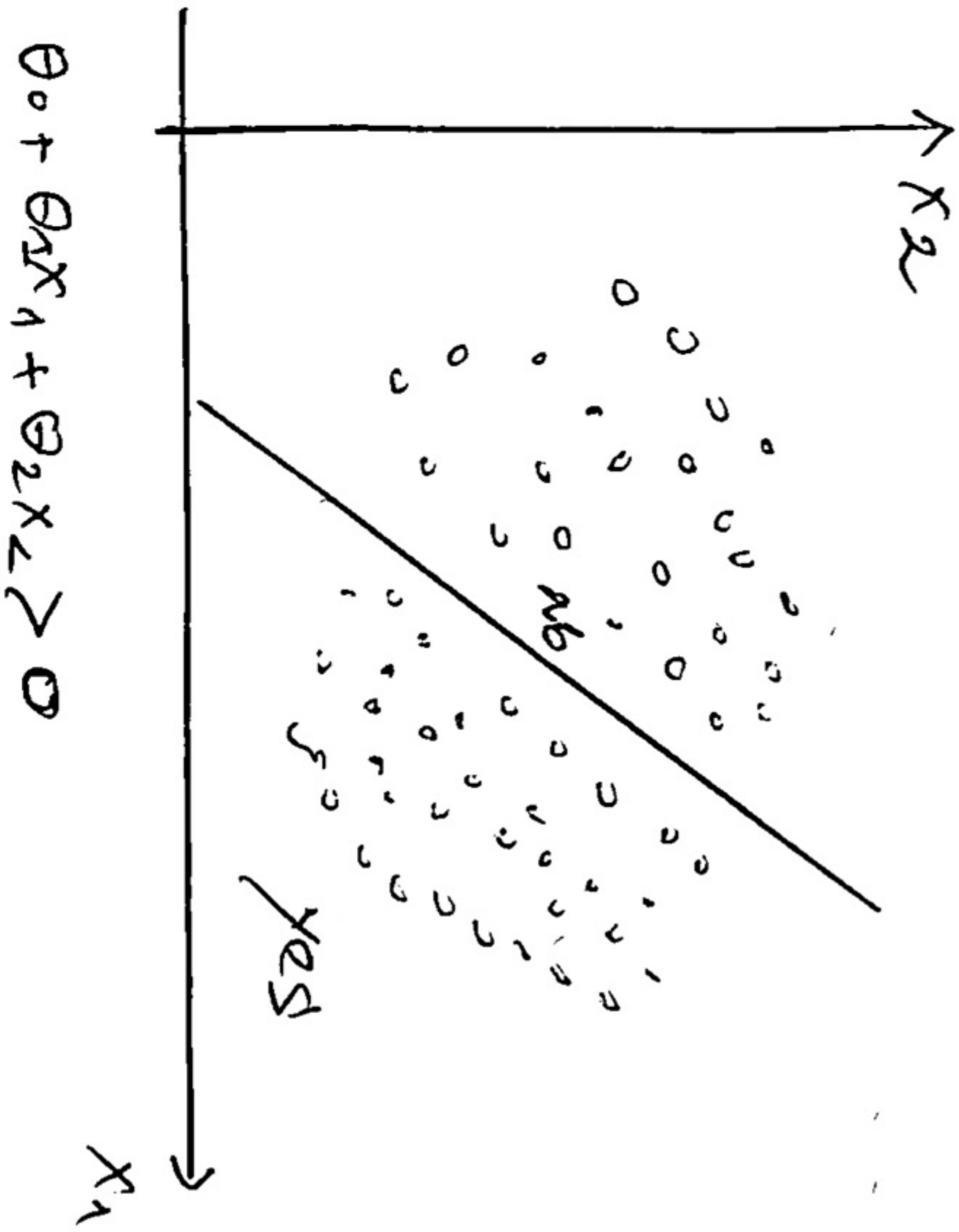


Logistic Regression \Rightarrow is a classification algorithm for categorical variables

1	yes
	no
	no
	1
	2

0	1
0	1



X : dependent variables
 Y : Independent

$X \in \mathbb{R}^{m \times n}$

$Y \in \{0, 1\}$

Prediction

$$\hat{y} = P(Y=1|x)$$

$$P(Y=0|x) = 1 - P(Y=1|x)$$

→ Logistic Regression Vs Linear Regression.

