

# 7CS108: Data Science and Data Mining

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# Classification

Classification is a technique or a function that classifies given data into one of the pre-defined categories/classes.



Apple



Pear

**Distinguishing feature - colour**





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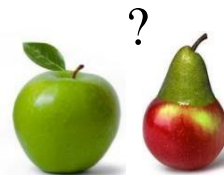


Apple



Pear

**Distinguishing feature - colour**



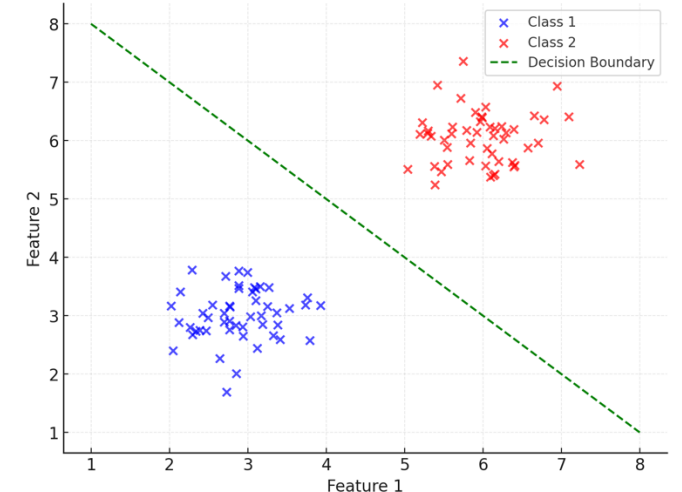
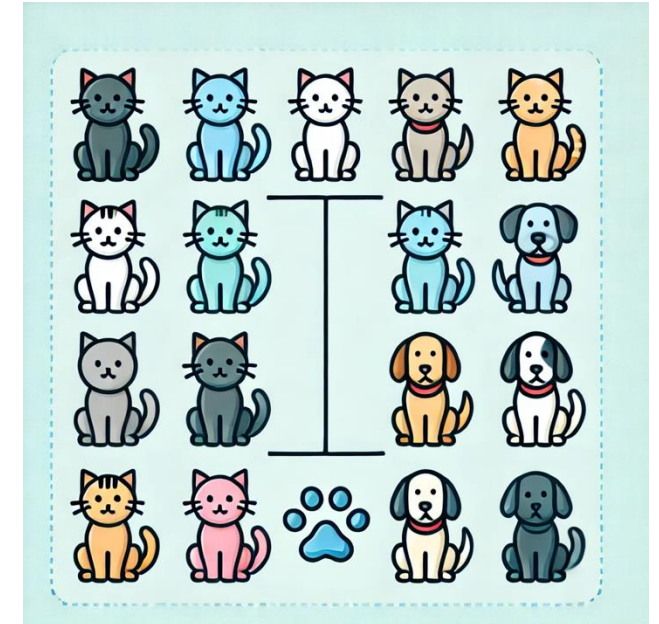
**Is colour enough?**

**It may lead to misclassification!**



# Classification

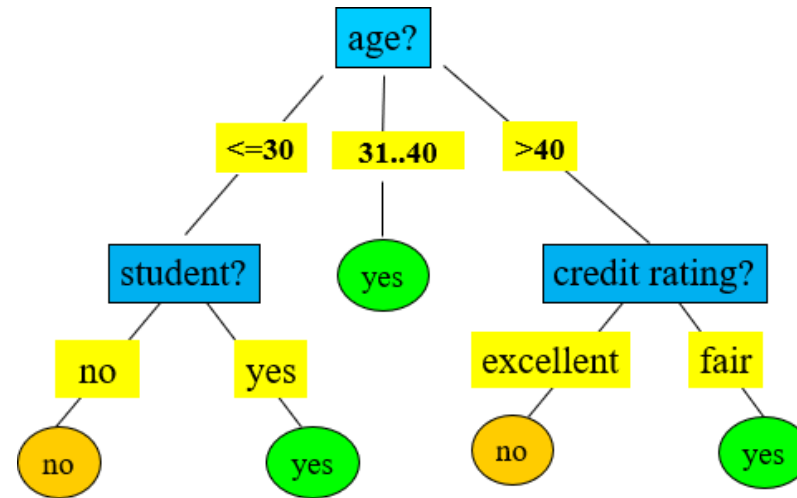
- Classification is a technique or a function that classifies given data into one of the pre-defined categories/classes.
- Classification is also known as “Supervised Learning”
  - There must be an “expert” (you) to “supervise” the computer.
  - In contrast, Clustering is known as “Unsupervised Learning”. We will discuss it in the later lectures.
  - There are semi-supervised and reinforcement learning too.
- From the data mining point of view...
  - Classification  $\approx$  Prediction  $\approx$  Forecasting
  - This is because the techniques are the same





