

K-Nearest Neighbour (K-NN) Classifier

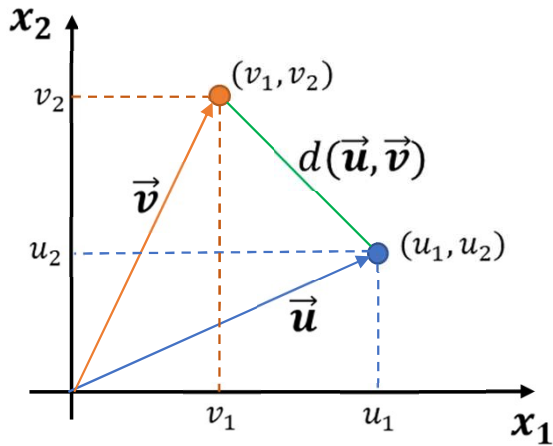
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OUTLINE

- Euclidean Distance
- Other Distance Metrics
- K-NN Classifier
- Decision Boundary of K -NN Classifier
- Choosing the value of K
- Merits and Demerits of K-NN classifier

EUCLIDEAN DISTANCE



- Consider the points in two-dimension. Each point in two-dimension can be represented by a vector of dimension two.
- The point (u_1, u_2) can be represented by the vector $\vec{u} = [u_1, u_2]^T$
- And the point (v_1, v_2) can be represented by the vector $\vec{v} = [v_1, v_2]^T$
- The Euclidean distance between the points is:

$$d(\vec{u}, \vec{v}) = \sqrt{(u_1 - v_1)^2 + (u_2 - v_2)^2}$$

- In general a point in n -dimensional space is represented as $\vec{u} = [u_1, u_2, u_3, \dots, u_n]^T$, a n -D vector
- Hence, the Euclidean distance between two points in n -dimensional space is represented as:

$$d(\vec{u}, \vec{v}) = \sqrt{(u_1 - v_1)^2 + (u_2 - v_2)^2 + (u_3 - v_3)^2 + \dots + (u_n - v_n)^2} = \sqrt{\sum_{i=1}^n (u_i - v_i)^2}$$

- In general in vector notation, the Euclidean distance is written as: $d(\vec{u}, \vec{v}) = \sqrt{(\vec{u} - \vec{v})^T (\vec{u} - \vec{v})}$

OTHER DISTANCE METRICS

- **Manhattan Distance:** For two data points denoted by \mathbf{x} and \mathbf{y} the Manhattan distance is defined as:

$$dist_{manhattan}(\mathbf{x}, \mathbf{y}) = \sum_{i=1}^n |x_i - y_i|$$

- **Minkowski Distance:** For two data points denoted by \mathbf{x} and \mathbf{y} the Minkowski distance is defined as:

$$dist_{minkowski}(\mathbf{x}, \mathbf{y}, h) = [\sum_{i=1}^n (x_i - y_i)^h]^{\frac{1}{h}}$$