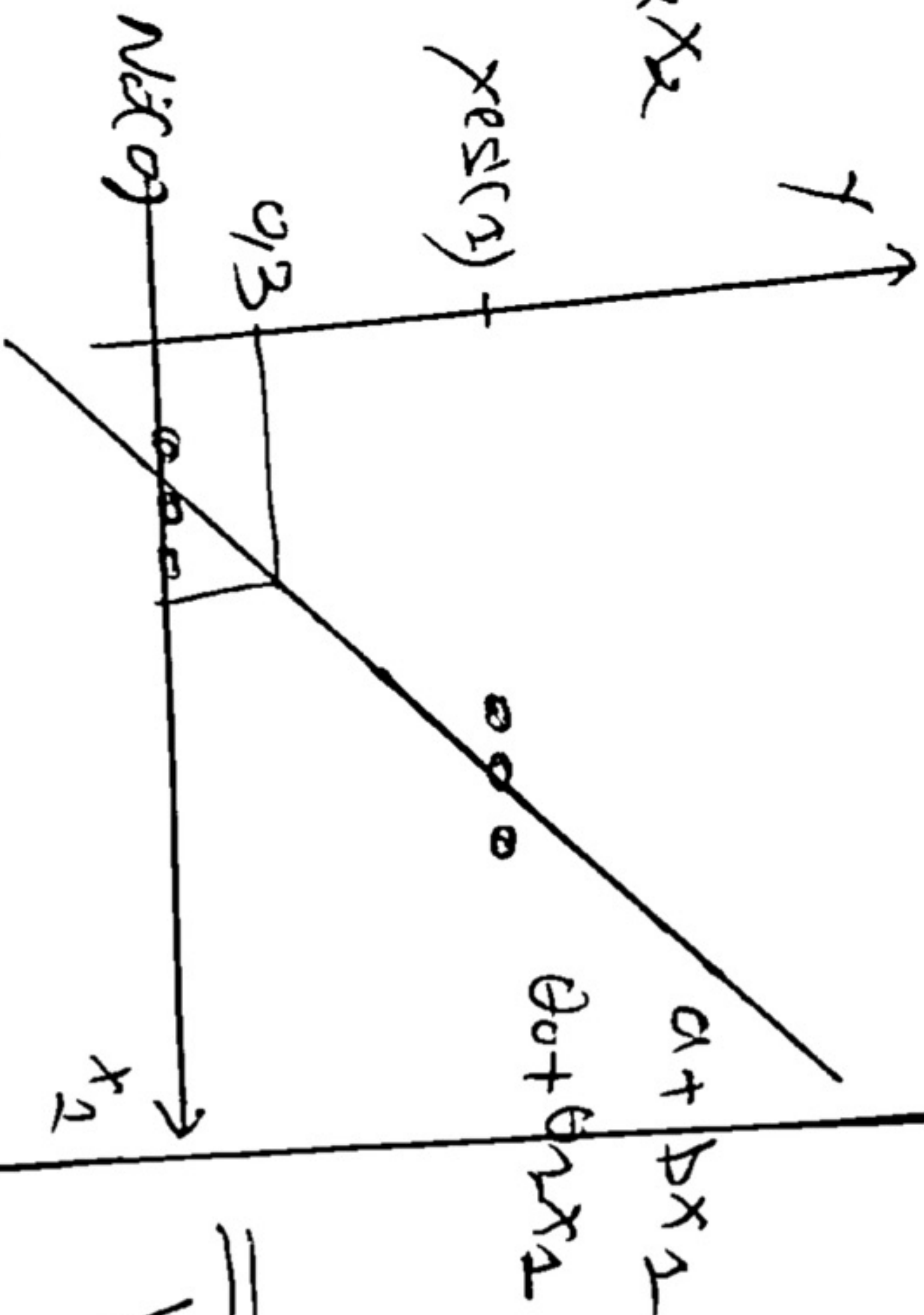
→ Linear Regression

$$\theta^T X = \theta_0 + \theta_1 x_1$$

$$\theta^T X = \theta_0 + \theta_1 x_1 + \dots + \theta_n x_n$$

$$\theta^T = [\theta_0, \theta_1, \dots, \theta_n]$$

$$X = \begin{bmatrix} x_1 \\ x_2 \\ \vdots \\ x_n \end{bmatrix}$$



Example

So if $\theta^T X = -1 + 0.1x$

$$y_1 = [13] \rightarrow \theta^T X = -1 + 0.1x_1$$

$$= -1 + 0.1x_1 \cdot 13$$

$$= 0.3$$

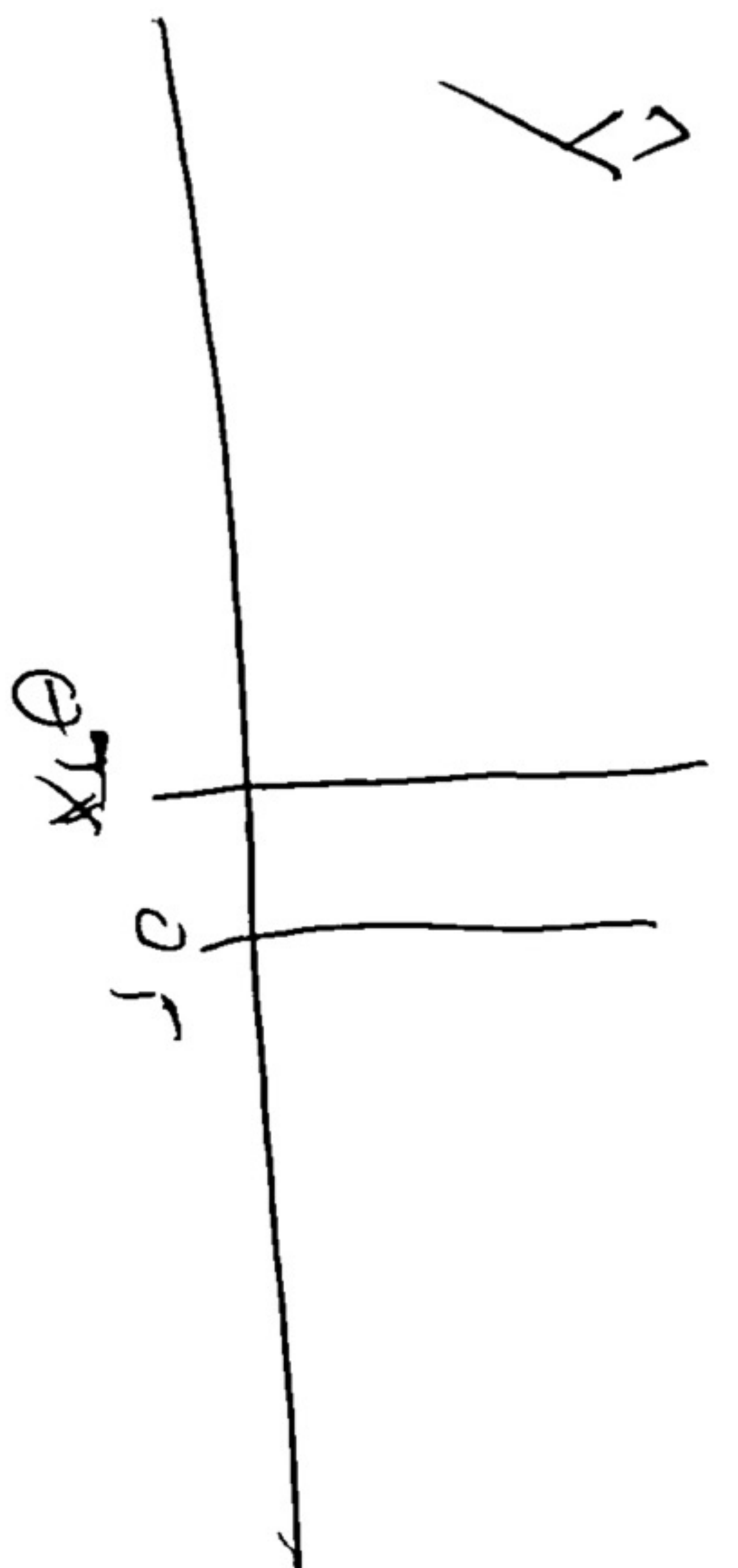
Decision

0 if $\theta^T X < 0.5 \rightarrow$ class 0

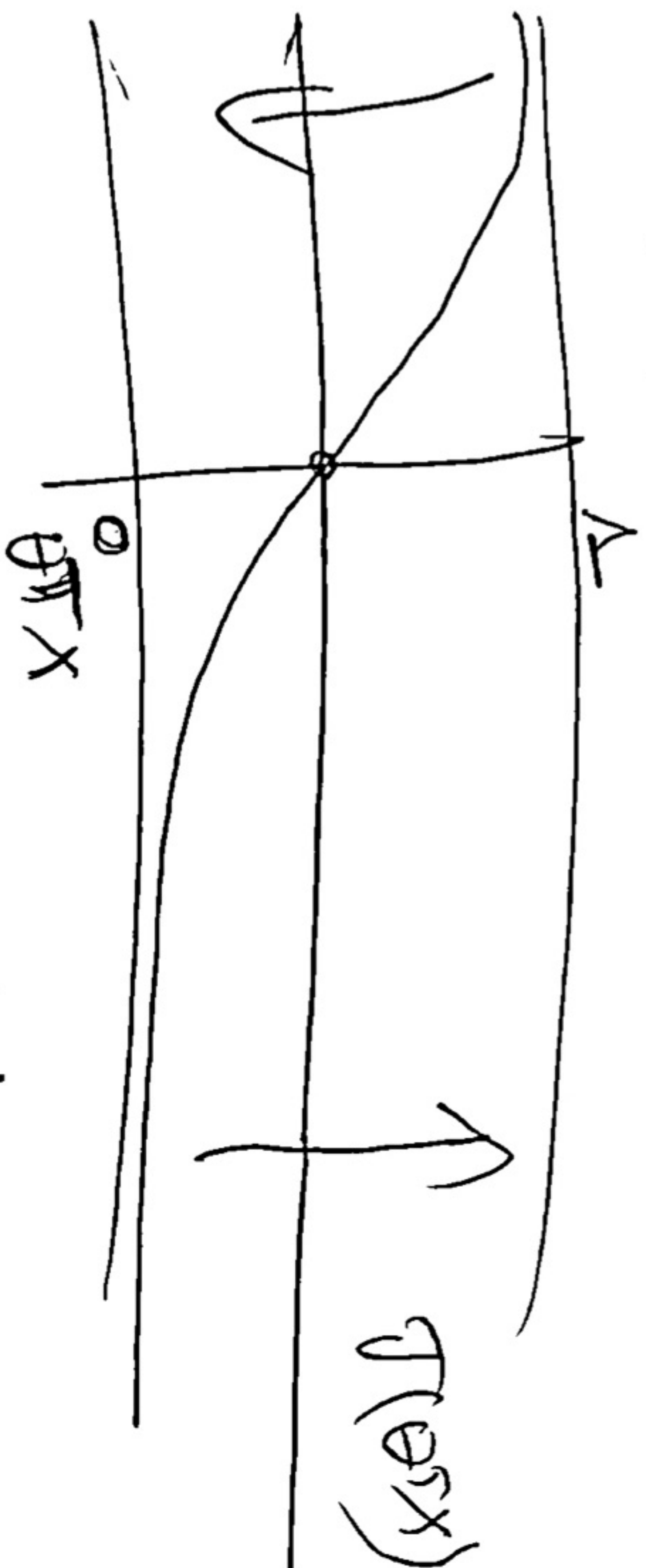