# Question?

Buys computer = yes ? no

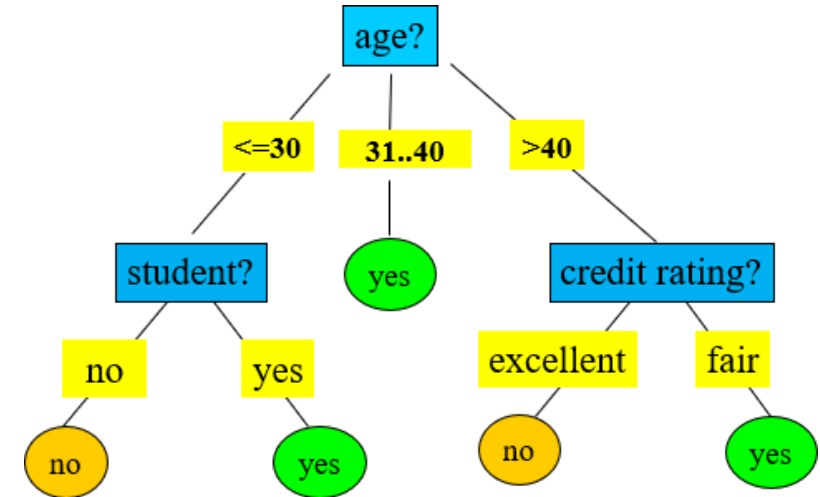


Age	Student	Credit rating	Buys computer
28	Yes	Fair	?



# What is Decision Tree?

- Flow-chart-like tree structure
  - Root node: a test on root node
    - Age < 30
  - Internal node: a test on an attribute
    - Student = yes ?
  - Leaf node: classes
    - Buys computer: yes or no?





# Training set

- Whether a customer is likely to buy a computer?

age	income	student	credit_rating	buys_computer
<=30	high	no	fair	no
<=30	high	no	excellent	no
31...40	high	no	fair	yes
>40	medium	no	fair	yes
>40	low	yes	fair	yes
>40	low	yes	excellent	no
31...40	low	yes	excellent	yes
<=30	medium	no	fair	no
<=30	low	yes	fair	yes
>40	medium	yes	fair	yes
<=30	medium	yes	excellent	yes
31...40	medium	no	excellent	yes
31...40	high	yes	fair	yes
>40	medium	no	excellent	no



