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Finishing School

# Getting started with Machine Learning



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# Getting started with Anaconda

<https://www.anaconda.com/products/individual>

Windows 

Python 3.8

☒ 64-Bit Graphical Installer (457 MB)

32-Bit Graphical Installer (403 MB)



# Machine Learning

- ✓ **What is Machine Learning**
  - ✓ **How does it contrast with other means of problem solving**
- ✓ **When to use Machine Learning**
- ✓ **How to use Machine Learning (Classification Problems)**



# Machine Learning

## Working with Real life problem – Spam Detection

- Classify emails as Spam or Ham
- Figure this out before the email is sent to the inbox



## Spam detection – Rule based approach

- **Define set of rules to identify spam**
- **Rules can be intuitive or logical**
- **Requires domain knowledge**
- **Needs access to historical data**



## Spam detection – Rule based approach (Technique)

- **Blacklist**
  - **Emails from specific IP address**
  - **Specific words and Phrases**
- **Whitelist**
  - **Emails form contacts of contacts of contacts**
  - **Certain domains**



## Spam detection – Rule based approach

From: Spammer@donottrust.com

Subject: You won a lottery !

Nirionline,

Congratulations! You won a 10 million lottery, we need the following details to credit the same to your account.



## Spam detection – Rule based approach

From: Spammer@donottrust.net  
Subject: look who won a lucky draw !

Niranjan,

You have won a lucky draw, we need the following details  
to credit the same to your account.





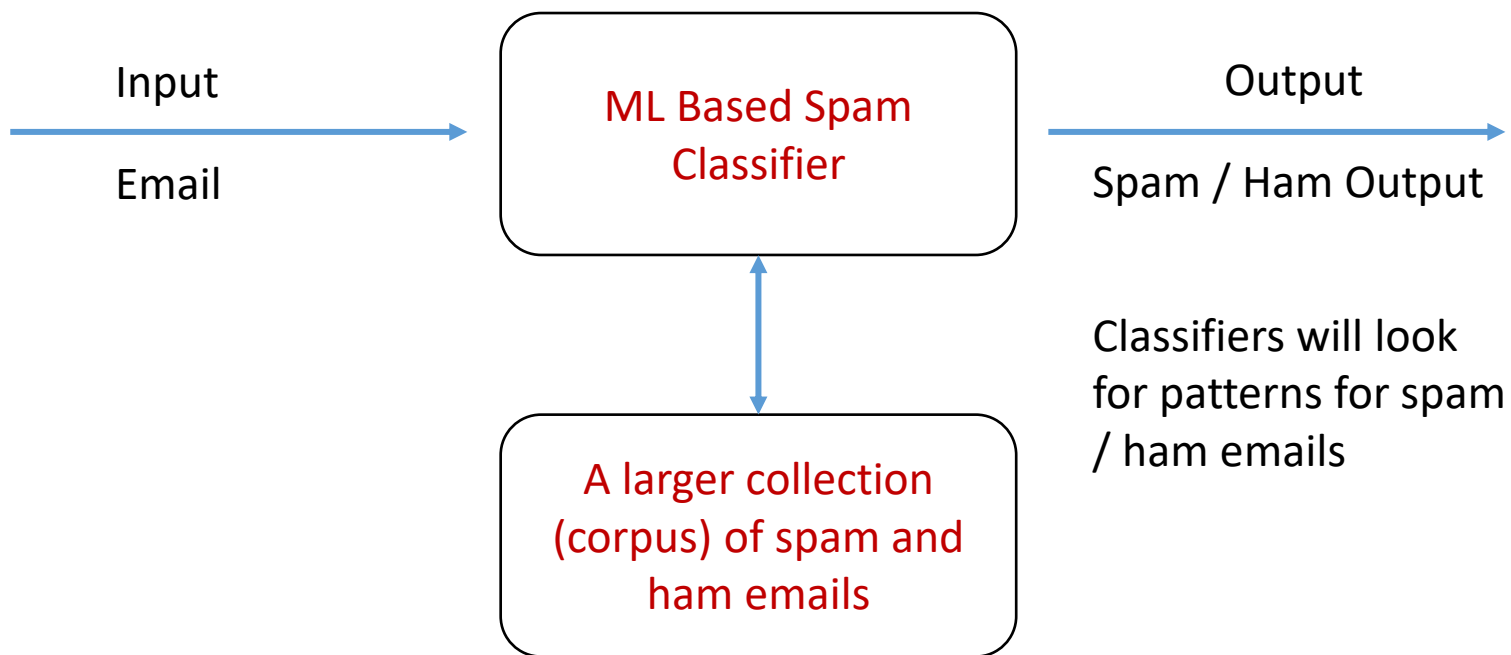
## Spam detection – Rule based approach (limitations)

- Rules get outdated
- Rules are static – Spammers are dynamic
- Hard to maintain many rules
- Rethinking rules is time consuming
- Easier for spammer to get around your rules than the time taken for you to build the rules.



# Machine Learning

## Spam detection – ML based approach





## Spam detection – ML based approach (Advantages)

- **Algorithm can vary approach based on data**
- **Corpus can be updated based on user feedback**
- **Can detect patterns not visible to humans**
- **May be less complex than rule based - approach**



# Machine Learning

## What is machine learning ?

**It's the ability of a program to read, understand and learn from the data and there by draw inferences about the data.**



## When to use machine learning

- **The alternative is a complex set of hard to maintain rules**
- **Dynamic environments with changing data**
- **Rule based solutions can not solve the problem**
- **To get insights on to your data**



# Machine Learning

## ML – based Vs Rule based

### ML Based

**Dynamic**  
**Experts optional**  
**Need corpus**  
**Training required**

### Rule Based

**Static**  
**Experts required**  
**Corpus required**  
**No training**



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# Machine Learning

## Types of Machine Learning



# Machine Learning

## Supervised Learning

- Learning under the supervision of a coach / Teacher
- ML algorithm given data along with labels / responses
- Learns from looking at questions and answers





# Machine Learning

## Supervised Learning

- **Classification**
  - **Spam or Ham**
  - **Mammal or fish or reptiles or birds or Amphibian**
- **Regression (predict numeric values)**
  - **Income of shopper**



# Machine Learning

## Supervised Learning



**Cause**



**Effect**

**X Cause Y**



# Machine Learning

## X Variables

- **Attributes which an ML algorithm focuses on are called features**
- **Each data point is a list (or vector) of such features**
- **Input to an ML algorithm is a feature vector**
- **Feature vectors are called x variables**
- **Also called independent variables or predictors**



# Machine Learning

## Y Variables

- **Attributes which an ML algorithm tries to predict are called labels**
- **Labels can be:**
  - **Categorical ( classification)**
  - **Continuous (regression)**
- **Labels are referred to as y variables**
- **Also called dependent variable**



# Machine Learning

## Data Set

**Attributes**

**Features**

**Target  
Variables**

| Sender          | Subject   | IP           | Body                                    | Result |
|-----------------|-----------|--------------|---|--------|
| Xyz@bank.com    | Loan      | 77.44.5.89   | Dear customer your loan application.... | Spam   |
| 123@builder.com | quotation | 101.22.76.81 | Dear Niranjana Your quotation.....      | Ham    |
| lmn@school.com  | reciept   | 301.99.23.45 | Hi sir your Sons school fee.....        | Ham    |

**Feature Vector**

**Labels**



# Machine Learning

## Supervised Learning ( Training Stage)

$$y = f(x)$$

Most machine learning algorithms aim to **learn** the function **f** which links the function **x** to label **y**



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# Machine Learning

## Supervised Learning models

- **Linear Regression**
- **Logistic Regression**



# Machine Learning

## Supervised Learning models

- **Linear Regression**
- **Logistic Regression**





# Machine Learning

## Linear Regression

$$Y = m (x) + C$$

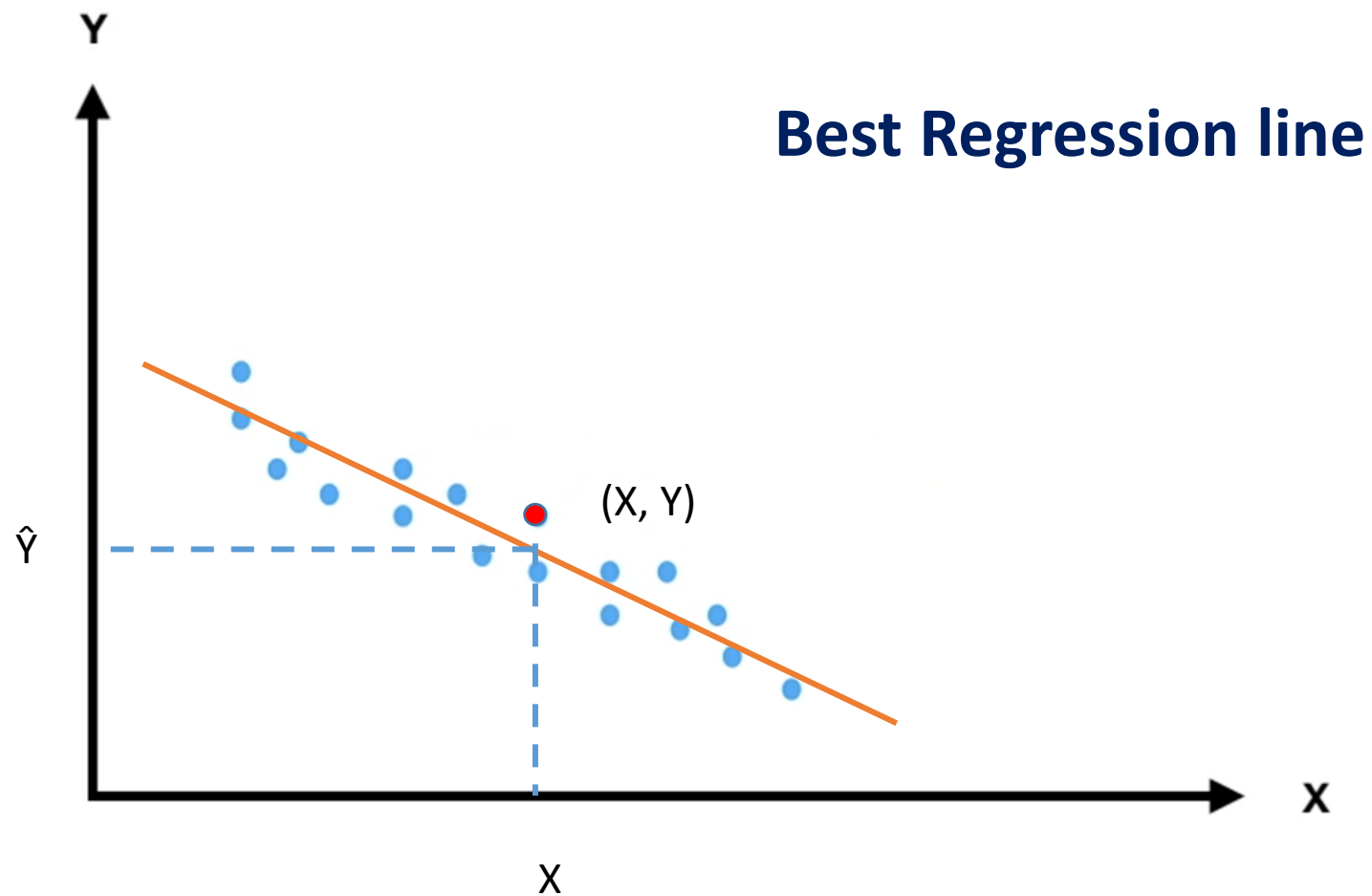
$$f(y) = m (x) + C$$

**Here Linear regression specifies upfront that the relationship between x & y represented by f is linear**



# Machine Learning

## Linear Regression





# Machine Learning

## Best regression line

$$\text{Residual} = e = Y - \hat{Y}$$

$$\text{Data} = (x_1, y_1), (x_2, y_2), (x_3, y_3) \text{ ----- } (x_n, y_n)$$

$$\text{Residual} = (Y_1 - \hat{Y}_1), (Y_2 - \hat{Y}_2), (Y_3 - \hat{Y}_3) \text{ ----- } (Y_n - \hat{Y}_n)$$

To get the best regression line our aim should be to get the minimum error for training data

Eg: least Square method

$$\text{Minise: } (Y_1 - \hat{Y}_1)^2 + (Y_2 - \hat{Y}_2)^2 + (Y_3 - \hat{Y}_3)^2 \text{ ----- } (Y_n - \hat{Y}_n)^2$$



# Machine Learning

## Cavities of regression line

- Can be used only when the errors are normally distributed
- Can also be used with multiple independent variables

# Logistic Regression

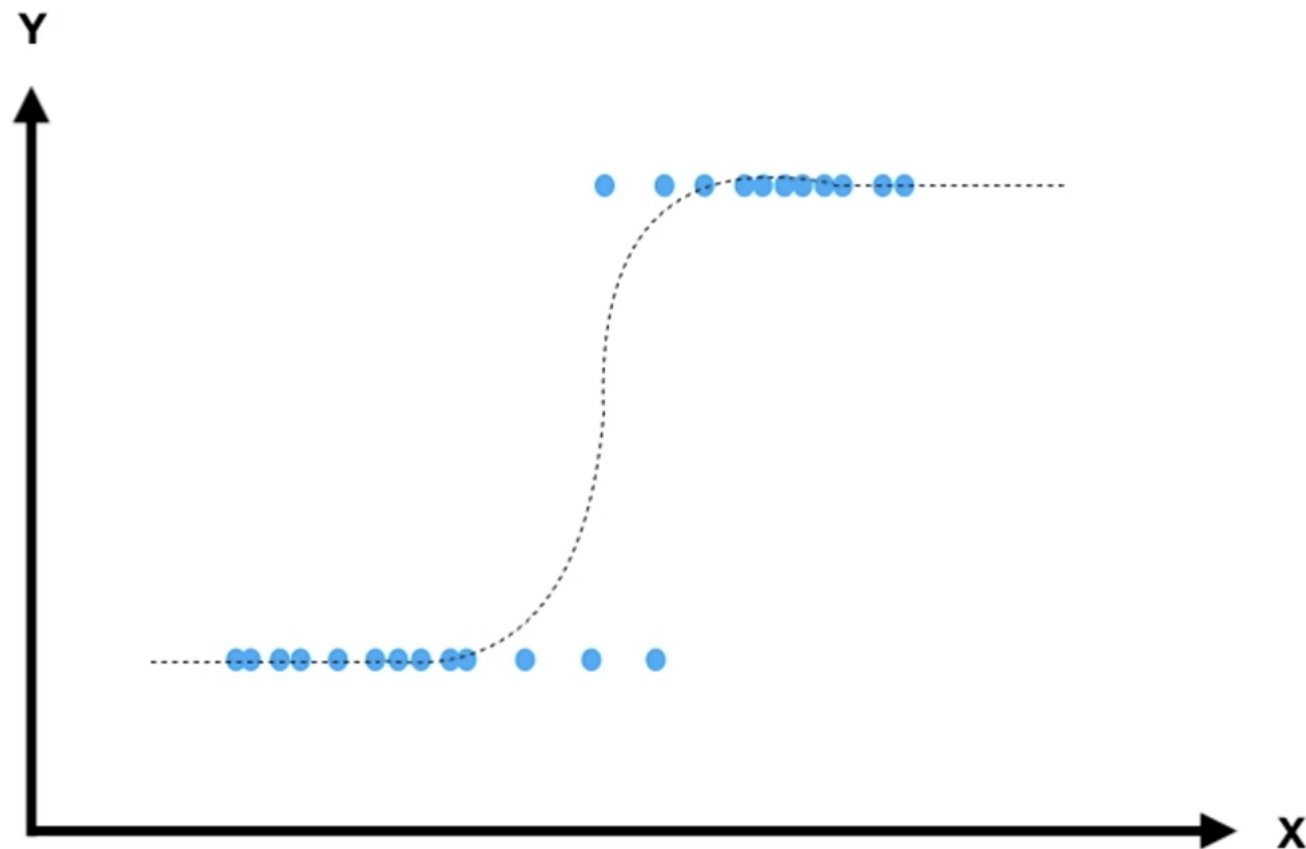
- Used when the dependent variable is categorical (y)
- Multiple Independent variables can be continuous or categorical (x)
- Predict probability of each outcome – assign result to category with highest probability



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# Machine Learning

## Logistic Regression





# Machine Learning

## Logistic Regression

### Objective function: Minimize cross entropy

Cross-entropy is a measure from the field of information theory, building upon entropy and generally **calculating the difference between two probability distributions**. It is closely related to but is different from KL divergence that calculates the relative entropy between two probability distributions, whereas cross-entropy can be thought to calculate the total entropy between the distributions.

Cross-entropy is also related to and often confused with logistic loss, called log loss. Although the two measures are derived from a different source, when used as loss functions for classification models, both measures calculate the same quantity and can be used interchangeably.



# Machine Learning

## Unsupervised Learning

- **There is only input data  $X$  – no output data**
- **Model the underlying structure to learn more about the data**
- **Algorithms self discover patterns of structure in the data**





# Machine Learning

## Unsupervised Learning - algorithms

- Autoencoding – Identify latent factors that drive data (eg: PCA)
- Clustering – Identify patterns in data items

## Looking within

- Be emotionally self sufficient
- Learn what matters (to you)
- Identify others who share them and those who don't
- Eliminate what does not matter
- Train yourself to navigate the outside world



# Machine Learning

## Why Look with in

### In Life

- Be emotionally self sufficient
- Learn what matters (to you)
- Identify others who share them
- and those who don't
- Eliminate what does not matter
- Train yourself to navigate the outside world

### In ML

- Make unlabeled data self sufficient
- Latent factor analysis (fields relevant for a cause)
- Clustering
- Anomaly detection
- Quantisation



# Machine Learning

## Why Look with in

### ML Technique

- Make unlabeled data self sufficient
- Latent factor analysis (fields relevant for a cause)
- Clustering
- Anomaly detection
- Quantisation

### In Life

- Identify photo of specific individual
- Find common drivers for 100 Stocks
- Find relevant document in a corpus
- Flag fraudulent credit card transactions
- Compress 24 bit true color to 8 bit



# Machine Learning

## Why Look with in

### What

- Make unlabeled data self sufficient
- Latent factor analysis (fields relevant for a cause)
- Clustering
- Anomaly detection
- Quantisation

### How

- Autoencoding
- Autoencoding
- Clustering
- Autoencoding
- Clustering



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# Machine Learning

## Support Vector Machines



# Machine Learning

## What is an SVM

- **SVMs are used to build binary classifiers**
- **Make Classification decision on basis of a “Linear function” of point’s co-ordinates**
- **Does not require prior knowledge of probability distribution of the points**
- **Involves an explicit training stage**



# Machine Learning

## What is an SVM

- **SVMs are supervised machine learning approach used to build linear, non probabilistic and binary classifiers**



# Machine Learning

## What is an SVM

The linear SVM classifier works by drawing a straight line between two classes. All the data points that fall on one side of the line will be labeled as one class and all the points that fall on the other side will be labeled as the second. Sounds simple enough, but there's an infinite amount of lines to choose from. How do we know which line will do the best job of classifying the data? This is where the LSVM algorithm comes in to play.

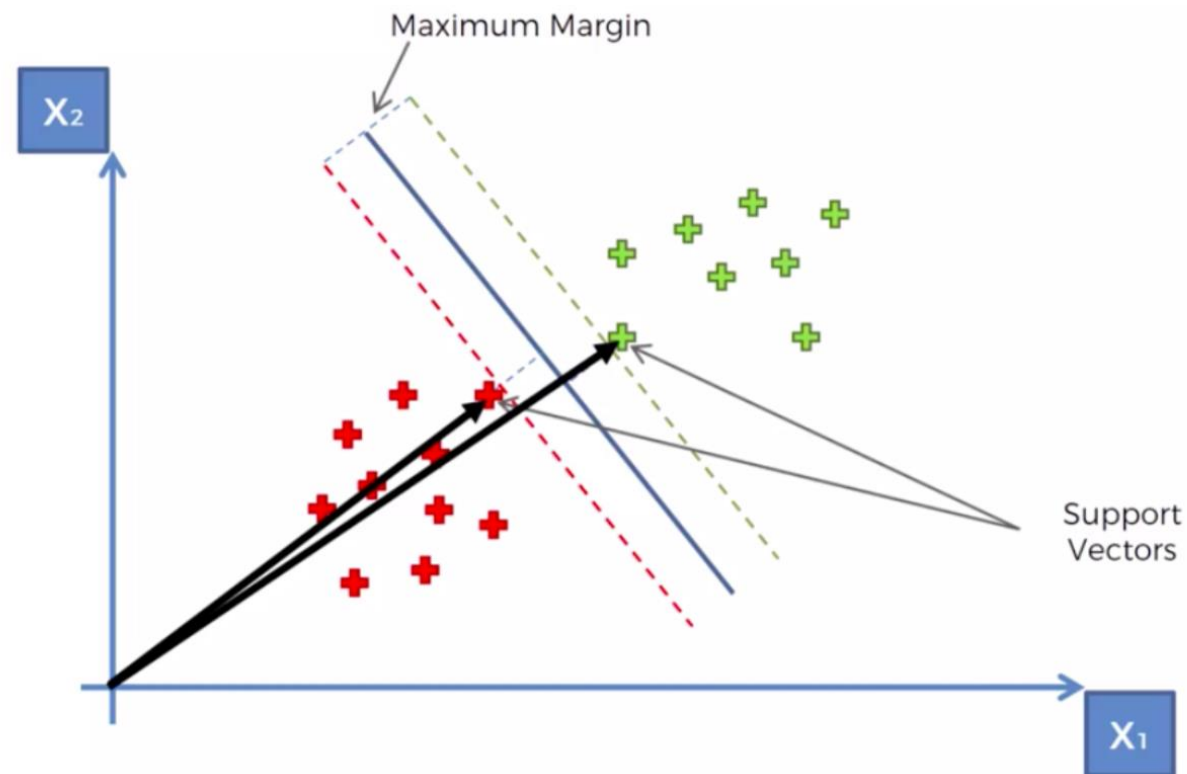




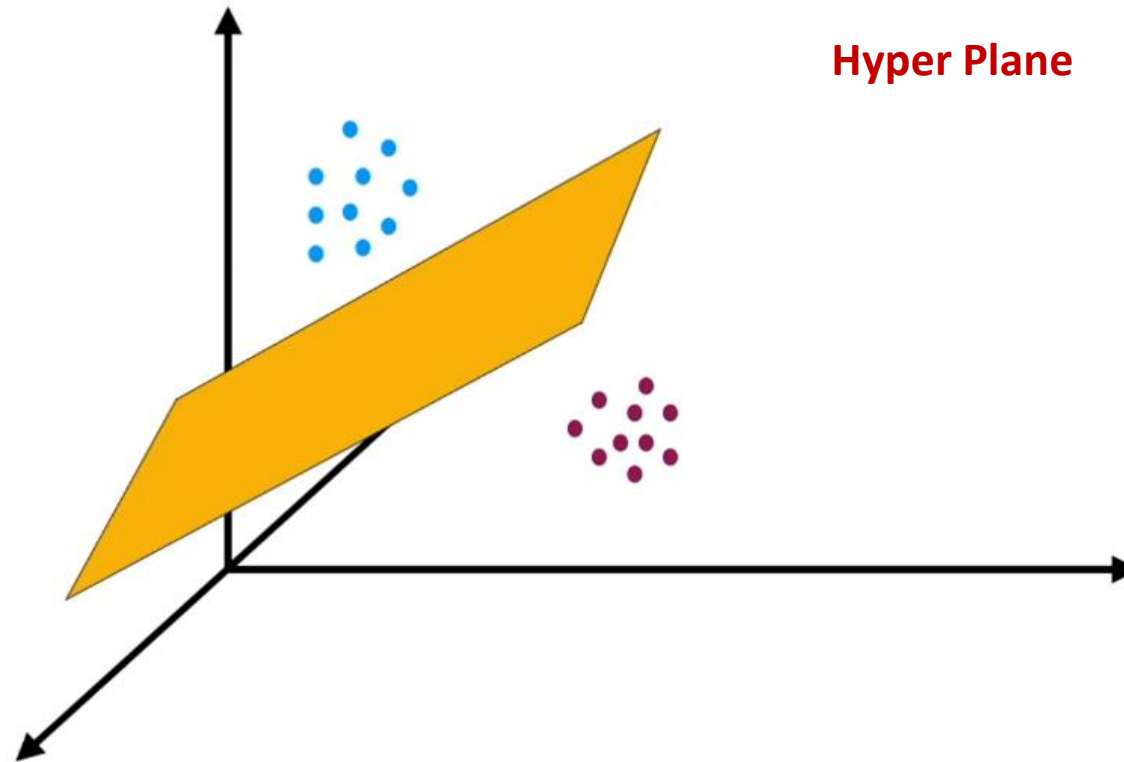
# Machine Learning

## What is an SVM

The LSVM algorithm will select a line that not only separates the two classes but stays as far away from the closest samples as possible. In fact, the “support vector” in “support vector machine” refers to two position vectors drawn from the origin to the points which dictate the decision boundary.