1 if $\theta^T X \geq 0.5 \rightarrow$ class 1

→ always number

classification not for



$$\sigma(\theta^T X) = \sigma(\theta_0 + \dots + \theta_n x_n)$$



$$y = \sigma(\theta^T X) \quad P(y=1|x)$$

Logistic function

$$\sigma(\theta^T X) = \frac{1 + e^{-\theta^T X}}{1 + e^{\theta^T X}}$$

$$\sigma(\theta^T X) = \frac{1}{1 + e^{-\theta^T X}}$$

FA - Score:

• Confusion Matrix

	1	0
1	TP	FN
0	FP	TN

• Precision = $\frac{TP}{TP + FP}$

• Recall = $\frac{TP}{TP + FN}$

• FA - Score = $\frac{2 \times (Prec \times Rec)}{Prec + Rec}$

FA - Score = 0,00

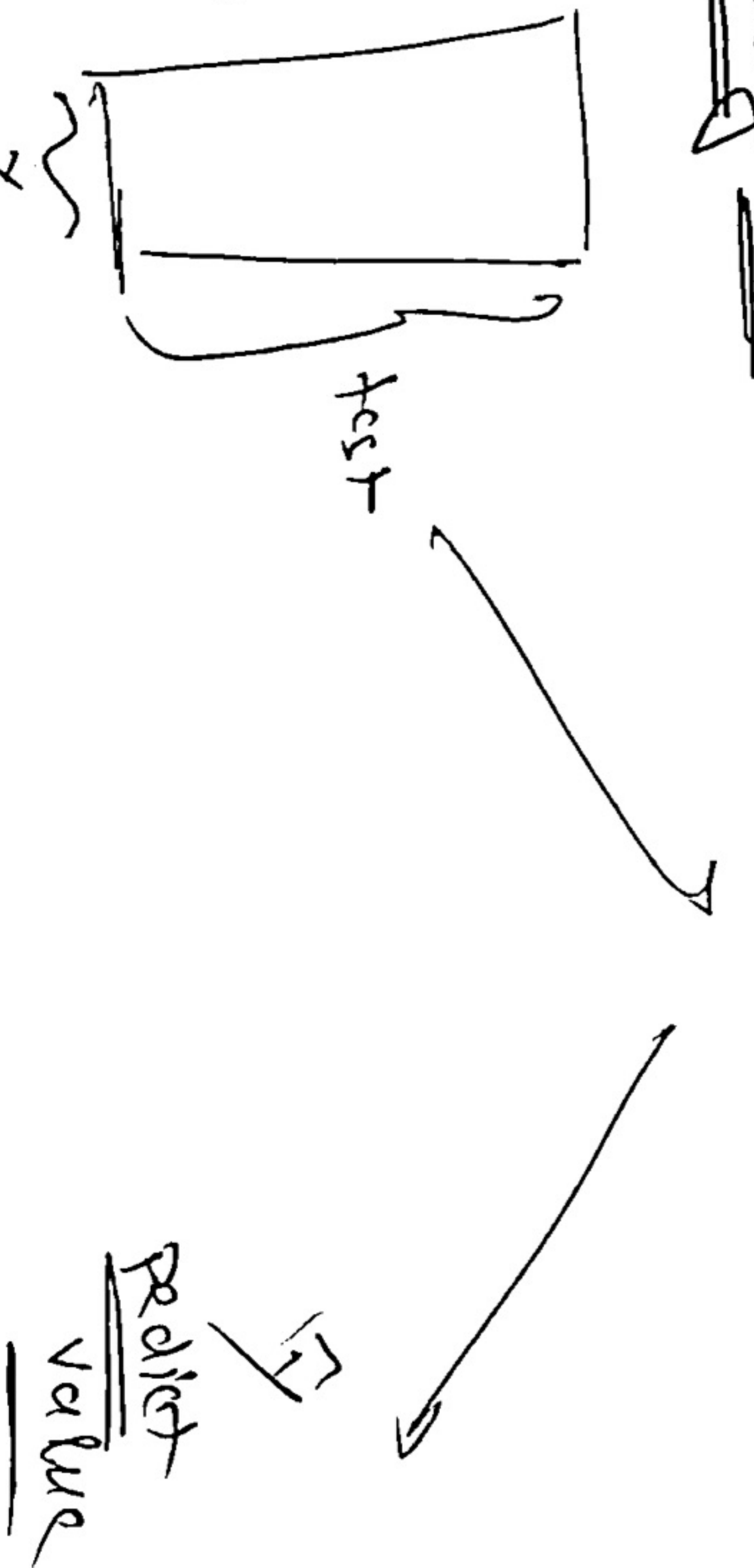
EX:

CRUN = 0
CRUN = 1

	Precision	Recall	High Accuracy / F1-Score
CRUN = 0	0,93	0,96	0,93
CRUN = 1	0,96	0,90	0,55

AVG = 969 69%

⊕ Log Loss



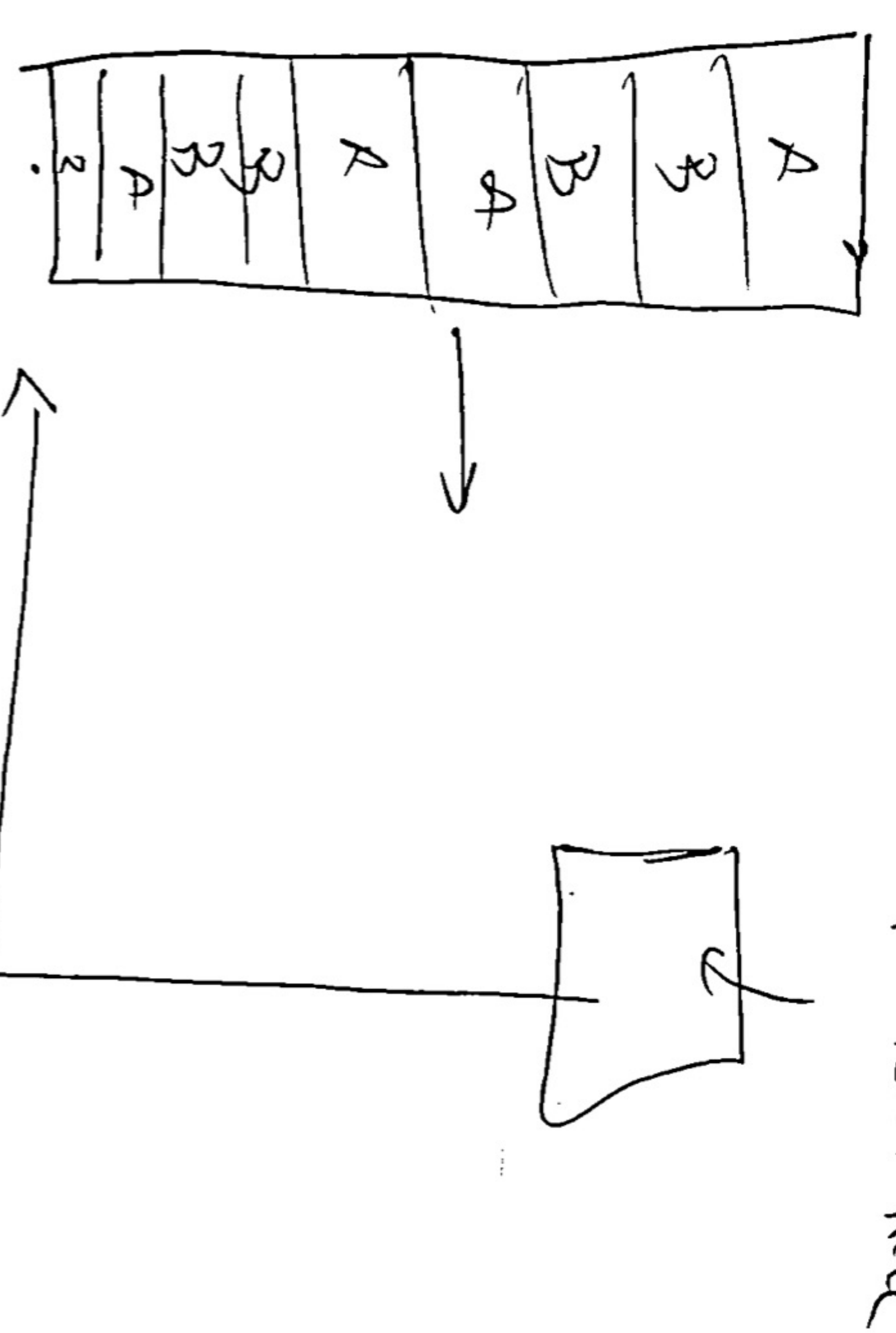
Log Loss: 0,00 → 1,00

High accuracy

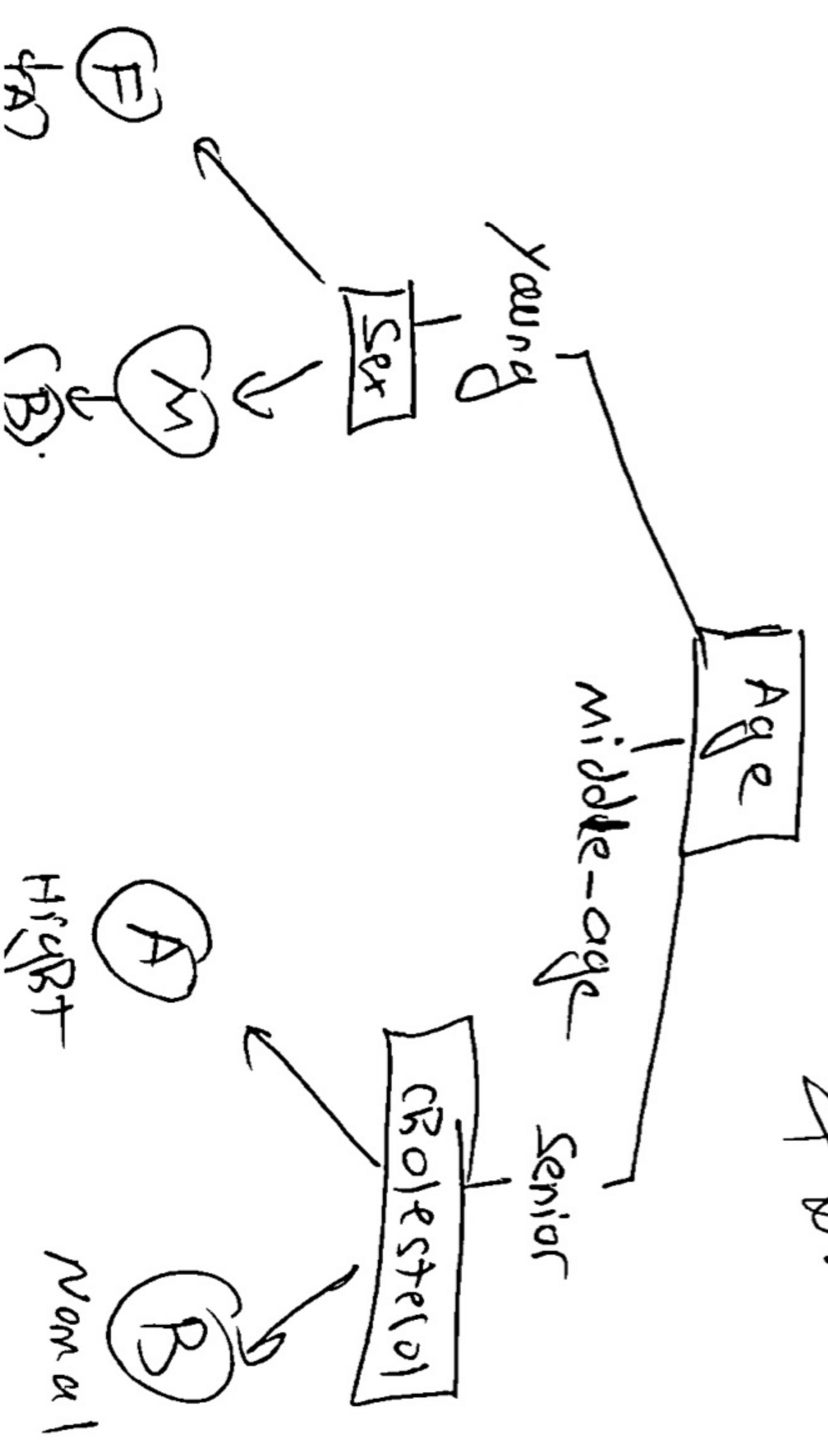
Log Loss = $-\frac{1}{n} \sum [y \log(\hat{y}) + (1-y) \log(1-\hat{y})]$

→ Introduction to:

Decision Trees



Example



Entropy:

• Measure of randomness or uncertainty: S

Example:

4 Purple A
7 Purple B

Entropy ↓

0 Purple A
2 Purple B

$S = 0$

5 Purple A
6 Purple B

Entropy ↑

$$S = -P(A) \log_2(P(A)) - P(B) \log_2(P(B))$$

Information Gain (+)