# Data Preparation

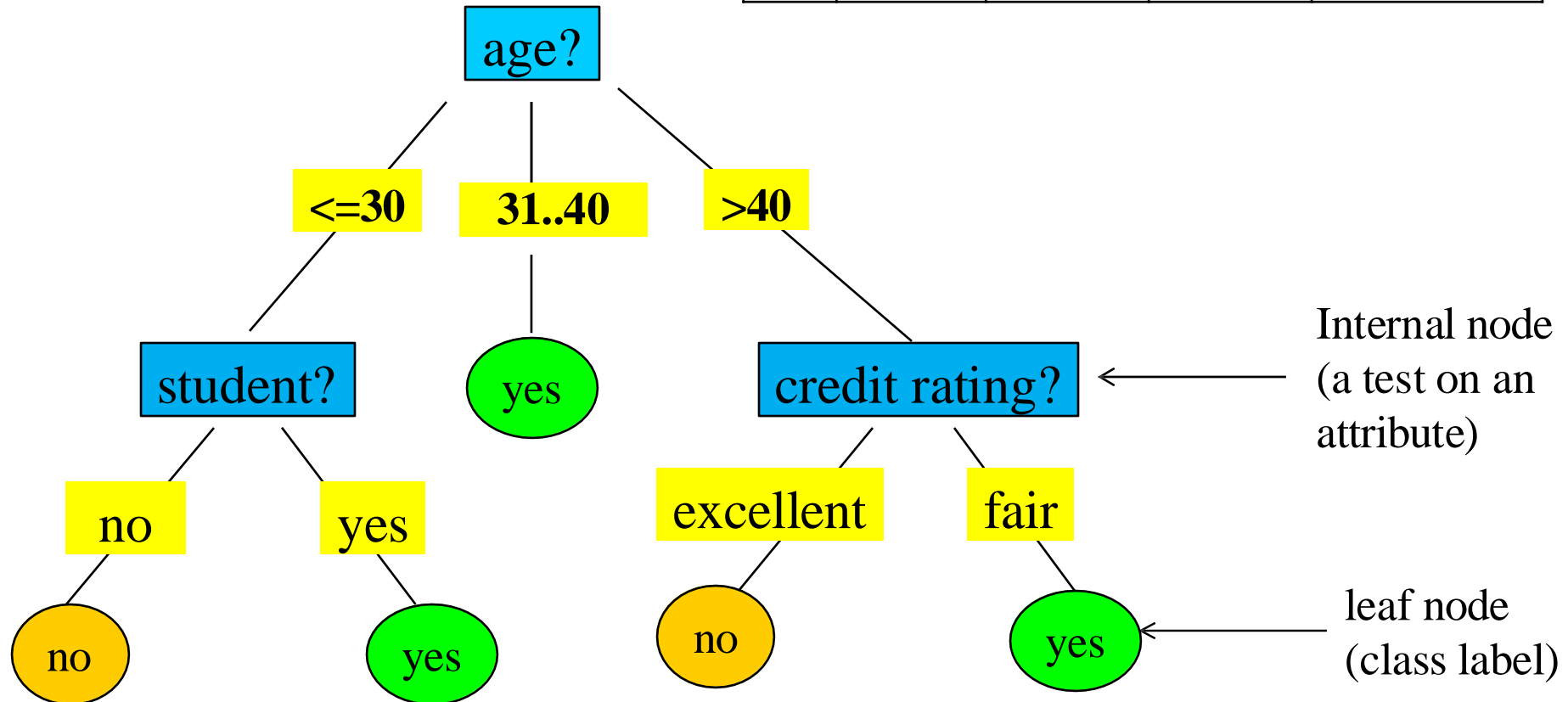
- Data Cleaning
  - Pre-process data in order to reduce noise and handle missing values
- Relevance analysis (feature selection)
  - Remove the irrelevant or redundant attributes
- Data transformation
  - Generalise and/or normalise data





# Decision Tree: “Buys Computer”

Age	Income	Student	Credit rating	Buys computer
25	Low	Student	Fair	?





# Decision Tree Induction

- Many algorithms:
  - Hunt's Algorithm
  - ID3 – late 1970's, J. Ross Quinlan
  - C4.5 – successor of ID3
  - CART – similar approach as C4.5
- General Structure
  - Attribute selection by calculating entropy and Information Gain
  - Decision tree induction
  - Rule generation



# A Simple Thought on Decision Tree

- A simple minded algorithm:
  - If Age = a and Income = b and Student = c and Credit Rating = d
  - then buys\_computer = ?
- This structure will yield a branch for each row, not eliminating superfluous attributes, and is unnecessarily complex.
- The complete classification space for m attributes is of size:  $\prod_{j=1}^m N_j$   
( $N_j$  is the number of values of attribute  $j$ )



# The Problem Statement

- Given a set of data described by **a set of attributes** and **an outcome class**, the problem is to find a **minimum decision tree** that will classify the values of the class based on the values of given attributes.



