  - -Let **C** be a set of clusters
  - -Initialize **C** to be a one-clustering of data
  - -While there exists a cluster c in C
    - remove *c* from **C**
    - partition c into 2 clusters (c₁ and c₂) using a flat clustering algorithm (e.g., k-means)
    - Add to  $c_1$  and  $c_2$  **C**
- Bisecting k-means











split using flat clustering







All observations start as their own cluster. Clusters meeting some criteria are merged. This process is repeated, growing clusters until some end point is reached.



#### **Hierarchical Agglomerative Clustering**

- Let **C** be a set of clusters
- Initialize C to all points/docs as separate clusters
- While **C** contains more than one cluster
  - -find  $c_1$  and  $c_2$  in **C** that are **closest together**
  - -remove  $c_1$  and  $c_2$  from **C**
  - merge  $c_1$  and  $c_2$  and add resulting cluster to **C**
- Merging history forms a binary tree or hierarchy
- Q: How to measure distance between clusters?



# **Single-link:** Similarity of the *most* similar (single-link)

![](_page_38_Figure_3.jpeg)

![](_page_39_Picture_0.jpeg)

# **Complete-link:** Similarity of the "furthest" points, the *least* similar

![](_page_39_Figure_3.jpeg)

![](_page_39_Figure_4.jpeg)

#### Weka: linkType=COMPLETE

![](_page_40_Picture_0.jpeg)

**Centroid:** Clusters whose centroids (centers of gravity) are the most similar

![](_page_40_Figure_3.jpeg)

$$\left\|\mu(L)-\mu(R)\right\|^2$$

Weka: linkType=CENTROID

![](_page_41_Picture_0.jpeg)

# **Average-link:** Average similarity between all pairs of elements

![](_page_41_Picture_3.jpeg)

 $\frac{1}{|L| \cdot |R|} \sum_{x \in L, y \in R} ||x - y||^2$  Weka: linkType=AVERAGE

	Weka Explorer
Preprocess Classify Cluster Associate Select att	ributes Visualize
Clusterer	
Chopse HierarchicalClusterer -N 3 -L SINGLE -P	-A "weka.core.EuclideanDistance -R first-last"
Cluster mode	Clusterer output
<ul> <li>Use training set</li> <li>Supplied test set</li> <li>Set</li> </ul>	Cluster 0 (((((((((((((((((((((((((((((())))))))
<ul> <li>Percentage split % 66</li> <li>Classes to clusters evaluation</li> <li>(Nom) class</li> </ul>	Cluster 2 ((((((((((((((((((((((((((((((((((((
✓ Store clusters for visualization	Time taken to build model (full training data) : 0.01 seconds
Ignore attributes Start Ignore attributes during clustering	<pre>g Clustered Instances 0 49 ( 33%)</pre>
Result list (right-click for options)	1 1 ( 1%) 2 100 ( 67%) Class attribute: class
	Classes to Clusters: 0 1 2 < assigned to cluster 49 1 0   Iris-setosa 0 0 50   Iris-versicolor 0 0 50   Iris-virginica Cluster 0 < Iris-setosa Cluster 1 < No class Cluster 2 < Iris-versicolor Incorrectly clustered instances : 51 0 24 %

Defaut **SINGLE** cluster distance gives poor results here

	Weka Explorer
Preprocess Classify Cluster Associate Select at	tributes Visualize
Clusterer	
Choose HierarchicalClusterer -N 3 -L AVERAGE	-P -A "weka.core.EuclideanDistance -R first-last"
Cluster mode	Clusterer output
<ul> <li>Use training set</li> <li>Supplied test set</li> <li>Percentage split</li> <li>Classes to clusters evaluation</li> </ul>	Cluster 1 ((((((((1.4:0.08775,(1.5:0.06508,1.5:0.06508):0.02267):0.04395,1.7:0.1317):0.01307,((1.5:0.0 Cluster 2 (((((2.5:0.12797,(2.3:0.10565,(2.4:0.06047,2.3:0.06047):0.04518):0.02232):0.06295,(((2.1:0.
(Nom) class       Image: Store clusters for visualization       Ignore attributes	Time taken to build model (full training data) : 0.01 seconds
Start     Stop       Result list (right-click for options)       10:09:16 - HierarchicalClusterer	Clustered Instances 0 50 (33%) 1 67 (45%) 2 33 (22%)
10:09:58 - HierarchicalClusterer	Class attribute: class Classes to Cluster: 0 1 2 < assigned to cluster 50 0 0   Iris-setosa 0 50 0   Iris-versicolor 0 17 33   Iris-virginica Cluster 0 < Iris-setosa Cluster 1 < Iris-versicolor Cluster 2 < Iris-virginica Incorrectly clustered instances : 17 0 11 2222 &

Using **AVERAGE** cluster distance measure improves results

### Knowing when to stop

- General issue is knowing when to stop merging/splitting a cluster
- We may have a problem specific desired range of clusters (e.g., 3-6)
- There are some general metrics for assessing quality of a cluster
- There are also domain specific heuristics for cluster quality

## (3) DBSCAN Algorithm

- Density-Based Spatial Clustering of Applications with Noise
- Clusters close points based on a distance and a minimum number of points
  - Key parameters: eps=maximum distance between two points; minPoints= minimal cluster size
- Marks as outliers points in low-density regions
- Needn't specify number of clusters expected
- Fast

![](_page_46_Picture_0.jpeg)

DBSCAN looks for densely packed observations and makes no assumptions about the number or shape of clusters.

- 1. A random observation, x;, is selected
- 2. If x; has a minimum of close neighbors, we consider it part of a cluster.
- 3. Step 2 is repeated recursively for all of x's neighbors, then heighbors' neighbors etc... These are the cluster's core members.
- 4. Once Step 3 runs out of observations, a new random point is chosen

Afterwards, observations not part of a core are assigned to a nearby cluster or marked as outliers.

ChrisAlbon

#### **Comparing Clustering algorithms**

![](_page_47_Figure_1.jpeg)

Scikit Learn — Plot Cluster Comparison

#### **DBSCAN Example**

![](_page_48_Figure_1.jpeg)

This gif (in ppt) shows how DBSCAN grows four clusters and identifies the remaining points as outliers

## Visualizing DBSCAN https://bit.ly/471dbscan

![](_page_49_Picture_1.jpeg)