

Basics: Machine Learning

Lecture 6

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Acknowledgement to all authors whose materials have been used

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Machine Learning

- The term **Machine Learning** was coined by **Arthur Samuel** in 1959:

“Machine Learning algorithms enable the computers to **learn from data**, and even **improve** themselves, **without being explicitly programmed**”.

- In 1997, Tom Mitchell gave a mathematical and relational definition that:

“A computer program is said to learn from **experience E** with respect to some **task T** and some **performance measure P**, if its performance on T, as measured by P, improves with experience E”.

ML Examples

Family-friendly hotels in Istanbul



Hotel Amira Istanbul

★★★★★ 4,017 Reviews

Istanbul, Turkey

"They were personable, funny, very helpful, and provided the **total package** in terms of the accommodation."



Muyan Suites

★★★★★ 1,149 Reviews

Istanbul, Turkey

"... card to travel to takism... Overall just perfect and awesome place thatz Muyan suites for a best **vacation** in istanbul.. Wish to go back again n stay only with thix small ill family of ours now.... Thank ..."



Hotel Yasmak Sultan

★★★★★ 1,842 Reviews

Istanbul, Turkey

"This hotel really is the **whole package** -- comfortable, clean, superbly located, has a lovely rooftop restaurant, and even has its own hamam."



White House Hotel Istanbul

★★★★★ 4,593 Reviews

Istanbul, Turkey

"Amazing and very clean Hotel !!! Have a great sleep and nicely **holiday**!!!! Amazing stuff very helpful when you need something !!! Amazing manager , very professional !!! Thank you very much to all team "

See all

Luxury hotels in Istanbul

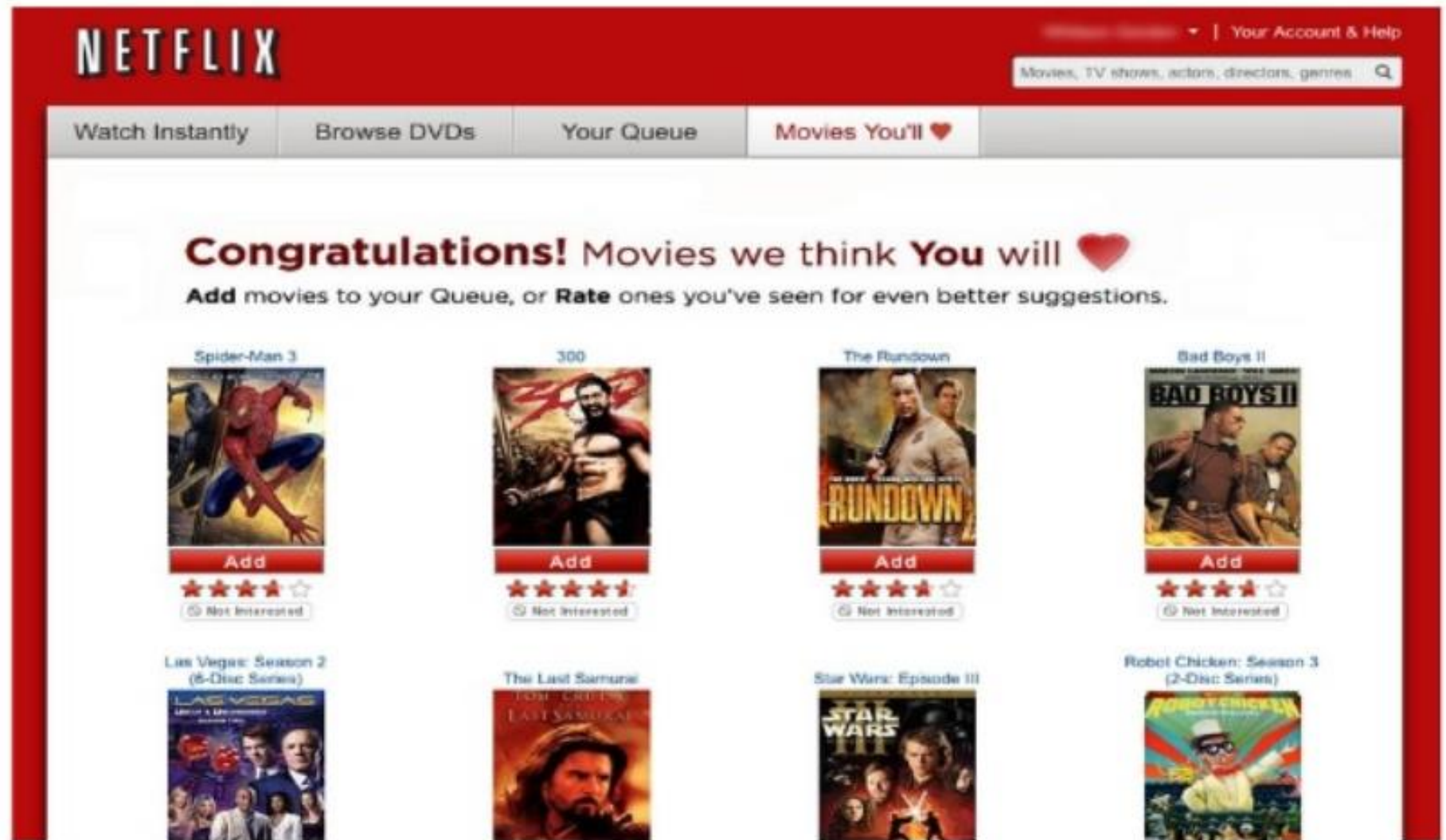


See all

Example 1:

- Suppose you decide to check out trip offers for a vacation
- You browse through the travel agency website and search for a hotel
- When you look at a specific hotel, just below the hotel description there is a section titled “**You might also like these hotels**”.
- This is a common **use case of Machine Learning** called “**Recommendation Engine**”
 - Many data points were used to train a model in order to predict what will be the best hotels to show you under that section, based on a lot of information they already know about you

Example: Netflix



ML Examples

Example 2:

- Program to **predict** traffic patterns at a busy intersection (task T)
- Run it through a machine learning algorithm with data (task T) about **past traffic patterns (experience E)** and, if it has successfully “**learned**”, it will then do better in **predicting future traffic patterns (performance measure P)**.

ML Examples

Example 3: Learn to detect SPAM

- T: Distinguish between SPAM and Non-SPAM
- P: % of emails correctly classified
- E: Labeled emails from your friend Abdullah

Examples of machine learning problems

- **Medical diagnoses:** ML is trained to recognize cancerous tissues
 - “Is this cancer?”
- **Graph Processing:**
 - “Which of these people are good friends with each other?”
- **Recommender Systems:**
 - “Will person X likes movie Y?”
- **Financial industry and trading** —fraud investigations and loan sanction
- **Speech Recognition:**
 - “Is this his/her voice?” (voice searches, voice dialing, call routing, and appliance control)

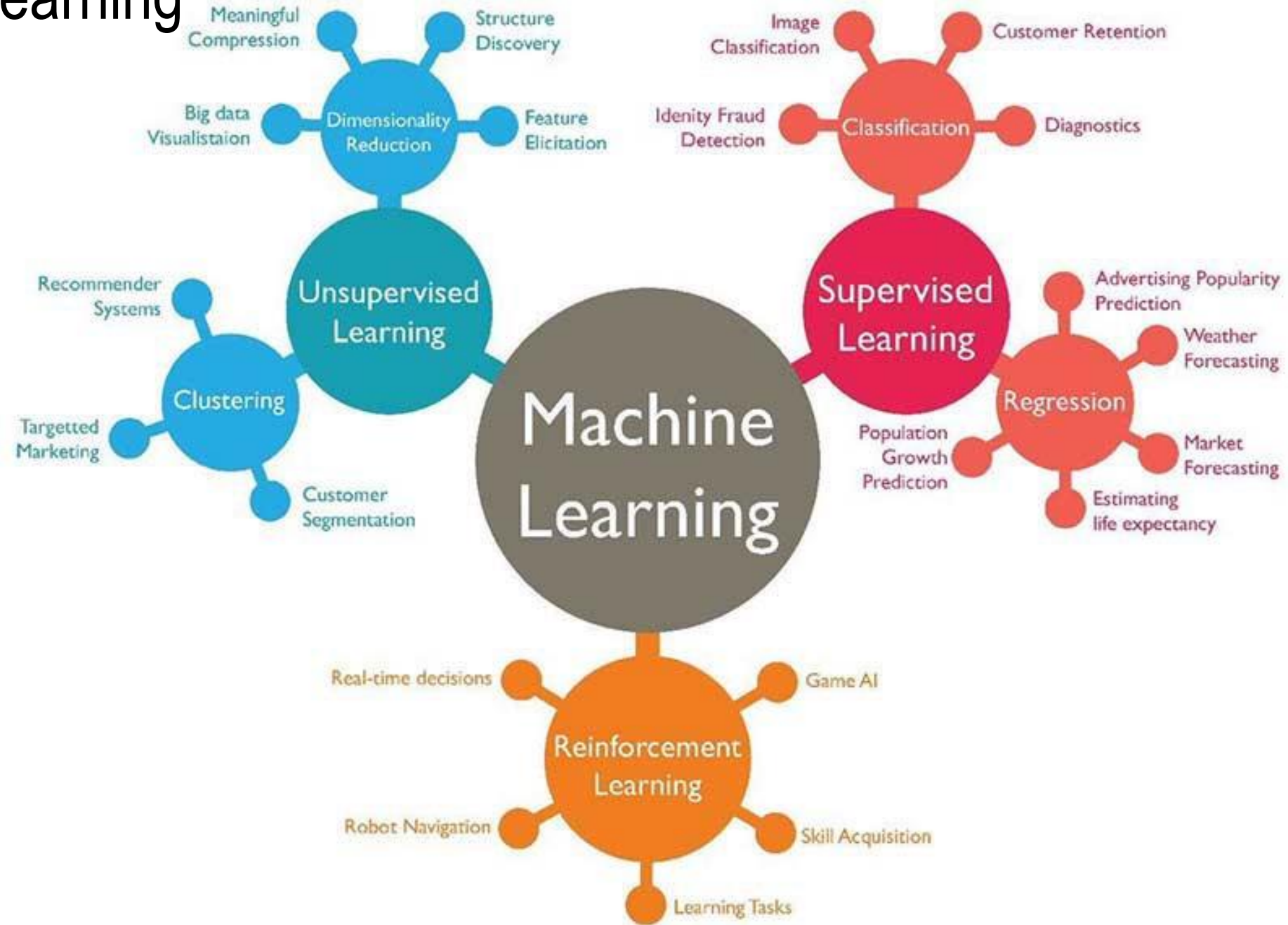
Such problems are excellent targets for Machine Learning, and in fact machine learning has been applied to such problems with great success.

What is a Model

- A **model** is a mathematical formula with **a number of parameters** that need to be **learned from the data**
 - Fitting a model to the data is a process known as **model training**
- **Example:** Consider a one feature/variable linear regression, where the goal is to fit a line (described by the equation $y = ax + b$) to a set of distributed data points.
- Once the model training is completed we get a model equation $y = 2x + 5$.
 - Then for a set of inputs $[1, 0, 7, 2, \dots]$ we would get a set of outputs $[7, 5, 19, 9, \dots]$.

Types of Machine Learning

Machine learning can be classified into 3 types of algorithms.