# The Problem Statement

- Given a set of data described by **a set of attributes** and **an outcome class**, the problem is to find a **minimum decision tree** that will classify the values of the class based on the values of given attributes.
- How do we select an attribute?
  - Amongst four attributes: **AGE, INCOME, STUDENT** and **CREDIT RATING**, **which attribute carries “more” information than others?**



# Information Theory

- Information Entropy  $\rightarrow$  uncertainty
- Information Gain  $\rightarrow$  reduce entropy

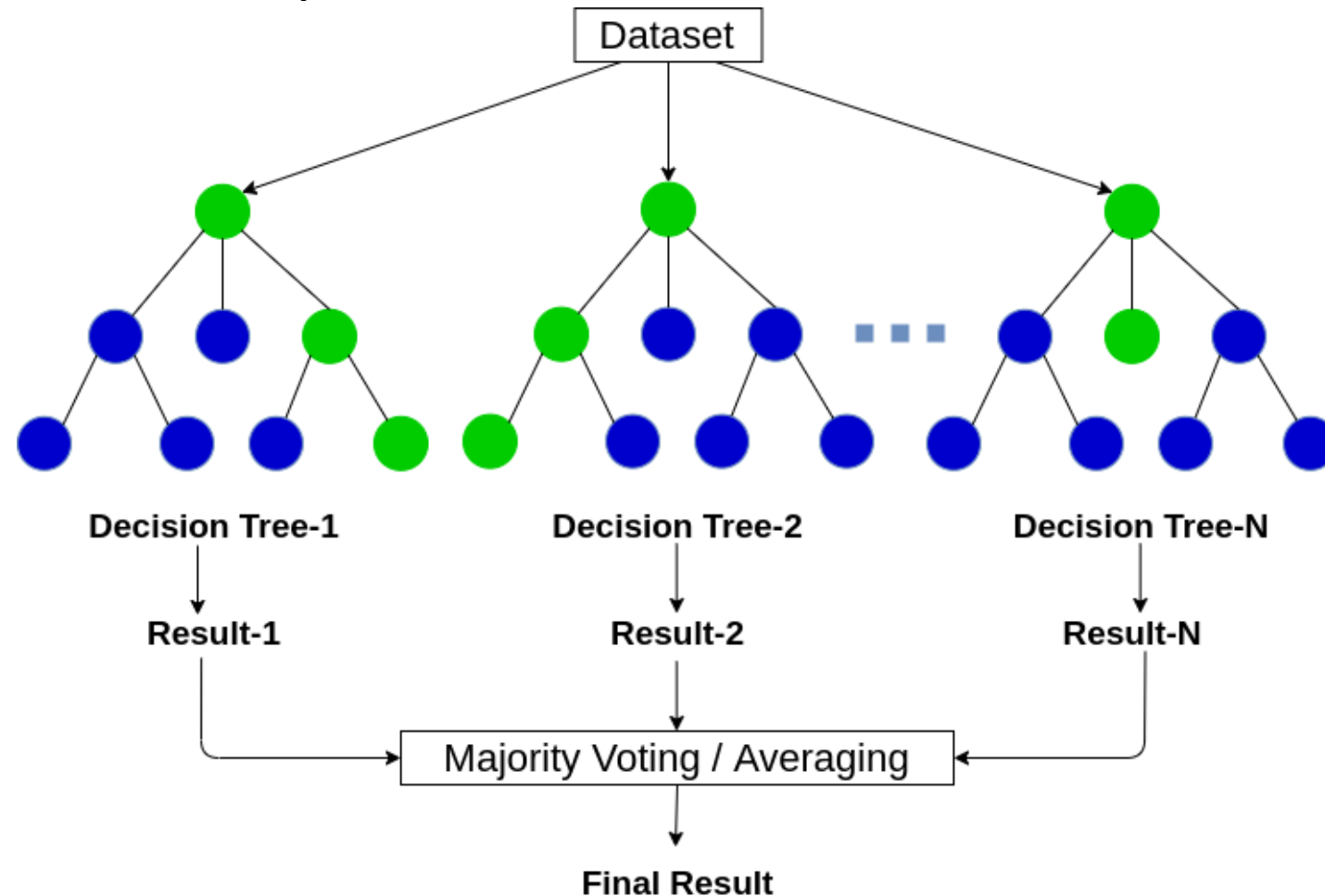
$$E = -\sum p_i \log_2(p_i)$$

