

# Machine Learning

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# Attribute Selection Measures

- **Information Gain** - Changes in entropy after the segmentation of a dataset based on an attribute

$$\text{Information Gain} = \text{Entropy}(S) - [(\text{Weighted Avg}) * \text{Entropy}(\text{each feature})]$$

**Entropy:** It measures the impurity in a given attribute

$$\text{Entropy}(s) = -P(\text{yes}) \log_2 P(\text{yes}) - P(\text{no}) \log_2 P(\text{no})$$

Where,  $S$  = Total number of samples,  $P(\text{yes})$  = probability of yes,  $P(\text{no})$  = probability of no

- **Gini Index** - Measure of impurity or purity used while creating a decision tree

$$\text{Gini Index} = 1 - \sum_j P_j^2$$



# Pruning -

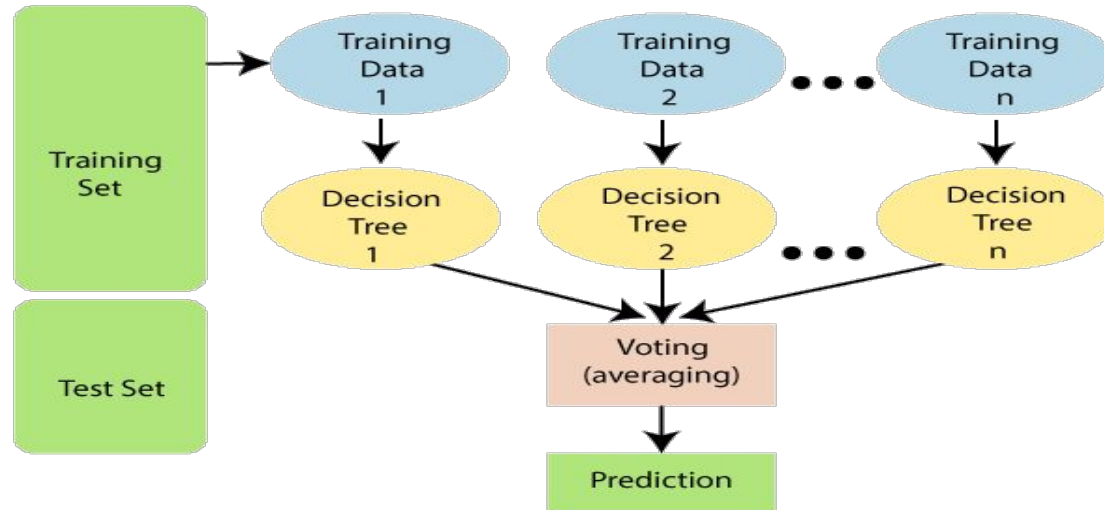
A process of deleting the unnecessary nodes from a tree in order to get the optimal decision tree.

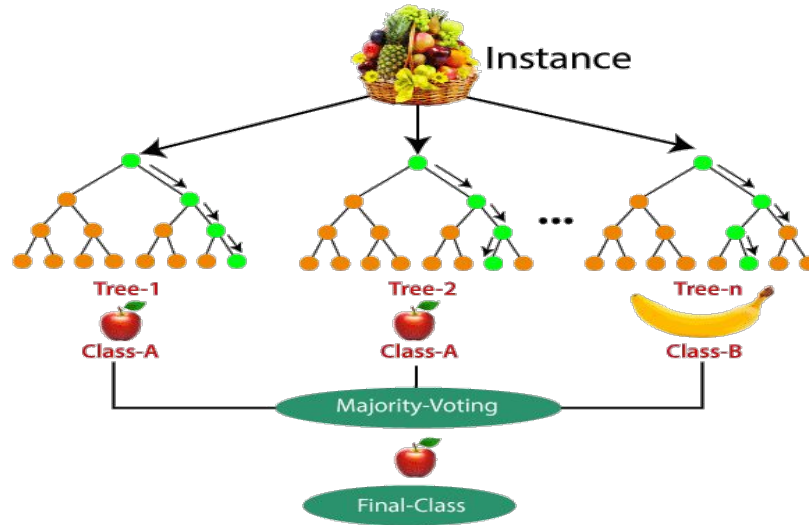
- **Cost Complexity Pruning**
- **Reduced Error Pruning.**



# Random Forest -

Random Forest is a classifier that contains a number of decision trees on various subsets of the given dataset and takes the average to improve the predictive accuracy of that dataset.