Supervised Learning

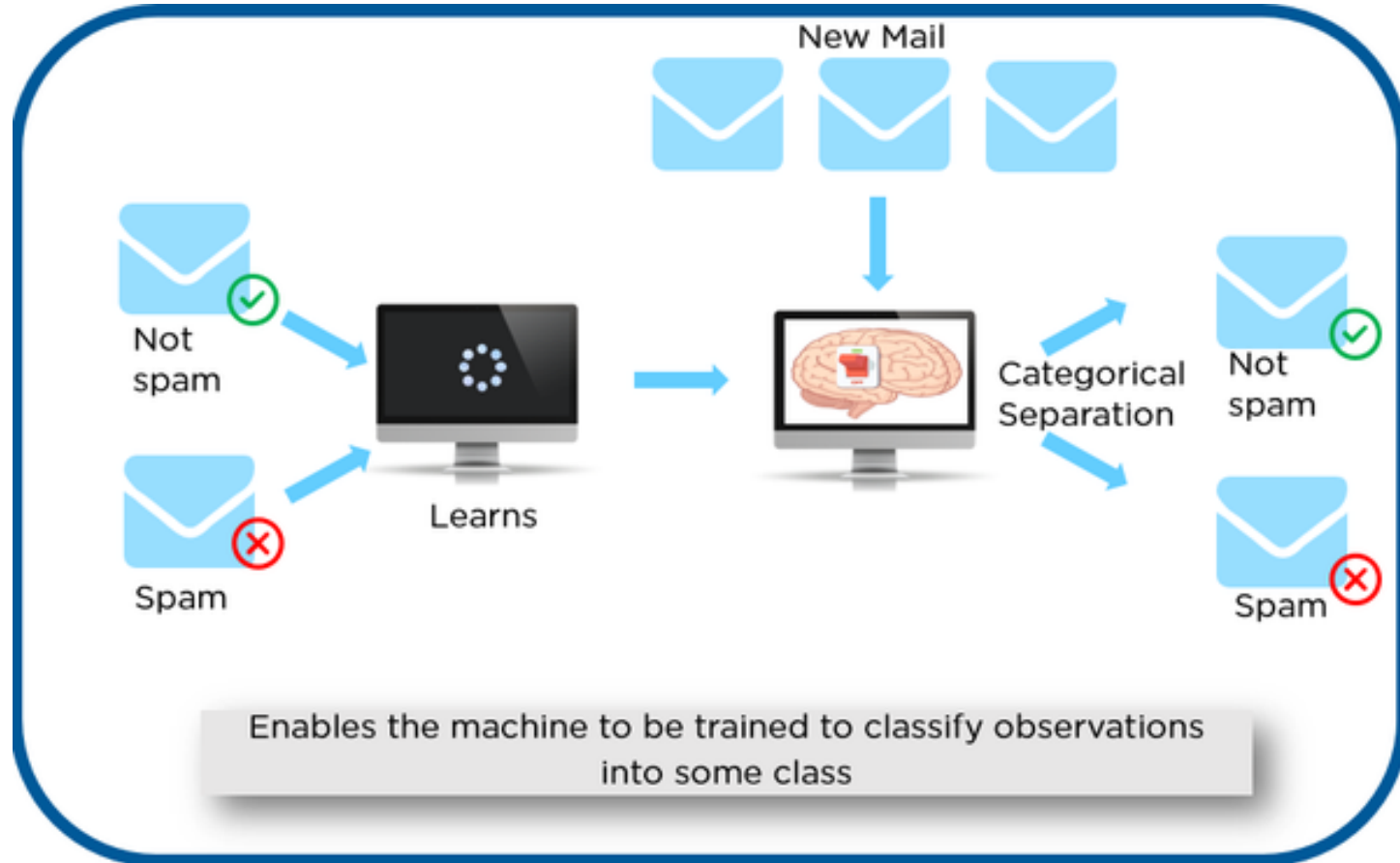
- Supervised learning is a learning model built to make prediction, given an unforeseen input instance.
- A supervised learning algorithm takes a known set of input dataset and its known responses to the data (output) to learn the model.
- A learning algorithm then trains a model to generate a prediction for the response to new data or the test dataset.
- Supervised learning uses classification algorithms and regression techniques to develop predictive models.
- The algorithms include linear regression, logistic regression, neural networks, decision tree, Support Vector Machine (SVM), random forest, naive Bayes, and k-nearest neighbor.

Supervised Learning

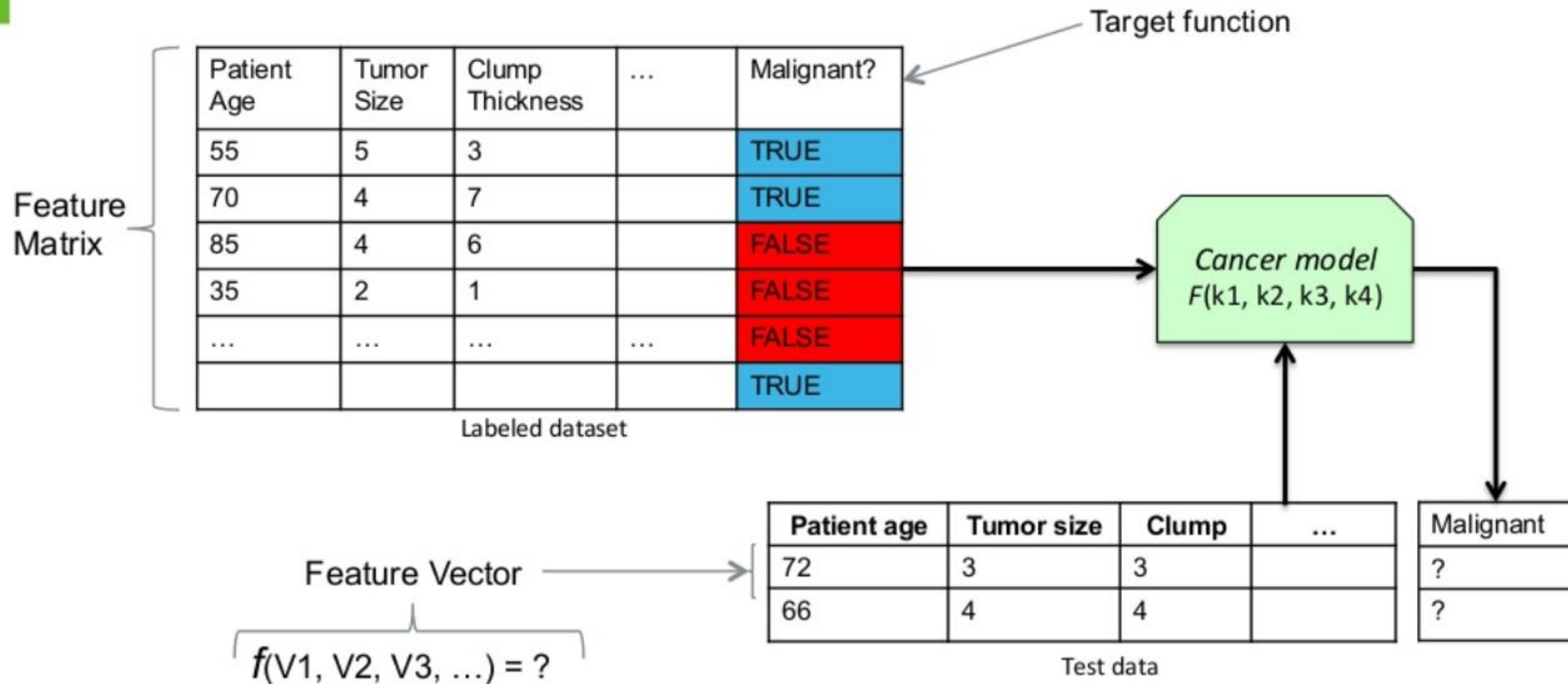
For example:

Spam filtering where Large number of email messages are labelled as either:

- spam
 - non-spam
-
- New email message will then be classified as spam or non-spam



Supervised Learning: learn from examples



Supervised Learning

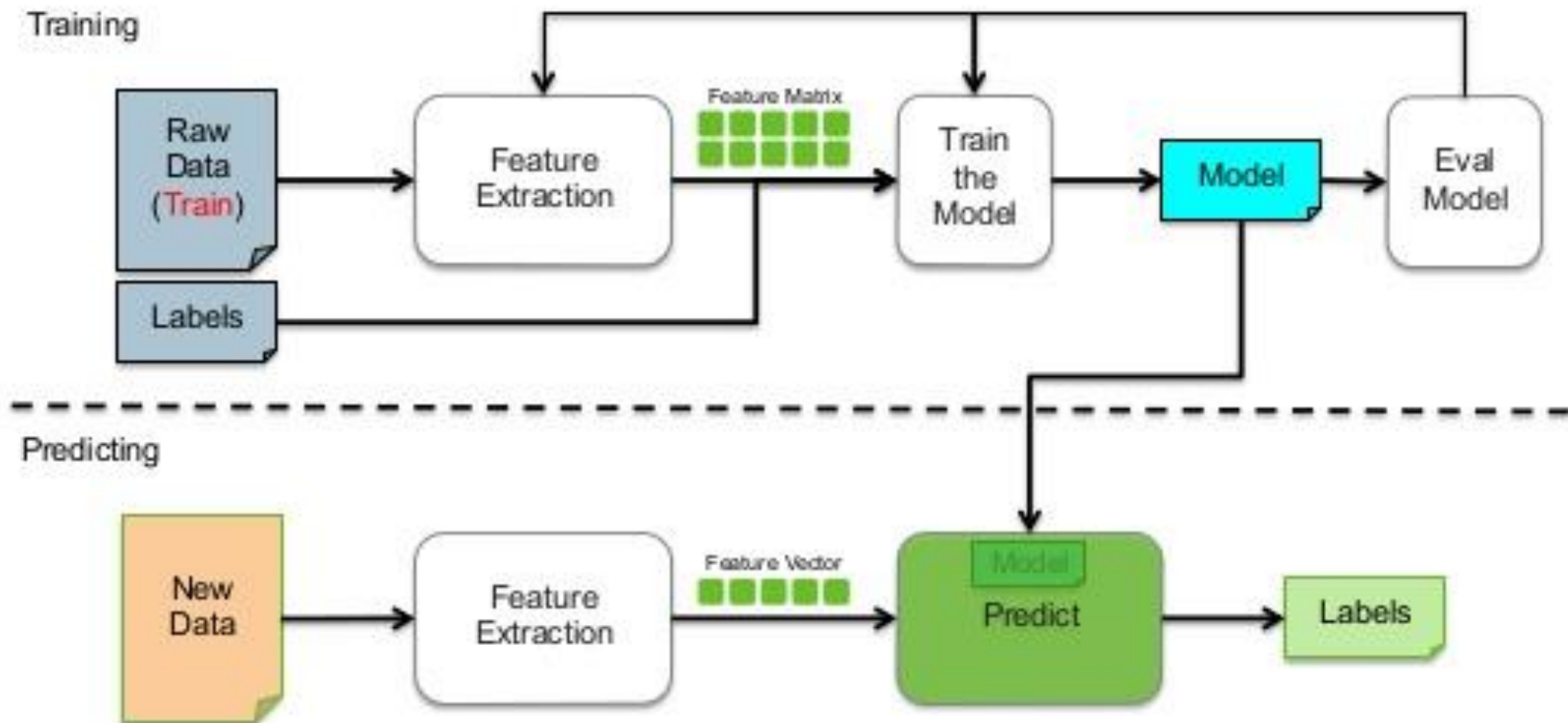


Training
Images



Test
Image

Supervised Learning Workflow



Classification

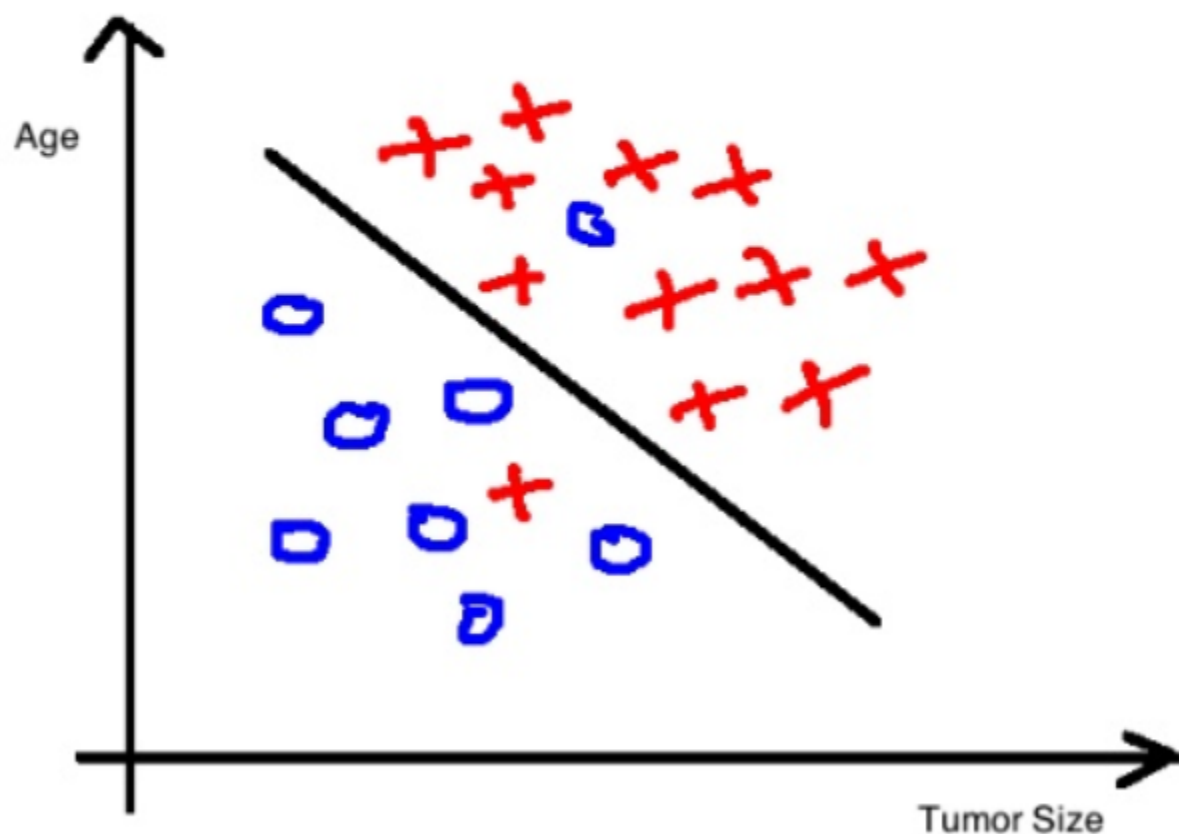
Classification learns from existing categorizations and then assigns unclassified items to the best category.