$$IG = E(Parent) - \sum w_i E(Child_i)$$



# Random Forest

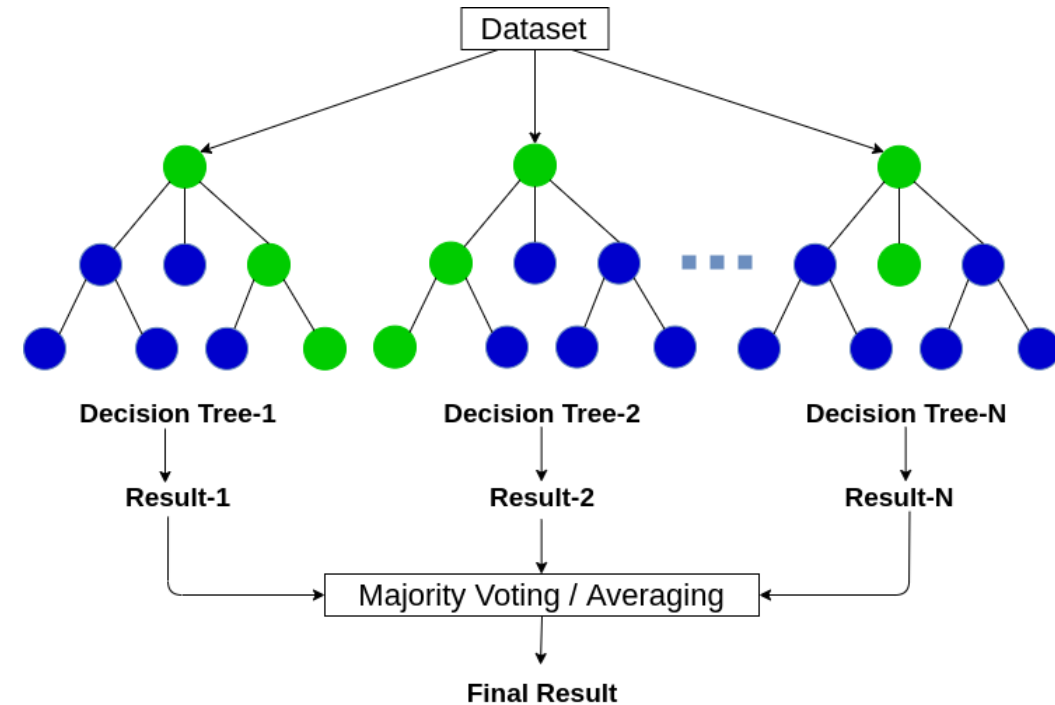
- Combines the output of multiple (randomly created) Decision Trees to generate the final output.





# Random Forest

- **Individual Trees:** Each Decision Tree is trained on a random subset of the data (with replacement, i.e., bootstrapping) and a random subset of features.
- **Tree Predictions:** Each tree makes an independent prediction based on its learned splits and conditions.
- **Aggregation by Majority Voting:** The Random Forest combines the predictions of all trees, and the class with the highest votes becomes the final classification.