## How SVMs work



In an  $n$  dimensional data/ space SVM finds an  $n-1$  hyperplane to separate points into two categories



# Machine Learning

## Hyper plane

In a vector space of  $n$  dimensions, a hyperplane is a geometrical shape with  $(n-1)$  dimensions and zero thickness in one dimension.

In our example  $Ax + By + Cz = D$

Points on one side of the hyperplane :

$$Ax + By + Cz > D$$

Points on other side of hyperplane:

$$Ax + By + Cz < D$$

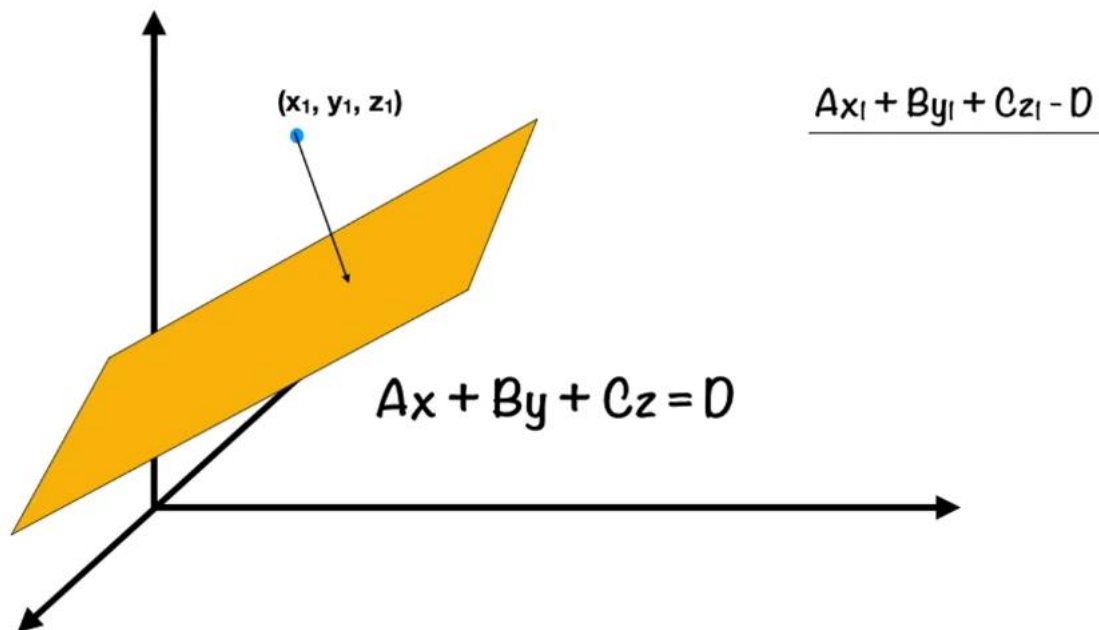
SVM tries to find such a linear equation



# Machine Learning

## Hyper plane

In a vector space of  $n$  dimensions, a hyperplane is a geometrical shape with  $(n-1)$  dimensions and zero thickness in one dimension.



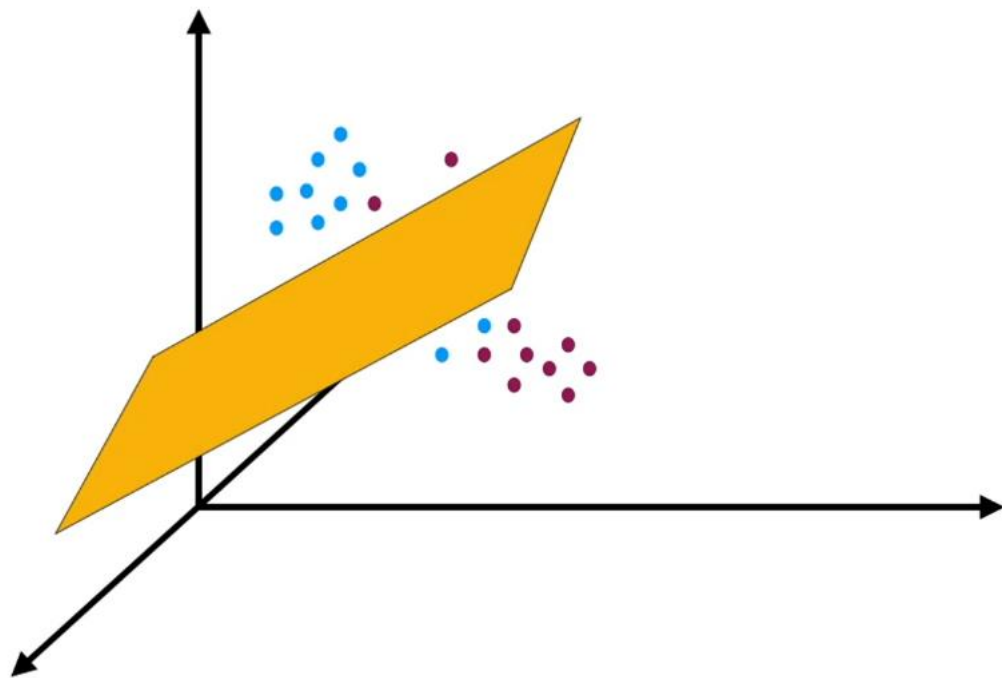
The best hyper plane is the one which maximizes the sum of distances of the nearest points on either side of the plane



# Machine Learning

## Hyper plane

**What if the points on the hyperplane are not linearly separable**



**The soft margin method finds a hyper plane which performs as clean a separation of points as possible**



# Machine Learning

## Non linear separation

**SVM is a linear classifier**

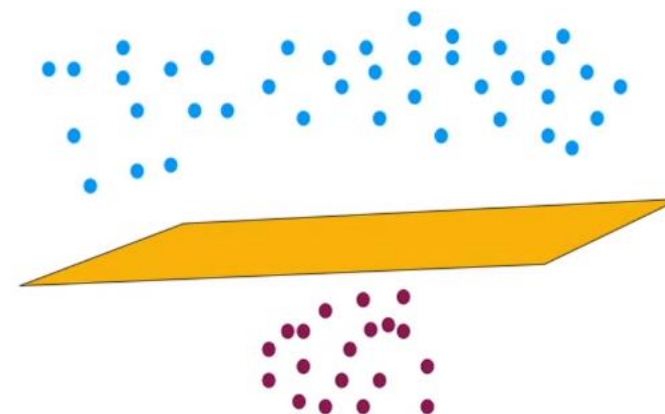
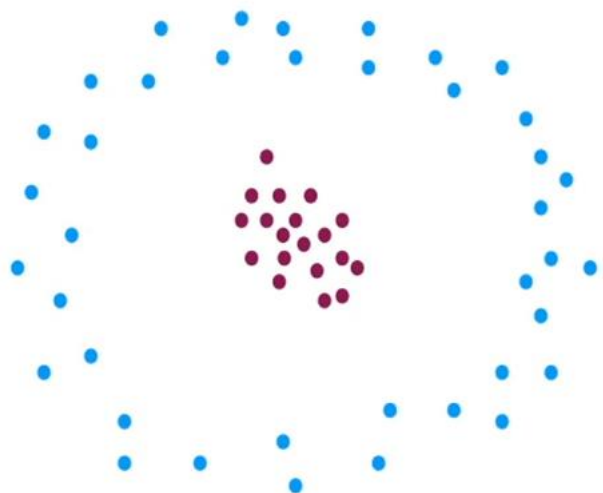
**But can be used to perform non linear separation**

**Achieved by using Kernel Trick**



# Machine Learning

## Kernel Trick



**Transformed / modified to be  
separated by a linear plane**

**Can be separated only by a circle (quadratic)**



# Machine Learning

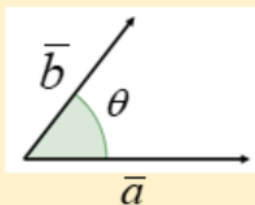
**Linear Classification does a dot product of vectors with many elements**

Algebraically, the dot product is **the sum of the products of the corresponding entries of the two sequences of numbers**. Geometrically, it is the product of the Euclidean magnitudes of the two vectors and the cosine of the angle between them. ... In modern geometry, Euclidean spaces are often defined by using vector spaces.

## Dot Product

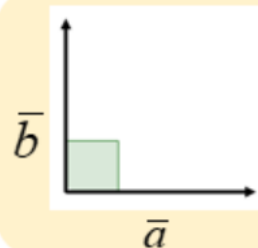
If  $\vec{a} = \langle a_1, a_2, a_3 \rangle$  and  $\vec{b} = \langle b_1, b_2, b_3 \rangle$   
then the dot product is

$$\vec{a} \cdot \vec{b} = a_1 b_1 + a_2 b_2 + a_3 b_3$$



If  $\theta$  is the angle between  $\vec{a}$  and  $\vec{b}$  then

$$\vec{a} \cdot \vec{b} = |\vec{a}| |\vec{b}| \cos \theta$$



$\vec{a} \cdot \vec{b}$  are orthogonal (perpendicular)

if and only if  $\vec{a} \cdot \vec{b} = 0$





# Machine Learning

## Non linear separation

Kernel function operates in feature space which may have which  
may have many more dimensions than original feature space

Finds the maximum margin in hyper plane in modified feature  
space

It's a linear function in the modified feature space

Kernel trick allows a way to solve problems where the data is not  
linearly separable by projecting such data into a higher  
dimensional space



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# Machine Learning

## SVM – Use case

<https://archive.ics.uci.edu/ml/machine-learning-databases/adult/adult.data>



# Machine Learning

## SVM – Use case

```
import numpy as np
```

```
import pandas as pd
```

```
original_data = pd.read_csv("adult.csv",
```

```
    names=['Age','Workclass', 'fnlwgt', 'Education', 'Education-Num',
```

```
    'Marital Status',
```

```
    'Occupation', 'Relationship', 'Race', 'Gender', 'Capital Gain',
```

```
    'Capital Loss',
```

```
    'Hours per week', 'Country', 'Target'], sep =r'\s*,\s*', engine
```

```
    ='python', na_values = "?")
```

```
original_data.head()
```



# Machine Learning

## SVM – Use case

```
import matplotlib.pyplot as plt
```

```
import math
```

```
%matplotlib inline
```

```
fig = plt.figure(figsize = (20,20))
```

```
cols = 3
```

```
rows = math.ceil(float(original_data.shape[1] / cols))
```



# Machine Learning

## SVM – Use case

```
for i, column in enumerate(['Age', 'Workclass', 'Education', 'Occupation', 'Race',  
'Gender']):
```

```
    ax = fig.add_subplot(rows, cols, i+1)
```

```
    ax.set_title(column)
```

```
    if original_data.dtypes[column] == np.object:
```

```
        original_data[column].value_counts().plot(kind='bar', axes = ax)
```

```
        plt.xticks(rotation='vertical')
```

```
plt.subplots_adjust(hspace=0.7, wspace=0.2)
```

```
plt.show()
```



# Machine Learning

## SVM – Use case

**# Use label encoder to convert text to numeric**

**import sklearn.preprocessing as preprocessing**

**le = preprocessing.LabelEncoder()**

**original\_data['Occupation'] =**

**le.fit\_transform(original\_data['Occupation'].astype(str))**

**original\_data.head()**

**()**



# Machine Learning

## SVM – Use case

```
original_data['Target'] =  
le.fit_transform(original_data['Target'].astype(str))  
original_data.tail()
```

```
original_data.Target.unique
```



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# Machine Learning

## SVM – Use case

<https://archive.ics.uci.edu/ml/machine-learning-databases/adult/adult.data>



## **SVM – Use case**

**<https://archive.ics.uci.edu/ml/machine-learning-databases/adult/adult.data>**



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# Machine Learning

## SVM – Use case

<https://archive.ics.uci.edu/ml/machine-learning-databases/adult/adult.data>



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# Machine Learning

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# Machine Learning

## SVM – Use case

<https://archive.ics.uci.edu/ml/machine-learning-databases/adult/adult.data>





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# Machine Learning

## Decision Tree



# Machine Learning (Decision Tree)

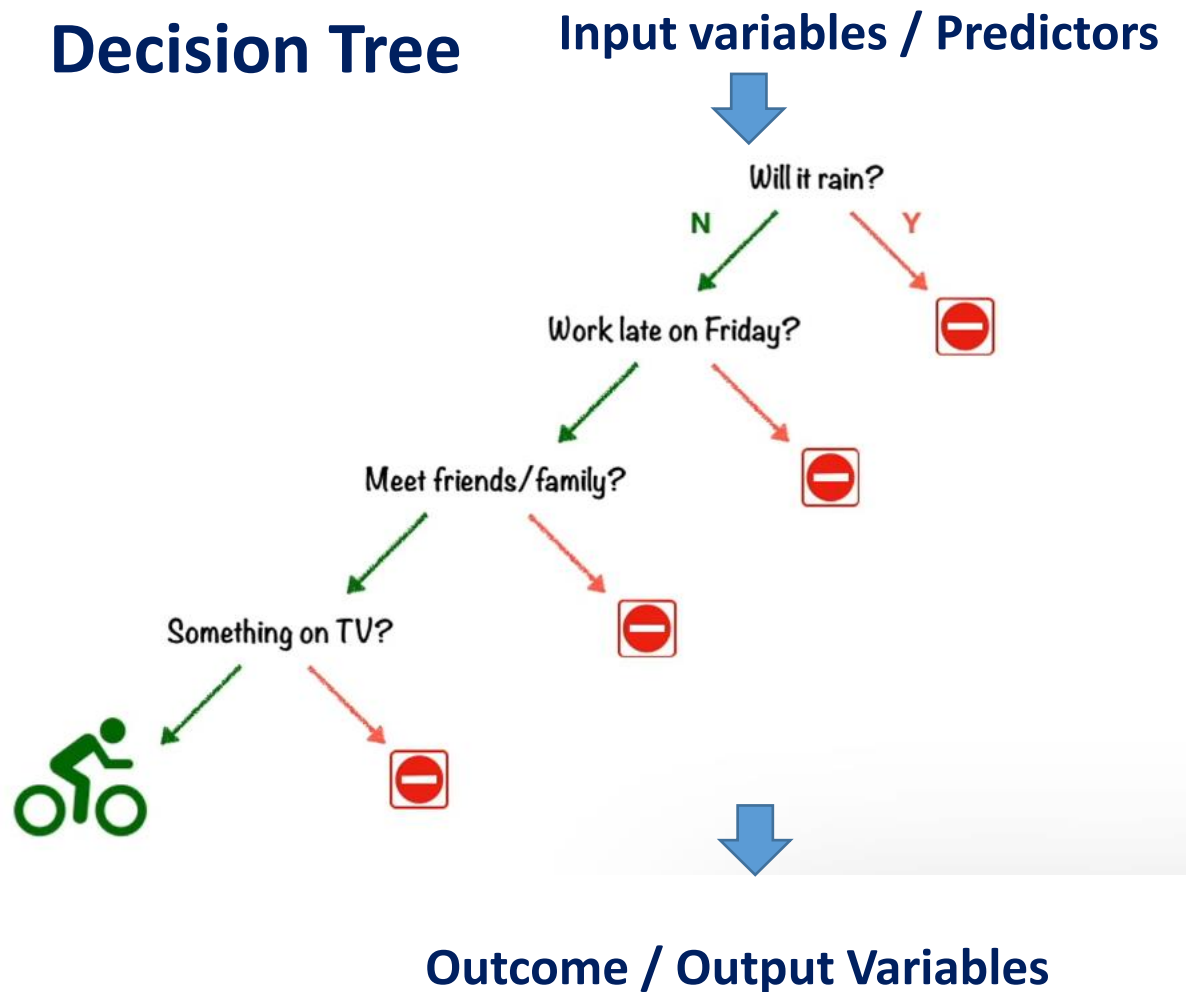
**Should I go biking on a Saturday morning**

- **Will it rain ?**
- **Will I work late on Friday ?**
- **Do I need to meet friends / family ?**
- **Is there something on TV ?**



# Machine Learning (Decision Tree)

## Decision Tree





# Machine Learning (Decision Tree)

## Decision Tree

- **Helps predict the outcome given a set of inputs**
- **In business, it represents visually how a decision is taken with inputs and consequences of each decision**



# Machine Learning (Decision Tree)

## Decision Tree

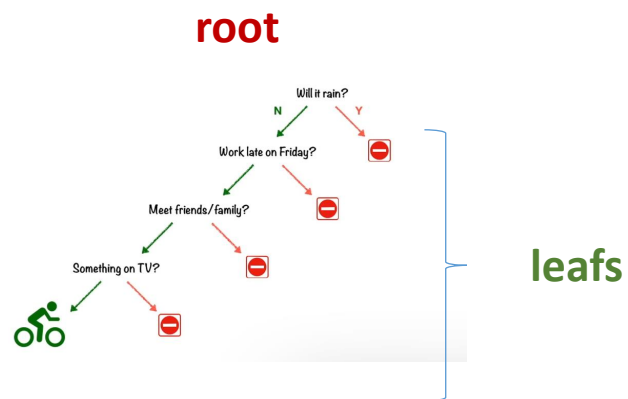
- Can be used for classification, just like SVM
- With SVM, can't understand the relationship between input variable and outcome
- Decision Trees are not a black box



# Machine Learning (Decision Tree)

## Decision Tree

- Inputs can be categorical or continuous
- Output can also be categorical or continuous
  - Called regression tree when outcome is continuous





# Machine Learning (Decision Tree)

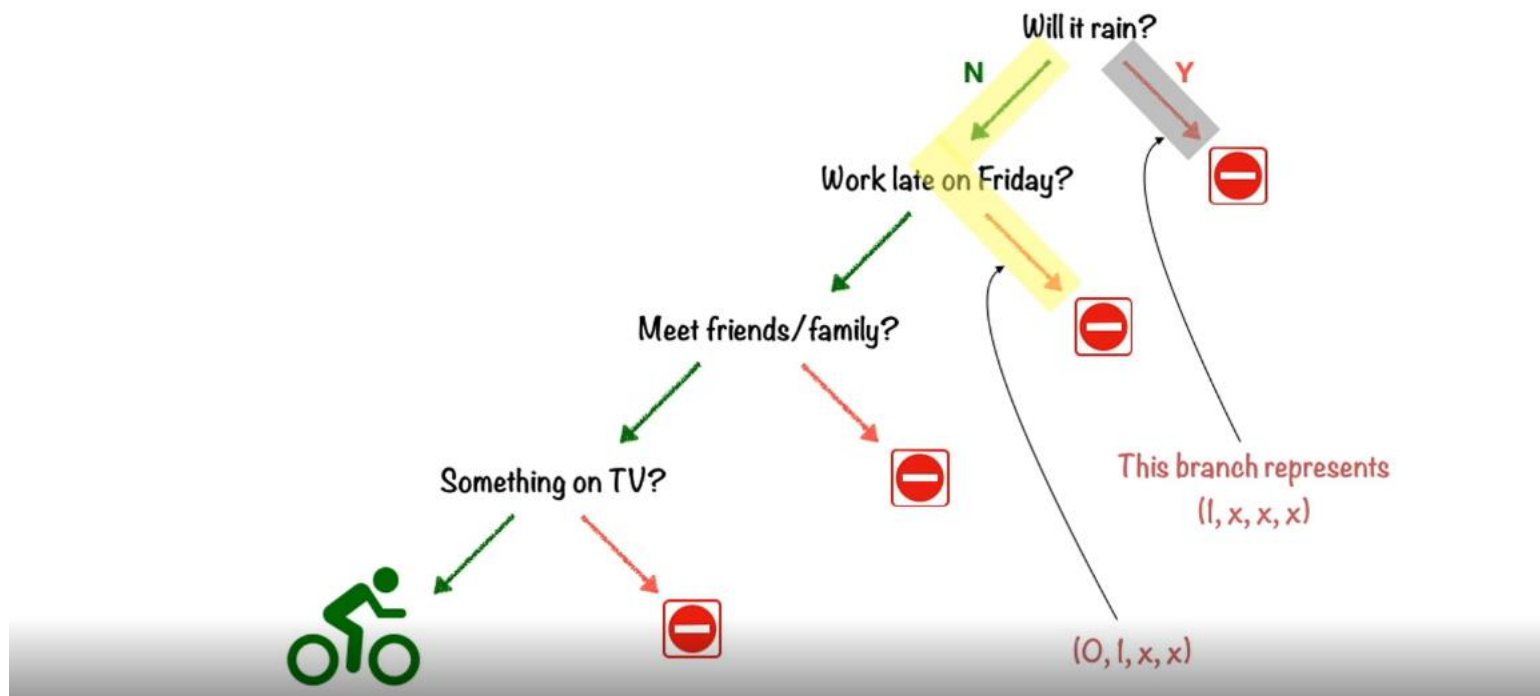
## Decision Tree

- The leaves of the tree are the outcomes
- In classification problem there are class labels
- The tree gives the more likely outcome given the values of the variables
- The branches are combinations of predictor values leading to an outcome



# Machine Learning (Decision Tree)

## Decision Tree







# Machine Learning (Decision Tree)

## Decision Tree Learning

- The process of creating / learning s decision tree from trining data
- Start with training data in the form of feature vector, label / outcome
  - (1, 0, 1, 1) No biking
  - (0, 0, 0, 0) Go biking
- Obtain a decision tree used to classify / predict a new instance
- A supervised learning approach