- Classification models classify input data into categories and **predicts discrete responses**
- Classification is recommended if the data can be **categorized, tagged, or separated into specific groups or classes**
- **Classification Examples:**
 - Bank credit scoring
 - Medical imaging
 - Speech recognition
 - To recognize letters and numbers in Handwriting
 - To check whether an email is genuine or spam
 - To detect whether a tumor is benign or cancerous
- **Classification Algorithms:**
 - k-nn, Decision Trees, Random Forest, SVM, Neural Network...

Classification Algorithms

- **Classification algorithms** attempt to estimate the mapping function (f) from the input variables (x) to discrete or categorical output variables (y).
 - In this case, y is a category that the mapping function predicts.
- **For example**, when provided with a dataset about houses, a classification algorithm predict whether the prices for the houses “sell more or less than the recommended retail price”
- **For example**, in a banking application, the customer who applies for a loan may be classified as a safe and risky according to his/her age and salary. The constructed model can be used to classify new data

Classification: predicting a category



Some techniques:

- Naïve Bayes
- Decision Tree
- Logistic Regression
- SGD
- Support Vector Machines
- Neural Network
- Ensembles

Basics: Regression Algorithms

Regression techniques predict continuous responses

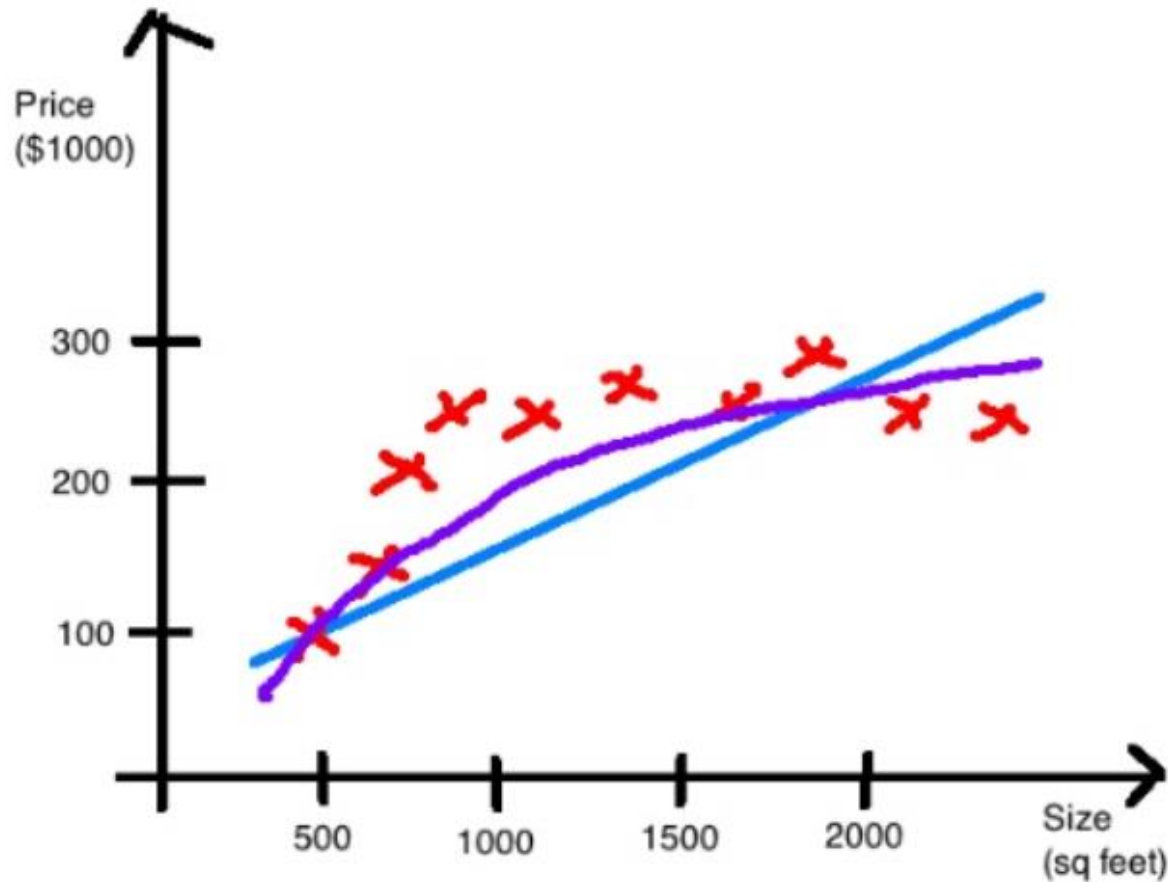
Regression techniques predict a continuous-valued attribute associated with an object

- **Regression algorithms** attempt to estimate the mapping function (f) from the input variables (x) to numerical or continuous output variables (y).
 - In this case, y is a real value, which can be an integer or a floating point value.
 - Therefore, regression prediction problems are usually quantities or sizes.
- For example, when provided with a dataset about houses, and you are asked to predict their prices, that is a regression task because price will be a continuous output.
- Regression algorithms include linear regression, Ensembles, Support Vector Regression (SVR), and regression trees.

Regression Examples:

- A linear regression attempts to model the relationship between two variables by fitting a linear equation to observed data
- For example, a data is collected about how happy people are after getting so many hours of sleep.
 - In this dataset, sleep and happy people are the variables.
- Other Examples:
 - Drug Response, Stock Price, ...

Regression: predict a continuous value



Some techniques:

- Linear Regression / GLM
- Decision Trees
- Support vector regression
- SGD
- Ensembles

Unsupervised Learning

In unsupervised learning the training data comprises examples of input vectors WITHOUT any corresponding target variables.

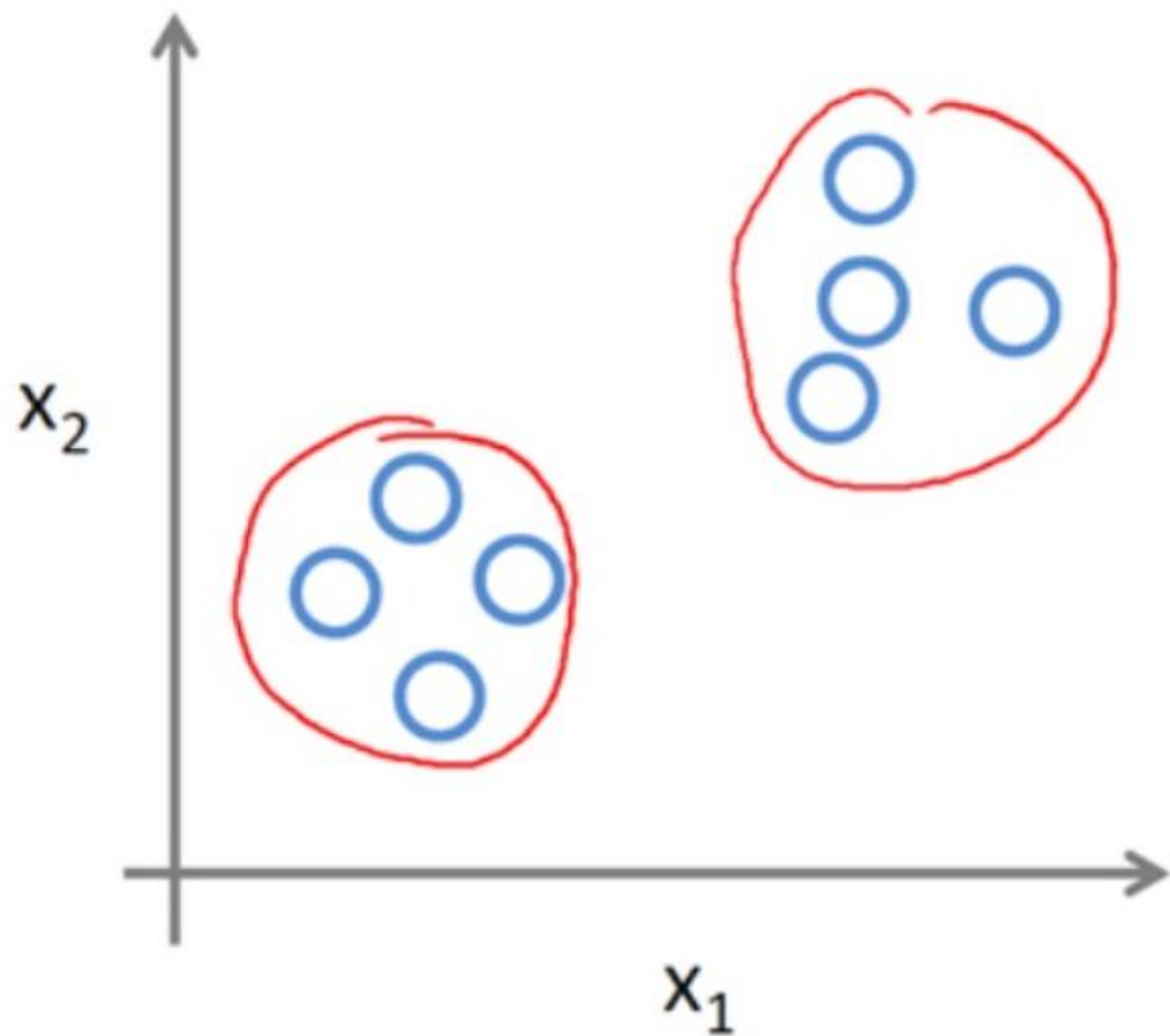
- In unsupervised learning, the algorithm **learns from plain examples without any associated response**, leaving to the **algorithm to determine the data patterns on its own**.
 - In an **unsupervised learning** you only have input data (X) and no corresponding output variables.
- The goal for unsupervised learning is to model the **underlying structure** or **distribution in the data** in order to learn more about the data.

Unsupervised Learning

Examples: Speech recognition, document clustering, and image compression.

- In **document clustering**, the aim is to group documents into various reports of politics, entertainment, sports, culture, heritage, art, and so on.
- **Fraud Detection**: Identify groups of motor insurance policy holders with a high average claim cost
- **Social Networks**: Recognize communities within large groups of people

Clustering: detect similar instance groupings



Some techniques:

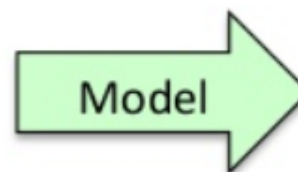
- k-means
- Spectral clustering
- DB-scan
- Hierarchical clustering

Unsupervised Learning: detect natural patterns

Age	State	Annual Income	Marital status
25	CA	\$80,000	M
45	NY	\$150,000	D
55	WA	\$100,500	M
18	TX	\$85,000	S
...