is the information that can increase the level of certainty after splitting.

$$\text{Information Gain} = (\text{Entropy before split}) - (\text{Entropy after split})$$

→ more accuracy

$$R^2 = 1 - KSE$$

$1 \Rightarrow 100\% \text{ accuracy}$
 $0 \Rightarrow \text{low}$

if $R^2 \in [0, 1] \approx \text{Good}$

→ Multiple Linear Regression

$$\hat{y} = \theta_0 + \theta_1 x_1 + \theta_2 x_2 \dots \theta_n x_n$$

$$\hat{y} = \theta^T X$$

$$\theta^T = [\theta_0, \theta_1, \dots, \theta_n] ; X = \begin{bmatrix} 1 & x_1 \\ 1 & x_2 \\ \vdots & \vdots \\ 1 & x_n \end{bmatrix}$$

↳ the same Error

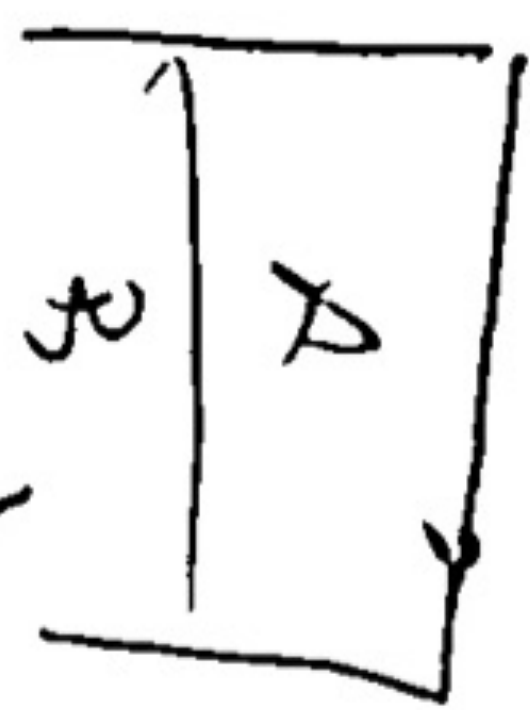
• How to estimate θ ?