# Machine Learning (Decision Tree)

## Decision Tree Learning

- Recursive partitioning is the most common strategy for decision tree learning
- Decision tree learning algorithms
  - CART
  - ID3
  - C4.5
  - CHAID



# **Machine Learning (Decision Tree)**

## **Decision Tree Learning**

- **Algorithm needs to tell us the order in which the predictors are evaluated**
- **If the predictors is continuous, the tree needs to split the variables into ranges**



# Machine Learning (Decision Tree)

## Greedy Algorithm for Learning a Decision Tree

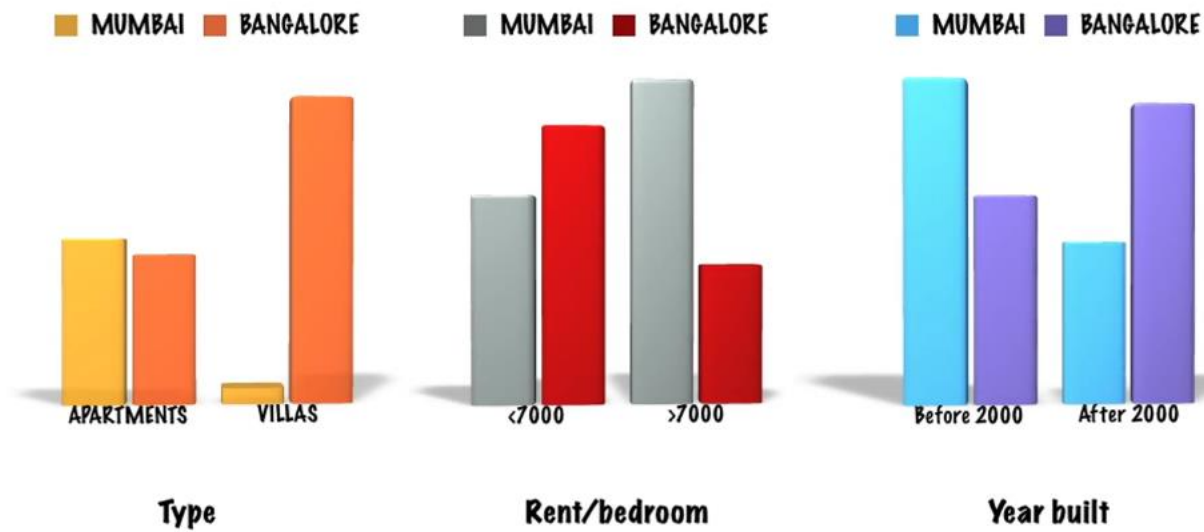
- **Given the type of housing, rent /bedroom and the year it was built**
- **Predict the city to which a residence belongs**
  - **Mumbai (or) Bangalore**



# Machine Learning (Decision Tree)

## Greedy Algorithm for Learning a Decision Tree

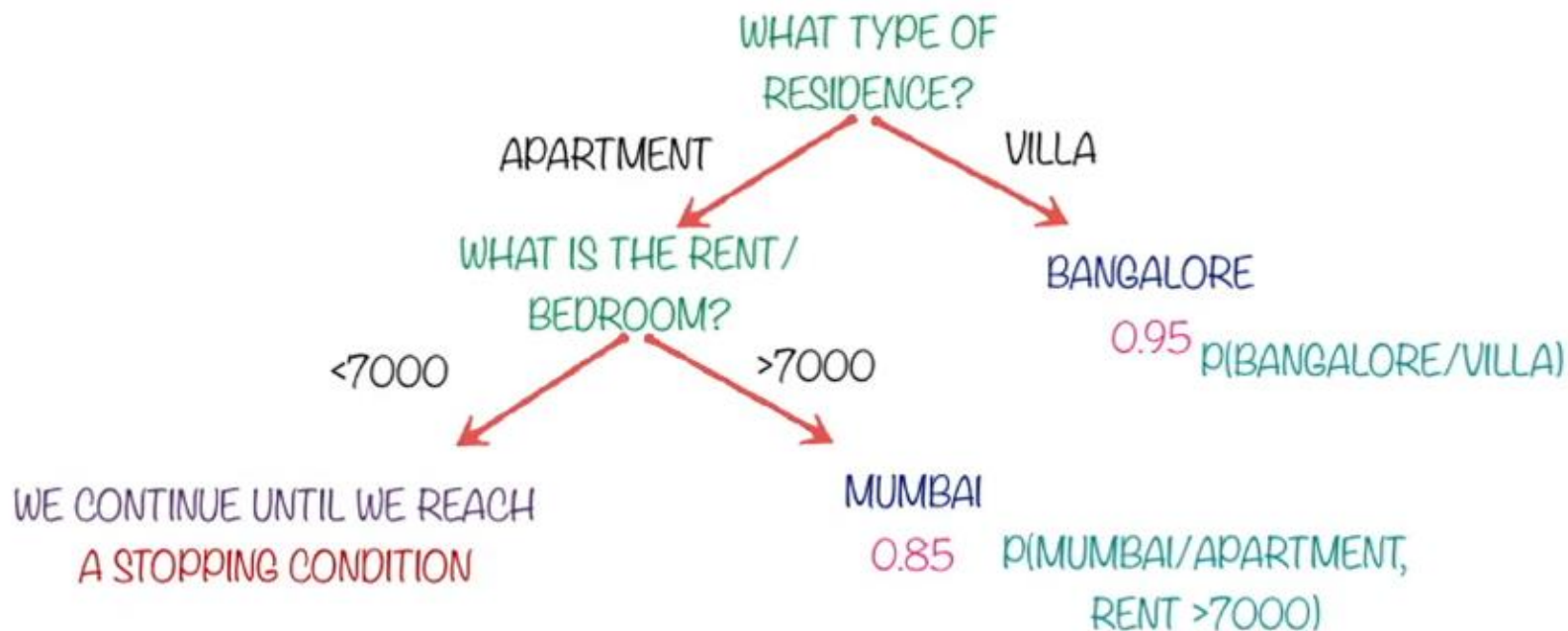
Draw a histogram for each attribute, for residences in each city





# Machine Learning (Decision Tree)

## Greedy Algorithm for Learning a Decision Tree



This is Recursive Partitioning



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# Machine Learning (Decision Tree)

The stopping condition could be

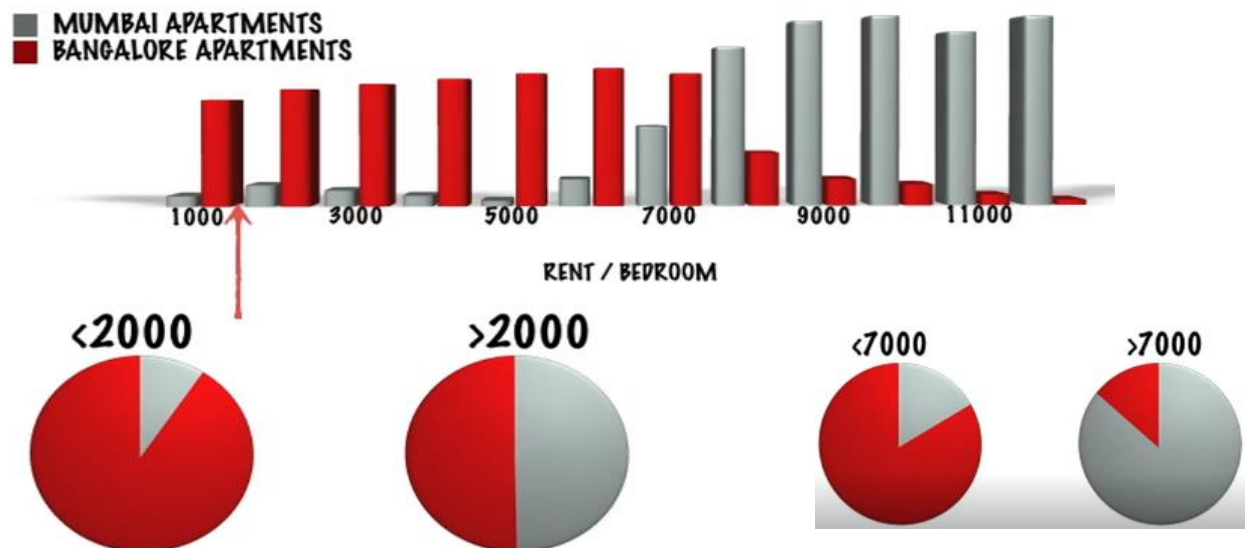
- All our subsets are mostly homogeneous
- We have run out of attributes
- The Tree is too large



# Machine Learning (Decision Tree)

The best split for a continuous input variable

- What is the best split for rent in our example ?
- Plot a histogram based on rent per bedroom



The best split is the point where the subsets we get are mostly homogeneous





# Machine Learning (Decision Tree)

## Information Gain

- Any statement, News or Message Contains Information
- Some have more information and some have less
- The Idea of information gain is to reduce entropy and maximize information



# Machine Learning (Decision Tree)

## Information Gain (closer look)



Let's say you have to classify an animal as a giraffe or a hippo

We are told the animal has 4 legs

This basically useless as both the animals have 4 legs

If we are told that the animal is 10 feet tall

This is a useful information, It tells us the animal is very likely a giraffe



So clearly – The values of some attributes give us more information than others  
&

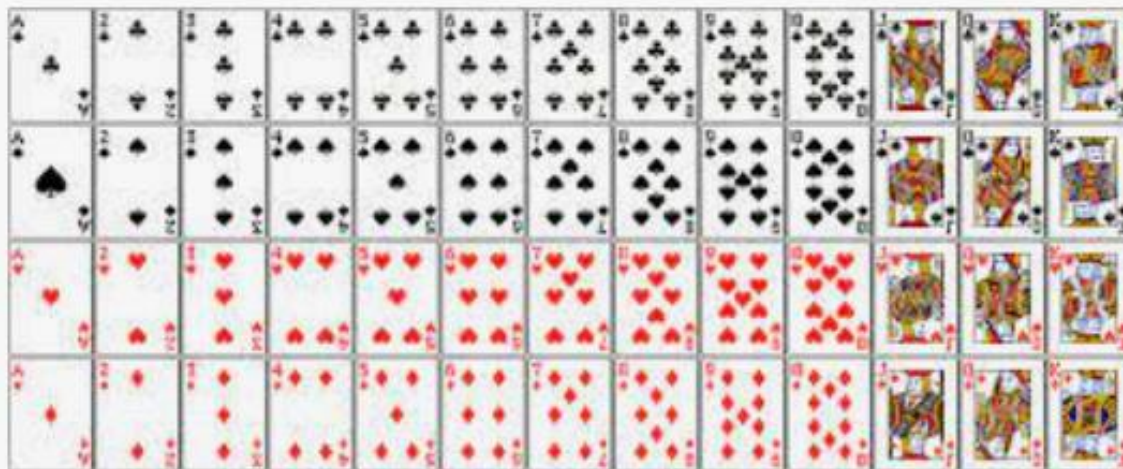
There are mathematical way to measure this information



# Machine Learning (Decision Tree)

## Information Gain

- Another example: Game of cards



Guess the card held by your opponent

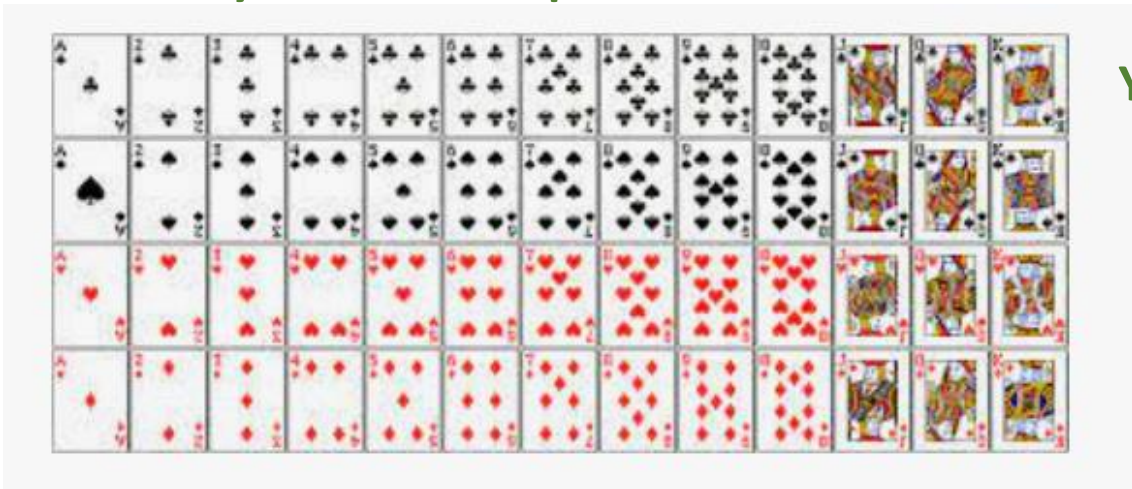
You are allowed to ask Yes or No  
questions

To begin with you have 52 possible outcomes



# Machine Learning (Decision Tree)

Initially there are 52 possible outcomes in all



Your question 1: Is the card an ACE ?

Yes

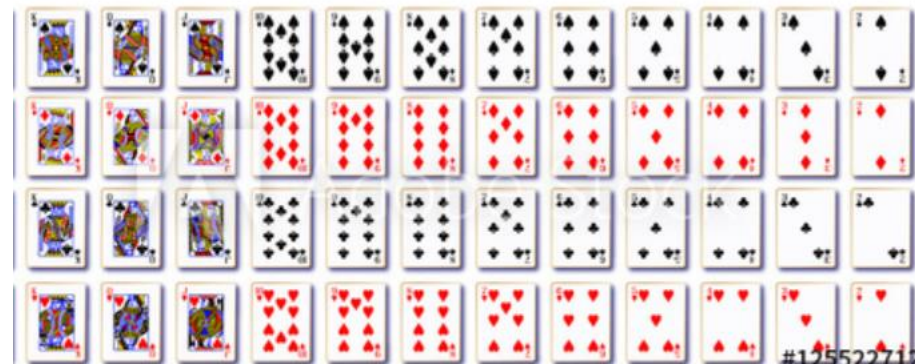
No



You are left  
with 4 possible  
outcomes

The answer Yes  
gives us **more**  
**information** then  
answer NO

You are left with  
48 possible  
outcomes



#125522710



# Machine Learning (Decision Tree)

Initially there are 52 possible outcomes in all

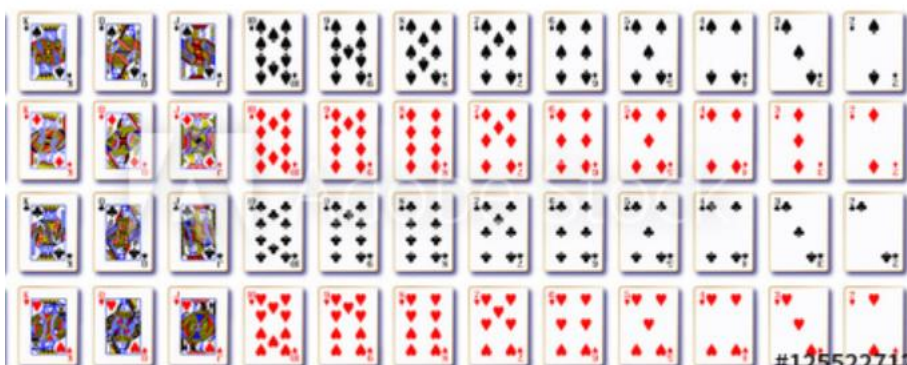
Yes



The answer “Yes” has a lower probability

$$P(\text{Yes}) = 4/52$$

No



The answer “No” has a Higher probability

$$P(\text{Yes}) = 48/52$$

The lower the probability of the answer, the more information you get

Your question 1: Is the card an ACE ?

The answer Yes gives us more information than answer NO

The probability of any occurrence is b/n 0-1 and the Log of value is always -ve regardless of the base of the log

The lower the probability the more -ve the log value will be, however since we are considering the -ve log for info content the lower the probability of the value Higher the information content

If X is the random variable that represents the answer to our question

$$\text{Information content of } (X = \text{Yes}) = -\text{Log } (P(X = \text{Yes}))$$

$$\text{Information content of } (X = \text{No}) = -\text{Log } (P(X = \text{No}))$$

$$\text{Information content of } (X = x) = -\text{Log } (P(X = x))$$



# Machine Learning (Decision Tree)

## Information Gain

- Any statement, News or Message Contains Information
- Some have more information and some have less
- The Idea of information gain is to reduce entropy and maximize information





# Machine Learning (Decision Tree)

## Information Gain

If  $X$  is the random variable that represents the answer to our question

Information content of  $(X = x) = -\text{Log} (P(X = x))$

Average value of the information content also called the expected value =  $\sum P(X = x) (-\text{Log} (P(X = x)))$

Entropy  $H(X)$

Entropy increases with:

- 1) Number of possible answers
- 2) The evenness of probability distribution

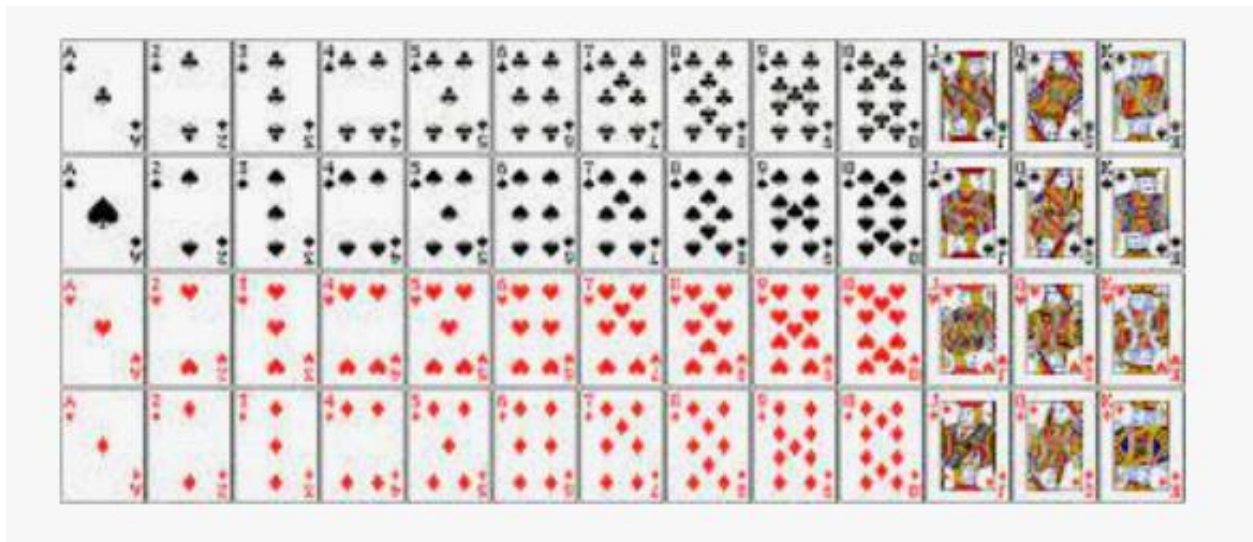
Entropy is the amount of uncertainty / Unpredictability there is in the answer

$P(\text{Yes}) = 0 \Rightarrow$  There is no uncertainty  $\Rightarrow$  entropy = 0  
Yes and No have equal probability  $\Rightarrow$  Very High entropy



# Machine Learning (Decision Tree)

Back to our Cards example



Initially there are 52  
possible outcomes

Each card has the same  
probability =  $1/52$

Information content of  $(X = x) = -\text{Log}(P(X = x))$

Entropy  $H(X)$  = Average value of the information  
content ( Also called as expected value  
 $= \sum P(X = x) (-\text{Log}(P(X = x)))$

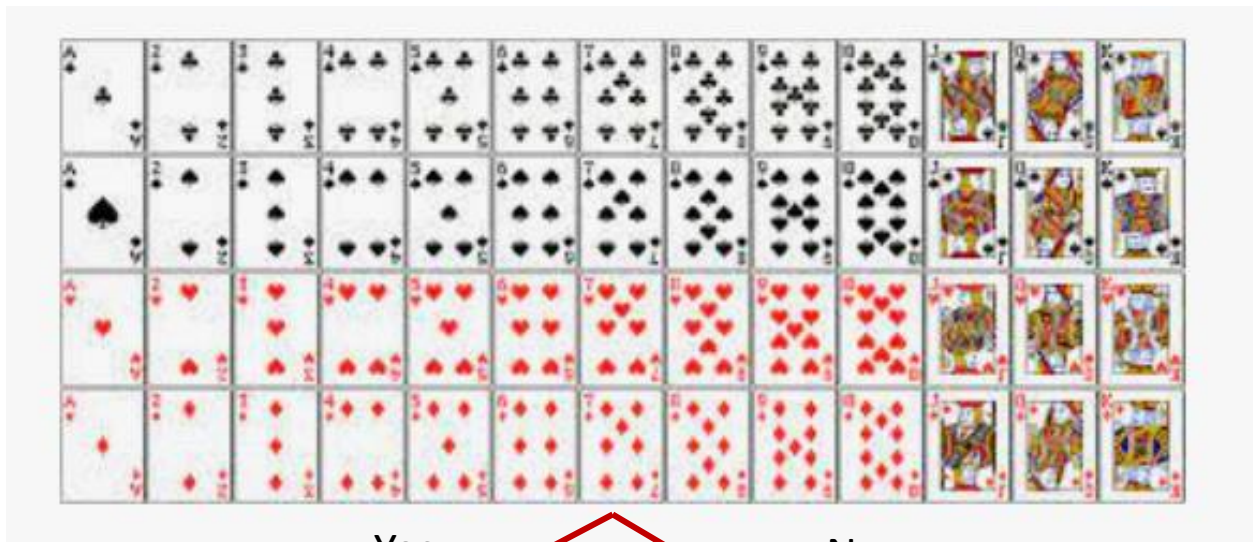
$$\begin{aligned}\text{Entropy} = H(X) &= \sum (1/52) (-\text{LOG}(1/52)) \\ &= \text{LOG}(52)\end{aligned}$$





# Machine Learning (Decision Tree)

Back to our Cards example



Yes

No

$$\text{Entropy} = H(X/Q1 = \text{Yes}) \\ = \text{Log}(4) \quad P(\text{Yes}) = 4/52$$

Within each group the  
entropy has decreased

$$\text{Entropy} = H(X/Q1 = \text{No}) = \\ \text{Log}(48) \quad P(\text{No}) = 48/52$$

The more homogenous  
each group is the lower the  
entropy

Entropy  $H(X)$  = Average value of the information  
content ( Also called as expected value  
 $= \sum P(X = x) ( - \text{Log} (P(X = x)))$

Before we have asked the question the uncertainty  
(entropy) in our guess is very high

$$\text{Entropy} = H(X) = \sum (1/52) (-\text{LOG}(1/52)) \\ = \text{LOG}(52)$$

However, Once we ask the question,  
Is the card an ACE ?