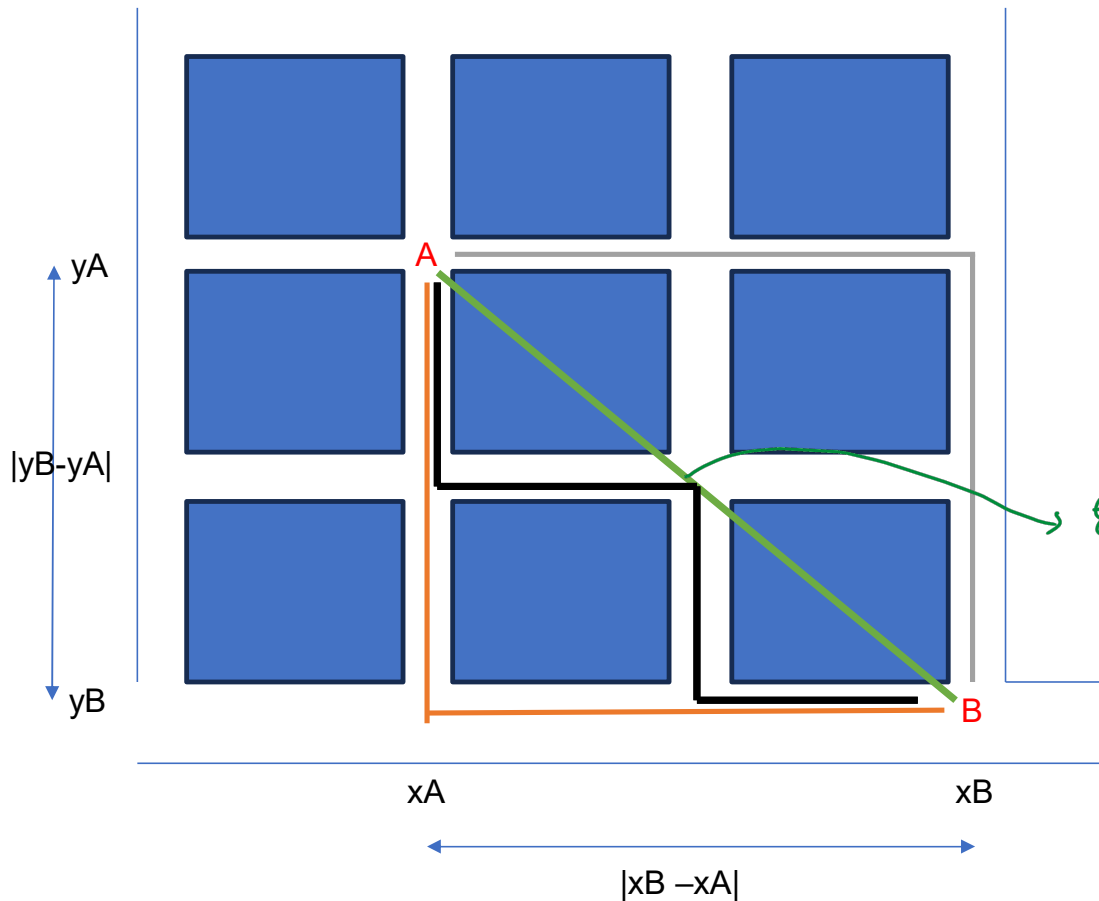
Note: for $h = 2$, Minkowski Distance is same as Euclidean Distance and for $h = 1$, it is Manhattan Distance

- **Chebyshev Distance:** For two data points in n -dimensional space it is defined as:

$$dist_{chebyshev}(\mathbf{x}, \mathbf{y}) = \max(|x_1 - y_1|, |x_2 - y_2|, |x_3 - y_3|, \dots, |x_n - y_n|)$$

There are other distance metric like: Mahalanobis Distance, Bhattacharya Distance etc. which are used for advanced statistical pattern recognition tasks.

Manhattan Distance or city-block distance



The shortest distance between A to B around the city using road

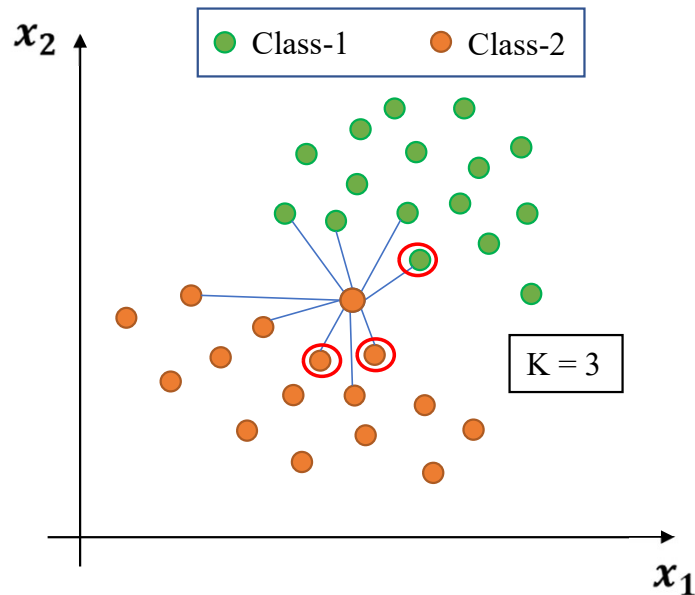
$$\underbrace{|x_B - x_A| + |y_B - y_A|}_{\text{City block distance}}$$

Manhattan "

Euclidean
distance
(x)

K-NN CLASSIFIER

- **Intuition:** One is known by the company one keeps.