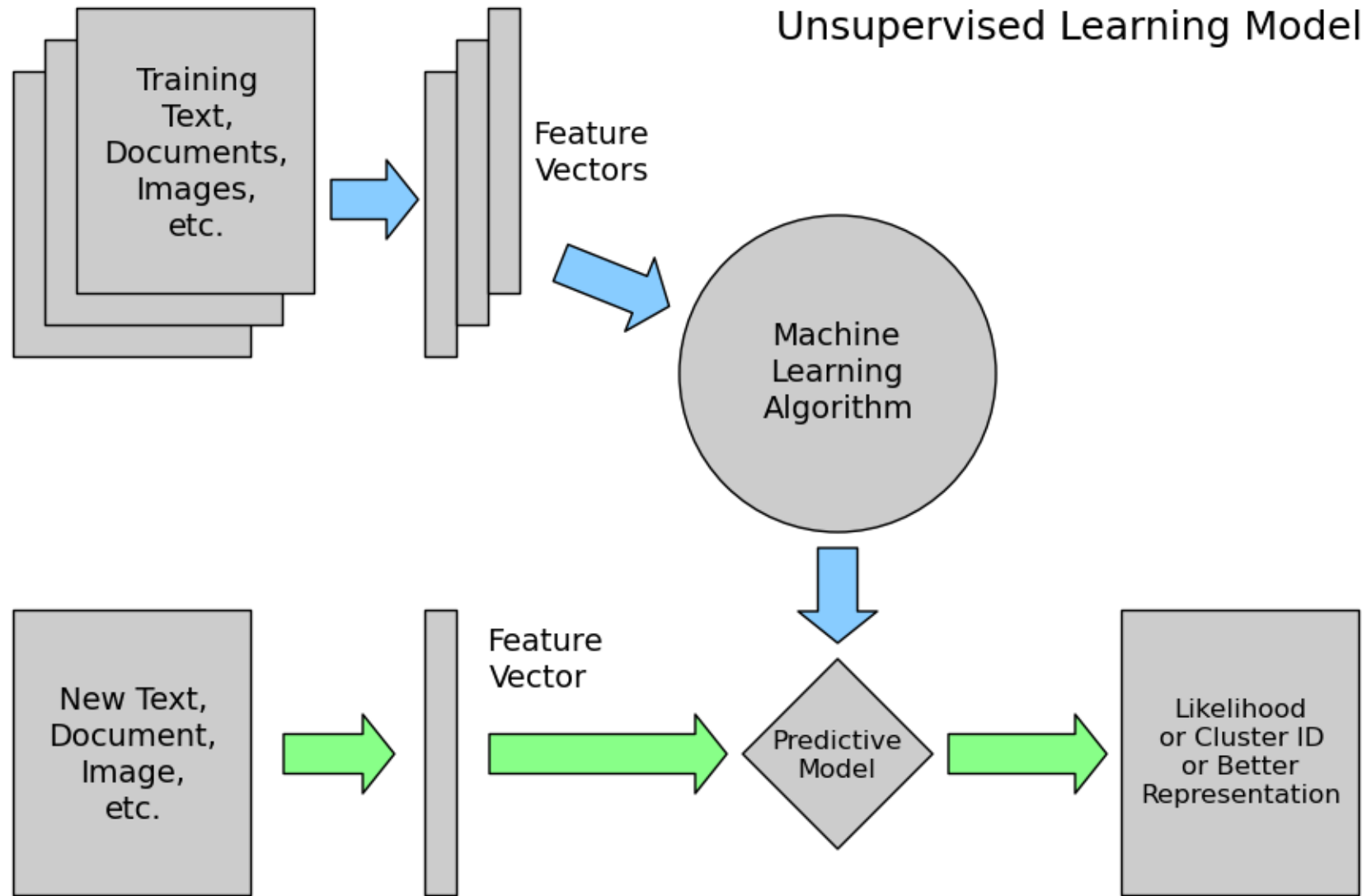


No labels



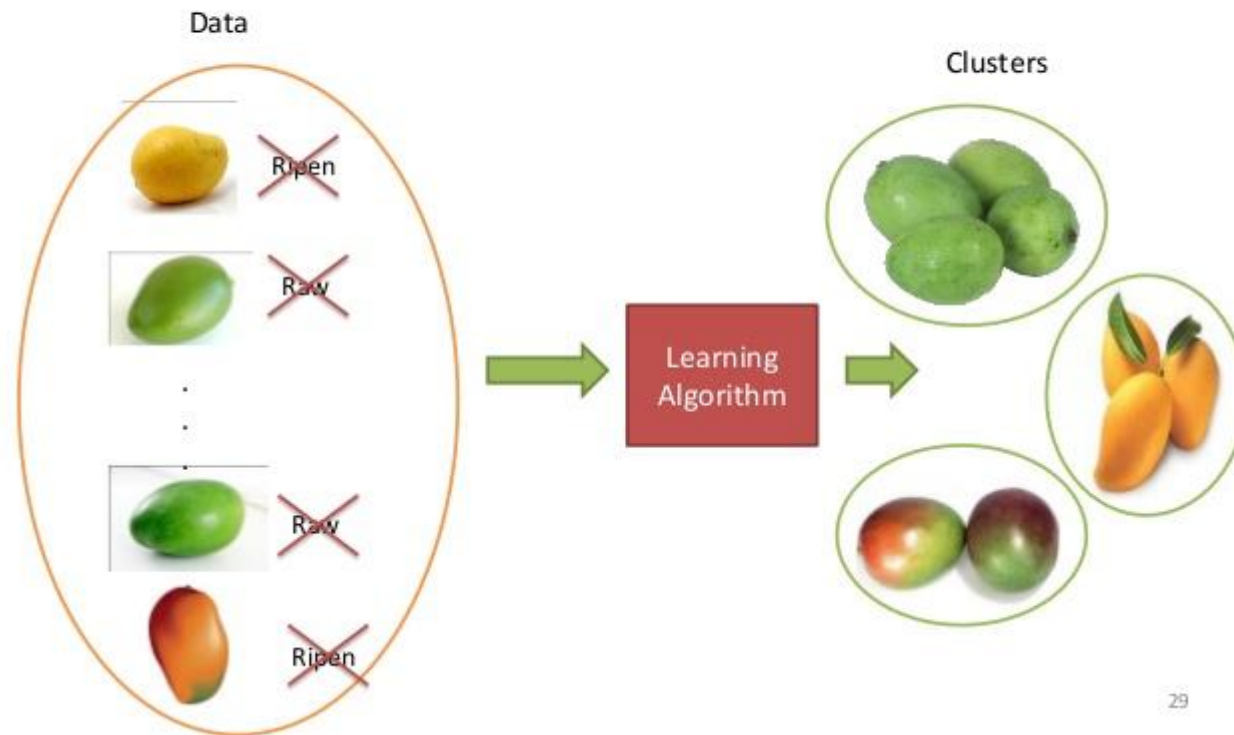
Naturally occurring
(hidden) structure

Unsupervised Learning



Unsupervised Learning

Unsupervised Learning



Unsupervised Learning



Objective is simply to divide above Images into N groups
Here ideally $N = 2$

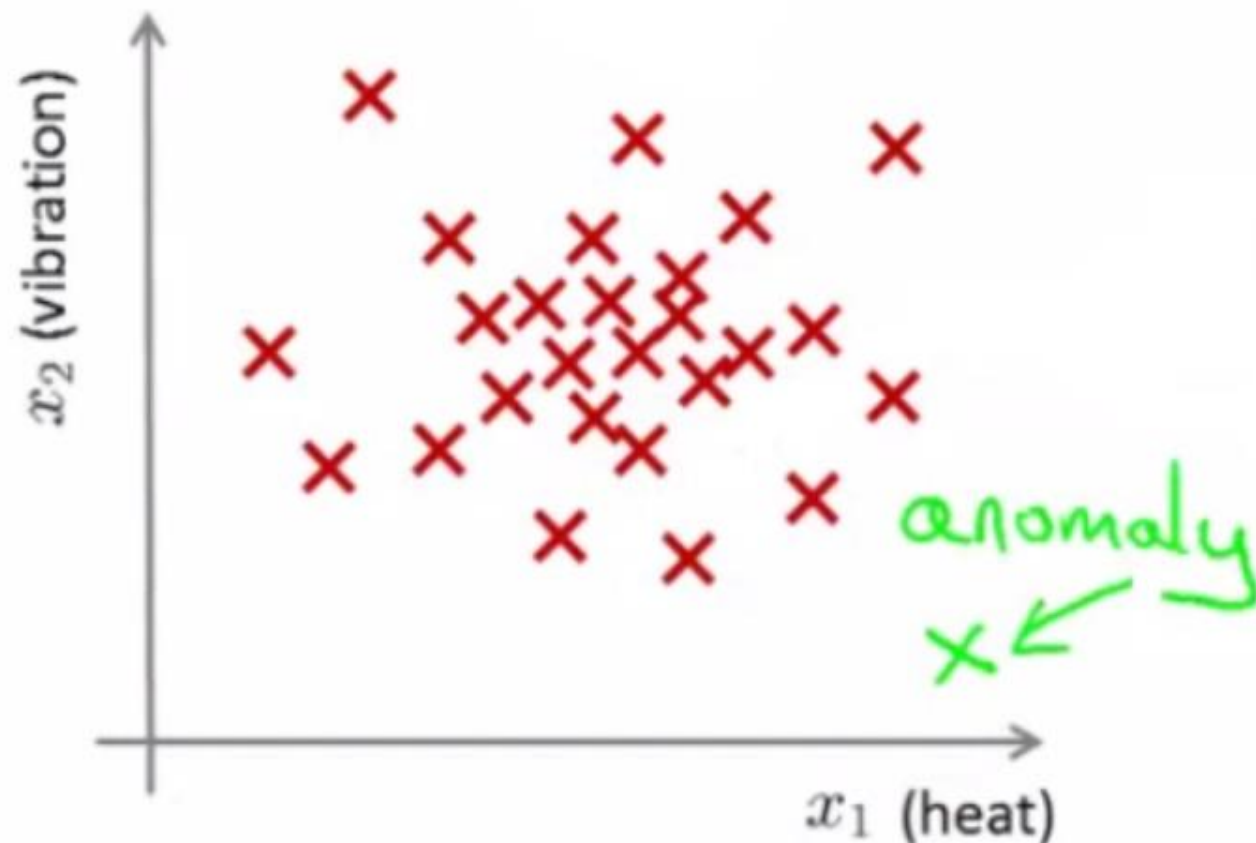
The method can also assume other groups e.g. Group with one object or group with multiple objects

Outlier Detection: identify abnormal patterns

Example: identify engine anomalies

Features:

- Heat generated
- Vibration of engine



Affinity Analysis: identifying frequent item sets

	Item 1	Item 2	Item 3	Item 4	Item 5	...
Tx 1	Y	N	N	Y	N	
Tx 2	Y	N	N	Y	N	
Tx 3	Y	Y	N	Y	N	
Tx 4	N	N	Y	Y	Y	
Tx 5						
...						

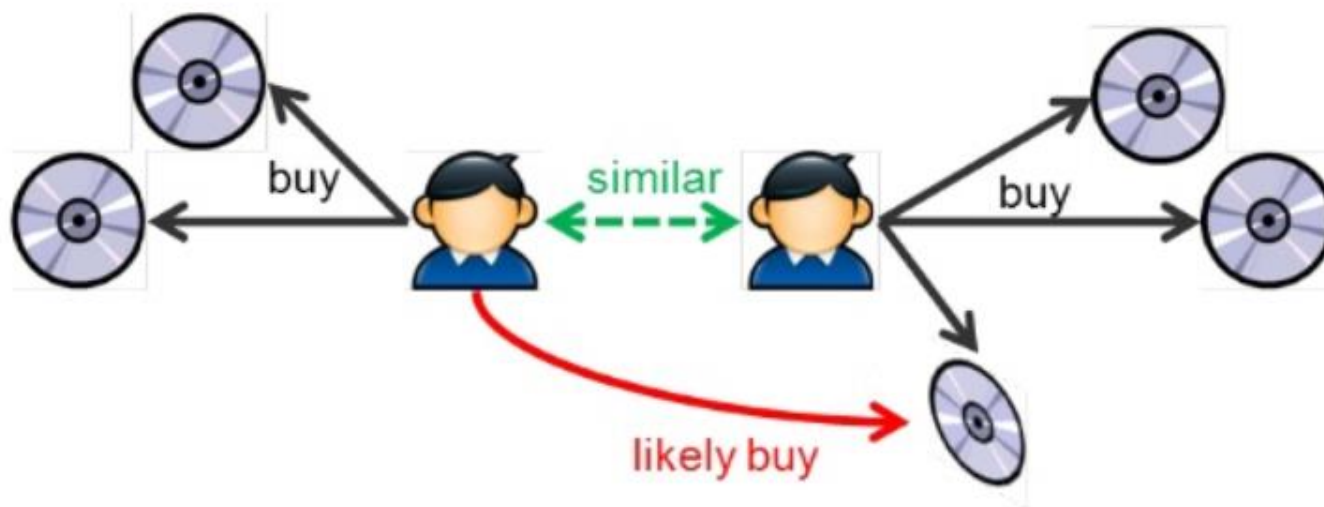


	Item 1	Item 2	Item 3	Item 4	Item 5	...
Tx 1	Y	N	N	Y	N	
Tx 2	Y	N	N	Y	N	
Tx 3	Y	Y	N	Y	N	
Tx 4	N	N	Y	Y	Y	
Tx 5						
...						