# Machine Learning (Decision Tree)

## Information Gain

Before we have asked any yes/no questions, the uncertainty (Entropy) in our guess is very high

$$\text{Entropy} = H(X)$$

After we ask the question, Is the card ACE ?

$$\text{Entropy after Q1} = H(X / Q1)$$

$$\text{Information Gain} = \text{Reduction in Entropy Overall} = H(X) - H(X / Q1)$$

As we saw whenever we ask a question, subsets are formed

When each of these subsets are homogenous, the information gain maximum



# Machine Learning (Decision Tree)

## Information Gain

$$\text{Entropy} = H(X)$$

$$\text{Entropy after } Q1 = H(X / Q1)$$

As we saw whenever we ask a question, subsets are formed

Is It an ACE ?

$$H(X/Q1) = 4/52 * \text{Log}(4) + 48 / 52 * \text{Log} (48)$$

$$IQ = H(X) - H(X/Q1) = 0.12$$

$$\text{Information Gain} = \text{Reduction in Entropy Overall} = H(X) - H ( X/Q1 )$$

When each of these subsets are homogenous, the information gain maximum

Is It a Black card ?

$$H(X/Q1) = \text{Log}(26)$$

$$IQ = H(X) - H(X/Q1) = \text{Log}(52) - \text{Log}(26) = 0.30$$

More Homogenous our subset the greater the information gain.....



# Machine Learning (Decision Tree)

## Decision Tree Learning

- Recursive partitioning is the most common strategy for decision tree learning

- Decision tree learning algorithms

|         |
|---------|
| • CART  |
| • ID3   |
| • C4.5  |
| • CHAID |

GINI Impurity

Information Gain

Chi Square Statistic

Each of these learning algorithms have slightly different way of arriving or measuring the homogeneity of a subset



# Machine Learning (Decision Tree)

## CART

CART is another decision tree learning method

**CART : Classification and Regression Trees**

**It uses a different way to choose an attribute**

**The idea is to Minimize the GINI Impurity at each step**

**Idea behind GINI impurity is to choose an attribute such that if you stop the decision tree with that attribute and go no further, **The probability of a false label is minimized****



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# Machine Learning (Decision Tree Lab)

## Data for sample case

<https://archive.ics.uci.edu/ml/machine-learning-databases/wine-quality/>

<https://www.kaggle.com/uciml/pima-indians-diabetes-database>



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# Machine Learning (Decision Tree)

```
(base) PS C:\Users\LENOVO> conda install python-graphviz
Collecting package metadata (current_repodata.json): done
Solving environment: done

## Package Plan ##
```



# Machine Learning (Decision Tree)

## Decision Tree Learning

- Recursive partitioning is the most common strategy for decision tree learning

- Decision tree learning algorithms

|         |
|---------|
| • CART  |
| • ID3   |
| • C4.5  |
| • CHAID |

GINI Impurity

Information Gain

Chi Square Statistic

Each of these learning algorithms have slightly different way of arriving or measuring the homogeneity of a subset





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# Machine Learning (Random Forest)

**Random Forest**



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# Machine Learning (Random Forest)

**Random Forest**

A single, smooth, red curved line that starts at the bottom left and arcs upwards towards the top right, spanning most of the width of the slide.



# Machine Learning (Overfitting)

**GFC -2007**

**Some ways of Mitigating this problem:**

- **Cross Validation**
- **Regularization**
- **Ensemble Learning**



# Machine Learning (Overfitting)

## Example: Restaurant rating

A Model should be simple and generic

Vs

Model is complex and more accurate on training data



# Machine Learning (Overfitting)

**Overfitting happens when model picks up random phenomena or noise present in the training set. Instead of the underlying relationship between the input and output**

**Why is Over fitting such a common problem:**

**The training set is only part of the much larger set**

**We are trying to find models that describe this much larger set**

**It's like trying to describe a photograph but you are shown only a small zoomed portion of the entire photograph**



# Machine Learning (Overfitting)

## Cross Validation

THE BELOW TABLE REPRESENTS THE  
ENTIRE TRAINING DATA SET

|       |       |       |       |       |       |       |       |          |          |          |          |          |          |          |          |
|-------|-------|-------|-------|-------|-------|-------|-------|----------|----------|----------|----------|----------|----------|----------|----------|
| $D_0$ |       |       |       |       |       |       |       | $D_1$    |          |          |          |          |          |          |          |
| $X_1$ | $X_2$ | $X_3$ | $X_4$ | $X_5$ | $X_6$ | $X_7$ | $X_8$ | $X_9$    | $X_{10}$ | $X_{11}$ | $X_{12}$ | $X_{13}$ | $X_{14}$ | $X_{15}$ | $X_{16}$ |
| $X_1$ | $X_2$ | $X_3$ | $X_4$ | $X_5$ | $X_6$ | $X_7$ | $X_8$ | $X_9$    | $X_{10}$ | $X_{11}$ | $X_{12}$ | $X_{13}$ | $X_{14}$ | $X_{15}$ | $X_{16}$ |
| TEST  |       |       |       |       |       |       |       | TRAINING |          |          |          |          |          |          |          |

THE BEST MODEL IS THE ONE WITH BEST AVERAGE  
PERFORMANCE

1. DIVIDE THE TRAINING SET RANDOMLY INTO TWO EQUAL  
PARTS -  $D_0$  AND  $D_1$

2. USE  $D_0$  TO TRAIN THE MODEL AND  $D_1$  TO TEST THE  
PERFORMANCE

3. THEN, USE  $D_1$  TO TRAIN  
THE MODEL AND  $D_0$  TO  
TEST THE  
PERFORMANCE

THIS TECHNIQUE IS CALLED

## 2-FOLD CROSS VALIDATION



# Machine Learning (Overfitting)

## Regularization

$$E'(f) = E(f) + \lambda R(f)$$





# Machine Learning (Overfitting)

## Ensemble Technique

### Bagging

- Random Forest

### Boosting

- ADABOOST
- Gradient Boosting
- XGBoosting





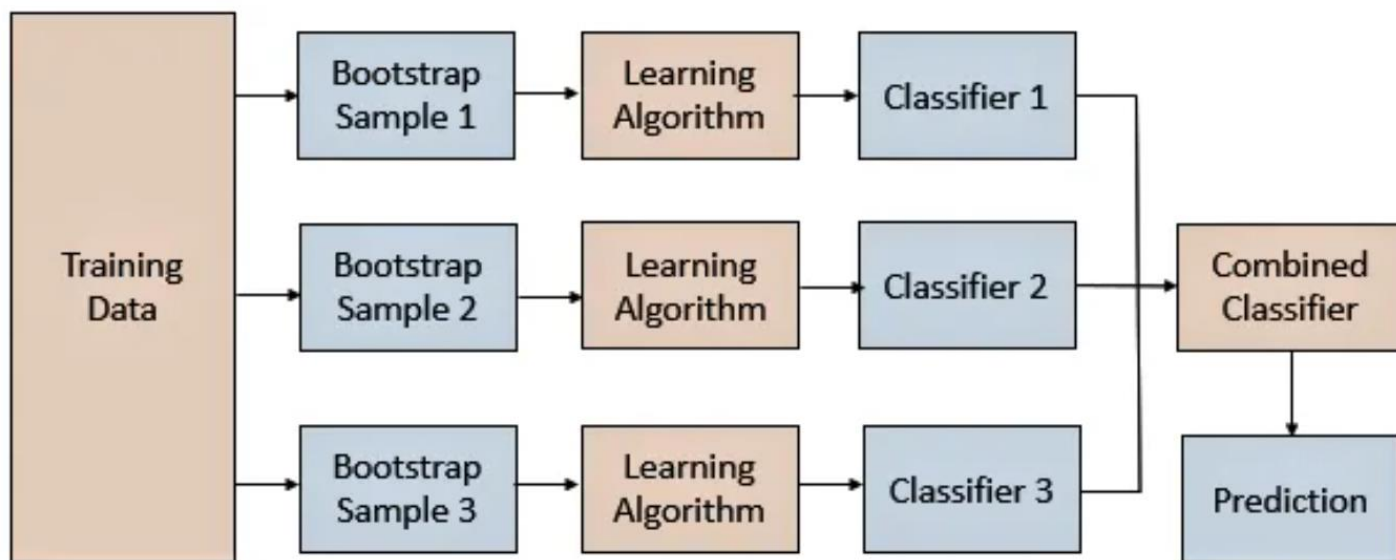
# Machine Learning (Decision Tree)

## Bagging

- Bagging is a technique used to reduce the variance of our predictions by combining the result of multiple classifiers models on different sub-samples of the same data set.

The steps followed in bagging are:

1. Create Multiple datasets
2. Build Multiple Classifiers
3. Combine Classifiers

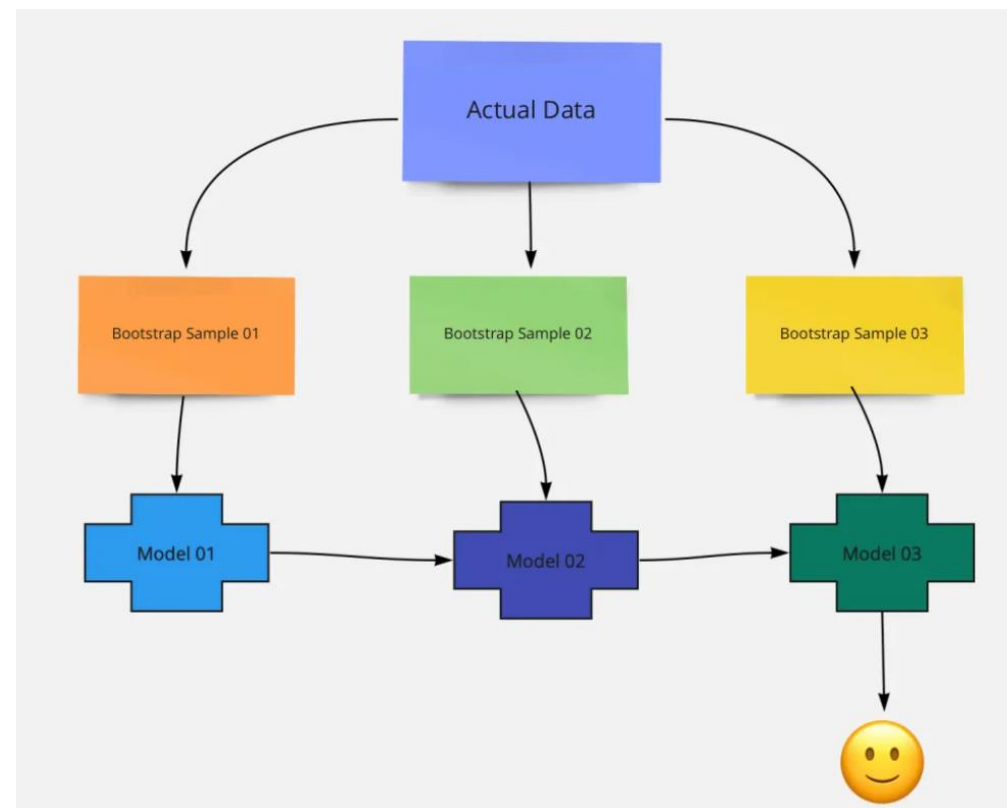




# Machine Learning (Decision Tree)

## Boosting

- In boosting method, all the individual models are built sequentially. Which means the outcome of the first model passes to the next model and so on.
- In bagging the models are built parallel so we don't know what the error of each model is. Whereas in boosting once the first model built we know the error of that model. So when we pass this first model to the next model the intention is to reduce the error further. In some boosting algorithm, each model has to reduce a minimum of 50% of error



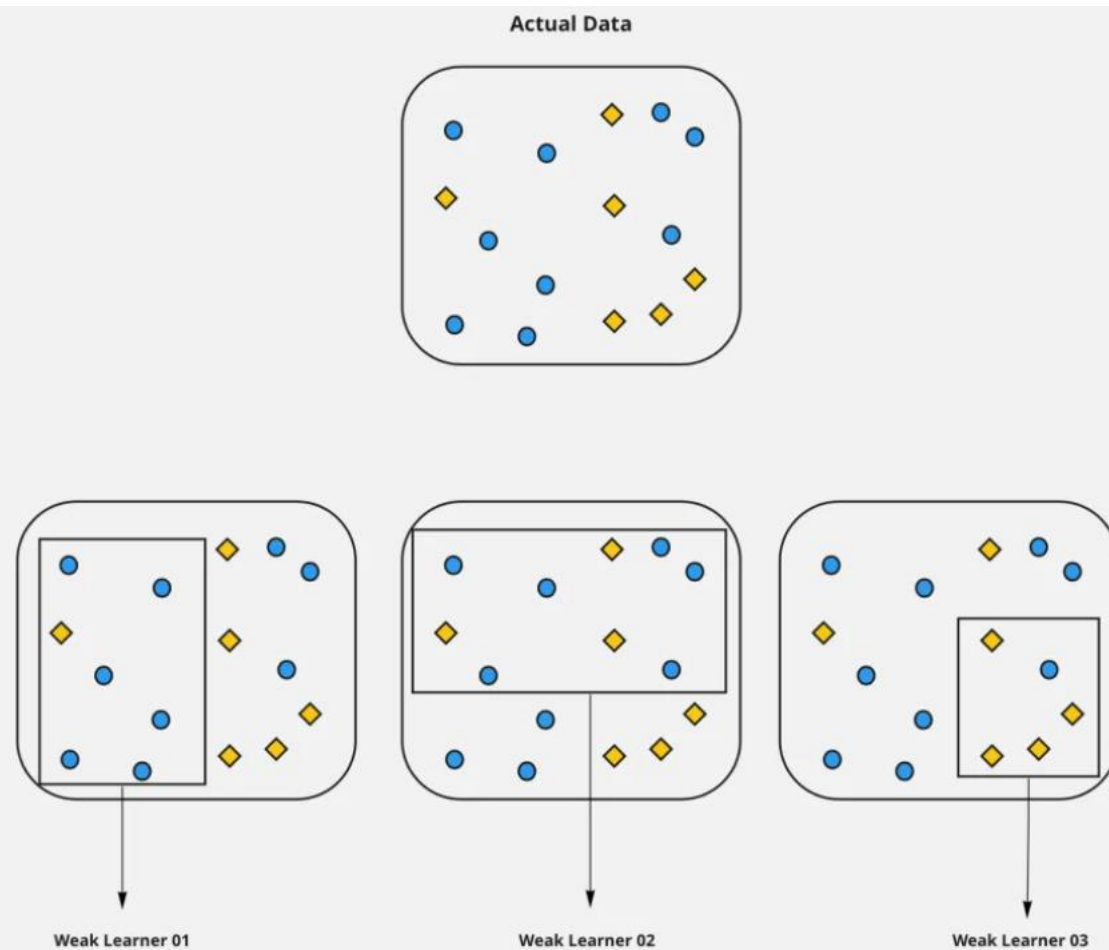


# Machine Learning (Decision Tree) Bootstrapping

For each model, we need to take a sample of data, but we need to be very careful while creating these samples of data. Because if we randomly take the data, in a single sample we will end up with only one target class or the target class distribution won't be the same. This will affect model performance.

To overcome this we need a smart way to create these samples, known as bootstrapping samples.

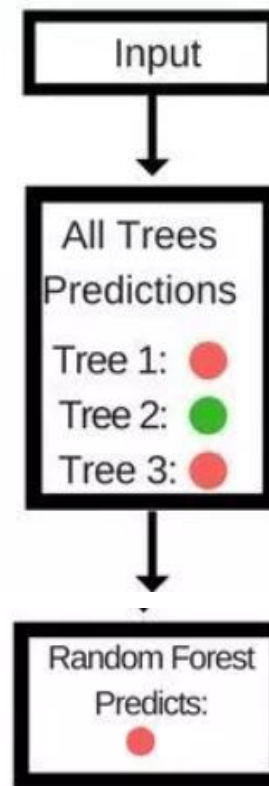
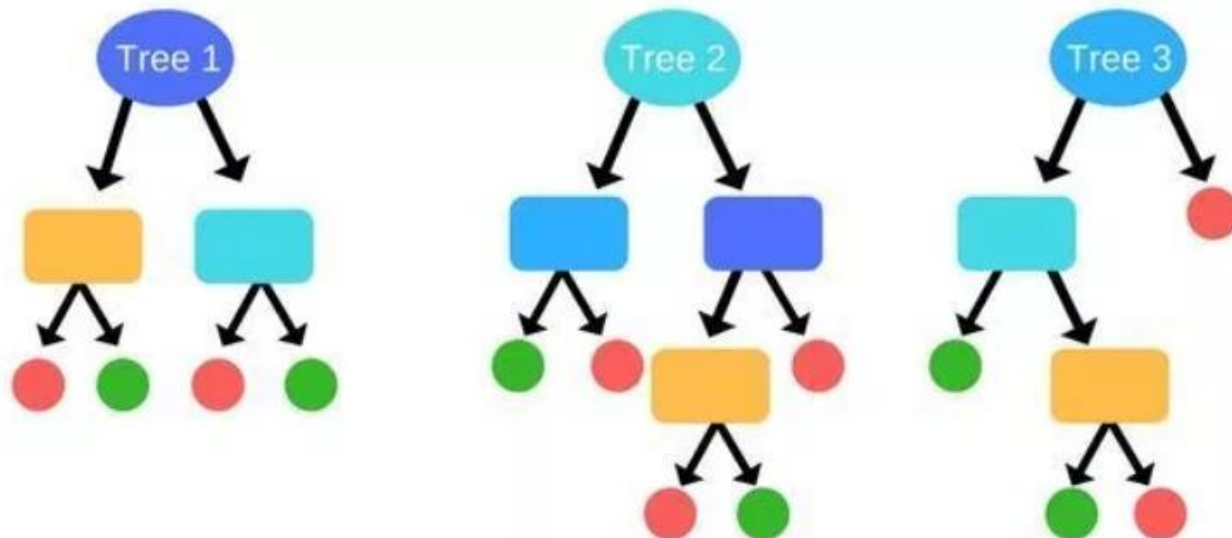
Bootstrapping is a statistical method to create sample data without leaving the properties of the actual dataset. The individual samples of data called bootstrap samples.





# Machine Learning (Decision Tree)

## Random Forest



- The same random forest algorithm or the random forest classifier can use for both classification and the regression task.
- Random forest classifier will handle the missing values.
- When we have more trees in the forest, a random forest classifier won't overfit the model.
- Can model the random forest classifier for categorical values also.



# Machine Learning (Decision Tree)

## Random Forest

### Random Forest pseudocode:

- Randomly select “k” features from total “m” features.
- Where  $k \ll m$
- Among the “k” features, calculate the node “d” using the best split point.
- Split the node into daughter nodes using the best split.
- Repeat 1 to 3 steps until “l” number of nodes has been reached.
- Build forest by repeating steps 1 to 4 for “n” number times to create “n” number of trees.

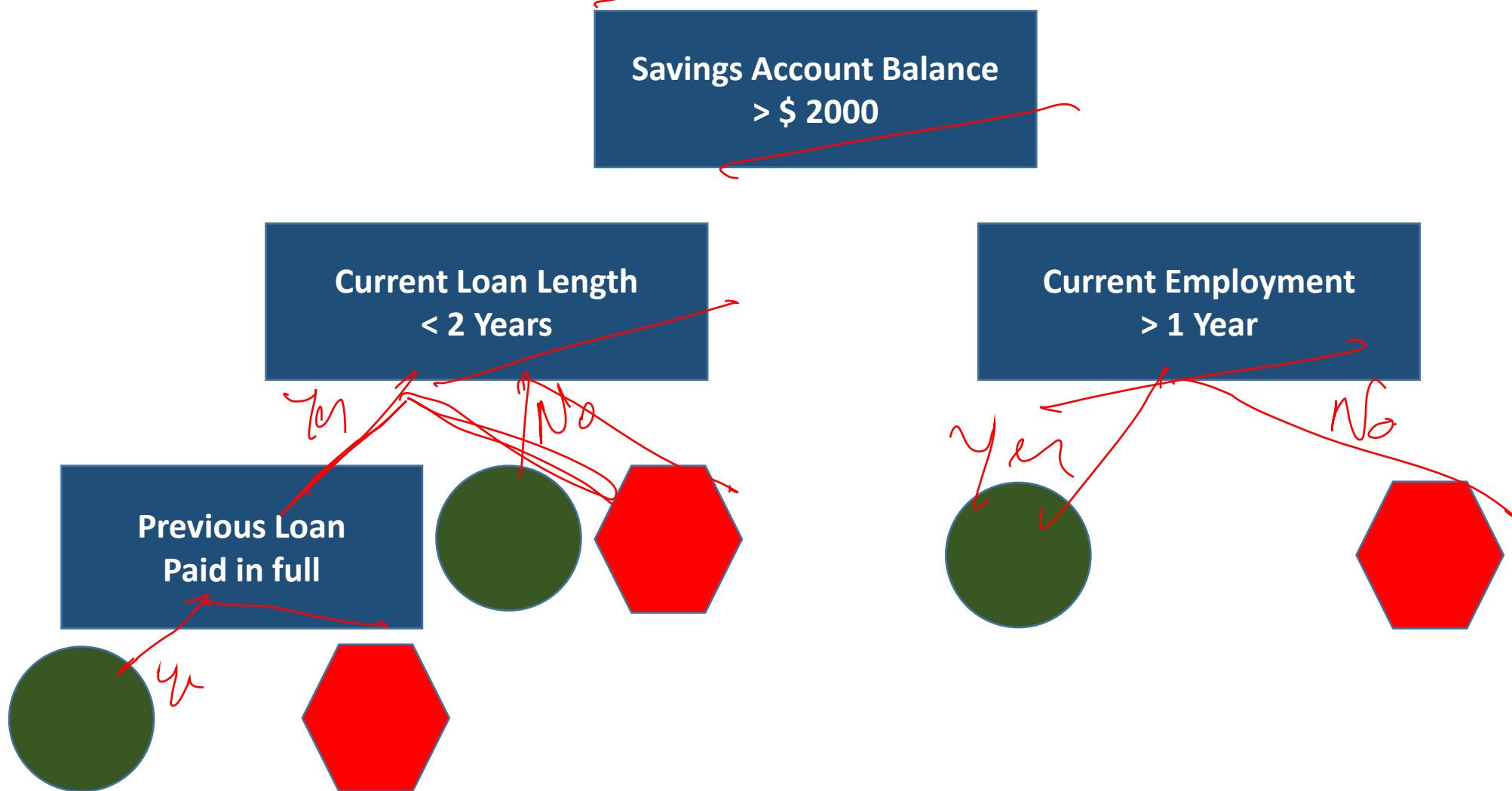
### Random forest prediction pseudocode:

- Takes the test features and use the rules of each randomly created decision tree to predict the outcome and stores the predicted outcome (target)
- Calculate the votes for each predicted target.
- Consider the high voted predicted target as the final prediction from the random forest algorithm.