- **K-NN Algorithm:**

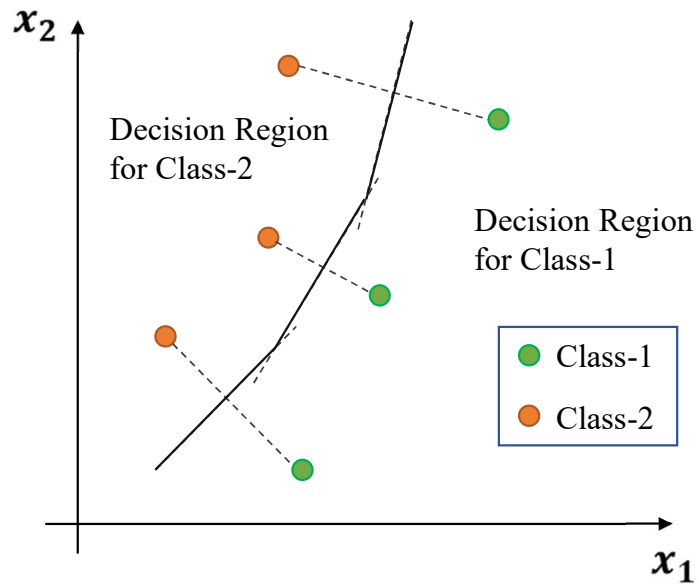
1. All the training samples/ points are available beforehand.
2. When a new test sample arrives calculate its **distance** from **all training points**.
3. Choose K-nearest neighbours based on the distance calculated. Usually the K is a positive odd integer and supplied by user.
4. Assign the class label of the test sample **based on majority**. i.e. for a test sample if most number of neighbours among those K-Nearest Neighbours belong to one particular class- c , then assign the class label of test sample as c .

- **Characteristics of K-NN Classifier:**

It doesn't create model based on the training patterns in advance. Rather, when a test instance comes for testing, runs the algorithm to get the class prediction of that particular testing instance. Hence, there is no learning in advance.

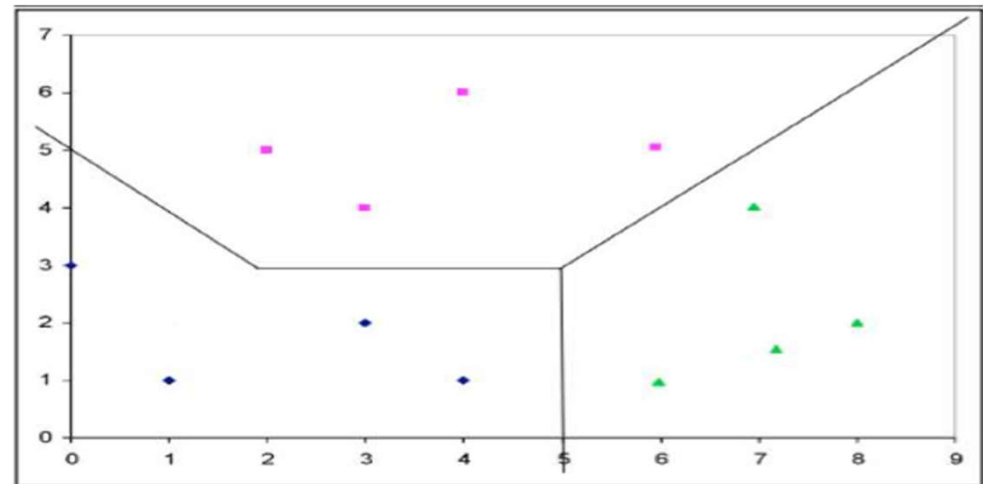
Hence, k-NN classifier is also known as **Lazy Learner**.

K-NN CLASSIFIER: DECISION BOUNDARY



- Boundary are the points those are equidistant between the points of Class-1 and Class-2
- Construct lines between closest pairs of points in different classes.
- Draw perpendicular bisectors. End bisectors at intersections.
- Note that locally the boundary is linear.
- Hence the decision boundary is piecewise linear curve.