- Linear algebra operation
- Gradient Descent Δ
- Proper approach Big Data

MatB please

→ scan

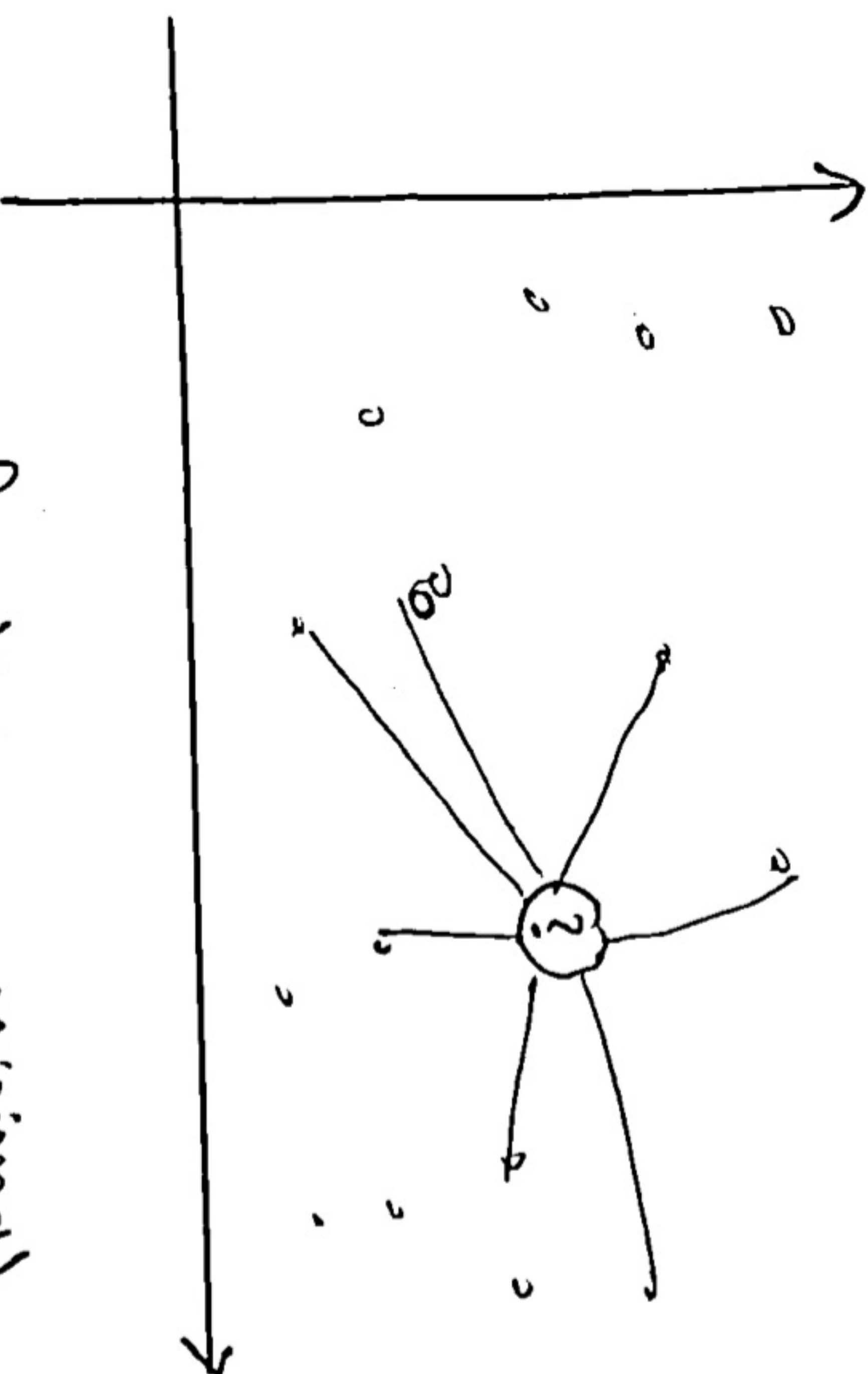
Precision tree



→ Introduction to classification

"", "A, B

① → K - Nearest Neighbors K^n



→ on noise the distance minimal

$$Dis(x_1, x_2) = \sqrt{\sum_{i=0}^n (x_{1i} - x_{2i})^2}$$

$K=0$ $n \in N$ $n \in class = [1, n]$

• Measure of randomness of uncertainty: Δ

Example:

\hat{y} : Actual labels
 \hat{y} : Predict labels

→ $E(\hat{y})$

→ Accuracy index

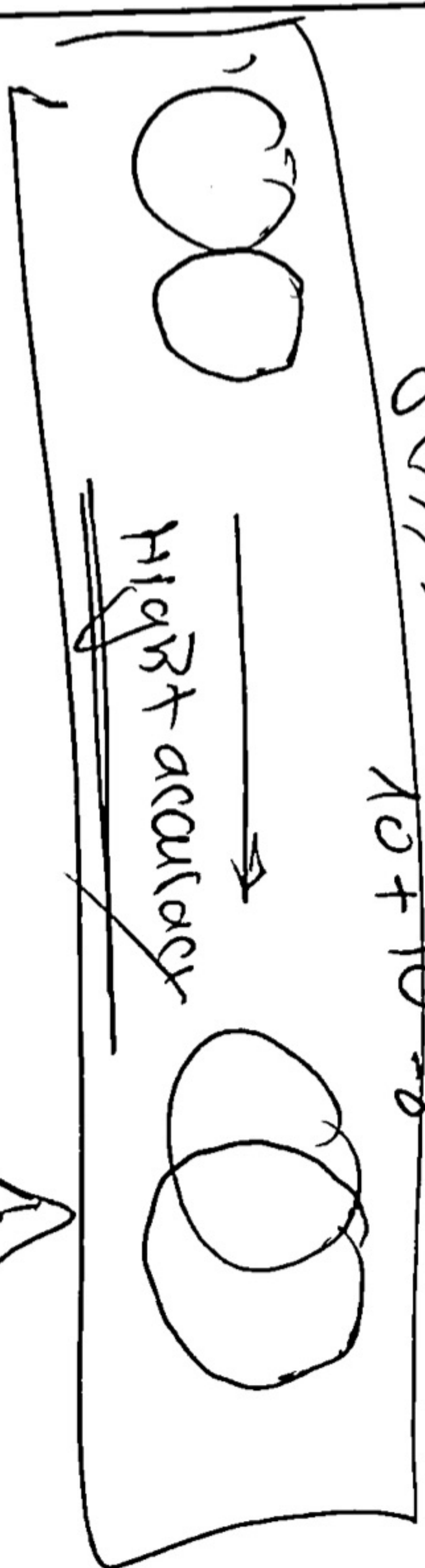
$$S(y, \hat{y}) = \frac{|y \cap \hat{y}|}{|\hat{y}|}$$

$$= \frac{|y \cap \hat{y}|}{|y| + |\hat{y}| - |y \cap \hat{y}|}$$

Example

$y = [0, 0, 0, 1, 1, 1, 1]$
 $\hat{y} = [1, 1, 0, 0, 1, 1, 1]$

$$S(y, \hat{y}) = \frac{8}{10 + 10 - 8} = 0.66 \Rightarrow 66\%$$



→ Redundant "Machine Learning"