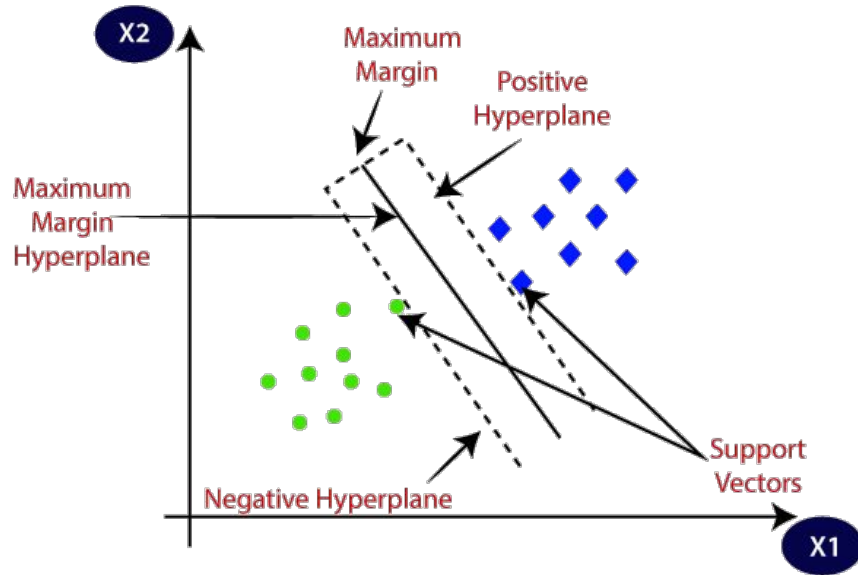


$$g(x) = f_0(x) + f_1(x) + f_2(x) + \dots$$

Random forest uses **Bagging or Bootstrap Aggregation** technique of ensemble learning in which aggregated decision tree runs in parallel and do not interact with each other.