# Machine Learning (Decision Tree)

## Credit Risk Prediction



# K Means



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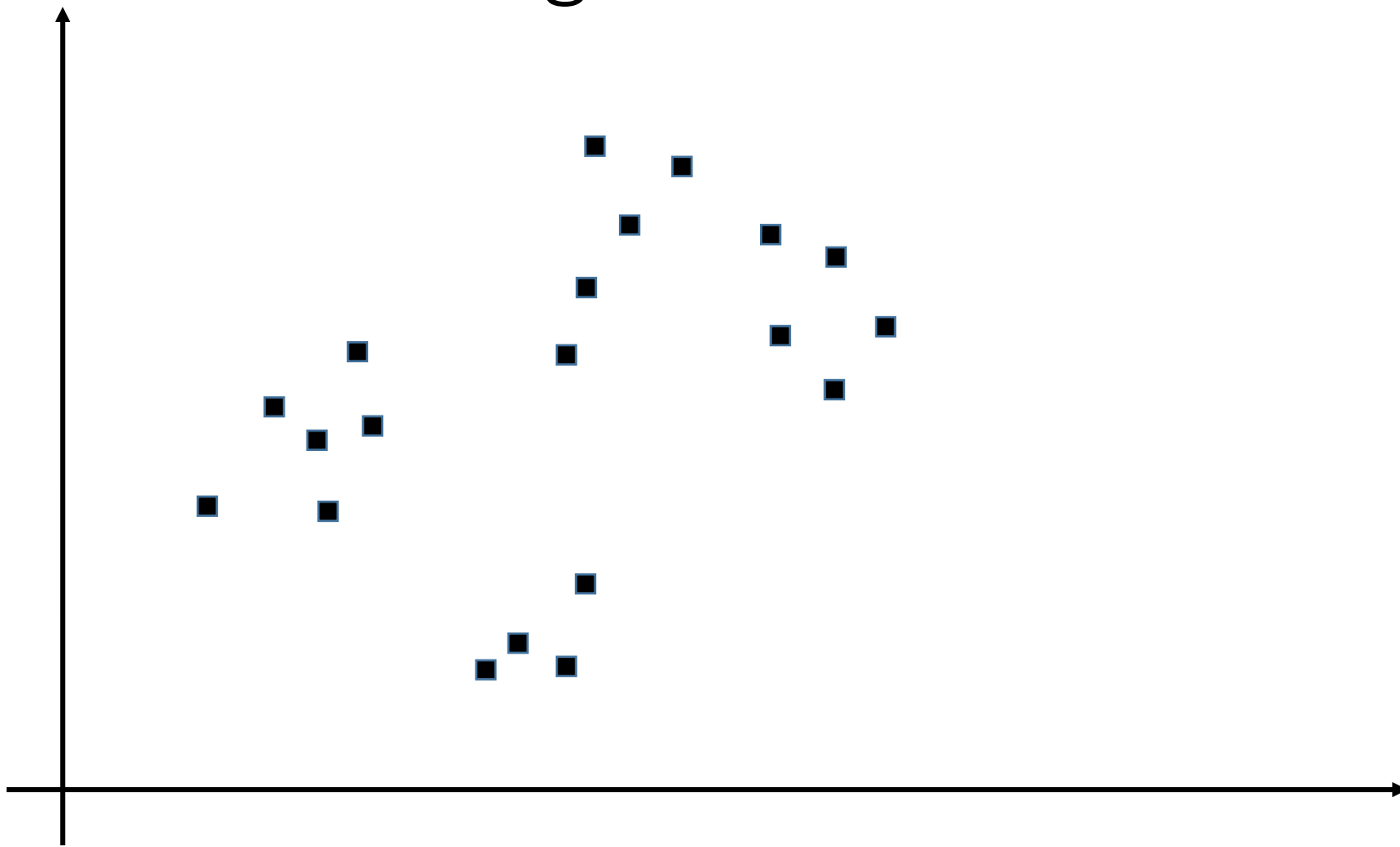
# Machine Learning

- Supervised Machine Learning
- Unsupervised Machine Learning
- Semi Supervised Machine Learning
- Reinforcement Machine Learning



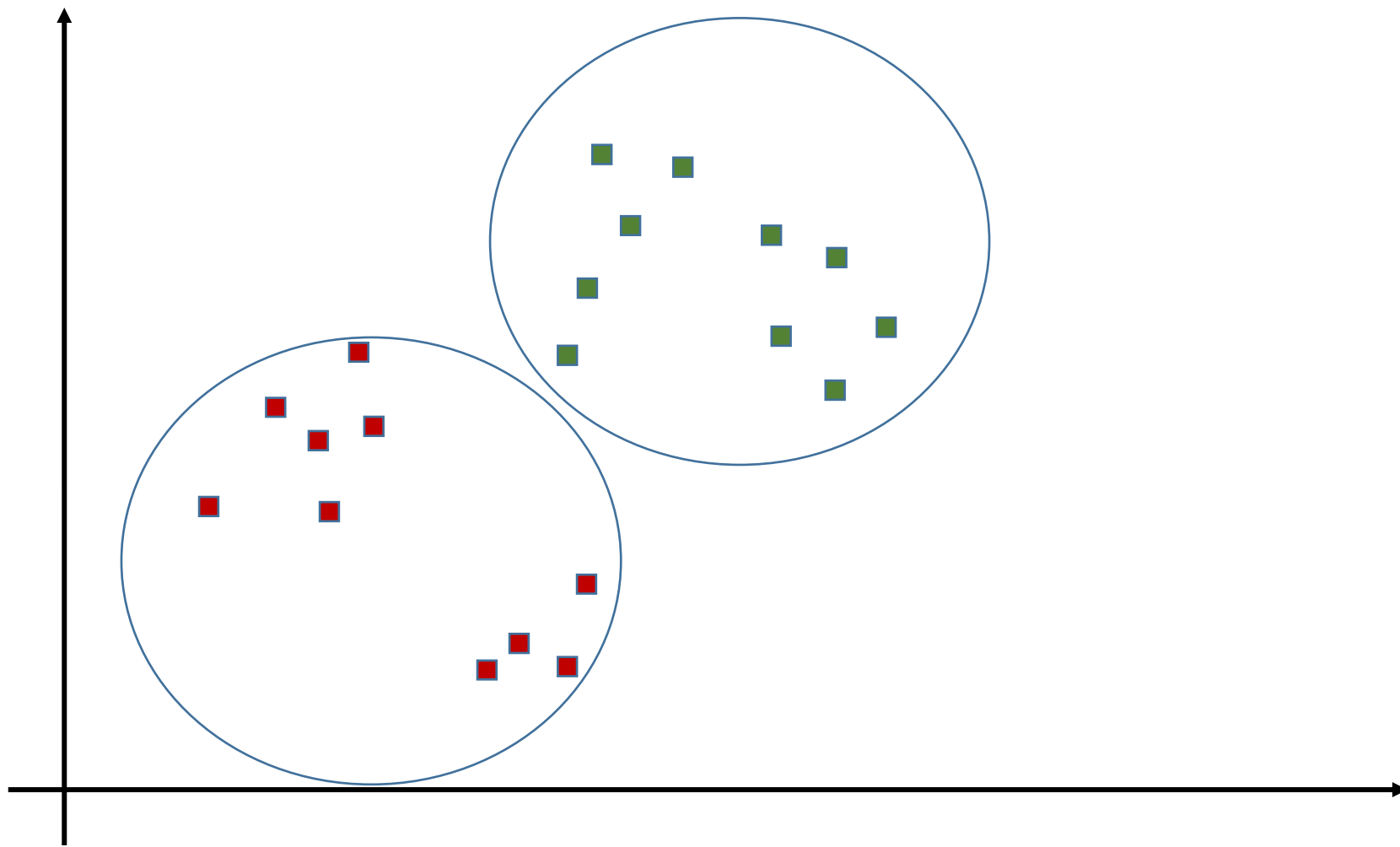


# What is clustering



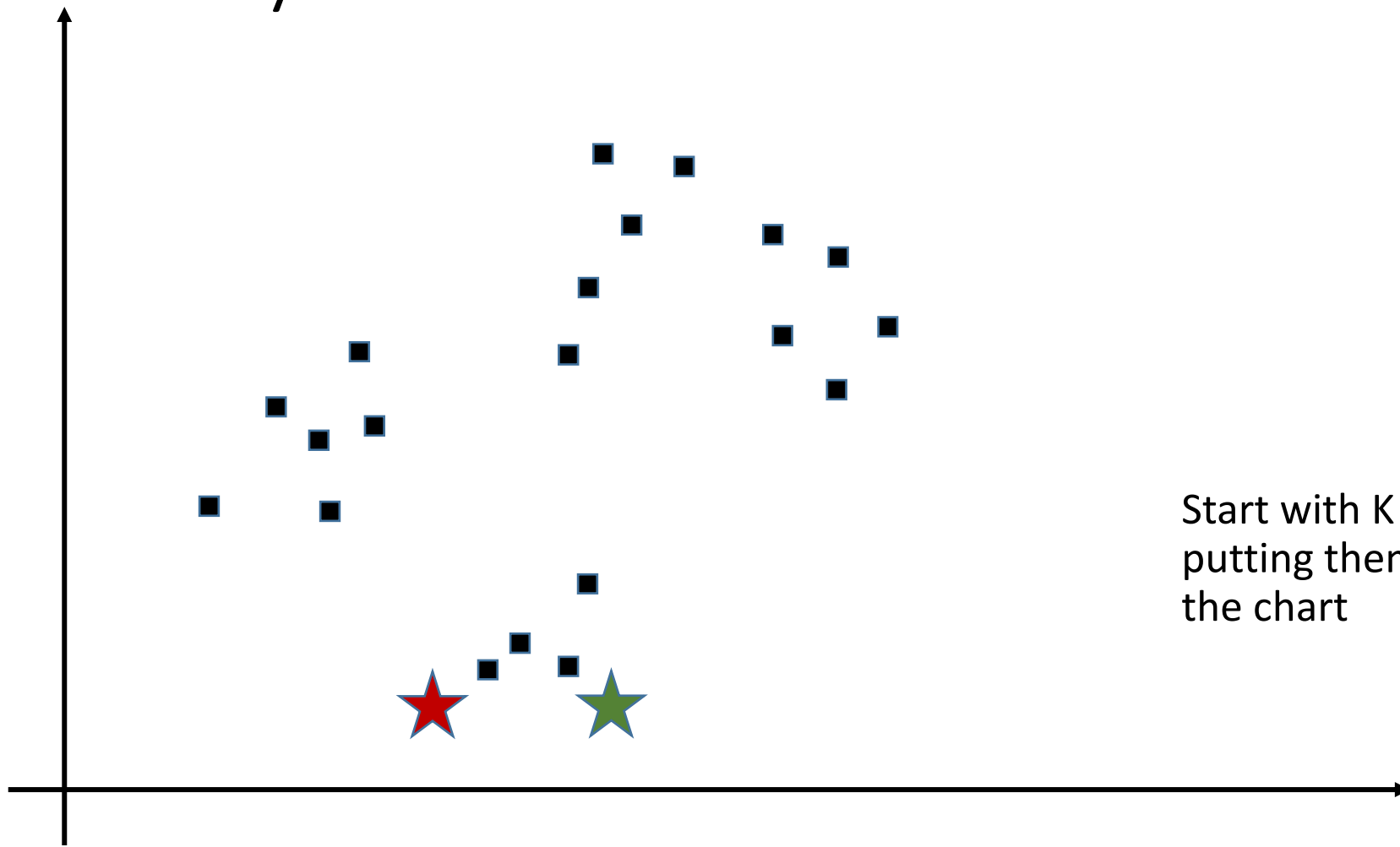


# How many clusters



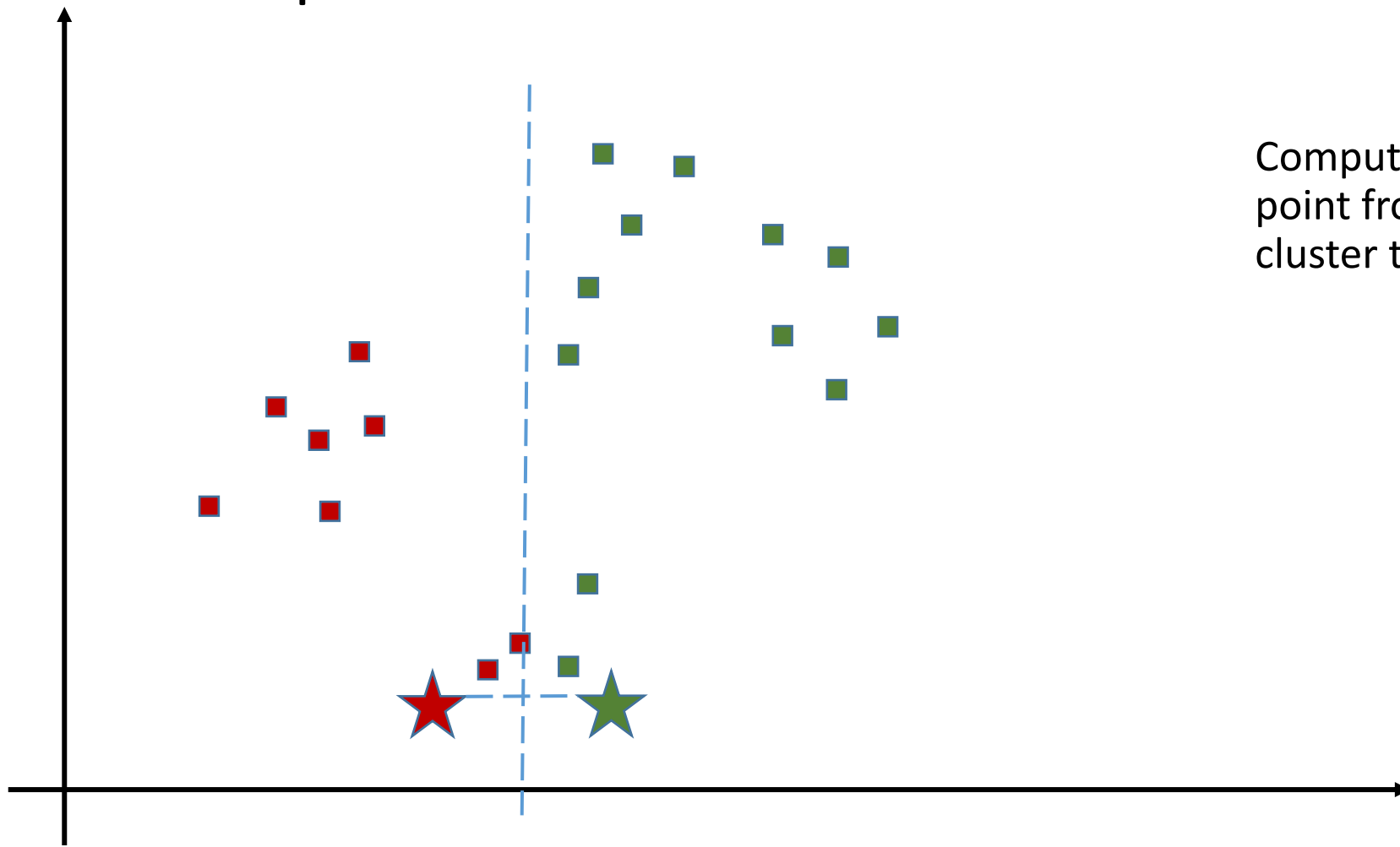


# How many centroids





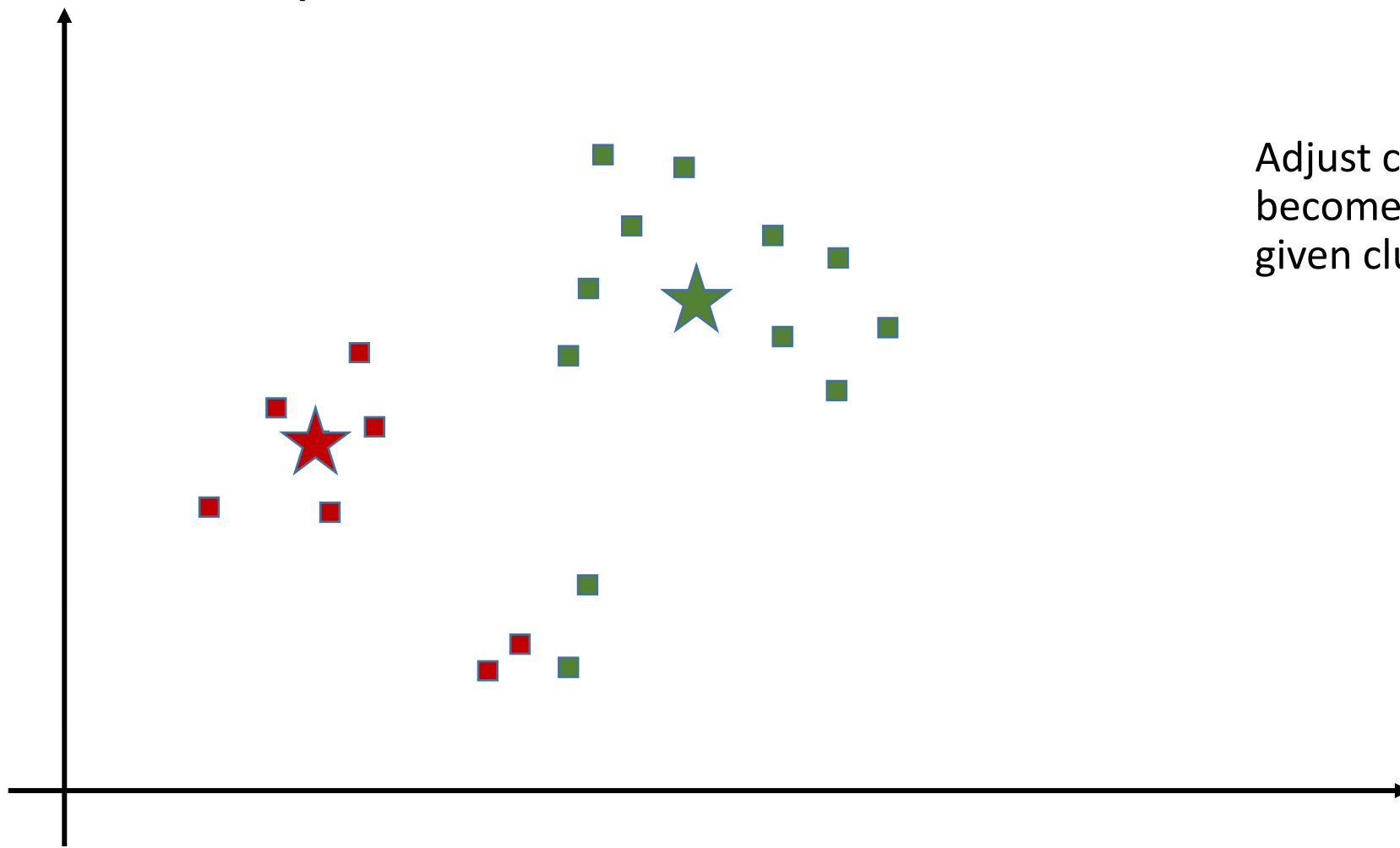
# How to improve



Compute distance of every point from centroid and cluster them accordingly



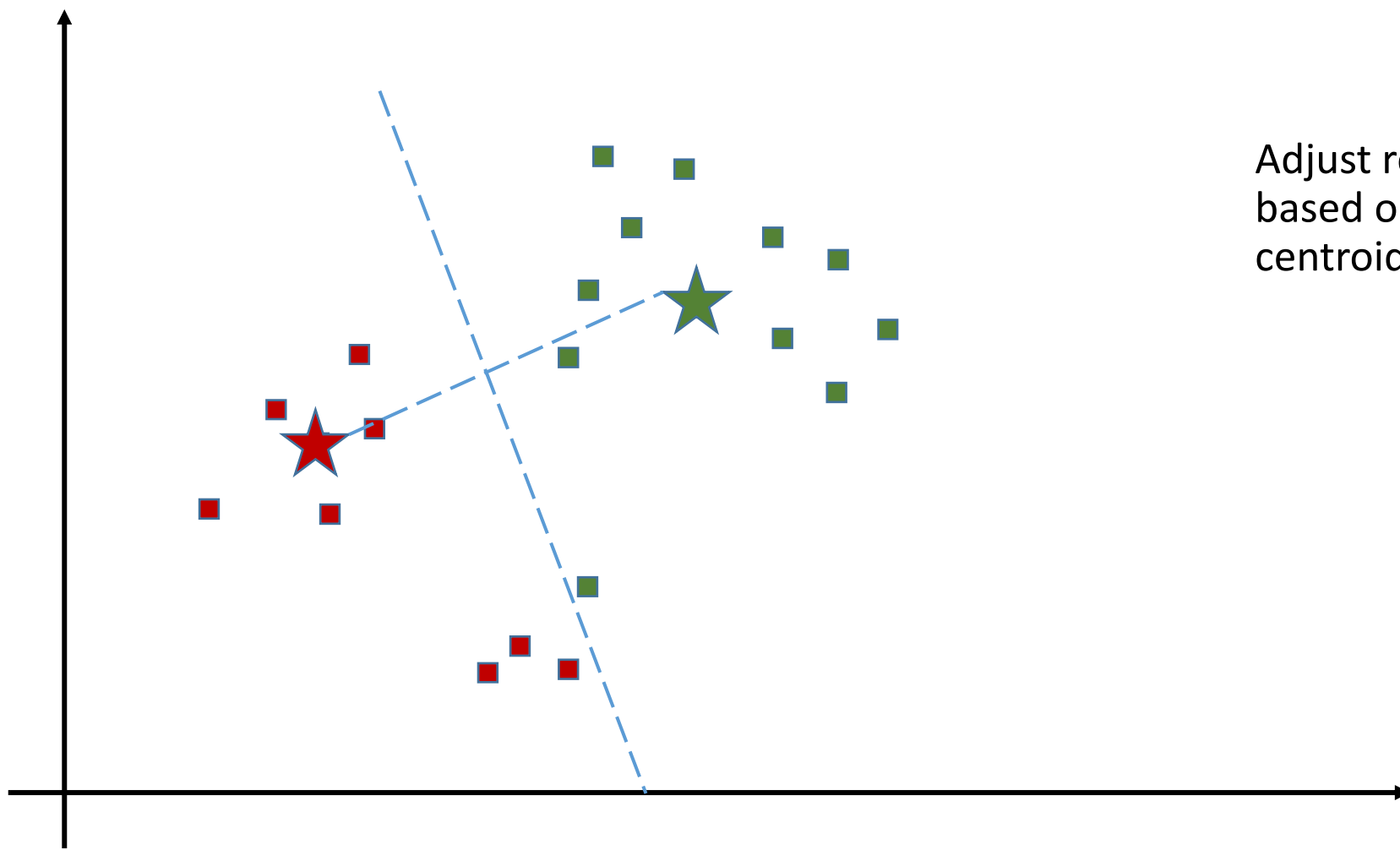
# How to improve



Adjust centroids so that they become center of gravity for given cluster



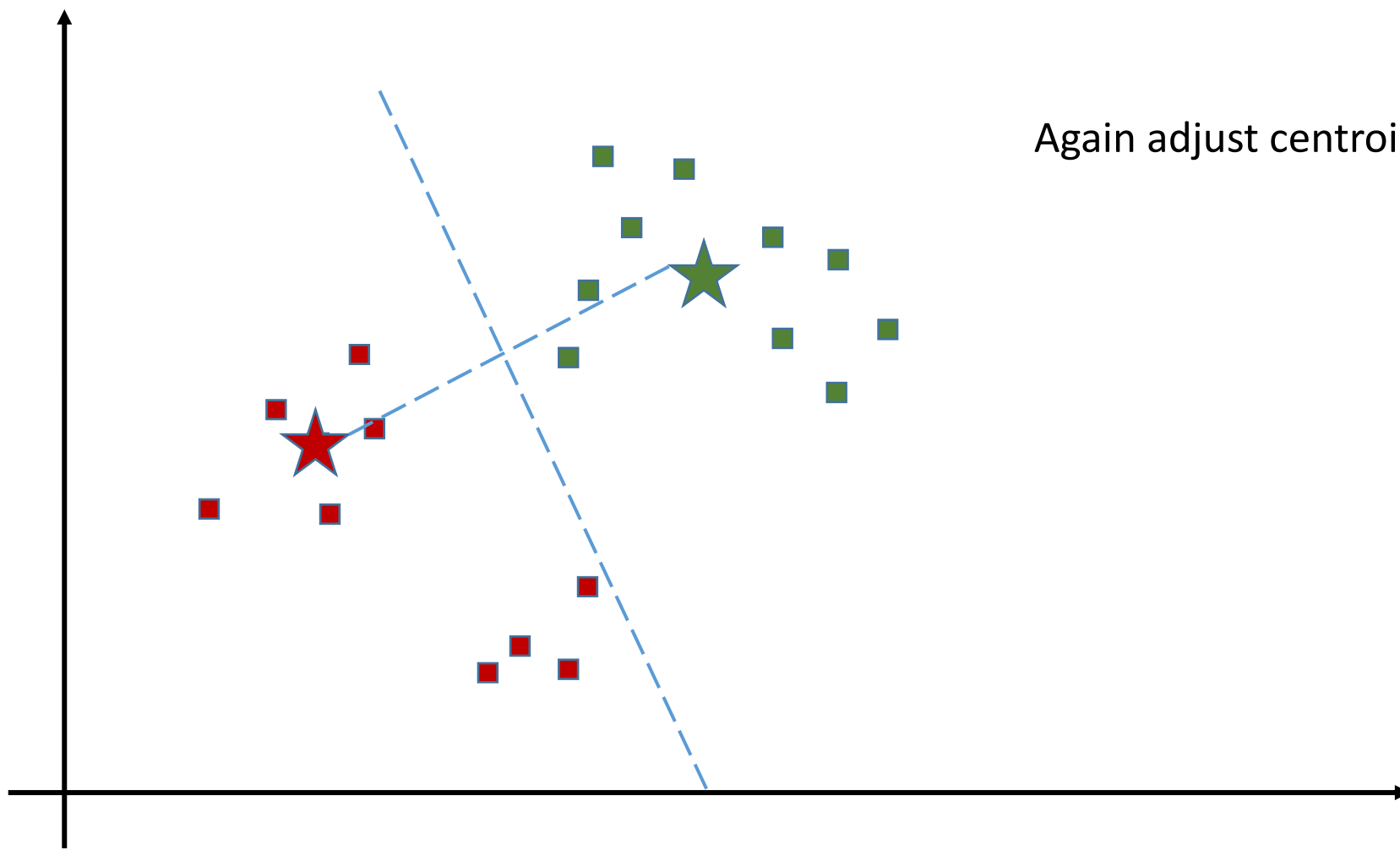
# Re-clustering



Adjust re-cluster every point based on their distance with centroid.