Learning

→ Supervised learning

→ Regression: Predicting continuous values; NP

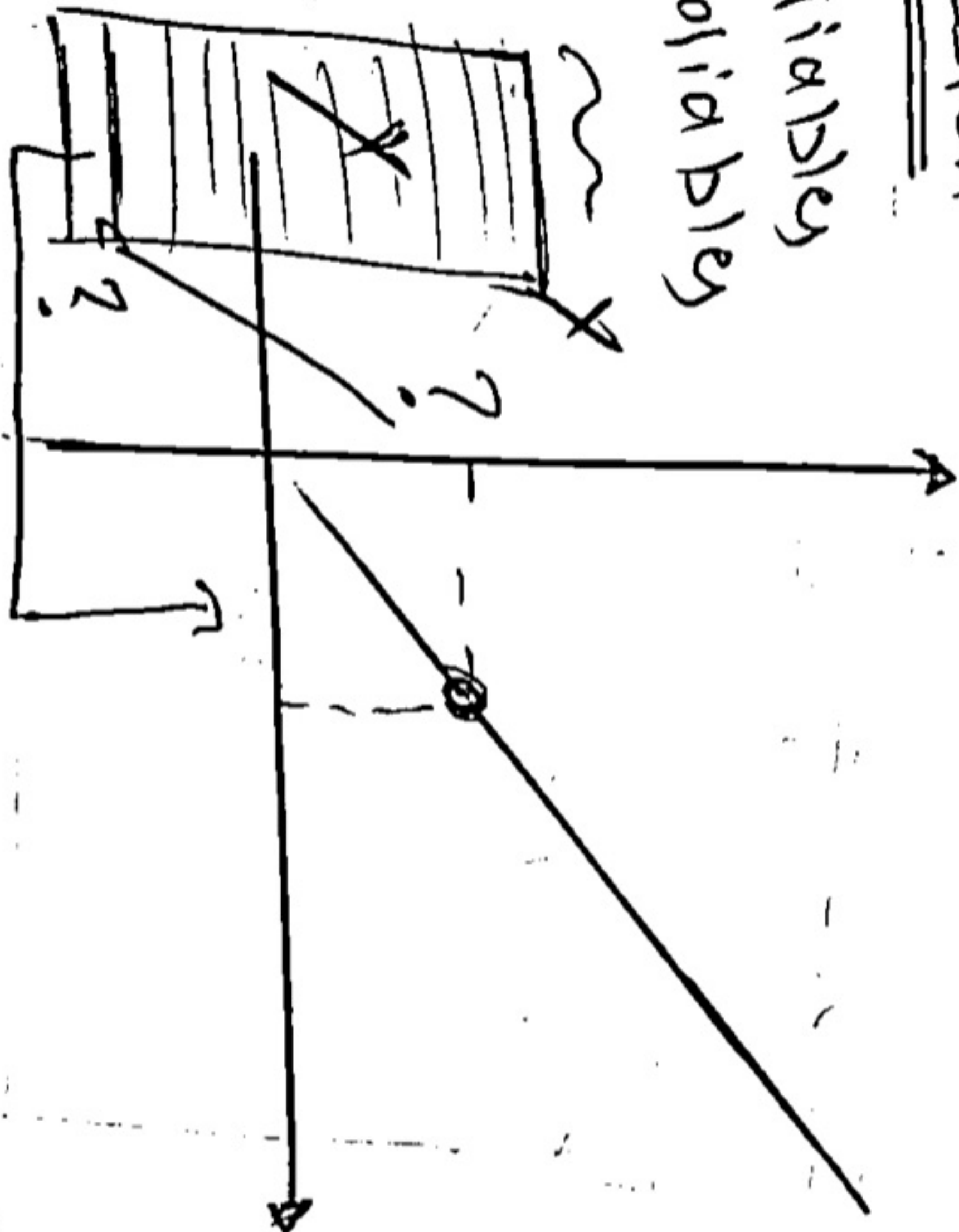
$n \in \mathbb{R}, \mathbb{N}, \mathbb{Z}, \mathbb{C}$