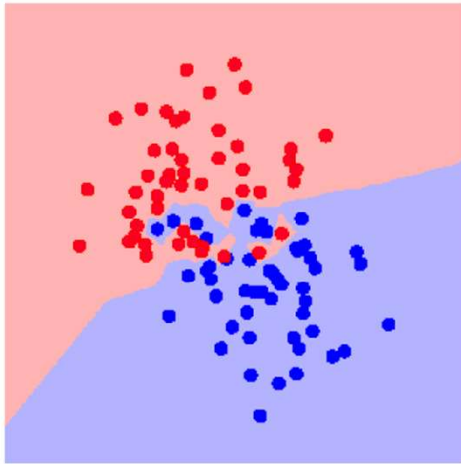
For multiclass classification also the same thing is done to find the decision boundary.



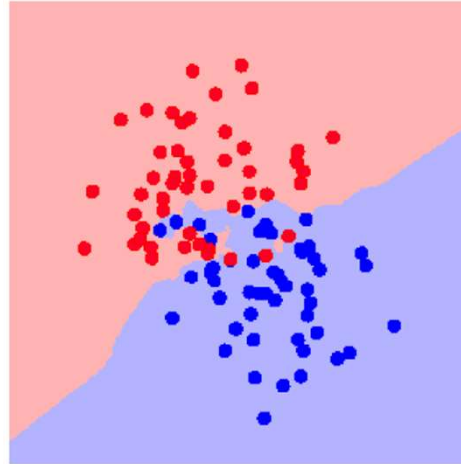
K-NN: CHOOSING THE VALUE OF K

- Increasing the 'K' simplifies the decision boundary. Because majority voting implies less emphasis on individual points

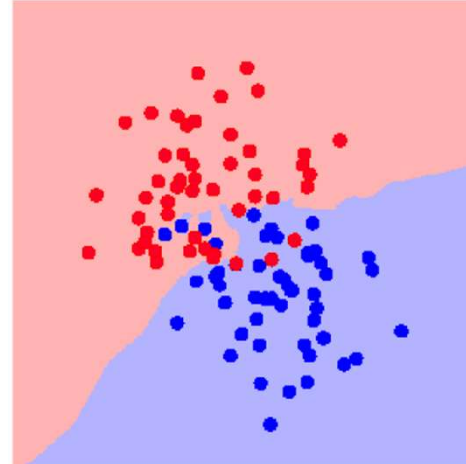
K = 1



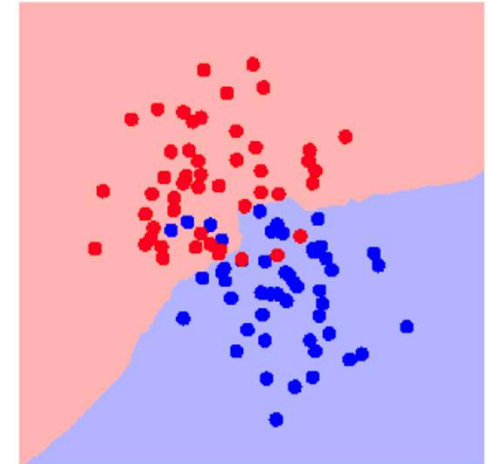
K = 3



K = 5



K = 7



- However increasing the K also increases computational cost.
- Hence, choosing K is an optimization between how much simplified decision boundary we want vs. how much computational cost we can afford.
- Usually K = 5, 7, 9, 11 works fine for most practical problems.

K-NN CLASSIFIER: MERITS AND DEMERITS

Merits:

- K-NN Classifier often works very well for practical problems.
- It is very easy to implement, as there is no complex learning algorithm involved.
- Robust to Noisy Data.

Demerits:

- Choosing the value of K may not be straightforward. Often the same training samples are used for different values of K, and we choose the most suitable value of K based on minimum misclassification errors on test samples.
- Doesn't work well for categorical attributes.
- Can encounter problem with sparse training data. (i.e. data points are located far away from each other)
- Can encounter problems in very high-dimensional spaces.
 - Most points are at corners.
 - Most points are at the edge of the space.

This problem is known as ***Curse of Dimensionality*** and affect many other Machine Learning algorithms.