Goal: identify frequent item set
Techniques: FP Growth, a priori




Product recommendation: predicting “preference”



Collaborative Filtering
Identify users with similar “taste”

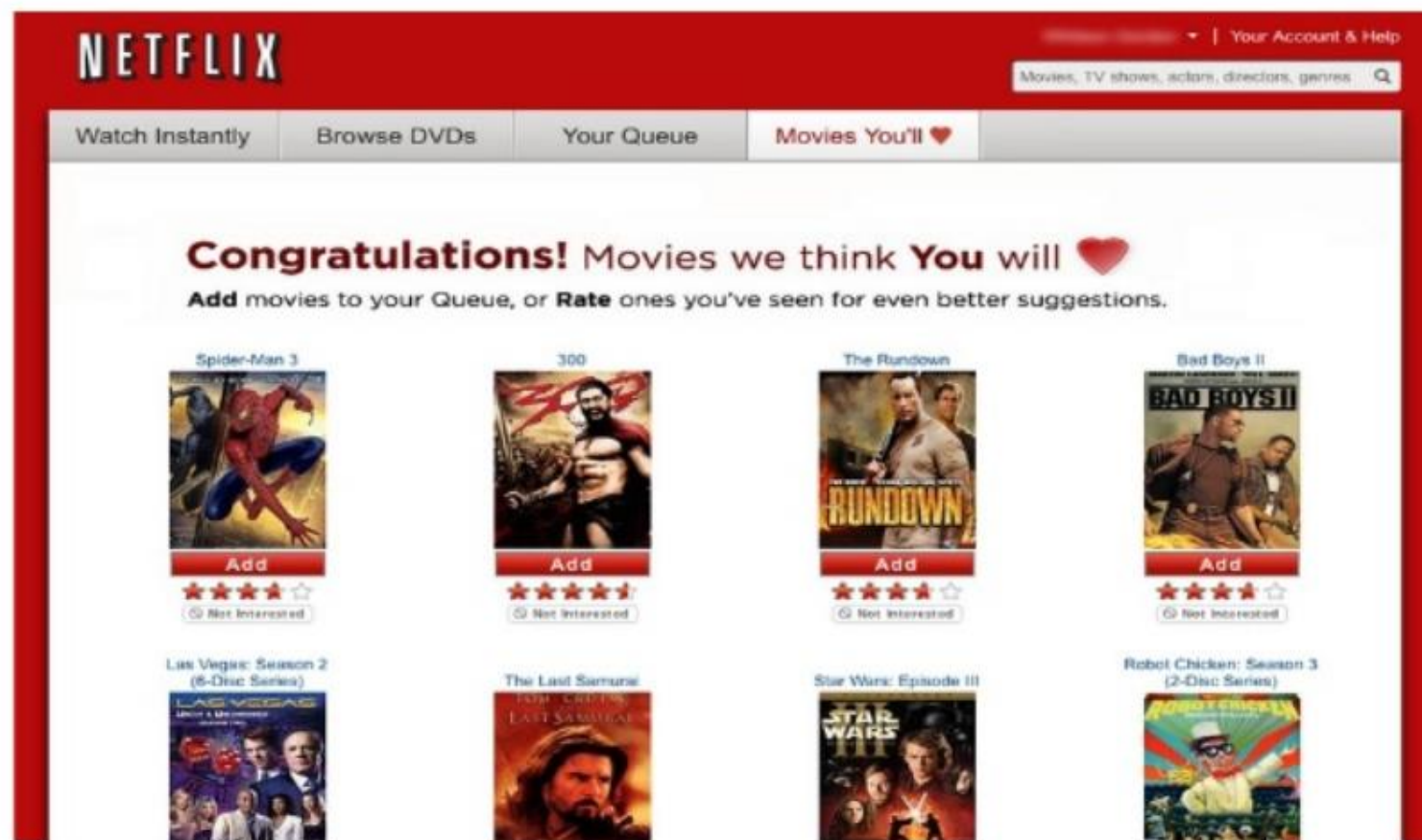
Collaborative filtering -> matrix completion

	Harry potter	X-Men	Hobbit	Argo	Pirates
101	5	2	4	?	?
102	?	?	5	2	?
103	1	2	?	?	3
104					
105					
...					



	Harry potter	X-Men	Hobbit	Argo	Pirates
101	5	2	4	1	3
102	4	1	5	2	3
103	1	2	4	1	3
104					
105					
...					

Example: Netflix



Example: market segmentation



Types of Unsupervised Learning

- In **Clustering** similar instances are grouped, based on their features or properties.
- **Association**: Association rules find associations amongst items within large commercial databases (e.g. **Collaborative filtering**)
 - Discover rules that describe large portions of data, such as people that buy X also tend to buy Y

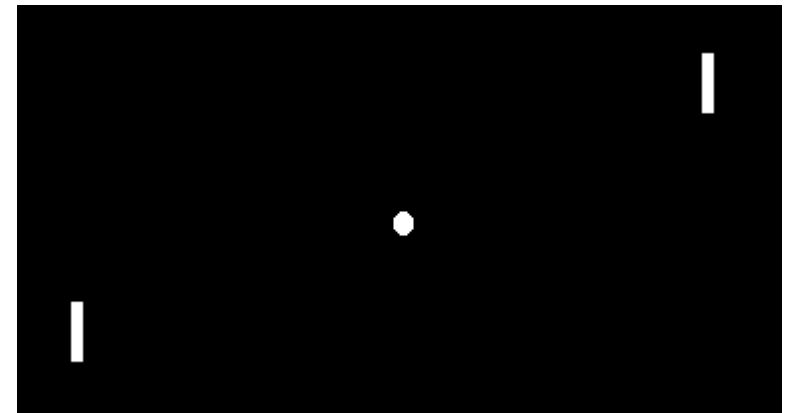
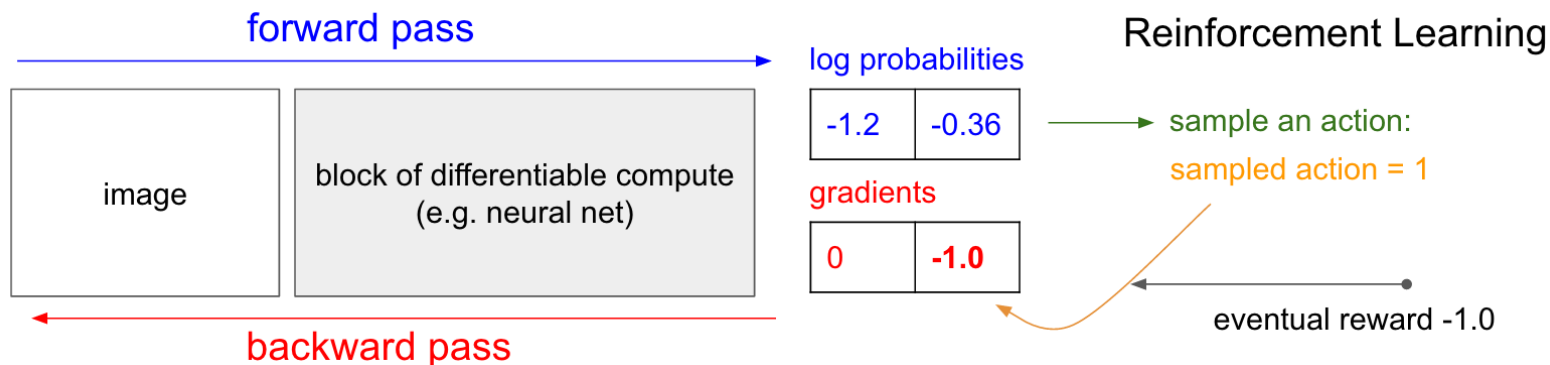
- The similarity between two objects is measured by the **similarity function**
 - The distance between those two object is measured.
 - Shorter the distance higher the similarity, conversely longer the distance higher the dissimilarity.

Reinforcement learning

- Reinforcement learning is a type of dynamic programming that trains algorithms using a system of reward and punishment.
- A reinforcement learning algorithm, or agent, learns by interacting with its environment without intervention from a human by maximizing its reward and minimizing its penalty.
 - The agent receives rewards by performing correctly and penalties for performing incorrectly.
- Reinforcement learning contrasts with other machine learning approaches in that the algorithm is not explicitly told how to perform a task, but works through the problem on its own.

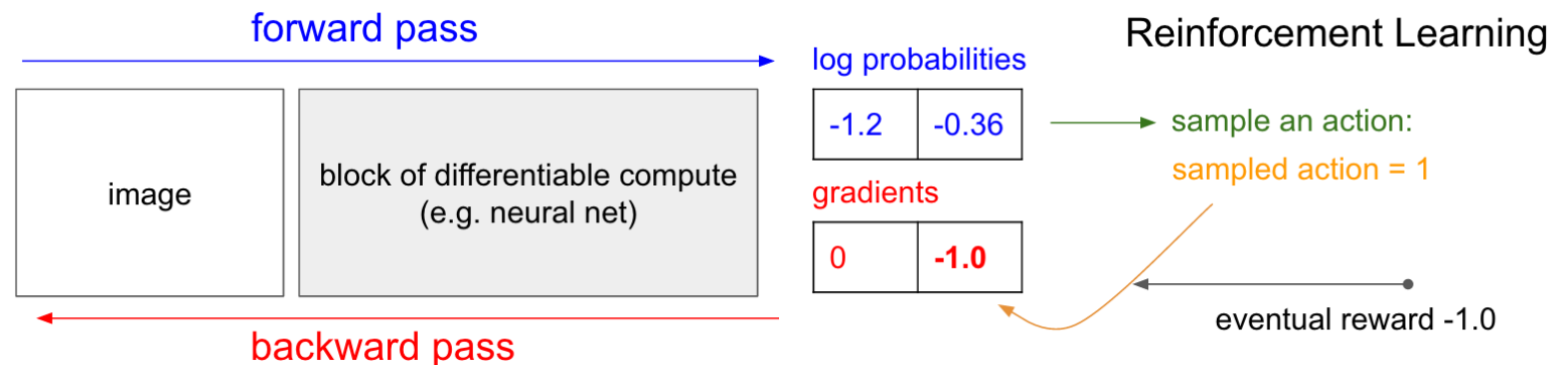
Reinforcement learning

- As an agent, which could be a self-driving car or a program playing chess, interacts with its environment, receives a **reward state** depending on how it performs, such as driving to destination safely or winning a game.
- Conversely, the agent receives a **penalty** for performing incorrectly, such as going off the road or being checkmated.
 - The agent over time makes decisions to **maximize its reward and minimize its penalty** using dynamic programming.
- The advantage of this approach to artificial intelligence is that it allows an AI program to learn without a programmer spelling out how an agent should perform the task.



Reinforcement learning

- The agent is supposed to find the **best possible path to reach the reward**.
- The goal of the robot is to get the reward that is the diamond and avoid the hurdles that is fire.
- The robot learns by trying all the possible paths and then choosing the path which gives him the reward with the least hurdles.
- **Each right step will give the robot a reward and each wrong step will subtract the reward of the robot.**
- The total reward will be calculated when it reaches the final reward that is the diamond.



The clustering Problem:

- Given an integer k and a set of n data points in \mathbb{R}^d , the goal is to choose k centers so as to minimize φ , the total squared distance between each point and its closest center.
- Solving this problem exactly is NP-hard, but Lloyd proposed a local search solution to this problem.
- k-means is the most popular clustering algorithm used in scientific and industrial applications” [3]

k-means Clustering

- K-means clustering is a **type of unsupervised learning**, which is used when you have unlabeled data (i.e., data without defined categories or groups).
- The goal of this algorithm is to find groups in the data, with the number of groups represented by the variable **k**.
- The algorithm works iteratively to assign each data point to one of **k** groups based on the features that are provided.
- Data points are clustered based on feature similarity.