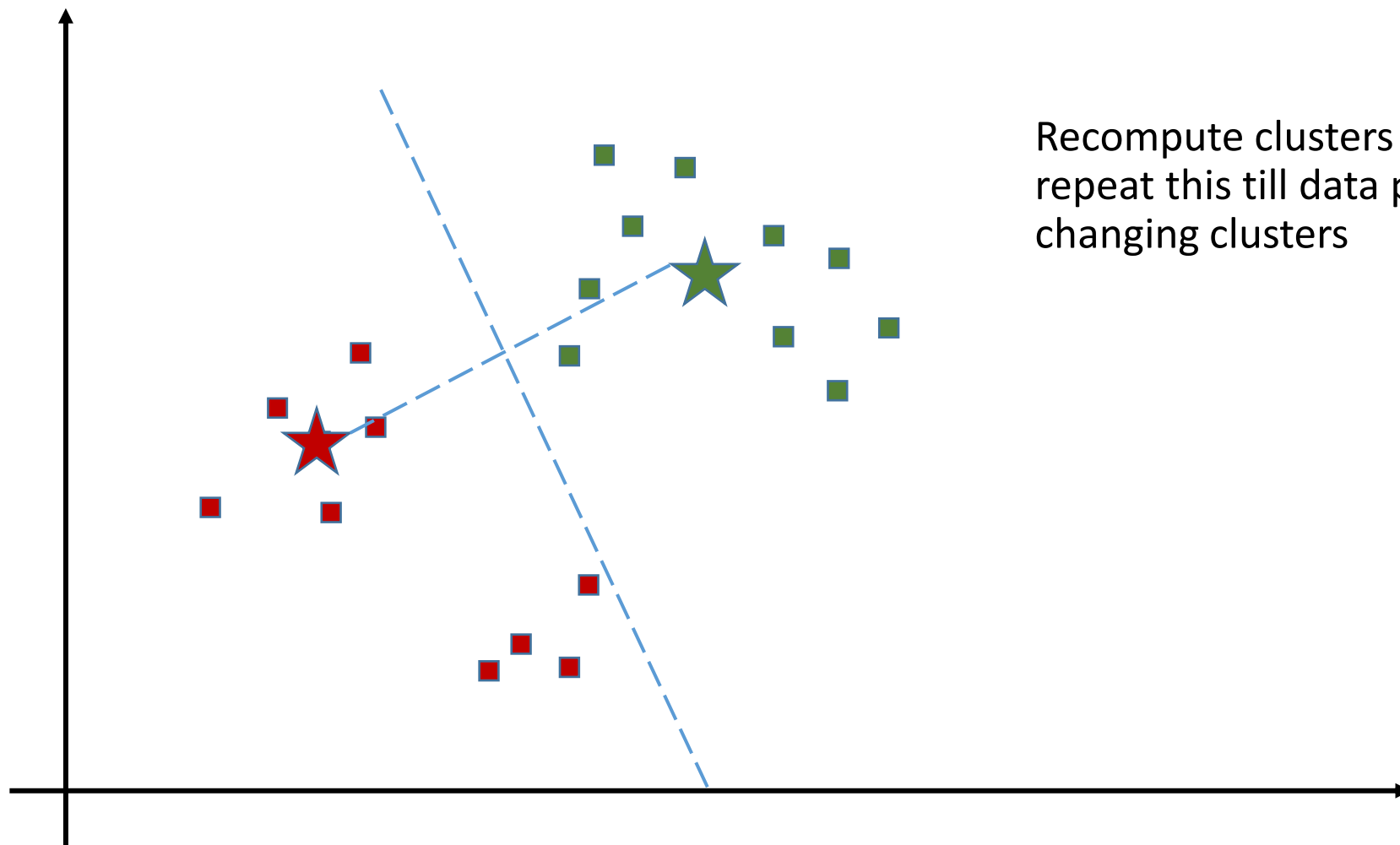


# Re-clustering





# Re-clustering

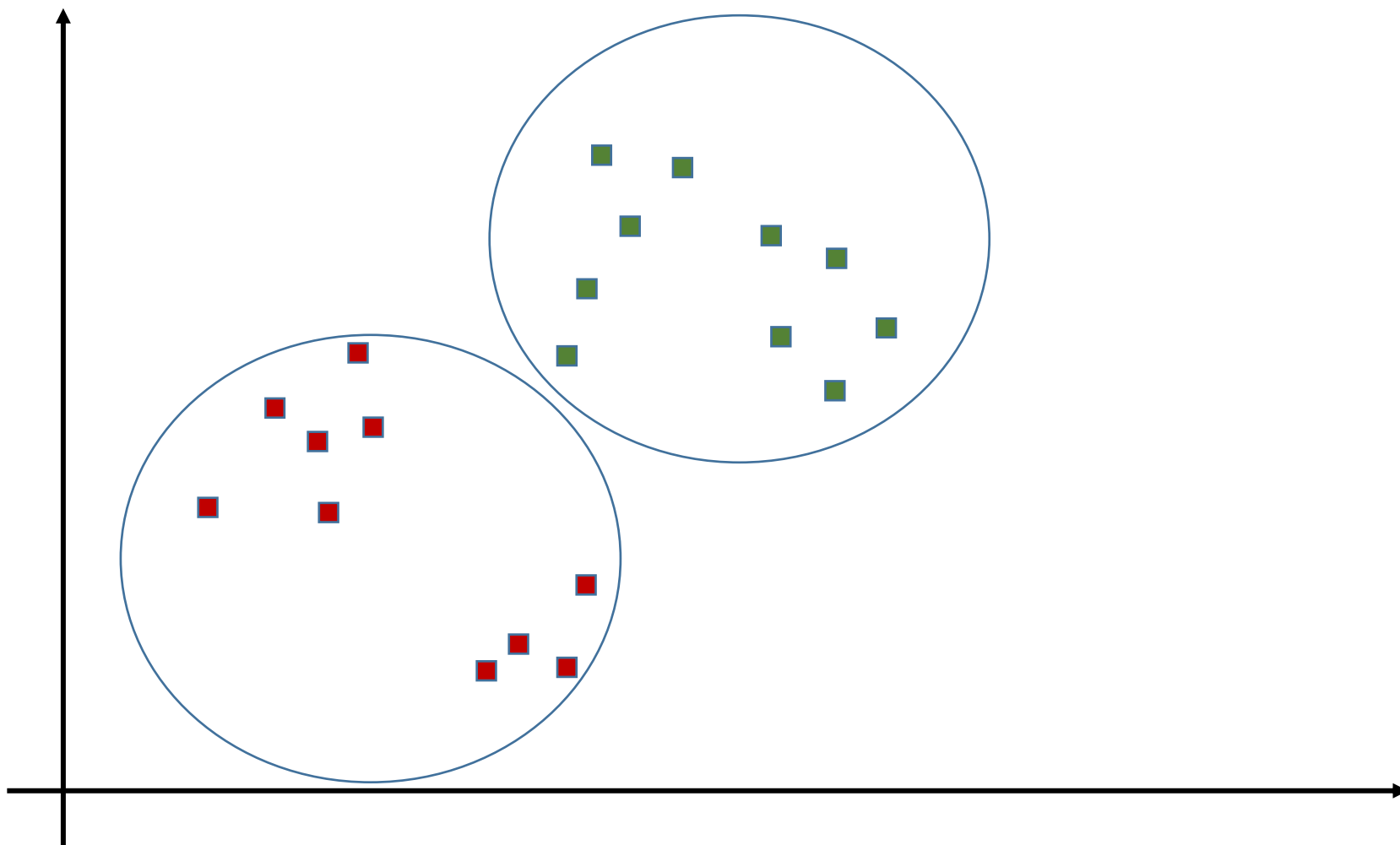


Recompute clusters and  
repeat this till data points stop  
changing clusters





# Re-clustering



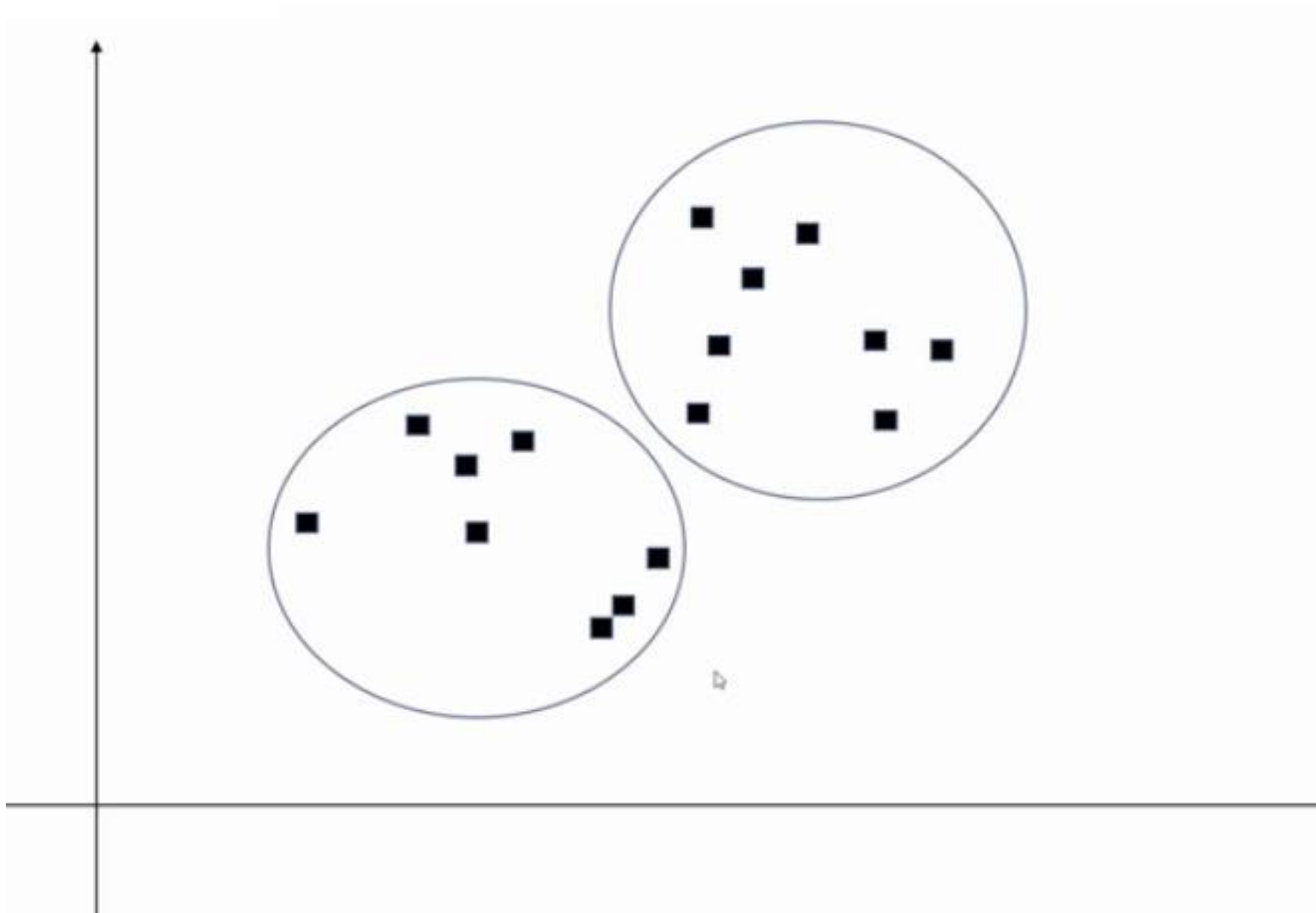


**How to determine correct  
number of clusters (k)?**



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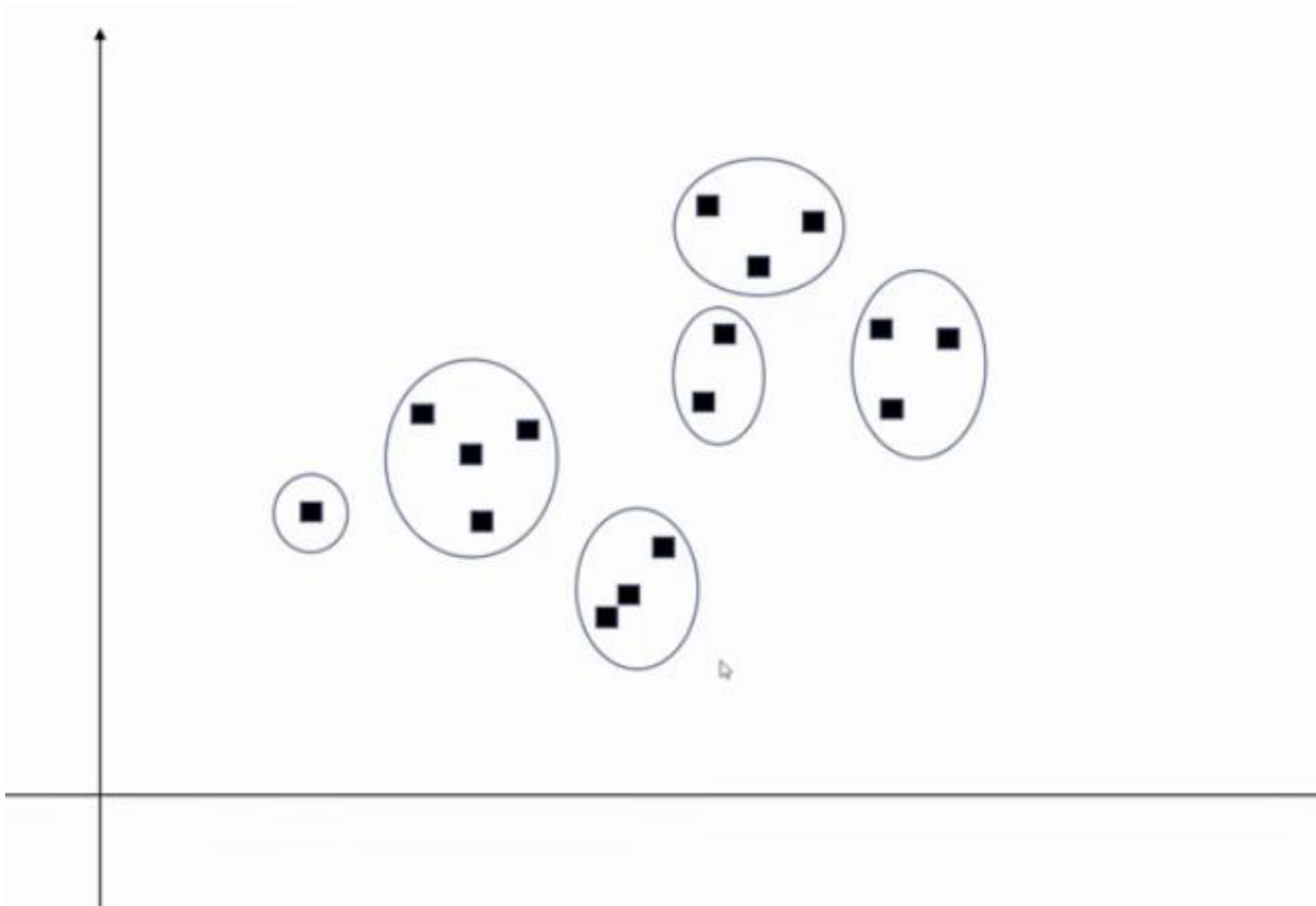
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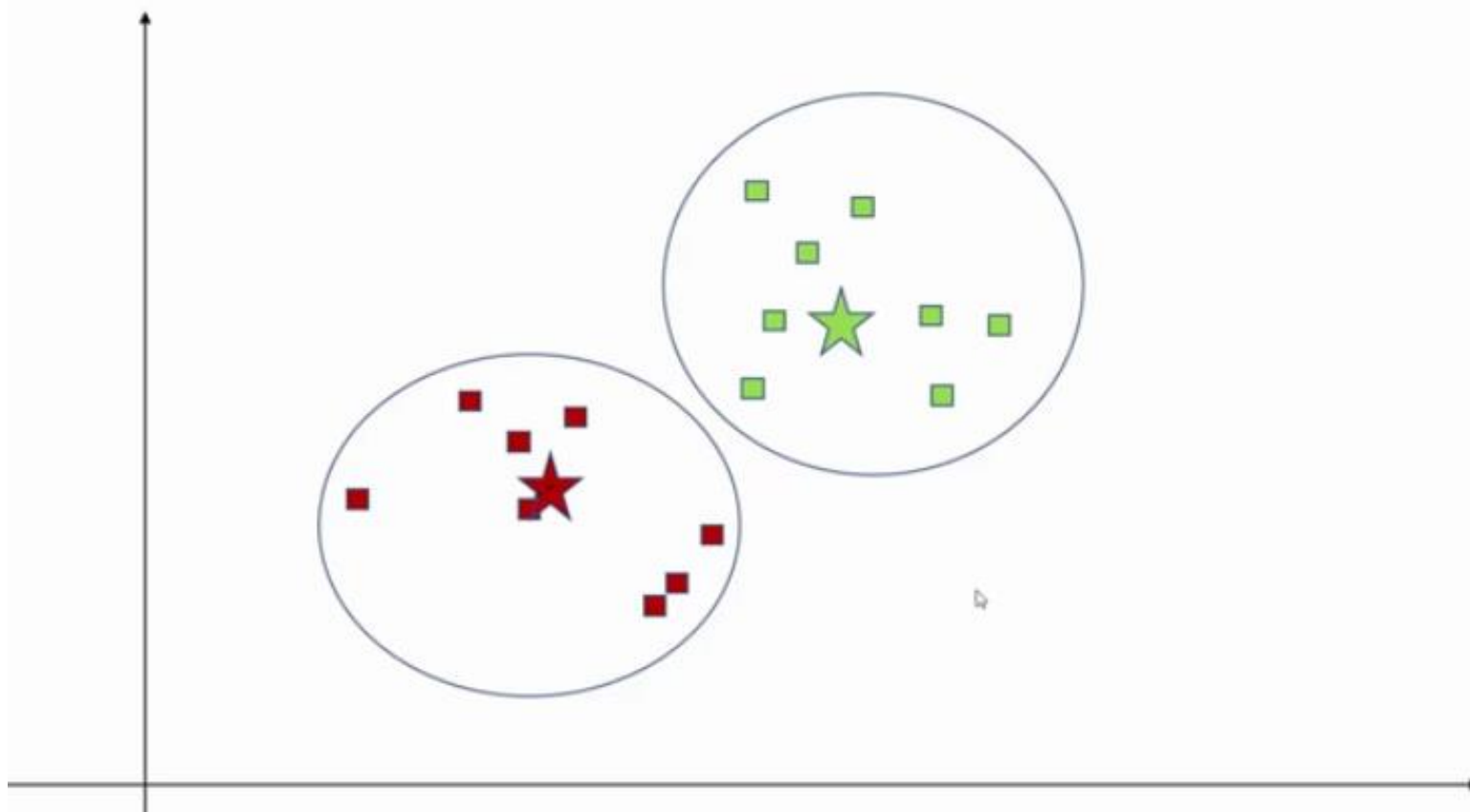
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# Sum of Squared Errors