1-D case:



define edge as: $[0, \epsilon] \cup [1-\epsilon, 1]$

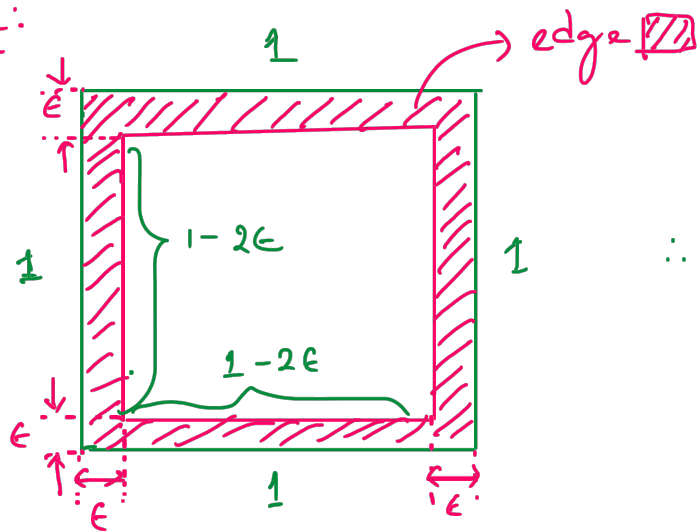
say $\epsilon = 0.05$

edge is $[0, 0.05] \cup [0.95, 1]$

x is uniformly distributed $[0, 1]$, what is the probability that x is at the edge of the line?

$$P(\text{edge}) = \frac{2\epsilon}{L} = 2 \times 0.05 = 0.1 \quad p = 1 - (1 - 2\epsilon)$$

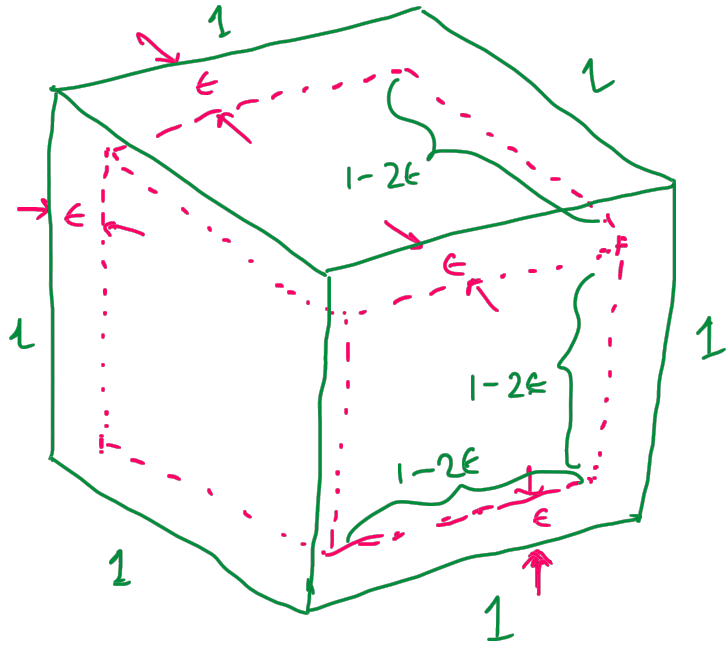
2-D case:



x is a point uniformly distributed in the unit square

$$\begin{aligned} \therefore P(\text{edge}) &= \text{Area of unit square} \\ &\quad - \text{area of the blank region} \\ &= 1 - (1 - 2\epsilon)^2 = 1 - 0.9^2 = 1 - 0.81 = 0.19 \end{aligned}$$

3D-case:



x is uniformly distributed in the cube

$$\therefore P(\text{edge}) = \text{Volume of unit cube} \\ - \text{volume of the inside cube}$$

$$= 1 - (1 - 2\epsilon)^3$$

$$\epsilon = 0.05, \quad \therefore P(\text{edge}) = 1 - 0.9^3 = 1 - 0.729 \\ = 0.271$$

For a n -dimensional hyper-cube

$$P(\text{edge}) = 1 - (1 - 2\epsilon)^n$$

$$\text{for } n=20, \epsilon=0.05 \quad P(\text{edge}) = 1 - (0.9)^{20} = 1 - 0.12 = 0.88$$

$$\text{for } n=50, \epsilon=0.01 \quad P(\text{edge}) = 1 - (0.98)^{50} = 1 - 0.36 = 0.64$$

Thank You