The results of the **k**-means clustering algorithm are:

- The centroids of the k clusters, which can be used to label new data
- Labels for the training data (each data point is assigned to a single cluster)

k-means Clustering

The k-means algorithm is a simple and fast algorithm for this problem, although it offers no approximation guarantees at all.

1. Arbitrarily choose an initial k centers $C = \{c_1, c_2, \dots, c_k\}$.
2. For each $i \in \{1, \dots, k\}$, set the cluster C_i to be the set of points in X that are closer to c_i than they are to c_j for all $j \neq i$.
3. For each $j \in \{1, \dots, k\}$, set c_j to be the center of mass of all points in C_j : $c_j = \frac{1}{|C_j|} \sum_{x \in C_j} x$.
4. Repeat Steps 2 and 3 until C no longer changes.

$$\text{e.g. } C_j = (170 + 168) / 2 \mid = (60 + 56) / 2$$

It is standard practice to choose the initial centers uniformly at random from X .

For Step 2, ties may be broken arbitrarily, as long as the method is consistent.

The idea here is that Steps 2 and 3 are both guaranteed to decrease ϕ , so the algorithm makes local improvements to an arbitrary clustering until it is no longer possible to do so.

Business Uses

- The *K*-means clustering algorithm is used to find groups which have not been explicitly labeled in the data.
- This can be used to confirm business assumptions about **what types of groups exist or to identify unknown groups in complex data sets.**

Some examples of use cases are:

1. Behavioral segmentation:

Segment by purchase history

Segment by activities on application, website, or platform

Define personas based on interests

Create profiles based on activity monitoring

2. Inventory categorization:

- Group inventory by sales activity
- Group inventory by manufacturing metrics

3. Sorting sensor measurements:

- Detect activity types in motion sensors
- Group images
- Separate audio
- Identify groups in health monitoring

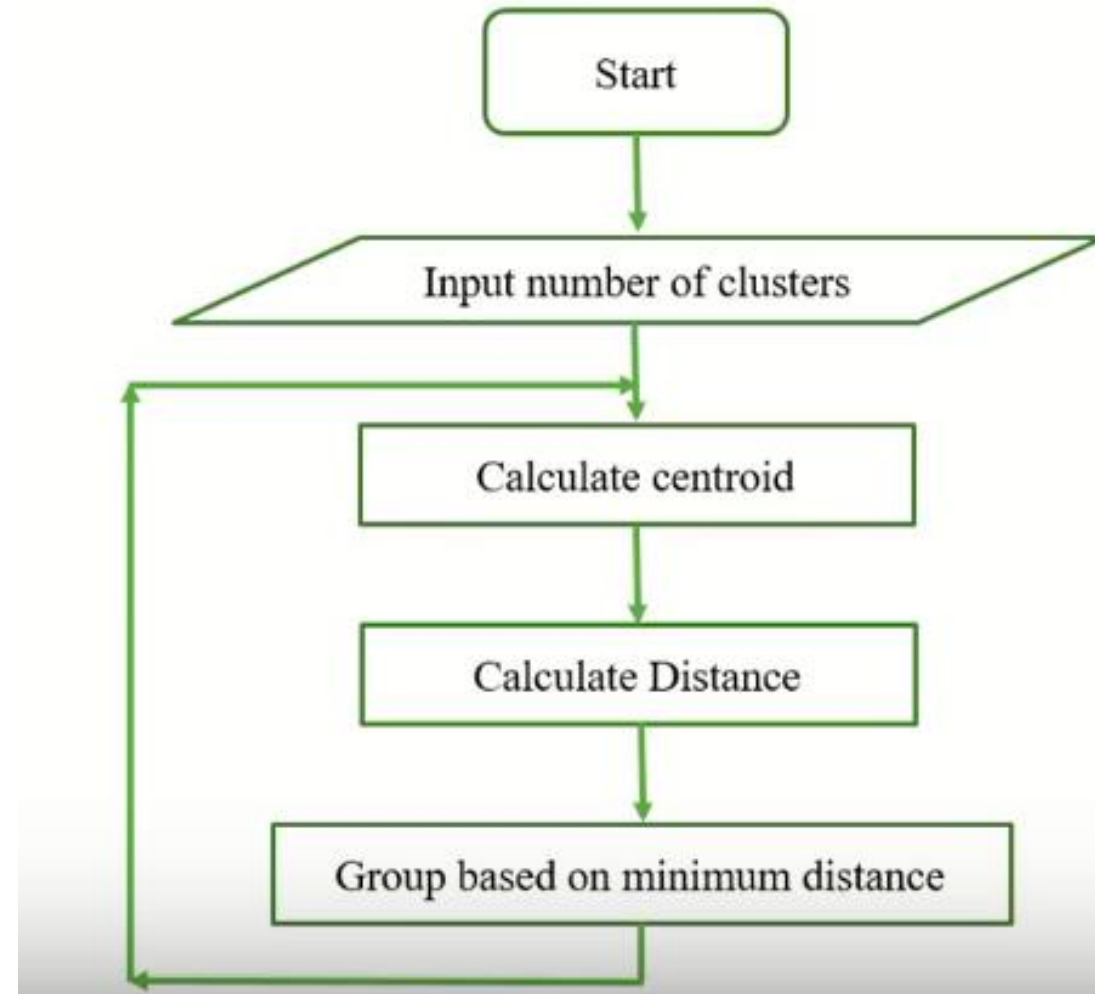
4. Detecting bots or anomalies:

- Separate valid activity groups from bots
- Group valid activity to clean up outlier detection

In addition, monitoring if a tracked data point switches between groups over time can be used **to detect meaningful changes in the data**.

K-means Clustering

- **D**etermines the centroid using the Euclidean method for distance calculation.
- **G**roups the objects based on minimum distance.



It analyze and explore the whole dataset.

K-means Clustering

Apply K-Mean Clustering for the following data sets for two clusters. Tabulate all the assignments.

Sample No	X	Y
1	185	72
2	170	56
3	168	60
4	179	68
5	182	72
6	188	77

k-means Clustering

- Given $k = 2$

Initial Centroid		
Cluster	X	Y
k1	185	72
k2	170	56

- Calculate Euclidean distance using the given equation.

$$\text{Distance } [(x,y), (a,b)] = \sqrt{(x - a)^2 + (y - b)^2}$$

$$\text{Cluster 1 (185,72)} = \sqrt{(185 - 185)^2 + (72 - 72)^2} = 0$$

$$\begin{aligned}\text{Distance from Cluster 2} &= \sqrt{(170 - 185)^2 + (56 - 72)^2} \\ (170,56) &= \sqrt{(-15)^2 + (-16)^2} \\ &= \sqrt{255 + 256} \\ &= \sqrt{481} \\ &= 21.93\end{aligned}$$

$$\begin{aligned}\text{Distance from Cluster 1} &= \sqrt{(185 - 170)^2 + (72 - 56)^2} \\ (185,72) &= \sqrt{(15)^2 + (16)^2} \\ &= \sqrt{255 + 256} \\ &= \sqrt{481} \\ &= 21.93\end{aligned}$$

$$\text{Cluster 2 (170,56)} = \sqrt{(170 - 170)^2 + (56 - 56)^2} = 0$$

Cluster	Centroid		
	X	Y	ASSIGNMENT
k1	0	21.93	1
k2	21.93	0	2

Initial Centroid

Cluster	X	Y
k1	185	72
k2	170	56

- Calculate Euclidean distance for the next dataset (168,60)

$$\text{Distance} [(x,y), (a,b)] = \sqrt{(x-a)^2 + (y-b)^2}$$

$$\begin{aligned} \text{Distance from Cluster 1} &= \sqrt{(168-185)^2 + (60-72)^2} \\ (185,72) &= \sqrt{(-17)^2 + (-12)^2} \\ &= \sqrt{289 + 144} \\ &= \sqrt{433} \\ &= 20.808 \end{aligned}$$

$$\begin{aligned} \text{Distance from Cluster 2} &= \sqrt{(168-170)^2 + (60-56)^2} \\ (170,56) &= \sqrt{(-2)^2 + (4)^2} \\ &= \sqrt{4 + 16} \\ &= \sqrt{20} \\ &= 4.472 \end{aligned}$$

k-means Clustering

Dataset	Euclidean Distance		
	Cluster 1	Cluster 2	ASSIGNMENT
(168,60)	20.808	4.472	2

- Update the cluster centroid.

Cluster	X	Y
k1	185	72
k2	$= (170 + 168) / 2$ $= 169$	$= (60 + 56) / 2$ $= 58$

- Calculate Euclidean distance for the next dataset (179,68)

$$\text{Distance } [(x,y), (a,b)] = \sqrt{(x-a)^2 + (y-b)^2}$$

$$\begin{aligned} \text{Distance from Cluster 1} &= \sqrt{(179-185)^2 + (68-72)^2} \\ (185,72) &= \sqrt{(-6)^2 + (-4)^2} \\ &= \sqrt{36 + 16} \\ &= \sqrt{52} \\ &= 7.211103 \end{aligned}$$

- Calculate Euclidean distance for the next dataset (179,68)

$$\text{Distance } [(x,y), (a,b)] = \sqrt{(x-a)^2 + (y-b)^2}$$

$$\begin{aligned} \text{Distance from Cluster 2} &= \sqrt{(179-169)^2 + (68-58)^2} \\ (169,58) &= \sqrt{(10)^2 + (10)^2} \\ &= \sqrt{100 + 100} \\ &= \sqrt{200} \\ &= 14.14214 \end{aligned}$$

The k-means Clustering

Dataset	Euclidean Distance		
	Cluster 1	Cluster 2	ASSIGNMENT
(179,68)	7.211103	14.14214	1

- Update the cluster centroid.

Cluster	X	Y
k1	$= 185+179/2$ =182	$= 72+68/2$ =70
k2	169	58

- Calculate Euclidean distance for the next dataset (182,72)

$$\text{Distance [(x,y), (a,b)]} = \sqrt{(x - a)^2 + (x - b)^2}$$

$$\begin{aligned} \text{Distance from Cluster 1} &= \sqrt{(182 - 182)^2 + (72 - 70)^2} \\ (182,70) &= \sqrt{(0)^2 + (2)^2} \\ &= \sqrt{0 + 4} \\ &= \sqrt{4} \\ &= 2 \end{aligned}$$

- Calculate Euclidean distance for the next dataset (182,72)

$$\text{Distance [(x,y), (a,b)]} = \sqrt{(x - a)^2 + (x - b)^2}$$

$$\begin{aligned} \text{Distance from Cluster 2} &= \sqrt{(182 - 169)^2 + (72 - 58)^2} \\ (169,58) &= \sqrt{(13)^2 + (14)^2} \\ &= \sqrt{169 + 196} \\ &= \sqrt{365} \\ &= 19.10 \end{aligned}$$



The k-means Clustering

Dataset	Euclidean Distance		
	Cluster 1	Cluster 2	ASSIGNMENT
(182,72)	2	19.10	1

- Update the cluster centroid.

Cluster	X	Y
k1	= 182+182/2 =182	= 70+72/2 = 71
k2	169	58

The k-means Clustering

■ Final Assignment

Dataset No	X	Y	Assignment
1	185	72	1
2	170	56	2
3	168	60	2
4	179	68	1
5	182	72	1
6	188	77	1

The k-means++ algorithm

We propose a specific way of choosing centers for the k-means algorithm. In particular, let $D(x)$ denote the shortest distance from a data point to the closest center we have already chosen. Then, we define the following algorithm, which we call k-means++.

- 1a. Take one center c_1 , chosen uniformly at random from \mathcal{X} .
- 1b. Take a new center c_i , choosing $x \in \mathcal{X}$ with probability $\frac{D(x)^2}{\sum_{x \in \mathcal{X}} D(x)^2}$.
- 1c. Repeat Step 1b. until we have taken k centers altogether.
- 2-4. Proceed as with the standard k-means algorithm.

We call the weighting used in Step 1b simply “ D^2 weighting”.