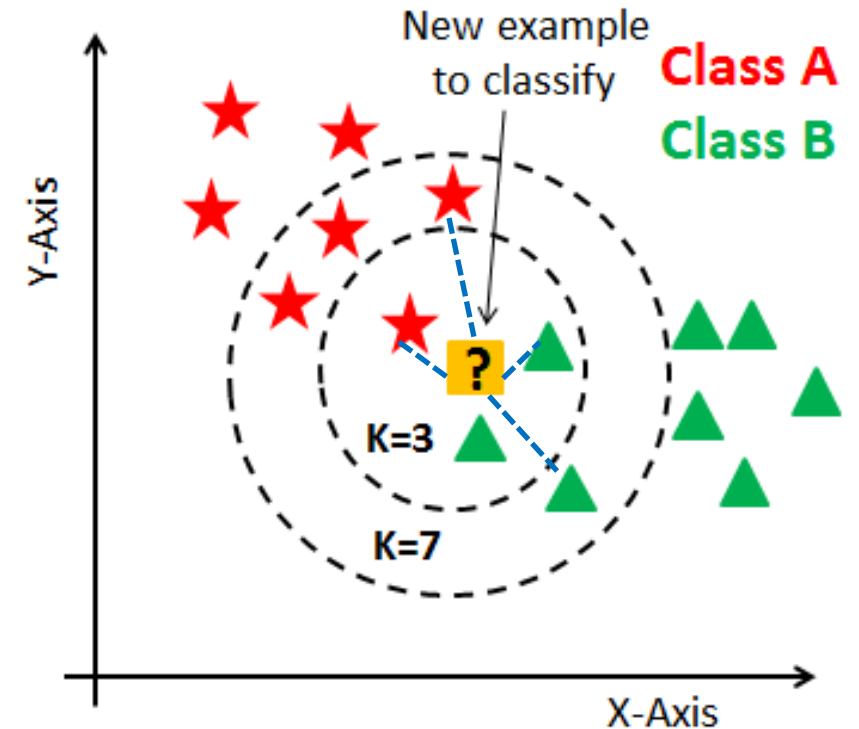






# K-Nearest Neighbours (kNN)

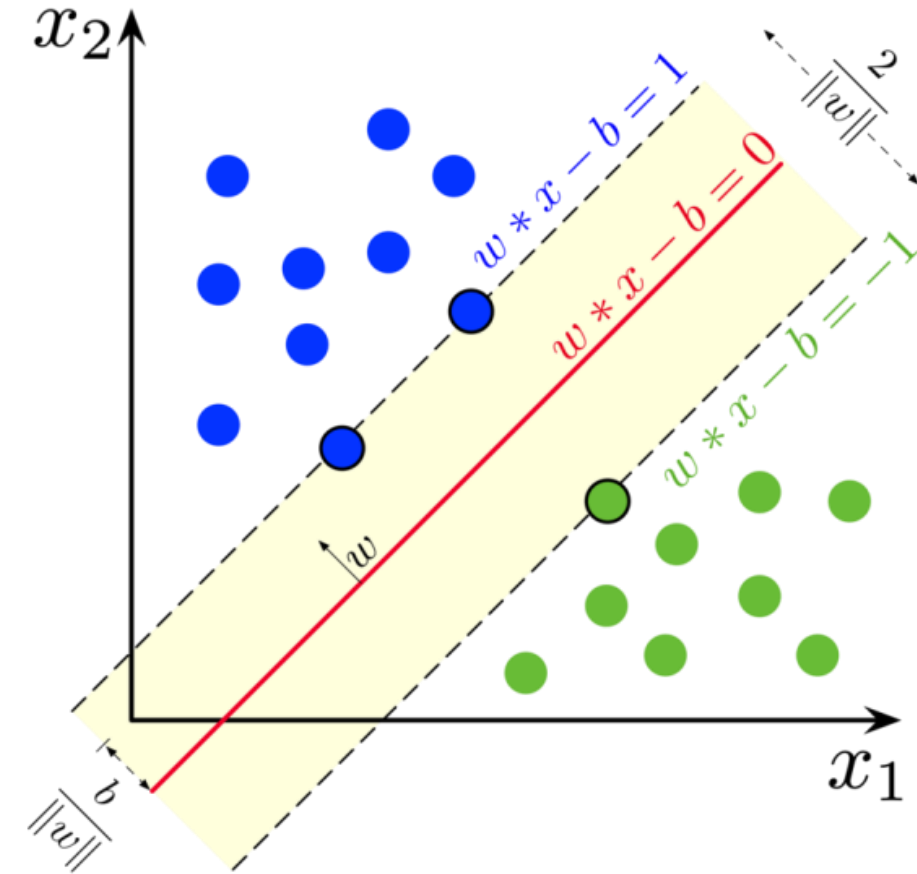
- Choose the value of  $k$  (number of neighbours).
- Compute the distance between the query point and all other data points.
- Identify the  $k$  nearest neighbours.
- Assign the class label based on the majority vote among  $k$  neighbours.
- In case of regression, take the mean of the  $k$  neighbours' values.