SSE = Sum of Squared Errors



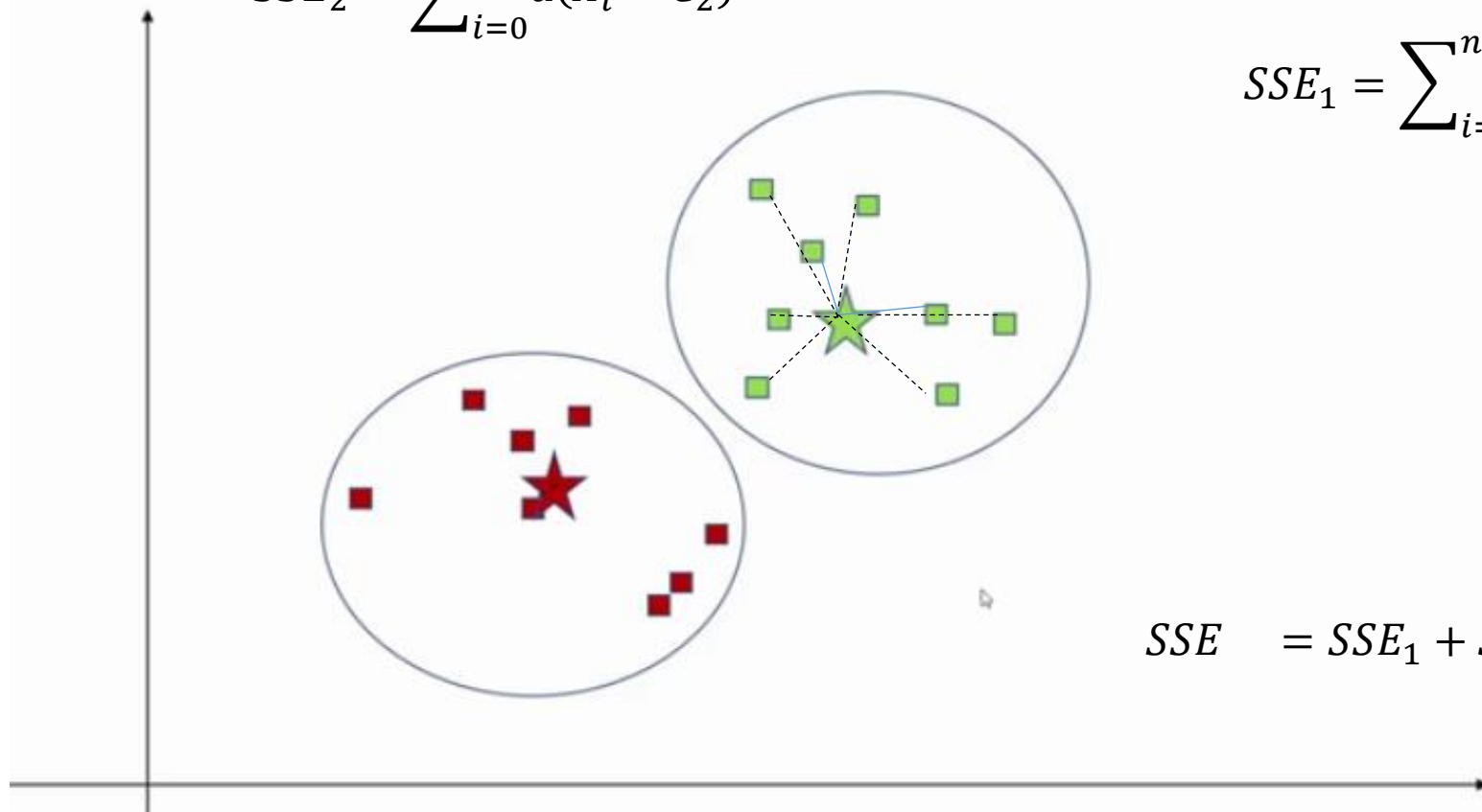


# Sum of Squared Errors

$$SSE_2 = \sum_{i=0}^m d(X_i - C_2)^2$$

SSE = Sum of Squared Errors

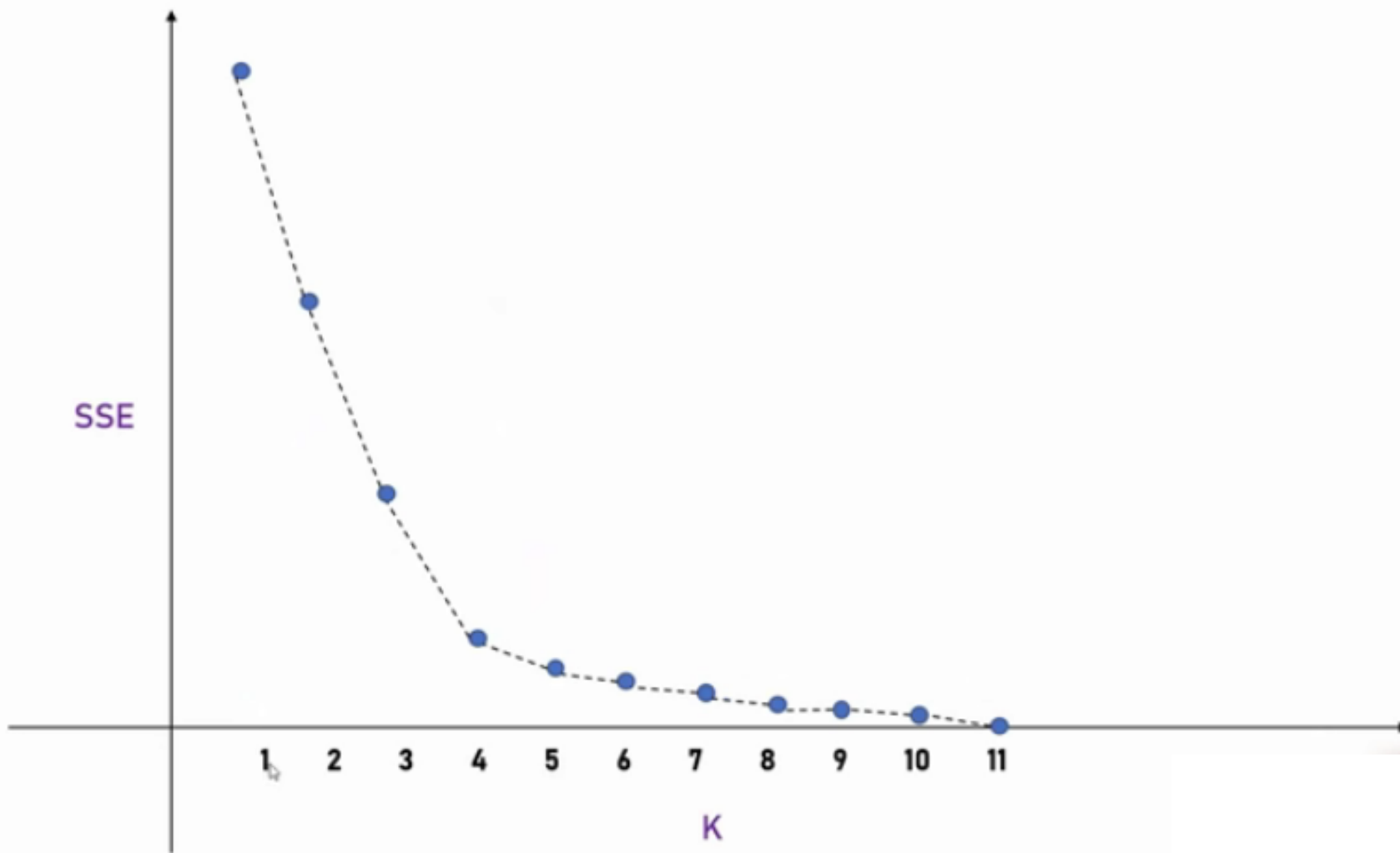
$$SSE_1 = \sum_{i=0}^n d(X_i - C_1)^2$$



$$SSE = SSE_1 + SSE_2 + SSE_3 + \dots + SSE_K$$



# Elbow Technique





# Code

```
import matplotlib.pyplot as plt
import pandas as pd
from sklearn.datasets import make_blobs
from sklearn.cluster import KMeans
from sklearn.metrics import silhouette_score
from sklearn.preprocessing import MinMaxScaler
%matplotlib inline
```





## Code

```
df = pd.read_csv("income.csv")  
df.head()
```

```
plt.scatter(df['Age'], df['Income($)'])
```

```
km = KMeans(n_clusters=3)
```

```
Km
```

```
y_predict = km.fit_predict(df[['Age', 'Income($)']])  
y_predict
```



# Code

```
df['cluster'] = y_predict  
df.tail()
```

```
df1 = df[df.cluster==0]  
df2 = df[df.cluster==1]  
df3 = df[df.cluster==2]
```

```
plt.scatter(df1['Age'], df1['Income($)', color = 'green')  
plt.scatter(df2['Age'], df2['Income($)', color = 'red')  
plt.scatter(df3['Age'], df3['Income($)', color = 'black')
```

```
plt.xlabel('Age')  
plt.ylabel('Income($)')  
plt.legend()
```



# Code

```
scaler = MinMaxScaler()  
features = ['Income($)']  
scaler.fit(df[features])  
df['Income($)'] = scaler.transform(df[features])
```

```
age = ['Age']  
scaler.fit(df[age])  
df.Age = scaler.transform(df[age])  
df
```



## Code

```
km = KMeans(n_clusters=3)
y_predict = km.fit_predict(df[['Age', 'Income($)']])
y_predict
```

```
df['cluster'] = y_predict
Df
```

```
km.cluster_centers_ # second run
```



# Code

```
df1 = df[df.cluster==0]
df2 = df[df.cluster==1]
df3 = df[df.cluster==2]

plt.scatter(df1['Age'], df1['Income($)', color = 'green')
plt.scatter(df2['Age'], df2['Income($)', color = 'red')
plt.scatter(df3['Age'], df3['Income($)', color = 'black')
plt.scatter(km.cluster_centers_[ :,0], km.cluster_centers_[ :,1], color = 'purple', marker='*',
label='centroid') # second run
plt.xlabel('Age')
plt.ylabel('Income($)')
plt.legend()
```



## Code

```
k_rng = range(1,10)
sse = []
for k in k_rng:
    km = KMeans(n_clusters = k)
    km.fit(df[['Age', 'Income($)']])
    sse.append(km.inertia_)
sse
```



# Code

```
plt.xlabel('x')  
plt.ylabel('sum of squared errors')  
plt.plot(k_rng,sse)
```

**<https://www.kaggle.com/patelprashant/employee-attrition>**





# Unsupervised Machine Learning (K-Means)

The k-means clustering method is an unsupervised machine learning technique used to identify clusters of data objects in a dataset. There are many different types of clustering methods, but k-means is one of the oldest and most approachable.

- What k-means clustering is
- When to use k-means clustering to analyze your data
- How to implement k-means clustering in Python with scikit-learn
- How to select a meaningful number of clusters



# Unsupervised Machine Learning (K-Means)

## What Is Clustering?

**Clustering is a set of techniques used to partition data into groups, or clusters. Clusters are loosely defined as groups of data objects that are more similar to other objects in their cluster than they are to data objects in other clusters.**

**Clustering helps identify two qualities of data:**

- **Meaningfulness**
- **Usefulness**



# Unsupervised Machine Learning (K-Means)

## What Is Clustering?

**Meaningful** clusters expand domain knowledge. For example, in the medical field, researchers applied clustering to gene expression experiments. The clustering results identified groups of patients who respond differently to medical treatments.

**Useful** clusters, on the other hand, serve as an intermediate step in a data pipeline. For example, businesses use clustering for customer segmentation. The clustering results segment customers into groups with similar purchase histories, which businesses can then use to create targeted advertising campaigns



# Unsupervised Machine Learning (K-Means)

## Overview of Clustering Techniques ?

Important factors that effects the selection of cluster algorithm include:

- Characteristics of the clusters
- Features of the dataset
- Number of outliers
- Number of data objects, etc

Three popular categories of clustering algorithms:

- Partitional clustering
- Hierarchical clustering
- Density-based clustering



# Unsupervised Machine Learning (K-Means)

## Overview of Clustering Techniques ?

*Non deterministic process*

### Partitional Clustering

**Partitional clustering** divides data objects into nonoverlapping groups. In other words, no object can be a member of more than one cluster, and every cluster must have at least one object.

These techniques require the user to specify the number of clusters, indicated by the variable  **$k$** . Many partitional clustering algorithms work through an iterative process to assign subsets of data points into  $k$  clusters. Two examples of partitional clustering algorithms are  **$k$ -means** and  **$k$ -medoids**.



# Unsupervised Machine Learning (K-Means)

## Overview of Clustering Techniques ?

### Partitional Clustering

#### Advantages:

They work well when clusters have a **spherical shape**.

They're **scalable** with respect to algorithm complexity.

#### Disadvantages:

They're not well suited for clusters with **complex shapes** and different sizes.

They break down when used with clusters of different **densities**.



# Unsupervised Machine Learning (K-Means)

## Overview of Clustering Techniques ?

*deterministic process*

### Hierarchical Clustering

**Hierarchical clustering** determines cluster assignments by building a hierarchy. This is implemented by either a bottom-up or a top-down approach:

These methods produce a tree-based hierarchy of points called a **dendrogram**. Similar to partitional clustering, in hierarchical clustering the number of clusters ( $k$ ) is often predetermined by the user. Clusters are assigned by cutting the dendrogram at a specified depth that results in  $k$  groups of smaller dendrograms.

**Agglomerative clustering** is the bottom-up approach. It merges the two points that are the most similar until all points have been merged into a single cluster.

**Divisive clustering** is the top-down approach. It starts with all points as one cluster and splits the least similar clusters at each step until only single data points remain.



# Unsupervised Machine Learning (K-Means)

## Overview of Clustering Techniques ?

### Hierarchical Clustering

#### Advantages

They often reveal the finer details about the **relationships** between data objects.  
They provide an **interpretable dendrogram**.

#### Disadvantages

They're **computationally expensive** with respect to algorithm complexity.  
They're sensitive to **noise** and **outliers**.





# Unsupervised Machine Learning (K-Means)

## Overview of Clustering Techniques ?

### Density-based clustering

*deterministic process*

**Density-based clustering** determines cluster assignments based on the density of data points in a region. Clusters are assigned where there are high densities of data points separated by low-density regions.

Unlike the other clustering categories, this approach doesn't require the user to specify the number of clusters. Instead, there is a distance-based parameter that acts as a tunable threshold. This threshold determines how close points must be to be considered a cluster member.

### Examples:

**DBSCAN:** Density-Based Spatial Clustering of Applications with Noise

**OPTICS:** Ordering Points To Identify the Clustering Structure



# Unsupervised Machine Learning (K-Means)

## Overview of Clustering Techniques ?

### Density-based clustering

#### Advantages

- They excel at identifying clusters of **nonspherical shapes**.
- They're resistant to **outliers**.

#### Disadvantages

- They aren't well suited for clustering in **high-dimensional spaces**.
- They have trouble identifying clusters of **varying densities**.



# Unsupervised Machine Learning (K-Means)

## Understanding the K-Means Algorithm

### Understanding the K-Means Algorithm

*Non deterministic process*

- Randomly select  $k$  centroids, where  $k$  is equal to the number of clusters you choose. (**Centroids** are data points representing the center of a cluster. )
- The main element of the algorithm works by a two-step process called **expectation-maximization**.
  - The **expectation** step assigns each data point to its nearest centroid.
  - The **maximization** step computes the mean of all the points for each cluster and sets the new centroid.



# Unsupervised Machine Learning (K-Means)

## Understanding the K-Means Algorithm

### Algorithm 1: K – means algorithm

- Specify the number  $k$  of clusters to assign
- Randomly initialize  $k$  centroids
- **Repeat:**
  - **Expectation:** Assign each point to its closest centroid.
  - **Maximization:** Compute the new centroid mean for each cluster
- **Untill:** The centroid positions do not change

*Non deterministic process*

The quality of the cluster assignments is determined by computing the **sum of the squared error (SSE)** after the centroids **converge**, or match the previous iteration's assignment. The SSE is defined as the sum of the squared Euclidean distances of each point to its closest centroid. Since this is a measure of error, the objective of  $k$ -means is to try to minimize this value.



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[file:///D:/Prime%20Intuit/centroids\\_iterations.webp](file:///D:/Prime%20Intuit/centroids_iterations.webp)



# Unsupervised Machine Learning (K-Means)

## K-Means Clustering Code in Python

- conda install -c conda-forge kneed

```
(base) PS C:\Users\LENO  
Collecting package meta  
Solving environment: do
```

```
## Package Plan ##
```

```
environment location: C:\Users\LENOVO\anaconda3
```

```
added / updated specs:  
- kneed
```

```
The following packages will be downloaded:
```

| package        | build          |        |             |
|----------------|----------------|--------|-------------|
| conda-4.10.3   | py38haa244fe_0 | 3.1 MB | conda-forge |
| kneed-0.7.0    | pyh9f0ad1d_0   | 12 KB  | conda-forge |
| python_abi-3.8 | 2_cp38         | 4 KB   | conda-forge |
| Total:         |                | 3.1 MB |             |

```
The following NEW packages will be INSTALLED:
```

|            |  |
|------------|--|
| kneed      | conda-forge/noarch::kneed-0.7.0-pyh9f0ad1d_0 |
| python_abi | conda-forge/win-64::python_abi-3.8-2_cp38    |

```
The following packages will be UPDATED:
```

|       |   |
|-------|---|
| conda | pkgs/main::conda-4.10.1-py38haa95532_1 --> conda-forge::conda-4.10.3-py38haa244fe_0 |
|-------|---|

```
Proceed ([y]/n)?
```



# Unsupervised Machine Learning (K-Means)

## Writing your K-Means Algorithm code

### Import the modules needed for the code

```
import matplotlib.pyplot as plt
from kneed import KneeLocator
from sklearn.datasets import make_blobs
from sklearn.cluster import KMeans
from sklearn.metrics import silhouette_score
from sklearn.preprocessing import StandardScaler
```



# Unsupervised Machine Learning (K-Means)

## Writing your K-Means Algorithm code

### Generate data using `make_blobs()` function

We can generate the data using `make_blobs()`, a convenience function in scikit-learn used to generate synthetic clusters. `make_blobs()` uses these parameters:

**`n_samples`** is the total number of samples to generate.

**`centers`** is the number of centers to generate.

**`cluster_std`** is the standard deviation.

**`make_blobs()`** returns a tuple of two values:

A two-dimensional NumPy array with the x- and y-values for each of the samples

A one-dimensional NumPy array containing the cluster labels for each sample





**PRIME INTUIT**

Finishing School

# Unsupervised Machine Learning (K-Means)

Writing your K-Means Algorithm code

# KNN Classifications



# K Nearest Neighbor Classifier

## Understanding the K Nearest Neighbour Algorithm

### **Purpose:**

The purpose of the k Nearest Neighbor (KNN) algorithm is to use a database in which the data points are separated into several separate classes to predict the classification of a new sample point.

### **Algorithm:**

The algorithm can be summarized as:

1. A positive integer  $k$  is specified, along with a new sample
2. We select the  $k$  entries in our database which are closest to the new sample
3. We find the most common classification of these entries
4. This is the classification we give to the new sample



# K Nearest Neighbor Classifier

## Understanding the K Nearest Neighbour Algorithm

### Algorithm 1: K – means algorithm

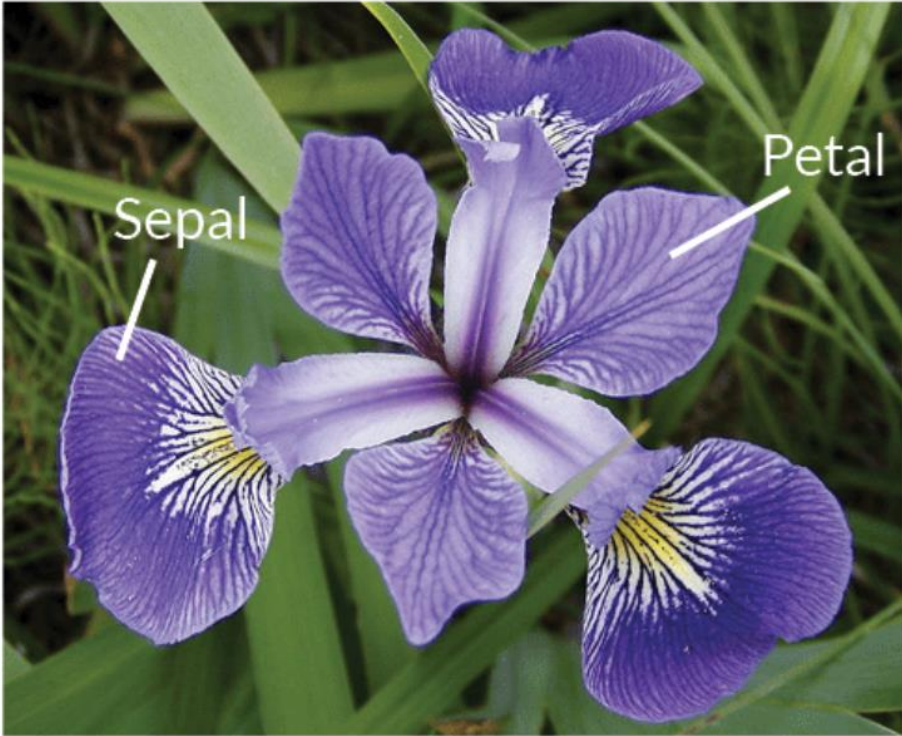
- Specify the number  $k$  of clusters to assign
- Randomly initialize  $k$  centroids
- **Repeat:**
  - **Expectation:** Assign each point to its closest centroid.
  - **Maximization:** Compute the new centroid mean for each cluster
- **Untill:** The centroid positions do not change

The quality of the cluster assignments is determined by computing the **sum of the squared error (SSE)** after the centroids **converge**, or match the previous iteration's assignment. The SSE is defined as the sum of the squared Euclidean distances of each point to its closest centroid. Since this is a measure of error, the objective of  $k$ -means is to try to minimize this value.