The k-means ++ Algorithm

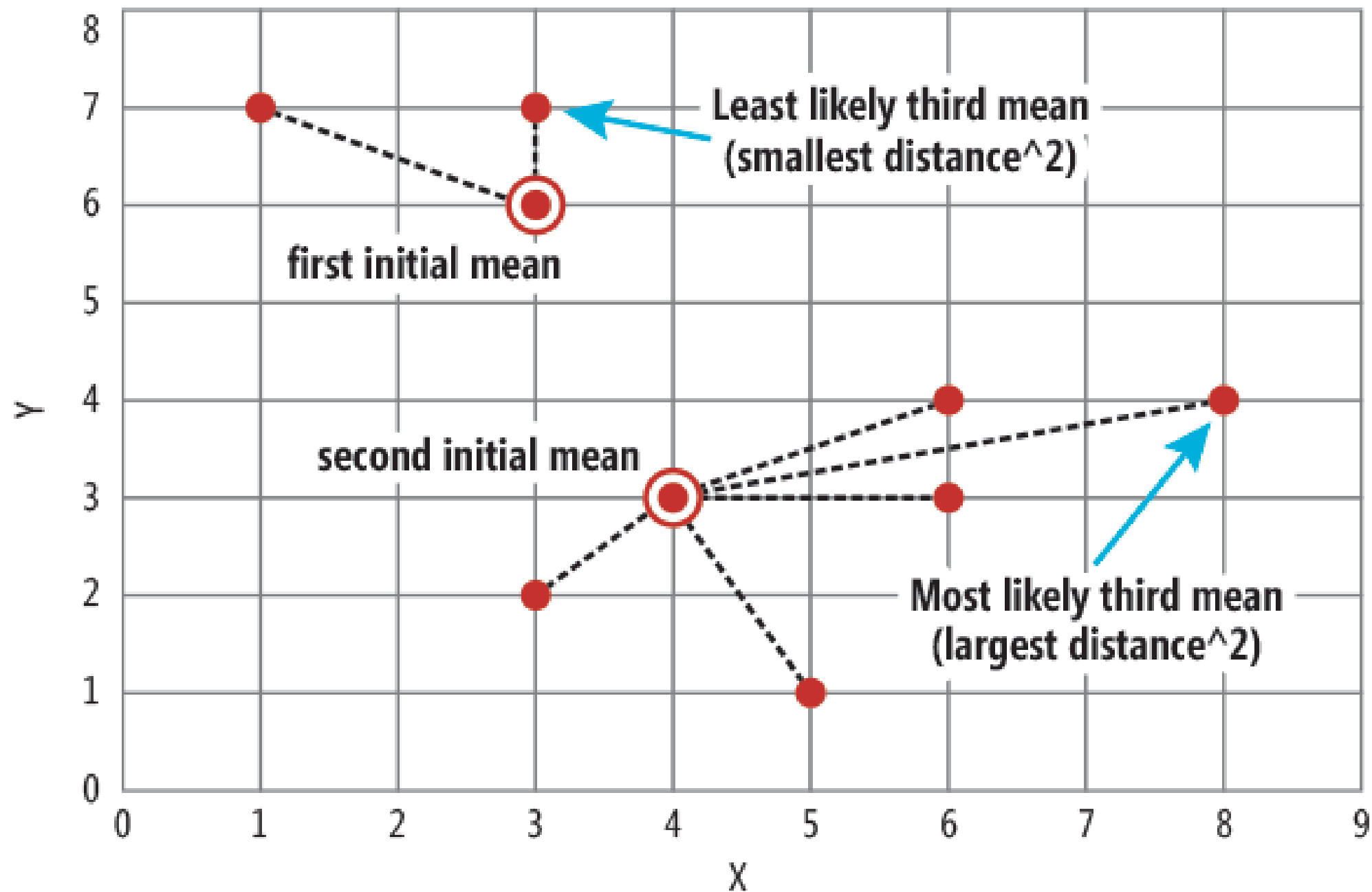
- The first step is to choose a data point at random. Call this point s_1 . Next, compute the squared distances

$$D_i^2 = ||Y_i - s_1||^2.$$

- Now choose a second point s_2 from the data. The probability of choosing Y_i is $D_i^2 / \sum_j D_j^2$
- Now recompute the distance as $D_i^2 = \min \left\{ ||Y_i - s_1||^2, ||Y_i - s_2||^2 \right\}$.
- Now choose a third point s_3 from the data where the probability of choosing Y_i is $D_i^2 / \sum_j D_j^2$.
- We continue until we have k points $s_1, s_2, s_3, \dots, s_k$.
- Finally, we run k-means clustering using $s_1, s_2, s_3, \dots, s_k$ as starting values. Call the resulting centers $c_1, c_2, c_3, \dots, c_k$.
- Arthur and Vassilvitskii proved that the expected value is over the randomness in the algorithm

$$\mathbb{E}[R(\hat{c}_1, \dots, \hat{c}_k)] \leq 8(\log k + 2) \min_{c_1, \dots, c_k} R(c_1, \dots, c_k).$$

Nine Data Items in Two-Dimension into Three Clusters



The k-means ++ Algorithm

- The first initial mean at (3, 6) was randomly selected.
- Then the **distance-squared from each of the other 8 data items to the first mean was computed**, and using that information, the second initial mean at (4, 3) was selected.
- To select a data item as the third initial mean, the squared distance from each data point to its closest mean is computed.
- The distances are shown as dashed lines.
- Using these squared distance values, the third mean will be selected so that data items with small squared distance values have a low probability of being selected, and data items with large squared distance values have a high probability of being selected.
- This technique is sometimes called proportional fitness selection.

The k-means ++ Algorithm: Proportional fitness selection using Roulette wheel selection

- Proportional fitness selection is the heart of the k-means++ initialization mechanism.
- There are several ways to implement proportional fitness selection.
- Here **we use Roulette wheel selection for proportional fitness selection.**
- Suppose there are four candidate items (0, 1, 2, 3) with associated values (20.0, 10.0, 40.0, 30.0).
- The sum of the values is $20.0 + 40.0 + 10.0 + 30.0 = 100.0$.
- Proportional fitness selection will pick item 0 with probability $20.0/100.0 = 0.20$; pick item 1 with probability $10.0/100.0 = 0.10$; pick item 2 with probability $40.0/100.0 = 0.40$; and pick item 3 with probability $30.0/100.0 = 0.30$.

The k-means ++ Algorithm: Proportional fitness selection using Roulette wheel selection

- If the probabilities of selection are stored in an array as (0.20, 0.10, 0.40, 0.30), the cumulative probabilities can be stored in an array with values (0.20, 0.30, 0.70, 1.00).
- Now, suppose a random p is generated with value 0.83.
- If i is an array index into the cumulative probabilities array, when $i = 0$, $\text{cum}[i] = 0.20$, which isn't greater than $p = 0.83$, so i increments to 1.
- Now $\text{cum}[i] = 0.30$, which is still not greater than p , so i increments to 2.
- Now $\text{cum}[i] = 0.70$, which is still not greater than p , so i increments to 3.
- Now $\text{cum}[i] = 1.00$, which is greater than p , so $i = 3$ is returned as the selected item.

K Nearest Neighbors - Classification

- K nearest neighbors algorithm stores all available cases and classifies new cases based on a similarity measure (e.g., distance functions).
- KNN has been used in [statistical estimation and pattern recognition](#) already in the beginning of 1970's as a non-parametric technique.
- A case is classified by a majority vote of its neighbors, with the case being assigned to the class most common amongst its k nearest neighbors measured by a distance function.
- If $k = 1$, then the case is simply assigned to the class of its nearest neighbor.

Distance functions

Euclidean

$$\sqrt{\sum_{i=1}^k (x_i - y_i)^2}$$

Manhattan

$$\sum_{i=1}^k |x_i - y_i|$$

Minkowski

$$\left(\sum_{i=1}^k (|x_i - y_i|)^q \right)^{1/q}$$

All three distance measures are only valid for continuous variables

K Nearest Neighbors - Classification

- In the instance of categorical variables the Hamming distance must be used.
- It also brings up the issue of standardization of the numerical variables between 0 and 1 when there is a mixture of numerical and categorical variables in the dataset.

Hamming Distance

$$D_H = \sum_{i=1}^k |x_i - y_i|$$

$$x = y \Rightarrow D = 0$$

$$x \neq y \Rightarrow D = 1$$

X	Y	Distance
Male	Male	0
Male	Female	1

- Choosing the optimal value for K is best done by first inspecting the data.
- In general, a large K value is more precise as it reduces the overall noise but there is no guarantee.
- Cross-validation is another way to retrospectively determine a good K value by using an independent dataset to validate the K value.
- Historically, the optimal K for most datasets has been between 3-10. That produces much better results than 1NN.

k Nearest Neighbors – Classification

Example:

- k-NN is a non-parametric method used for classification
- Prediction for the test data is done on the basis of its neighbor
- k is an integer(small), if k=1, k is assigned to the class of single nearest neighbor

Name	Acid Durability	Strength	Class
Type-1	7	7	Bad
Type-2	7	4	Bad
Type-3	3	4	Good
Type-4	1	4	Good
Assume the Test Data is: Acid Durability=3, and Strength=7. What is the class?			

k Nearest Neighbors – Classification

The similarity is calculated using distance measure like Euclidean

$$d(\mathbf{p}, \mathbf{q}) = d(\mathbf{q}, \mathbf{p}) = \sqrt{(q_1 - p_1)^2 + (q_2 - p_2)^2 + \cdots + (q_n - p_n)^2}$$
$$= \sqrt{\sum_{i=1}^n (q_i - p_i)^2}.$$

Name	Acid Durability	Strength	Class	Distance
Type-1	7	7	Bad	Sqrt((7-3) ² + (7-3) ²)=4
Type-2	7	4	Bad	5
Type-3	3	4	Good	3
Type-4	1	4	Good	3.6

k Nearest Neighbors – Classification

Rank these Attributes

Name	Acid Durability	Strength	Class	Distance	Rank
Type-1	7	7	Bad	4	3
Type-2	7	4	Bad	5	4
Type-3	3	4	Good	3	1
Type-4	1	4	Good	3.6	2

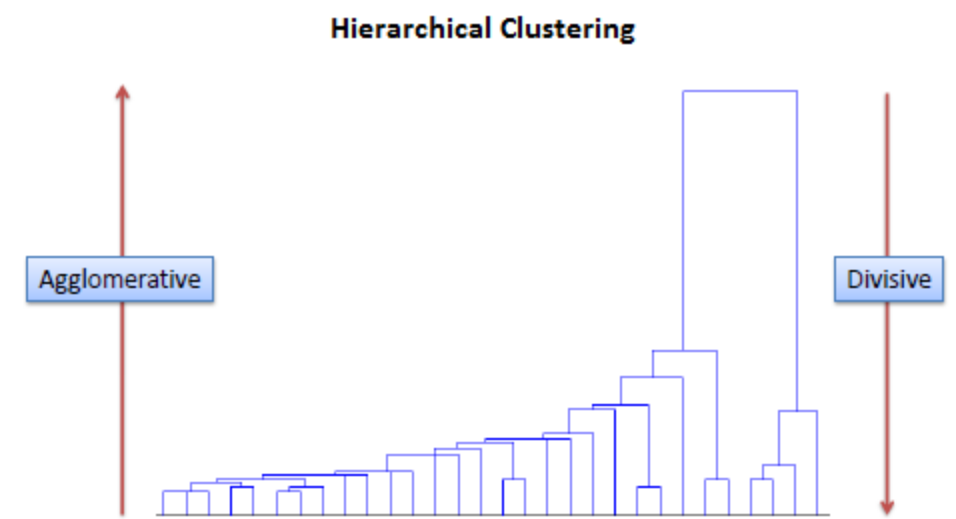
k Nearest Neighbors - Classification

k= 1

Name	Acid Durability	Strength	Class	Distance	Rank
Type-1	7	7	Bad	4	3
Type-2	7	4	Bad	5	4
Type-3	3	4	Good	3	1
Type-4	1	4	Good	3.6	2

Acceptance level is good in the two neighbors

Hierarchical clustering



Hierarchical clustering involves creating clusters that have a predetermined ordering from top to bottom.

Divisive method

- In divisive or top-down clustering method we assign all of the observations to a single cluster and then partition the cluster to two least similar clusters.
- Finally, we proceed recursively on each cluster until there is one cluster for each observation.
- There is evidence that divisive algorithms produce more accurate hierarchies than agglomerative algorithms in some circumstances but is conceptually more complex.