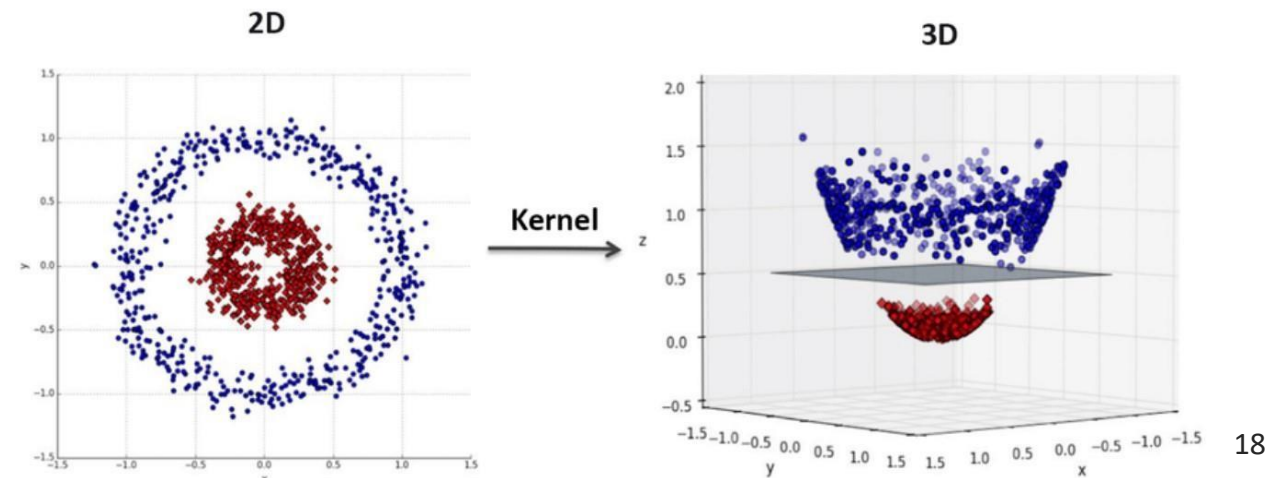
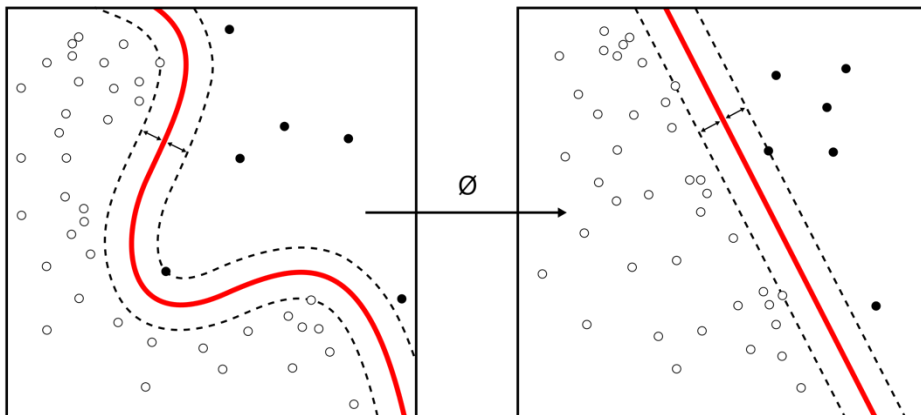


# Support Vector Machine (SVM)

- The objective is to find a hyperplane in an n-dimensional space that separates the data points to their potential classes. The hyperplane should be positioned with the maximum distance to the data points.
- The data points with the minimum distance to the hyperplane are called Support Vectors



- The original maximum-margin hyperplane algorithm proposed by Vapnik in 1963 constructed a linear classifier that works when data are linearly separable (using a line or hyperplane)
- For data that are not linearly separable, we need to use a kernel function and transform data to another space where they are linearly separable. This is called the kernel trick.
- There are different kernel functions including linear function, Polynomial function, Radial basis function (RBF), and Sigmoid function.





# Evaluation

- Confusion matrix
- Accuracy
- Precision
- Recall



# Evaluation

## Confusion Matrix

	Actual Positive (Apple)	Actual Negative (Non-apple)
Predicted Positive (Apple)	TP	FP
Predicted Negative (Non-apple)	FN	TN

TP: true positive

FP: false positive

FN: false negative

TN: true negative

$$\text{Accuracy} = \frac{TP+TN}{TP+TN+FP+FN}$$

**NOTE:** Accuracy is NOT always a good measure!



# Confusion Matrix

**Total fruits: 20**

	Actual Positive (Apple)	Actual Negative (Non-apple)
Predicted Positive (Apple)	10 TP	5 FP
Predicted Negative (Non-apple)	2 FN	3 TN

TP: true positive  
FN: false negative

FP: false positive  
TN: true negative



# Confusion Matrix

**Total fruits: 20**

	Actual Positive (Apple)	Actual Negative (Non-apple)
Predicted Positive (Apple)	10 TP	5 FP
Predicted Negative (Non-apple)	2 FN	3 TN

**True Positive (TP):** the classifier correctly makes positive decisions

**True Negative (TN):** the classifier correctly makes negative decision

**False Positive (FP):** the classifier mistakenly makes positive decisions

**False Negative (FN):** the classifier mistakenly makes negative decisions



# Limitation

Suppose

- Total number of fruits in the testing examples = 10,000
- Number of Non-apple = 9990
- Number of Apple = 10

Can you classify Apple?