# K Nearest Neighbor Classifier

## Understanding the K Nearest Neighbour Algorithm



**Iris Versicolor**



**Iris Setosa**

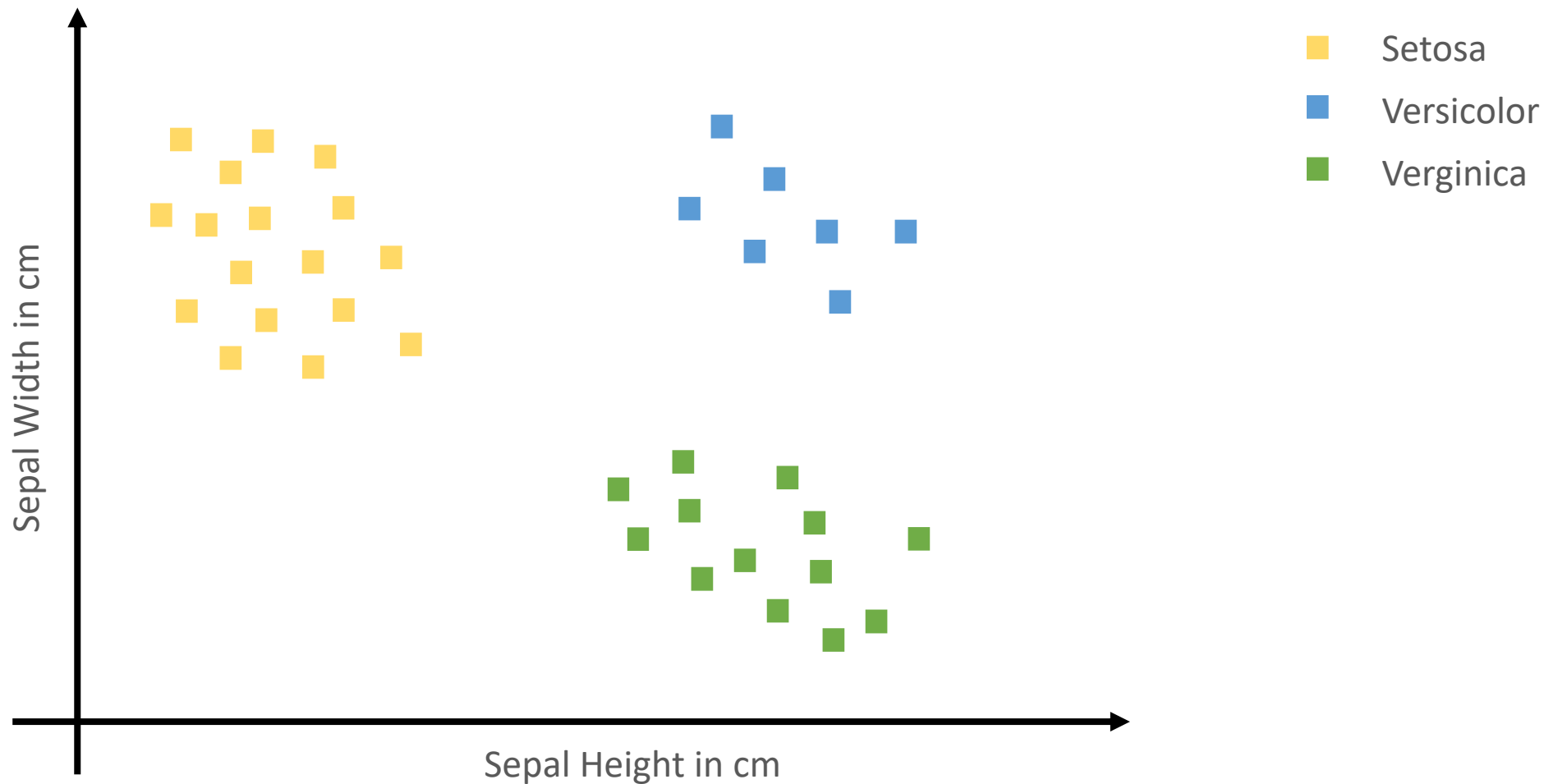


**Iris Virginica**



# K Nearest Neighbor Classifier

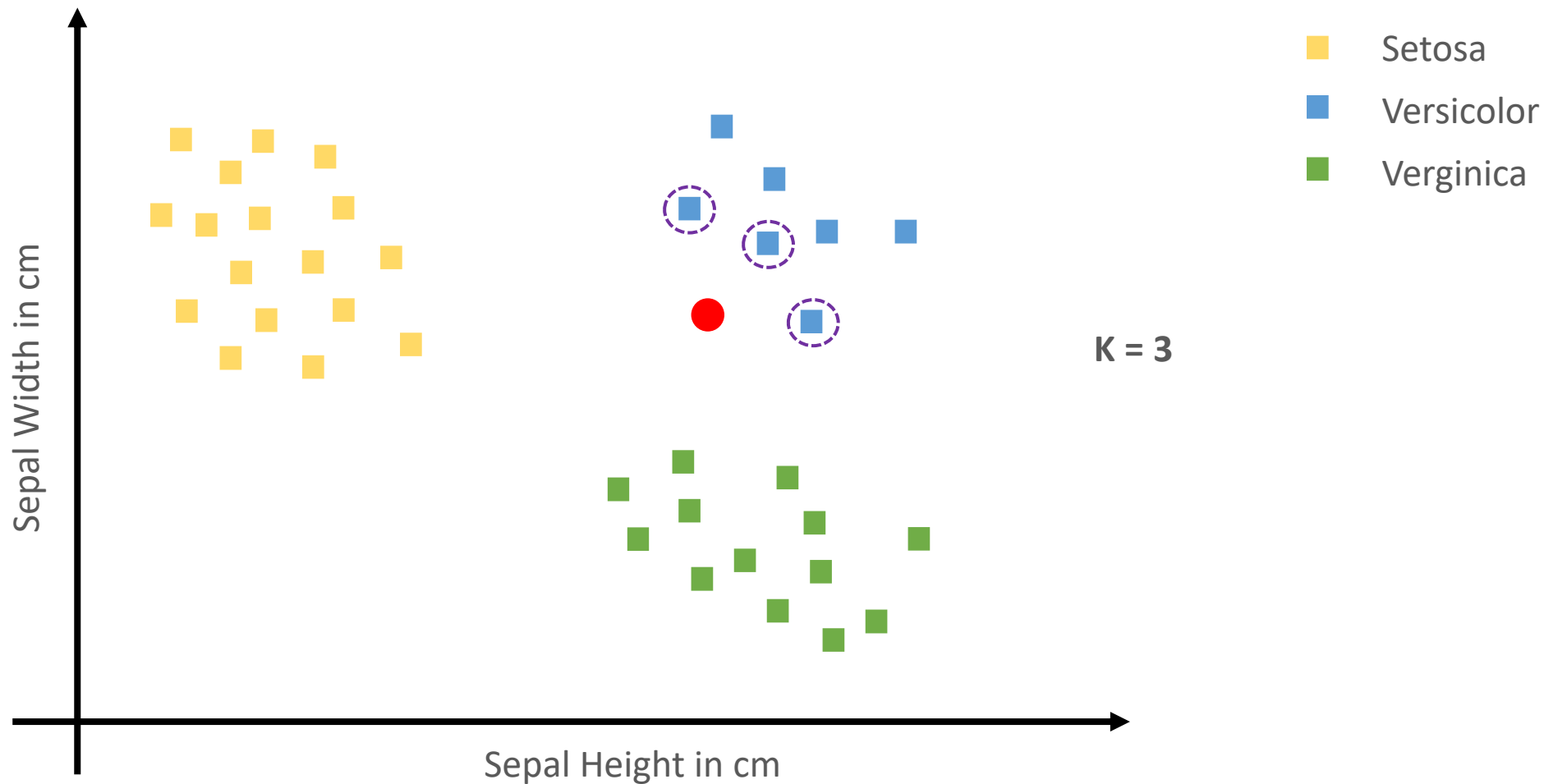
## Understanding the K Nearest Neighbour Algorithm





# K Nearest Neighbor Classifier

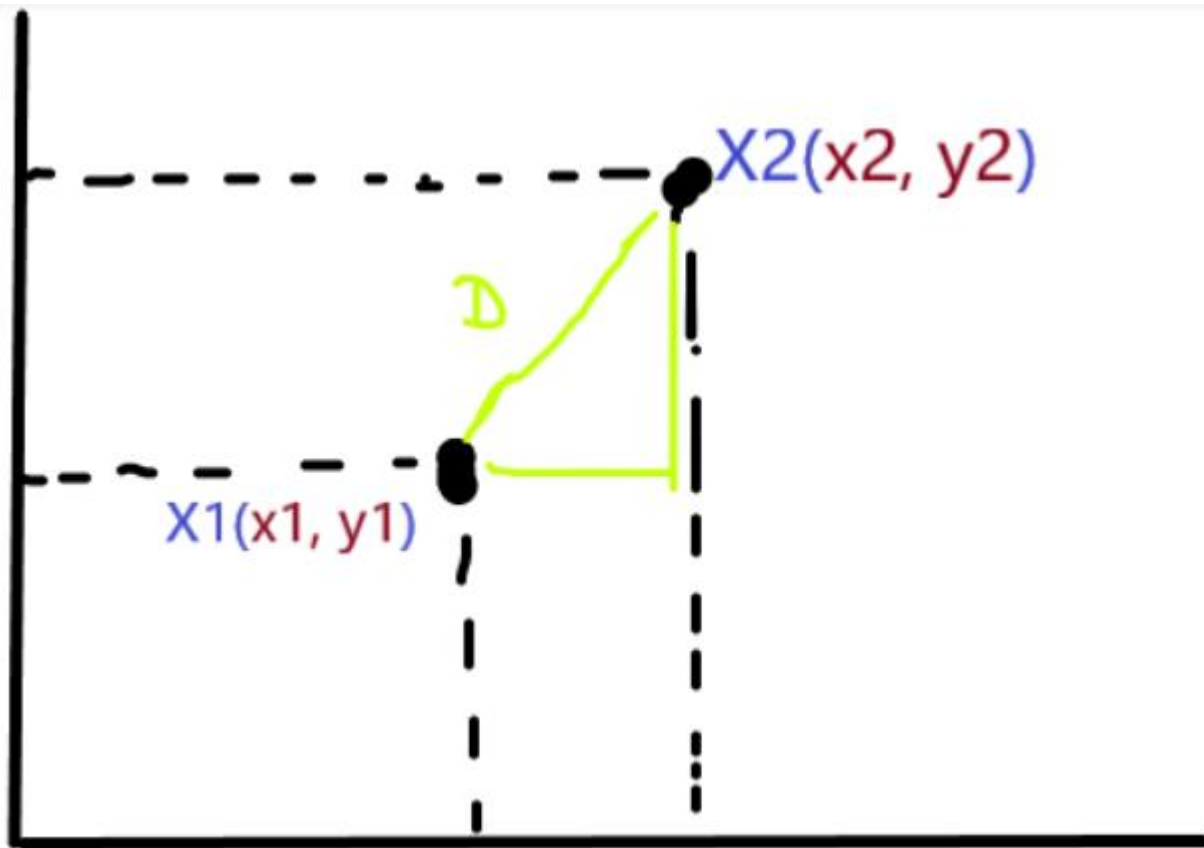
## Understanding the K Nearest Neighbour Algorithm





# K Nearest Neighbor Classifier

## Understanding the K Nearest Neighbour Algorithm - Euclidean Distance



There are 2-Dim data X1 and X2 placed at certain coordinates in 2 dimensions, suppose X1 is at (x1,y1) coordinates and X2 is at (x2,y2) coordinates. We have 2-dim data so we considered F1 and F2 two features and D is considered as the shortest line from X1 and X2, If we find the distance between these data points that can be found by the Pythagoras theorem and can be written as:

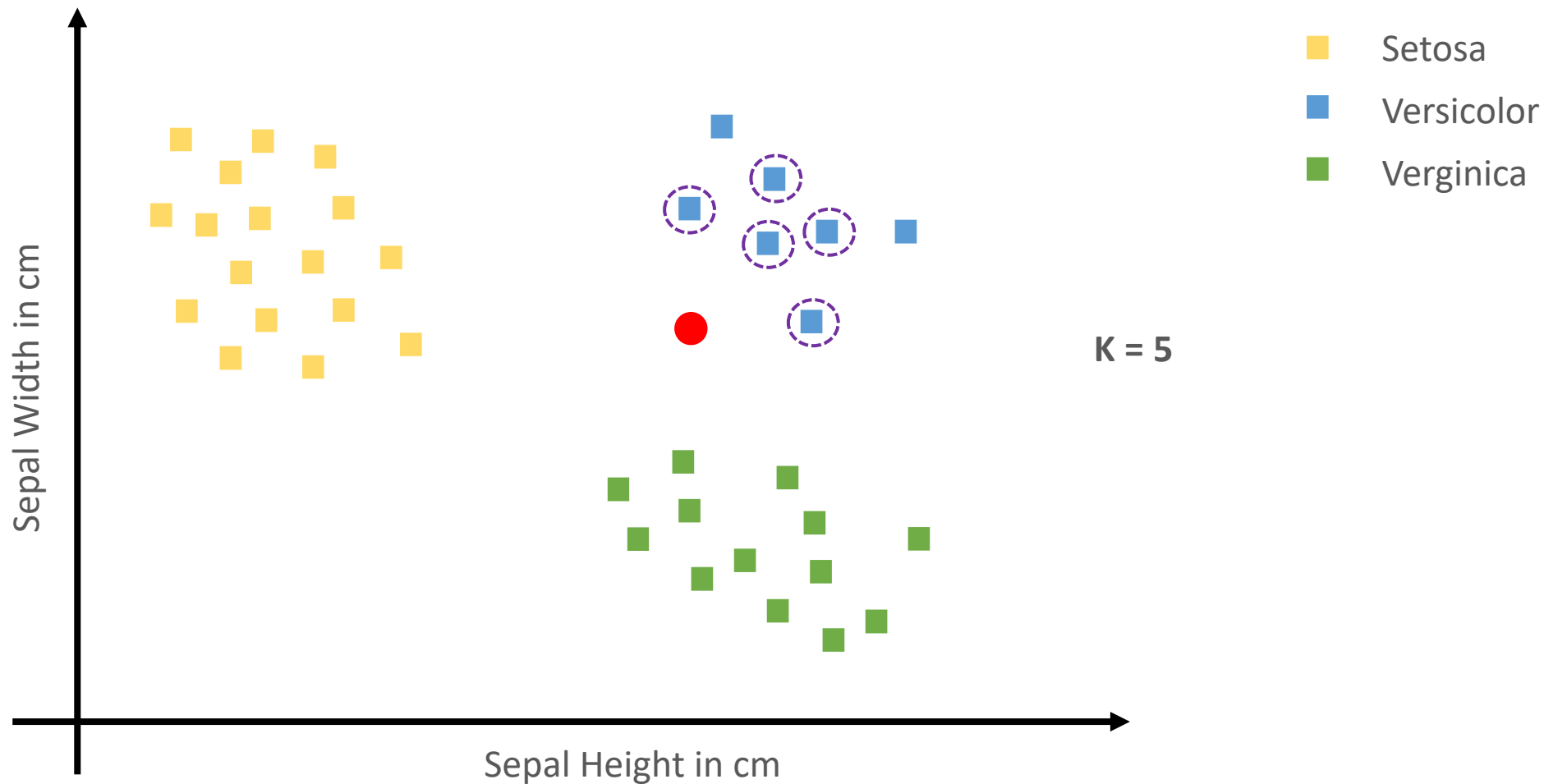
$$d = \sqrt{(x_2 - x_1)^2 + (y_2 - y_1)^2}$$





# K Nearest Neighbor Classifier

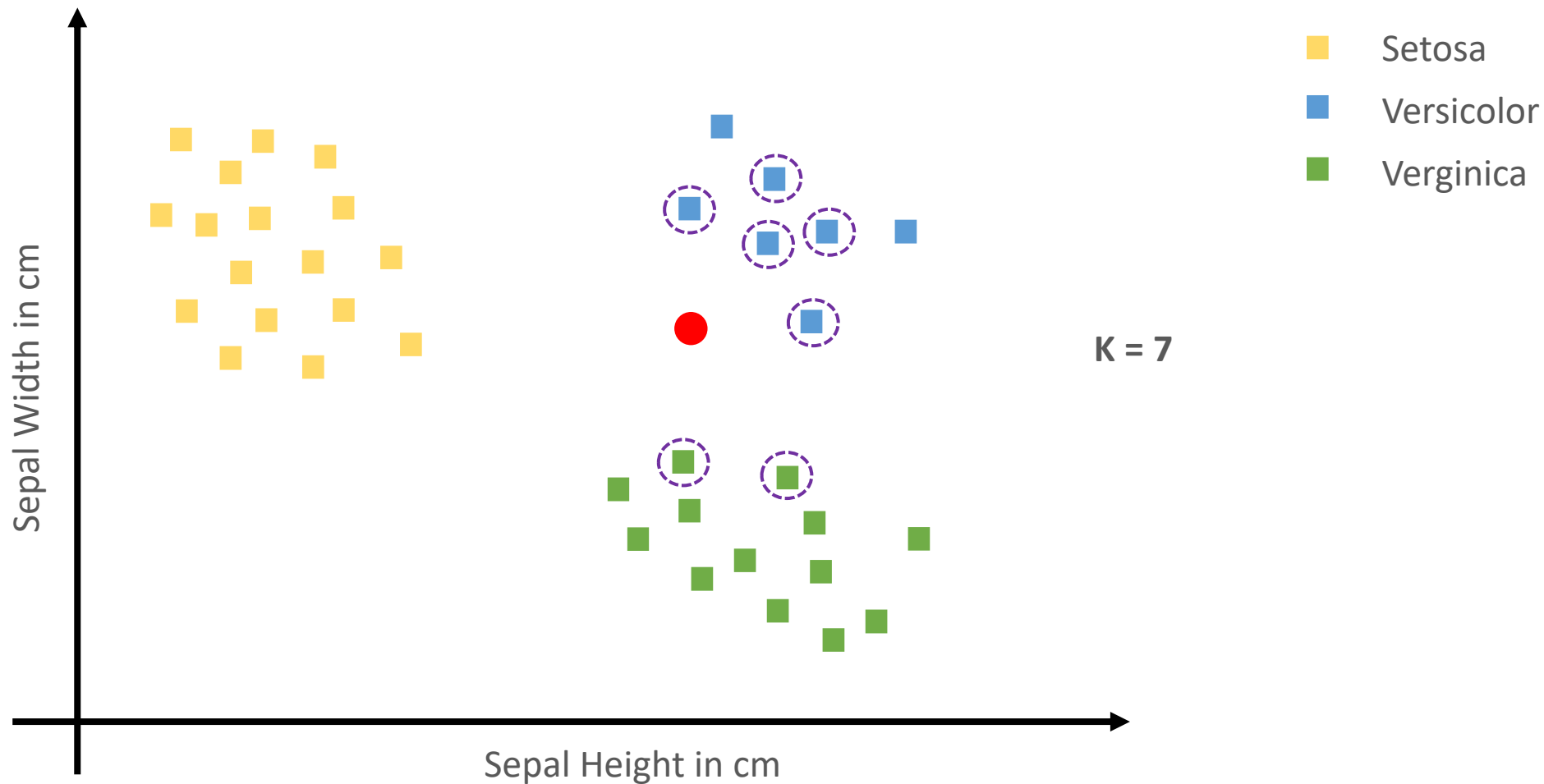
## Understanding the K Nearest Neighbour Algorithm





# K Nearest Neighbor Classifier

## Understanding the K Nearest Neighbour Algorithm





# K Nearest Neighbor Classifier

## Working with Iris data set

```
import numpy as np
import pandas as pd
df = pd.read_csv('Iris.csv')
df.head()
```

```
df.shape
df.describe()
df.groupby('Species').size()
```

# Dividing data into features and labels

```
feature_columns = ['SepalLengthCm', 'SepalWidthCm', 'PetalLengthCm', 'PetalWidthCm']
X = df[feature_columns].values
y = df['Species'].values
```



# K Nearest Neighbor Classifier

## Working with Iris data set

```
# Change Species coloumn to numbers
```

```
from sklearn.preprocessing import LabelEncoder  
le = LabelEncoder()  
y = le.fit_transform(y)
```

```
# Split data into training and test
```

```
from sklearn.model_selection import train_test_split  
X_train, X_test, y_train, y_test = train_test_split(X, y, test_size = 0.2, random_state = 0)
```



# K Nearest Neighbor Classifier

## Working with Iris data set

# Data Visualisation

```
import matplotlib.pyplot as plt  
import seaborn as sns  
%matplotlib inline
```

# Pair Plot

```
plt.figure()  
sns.pairplot(df.drop("Id", axis=1), hue = "Species", size=3, markers=["o", "s", "D"])  
plt.show()
```



# K Nearest Neighbor Classifier

## Working with Iris data set

# Box Plots

```
plt.figure()  
df.drop("Id", axis=1).boxplot(by="Species", figsize=(15, 10))  
plt.show()
```

# Fitting classifier to the Training set

# Loading libraries

```
from sklearn.neighbors import KNeighborsClassifier  
from sklearn.metrics import confusion_matrix, accuracy_score  
from sklearn.model_selection import cross_val_score
```



# K Nearest Neighbor Classifier

## Working with Iris data set

```
# Instantiate learning model (k = 3)
classifier = KNeighborsClassifier(n_neighbors=3)

# Fitting the model
classifier.fit(X_train, y_train)

# Predicting the Test set results
y_pred = classifier.predict(X_test)

print(y_pred)
```



# K Nearest Neighbor Classifier

## Working with Iris data set

# Define a function module to print results of ML Classifier Score

```
from sklearn.metrics import accuracy_score, confusion_matrix, precision_score, recall_score, f1_score
```

```
def print_score(clf, x_train, y_train, x_test, y_test, train = True):
```

```
    if train:
```

```
        pred = clf.predict(x_train)
```

```
        print("Train Result:\n =====")
```

```
        print(f"accuracy score: {accuracy_score(y_train, pred):.4f}\n")
```

```
        print("Classification Data:")
```

```
        print(f"Precision: {precision_score(y_train, pred, average=None, zero_division=1)}\n")
```

```
        print(f"Recall Score: {recall_score(y_train, pred, average=None, zero_division=1)}\n")
```

```
        print(f"Confusion_matrix:\n {confusion_matrix(y_train, clf.predict(x_train))}\n")
```

```
    elif train == False:
```

```
        pred = clf.predict(x_test)
```

```
        print("Test Result:\n =====")
```

```
        print(f"accuracy score: {accuracy_score(y_test, pred)}\n")
```

```
        print("Classification Data:")
```

```
        print(f"Precision: {precision_score(y_test, pred, average=None, zero_division=1)}\n")
```

```
        print(f"Recall Score: {recall_score(y_test, pred, average=None, zero_division=1)}\n")
```

```
        print(f"Confusion_matrix:\n {confusion_matrix(y_test, clf.predict(x_test))}\n")
```





# K Nearest Neighbor Classifier

## Working with Iris data set

# Print results

```
print_score(classifier, X_train, y_train, X_test, y_test, train = True)
```

```
print_score(classifier, X_train, y_train, X_test, y_test, train = False)
```

# Using cross-validation for parameter tuning:

# creating list of K for KNN

```
k_list = list(range(1,50,2))
```

# creating list of cv scores

```
cv_scores = []
```

# perform 10-fold cross validation

for k in k\_list:

```
    knn = KNeighborsClassifier(n_neighbors=k)
```

```
    scores = cross_val_score(knn, X_train, y_train, cv=10, scoring='accuracy')
```

```
    cv_scores.append(scores.mean())
```

```
print(cv_scores)
```



# K Nearest Neighbor Classifier

## Working with Iris data set

```
# changing to misclassification error
```

```
MSE = [1 - x for x in cv_scores]
```

```
plt.figure()
```

```
plt.figure(figsize=(15,10))
```

```
plt.title('The optimal number of neighbors', fontsize=20, fontweight='bold')
```

```
plt.xlabel('Number of Neighbors K', fontsize=15)
```

```
plt.ylabel('Misclassification Error', fontsize=15)
```

```
sns.set_style("whitegrid")
```

```
plt.plot(k_list, MSE)
```

```
plt.show()
```

```
# finding best k
```

```
best_k = k_list[MSE.index(min(MSE))]
```

```
print("The optimal number of neighbors is %d." % best_k)
```



# K Nearest Neighbor Classifier

## Working with Iris data set

```
# Instantiate learning model (k = 9)
classifier = KNeighborsClassifier(n_neighbors=9)

# Fitting the model
classifier.fit(X_train, y_train)

print_score(classifier, X_train, y_train, X_test, y_test, train = True)
print_score(classifier, X_train, y_train, X_test, y_test, train = False)
```