Hierarchical clustering

Agglomerative method

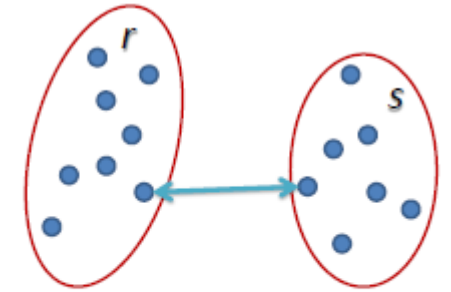
- In agglomerative or bottom-up clustering method we assign each observation to its own cluster.
- Then, compute the similarity (e.g., distance) between each of the clusters and join the two most similar clusters, until there is only a single cluster left.

Hierarchical clustering

	BA	FI	MI	NA	RM	TO
BA	0	662	877	255	412	996
FI	662	0	295	468	268	400
MI	877	295	0	754	564	138
NA	255	468	754	0	219	869
RM	412	268	564	219	0	669
TO	996	400	138	869	669	0

Before any clustering is performed, it is required to determine the proximity matrix containing the distance between each point using a distance function.

Then, the matrix is updated to display the distance between each cluster. The following three methods differ in **how the distance between each cluster is measured**.



$$L(r, s) = \min(D(x_{ri}, x_{sj}))$$

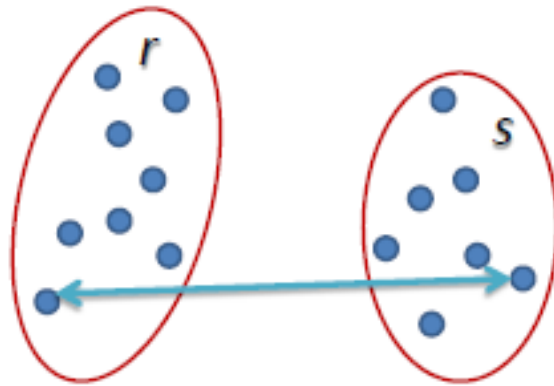
Single Linkage

- In single linkage hierarchical clustering, the distance between two clusters is defined as the shortest distance between two points in each cluster.
- For example, the distance between clusters “r” and “s” to the left is equal to the length of the arrow between their two closest points.

Hierarchical clustering

Complete Linkage

- In complete linkage hierarchical clustering, the distance between two clusters is defined as the longest distance between two points in each cluster.
- For example, the distance between clusters “r” and “s” to the left is equal to the length of the arrow between their two furthest points.

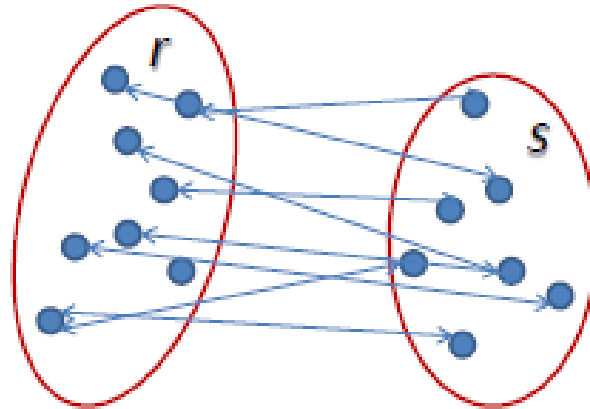


$$L(r, s) = \max(D(x_{ri}, x_{sj}))$$

Hierarchical clustering

Average Linkage

- In average linkage hierarchical clustering, the distance between two clusters is defined as the average distance between each point in one cluster to every point in the other cluster.
- For example, the distance between clusters “r” and “s” to the left is equal to the average length each arrow between connecting the points of one cluster to the other.



$$L(r, s) = \frac{1}{n_r n_s} \sum_{i=1}^{n_r} \sum_{j=1}^{n_s} D(x_{ri}, x_{sj})$$

Hierarchical clustering

- Begin with the disjoint clustering having level $L(0) = 0$ and sequence number $m = 0$.
- Find the least dissimilar pair of clusters in the current clustering, say pair $(r), (s)$, according to
$$d[(r),(s)] = \min d[(i),(j)]$$
 where the minimum is over all pairs of clusters in the current clustering.
- Increment the sequence number : $m = m + 1$. Merge clusters (r) and (s) into a single cluster to form the next clustering m . Set the level of this clustering to
$$L(m) = d[(r),(s)]$$
- Update the proximity matrix, D , by deleting the rows and columns corresponding to clusters (r) and (s) and adding a row and column corresponding to the newly formed cluster. The proximity between the new cluster, denoted (r,s) and old cluster (k) is defined in this way:
$$d[(k), (r,s)] = \min d[(k),(r)], d[(k),(s)]$$
- If all objects are in one cluster, stop. Else, go to step 2.

Hierarchical Clustering

	BA	FI	MI	NA	RM	TO
BA	0	662	877	255	412	996
FI	662	0	295	468	268	400
MI	877	295	0	754	564	138
NA	255	468	754	0	219	869
RM	412	268	564	219	0	669
TO	996	400	138	869	669	0



Hierarchical Clustering

	BA	FI	MI/TO	NA	RM
BA	0	662	877	255	412
FI	662	0	295	468	268
MI/TO	877	295	0	754	564
NA	255	468	754	0	219
RM	412	268	564	219	0



Hierarchical Clustering

	BA	FI	MI/TO	NA/RM
BA	0	662	877	255
FI	662	0	295	268
MI/TO	877	295	0	564
NA/RM	255	268	564	0



Hierarchical Clustering

	BA/NA/RM	FI	MI/TO
BA/NA/RM	0	268	564
FI	268	0	295
MI/TO	564	295	0



$\min d(i,j) = d(\text{BA/NA/RM}, \text{FI}) = 268 \Rightarrow$ merge BA/NA/RM and FI into a new cluster called BA/FI/NA/RM

$L(\text{BA/FI/NA/RM}) = 268$

$m = 4$

Hierarchical Clustering

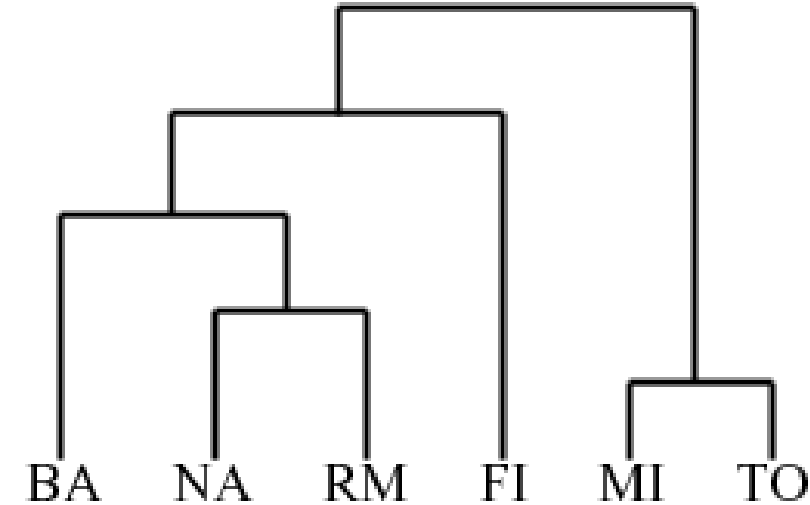
	BA/FI/NA/RM	MI/TO
BA/FI/NA/RM	0	295
MI/TO	295	0



Hierarchical Clustering

Finally, we merge the last two clusters at level 295.

The process is summarized by the following hierarchical tree:



Problems with the Hierarchical Clustering:

The main weaknesses of agglomerative clustering methods are:

- they do not scale well: time complexity of at least $O(n^2)$, where n is the number of total objects;
- they can never undo what was done previously.

Naive Bayes Classifier

What is a classifier?

- A classifier is a machine learning model that is used to discriminate different objects based on certain features.

Principle of Naive Bayes Classifier:

- A Naive Bayes classifier is a probabilistic machine learning model that's used for classification task. The classifier is based on the Bayes theorem.
- We can find the probability of **A** happening, given that **B** has occurred.
- Here, **B** is the evidence and **A** is the hypothesis.
- The assumption made here is that the predictors/features are independent. That is presence of one particular feature does not affect the other. Hence it is called naive.

$$P(A|B) = \frac{P(B|A)P(A)}{P(B)}$$

Consider the problem of playing golf. The dataset is represented as below.

We classify whether the day is suitable for playing golf, given the features of the day. According to this example, Bayes theorem can be written as:

$$P(y|X) = \frac{P(X|y)P(y)}{P(X)}$$

The variable **y** is the class variable(play golf), which represents if it is suitable to play golf or not given the conditions. Variable **X** represent the parameters/features. **X** is given as.

$$X = (x_1, x_2, x_3,, x_n)$$

$$P(y|x_1, ..., x_n) = \frac{P(x_1|y)P(x_2|y)...P(x_n|y)P(y)}{P(x_1)P(x_2)...P(x_n)}$$

	OUTLOOK	TEMPERATURE	HUMIDITY	WINDY	PLAY GOLF
0	Rainy	Hot	High	False	No
1	Rainy	Hot	High	True	No
2	Overcast	Hot	High	False	Yes
3	Sunny	Mild	High	False	Yes
4	Sunny	Cool	Normal	False	Yes
5	Sunny	Cool	Normal	True	No
6	Overcast	Cool	Normal	True	Yes
7	Rainy	Mild	High	False	No
8	Rainy	Cool	Normal	False	Yes
9	Sunny	Mild	Normal	False	Yes
10	Rainy	Mild	Normal	True	Yes
11	Overcast	Mild	High	True	Yes
12	Overcast	Hot	Normal	False	Yes
13	Sunny	Mild	High	True	No

Question on NAIVE BAYE'S Algorithm:

Ques:1) For the given dataset, Apply Naive-Bayes Algorithm and Predict the outcome for a Car = { Red, Domestic, SUV }

Color	Type	Origin	Stolen
Red	Sports	Domestic	Yes
Red	Sports	Domestic	NO
Red	Sports	Domestic	Yes
Yellow	Sports	Domestic	NO
Yellow	Sports	Imported	Yes
Yellow	SUV	Imported	NO
Yellow	SUV	Imported	Yes
Yellow	SUV	Domestic	NO
Red	SUV	Imported	NO
Red	Sports	Imported	Yes

Posterior $P(A|B) = \frac{P(B|A) \cdot P(A)}{P(B)}$

Prob. of B when A is True (Likelihood) → Proposition
 Prob. (A) when B is true → evidence

NO ANS..

$X = [\text{Red, Domestic, SUV}] = P(X|Yes) = \frac{3}{5} \cdot \frac{2}{5} \cdot \frac{1}{5} = \frac{6}{125} = 0.024$ Yes

ii) $P(\text{Red}|Yes) = \frac{P(Yes|\text{Red}) \cdot P(\text{Red})}{P(Yes)} = \frac{\frac{3}{5} \cdot \frac{5}{10}}{\frac{5}{10}} = \frac{3}{5}$

iii) $P(\text{Domestic}|Yes) = \frac{2}{5}$ iii) $P(\text{SUV}|Yes) = \frac{1}{5}$

$P(\text{Red}|No) = \frac{P(No|\text{Red}) \cdot P(\text{Red})}{P(No)} = \frac{\frac{2}{5} \cdot \frac{5}{10}}{\frac{5}{10}} = \frac{2}{5}$

$P(\text{Domestic}|No) = \frac{3}{5}$, $P(\text{SUV}|No) = \frac{2}{5}$ ^{5/10} = $\frac{2}{5} \times \frac{3}{5} \times \frac{2}{5} = 0.072$ **NO**

Sample Dataset

500.0	4.0	1.8	15.6	349.99
250.0	4.0	2.0	13.3	80.0
80.0	4.0	1.66		144.0
160.0	1.0	1.6	10.0	31.0
80.0	4.0	1.8		129.95
80.0	4.0	1.66		144.0
80.0	4.0			136.96
500.0	4.0	1.7	15.6	255.0
			11.6	60.0
				100.0
80.0	2.0	1.86	14.1	118.08
80.0	4.0			136.96
1024.0	12.0	1.7	15.6	529.99
160.0	1.0	1.6	10.0	31.0
1024.0	4.0	2.0	15.6	249.99
80.0	4.0	1.8		129.95
500.0	4.0	2.53	15.6	102.5
			11.6	60.0
1024.0	8.0	2.0	15.6	469.99
			10.6	40.0
1024.0	4.0	2.0	15.6	249.99
500.0	4.0	2.53	15.6	102.5
				100.0