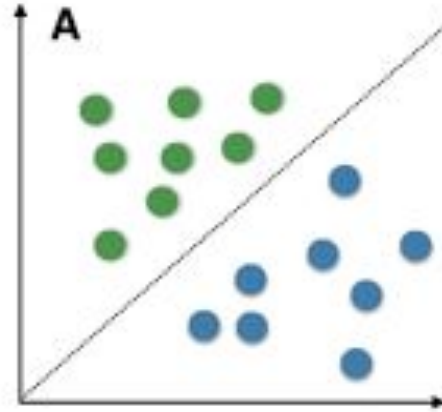


# Support Vector Machine -

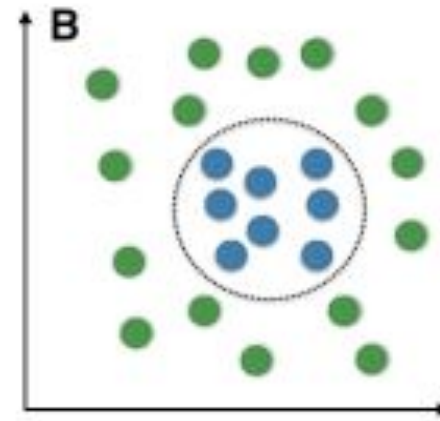


- Kernel
- Hyperplane
- Boundary line
- Support vectors





**Linear SVM**

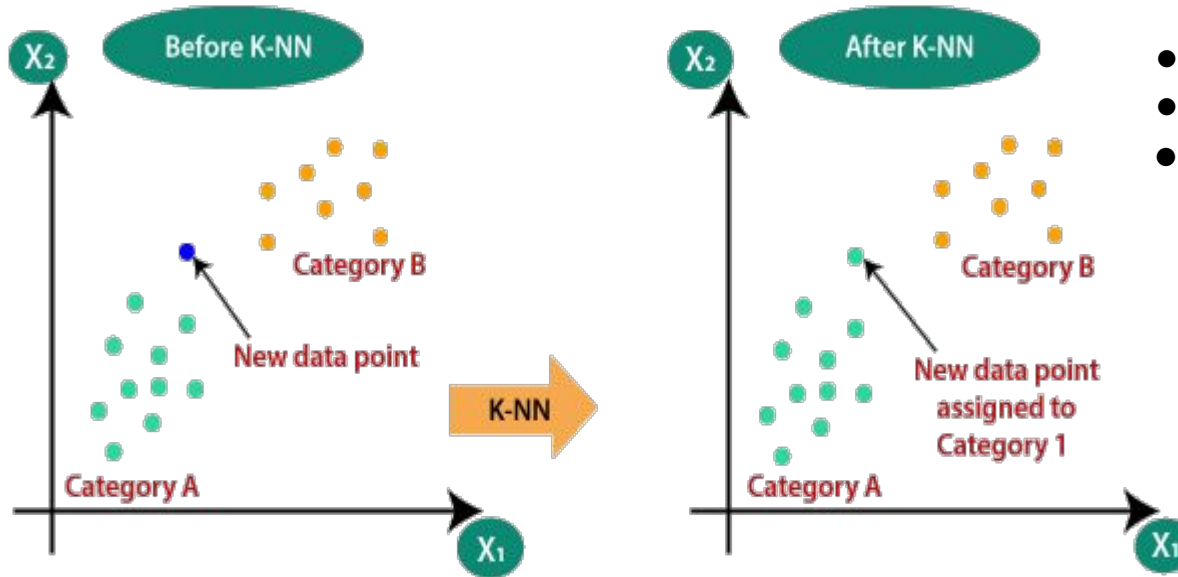


**Non-Linear SVM**

The main goal of SVR is to consider the maximum data points within the boundary lines and the hyperplane (best-fit line) must contain a maximum number of datapoints



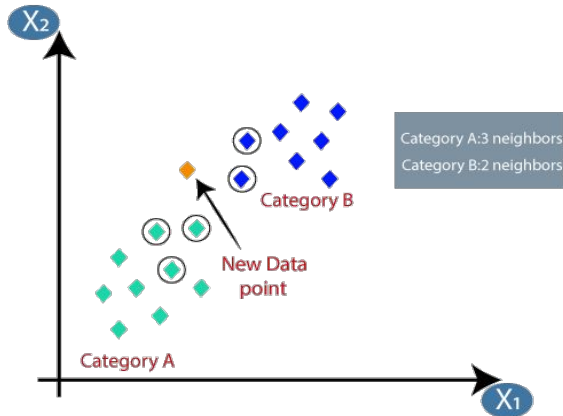
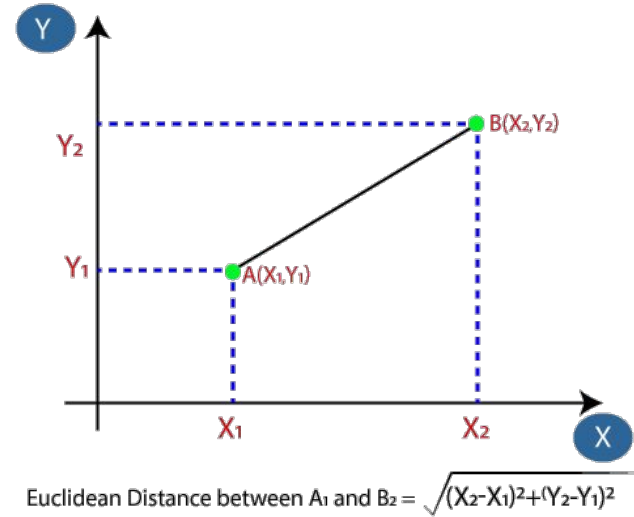
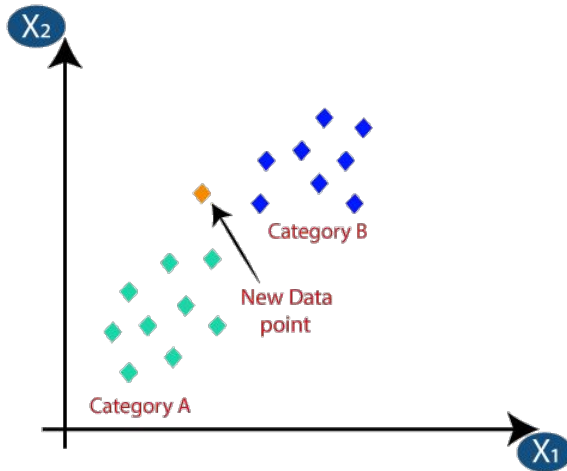
# K-Nearest Neighbor(KNN) -



- lazy learner algorithm
- non-parametric algorithm
- Based on the similarity







K is the number of nearest neighbors

Here  $k=5$ , as new datapoint is having 3 nearest neighbors from category A so it belongs to Category A



# Naïve Bayes -

- Based on **Bayes theorem**
- It is mainly used in *text classification* that includes a high-dimensional training dataset.
- It is a probabilistic classifier, which means it predicts on the basis of the probability of an object.

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

Where, **P(A|B)** is **Posterior probability**: Probability of hypothesis A on the observed event B, **P(B|A)** is **Likelihood probability**: Probability of the evidence given that the probability of a hypothesis is true, **P(A)** is **Prior Probability**: Probability of hypothesis before observing the evidence, **P(B)** is **Marginal Probability**: Probability of Evidence.



## Steps -

1. Convert the given dataset into frequency tables.
2. Generate Likelihood table by finding the probabilities of given features.
3. Now, use Bayes theorem to calculate the posterior probability.

## Types -

- **Gaussian**
- **Multinomial**
- **Bernoulli**