If model predicts everything to be of class non-apple, the accuracy is  $9990/10000 = 99.9\%$  !!!

Here, accuracy is misleading because model cannot detect any Apple at all, still achieving high accuracy.





# Evaluation

	Actual Positive (Apple)	Actual Negative (Non-apple)
Predicted Positive (Apple)	TP	FP
Predicted Negative (Non-apple)	FN	TN

$$Precision = \frac{TP}{TP + FP}$$

$$Recall = \frac{TP}{TP + FN}$$

$$F_1 = 2 * \frac{Precision * Recall}{Precision + Recall}$$



# Evaluation

	Actual Positive (Apple)	Actual Negative (Non-apple)
Predicted Positive (Apple)	TP	FP
Predicted Negative (Non-apple)	FN	TN

$$Precision = \frac{TP}{TP + FP} = \frac{TP}{\text{predicted positive}}$$

**Precision** is the total number of **correctly identified actual apple cases out of retrieved apple**

$$Recall = \frac{TP}{TP + FN} = \frac{TP}{\text{actual positive}}$$

**Recall** is the number of **correctly identified apple cases from the total number of actual apple cases**



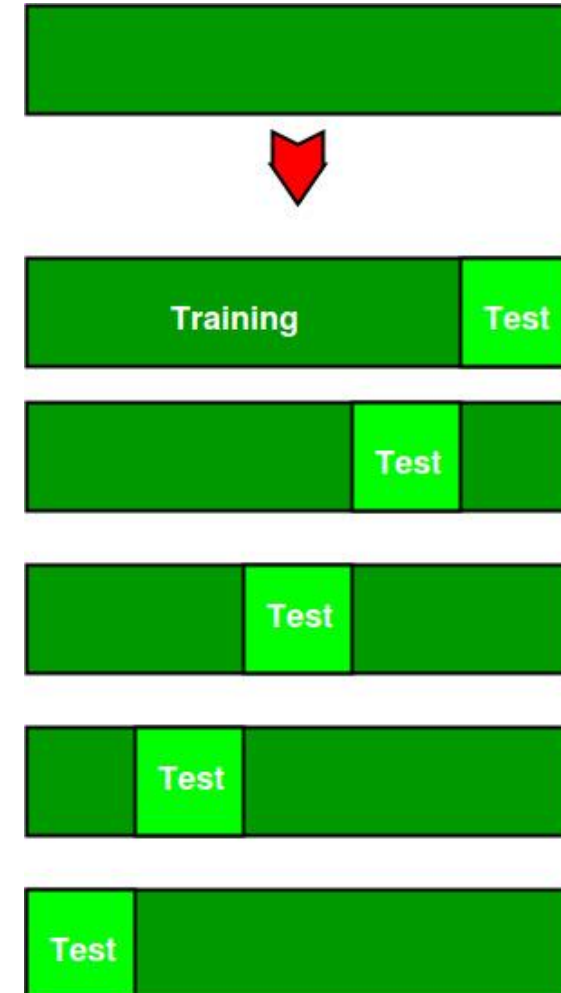
# Evaluation

- Methods for Performance Evaluation
  - How to obtain reliable estimates?
  - i.e. how to partitioning the data?
- N-Fold cross validation
- Training and Testing sets



# Cross Validation

- Ensures generalization to unseen data.
- Prevents overfitting.
- Provides reliable performance evaluation.
- Improves model robustness and reduces evaluation bias.
- Popular approaches:
  - K-Fold Cross Validation
  - Leave-One-Out





# train\_test\_split()

```
x_train, x_test, y_train, y_test = train_test_split(features, targets, test_size= .30,  
random_state = 50)
```