① → Linear regression

X : independent variables

Y : dependent variables

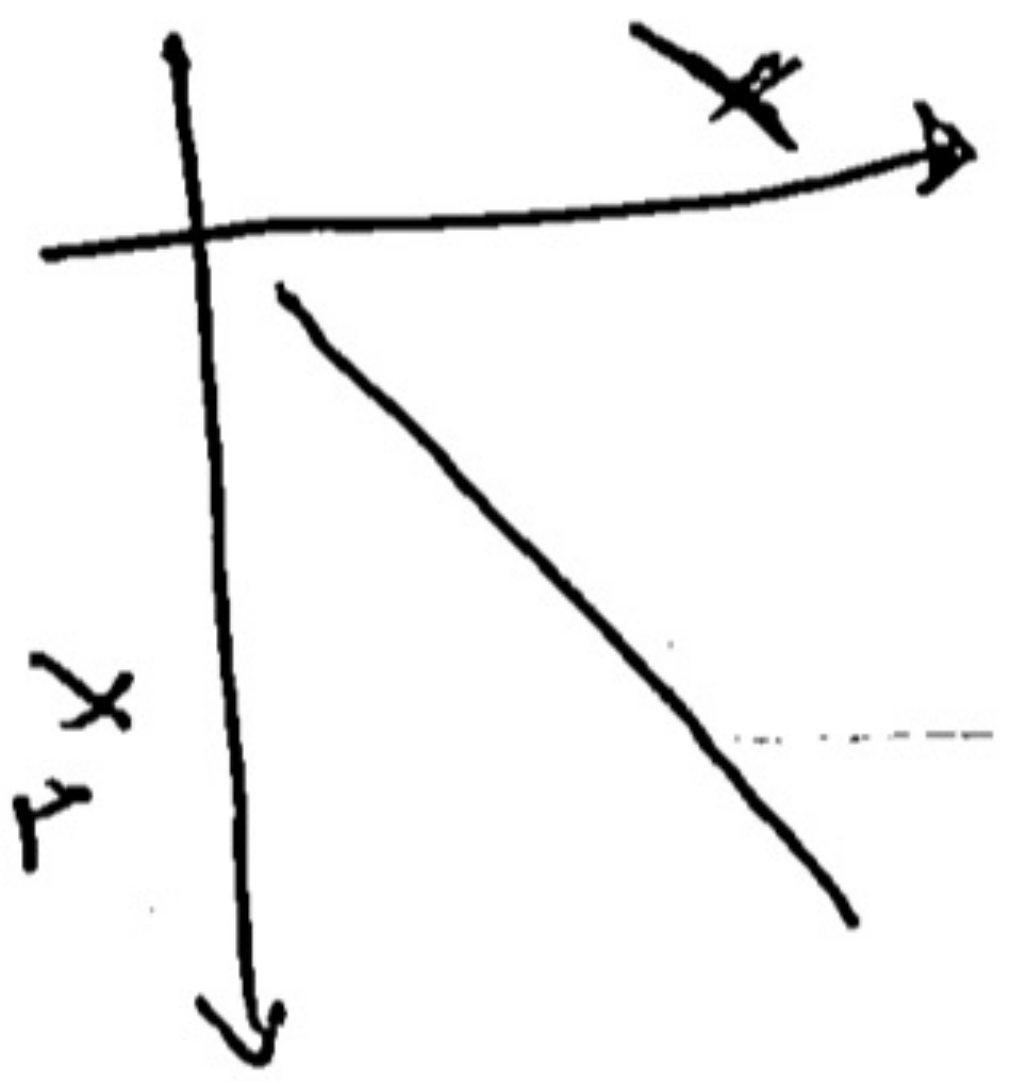
misalignment



Math (Solve) Prove

$$\hat{y} = \theta_0 + \theta_1 x_1$$

we know

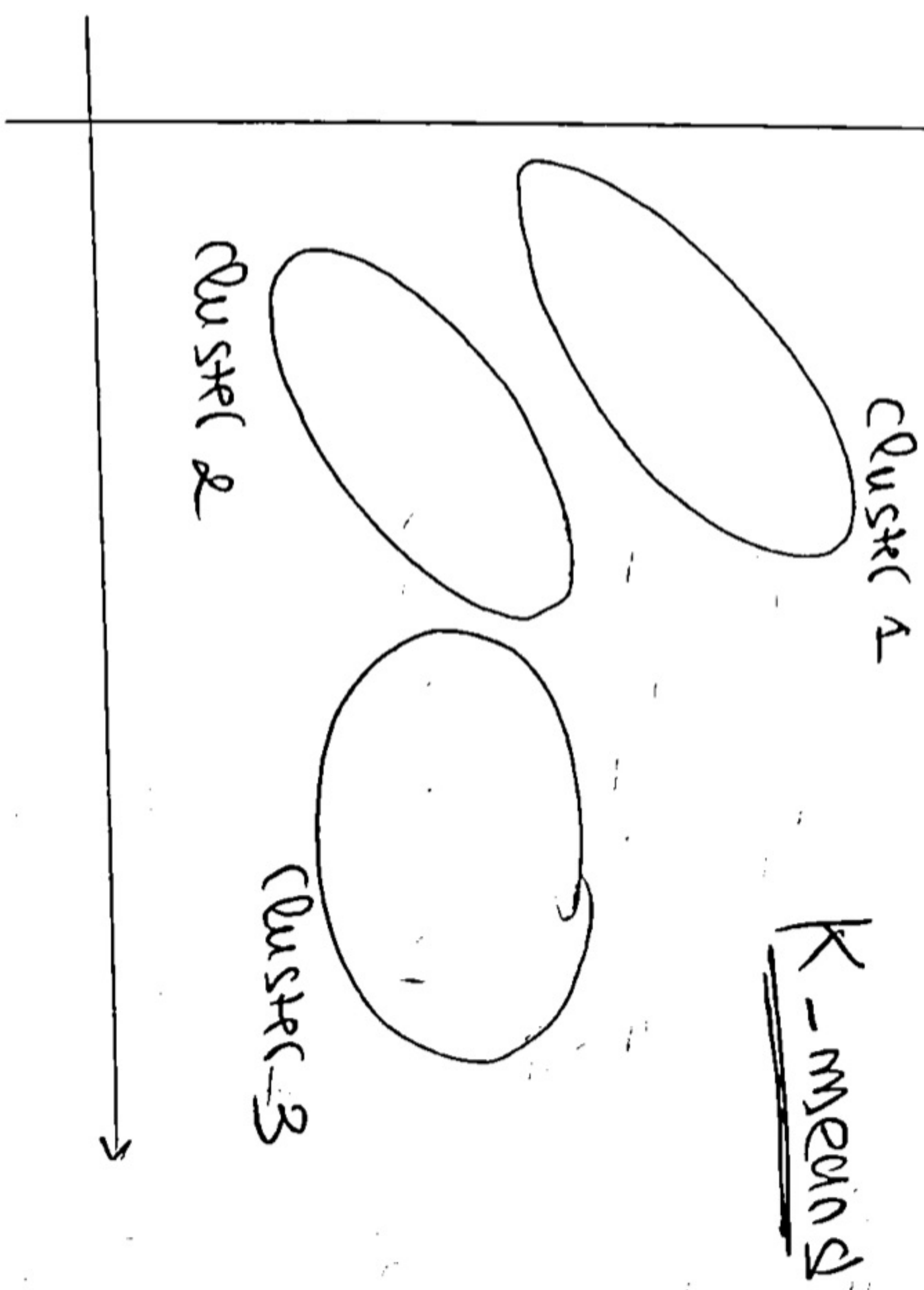


→ unsupervised learning

→ model discovery info

→ Dimension reduction, Density estimation

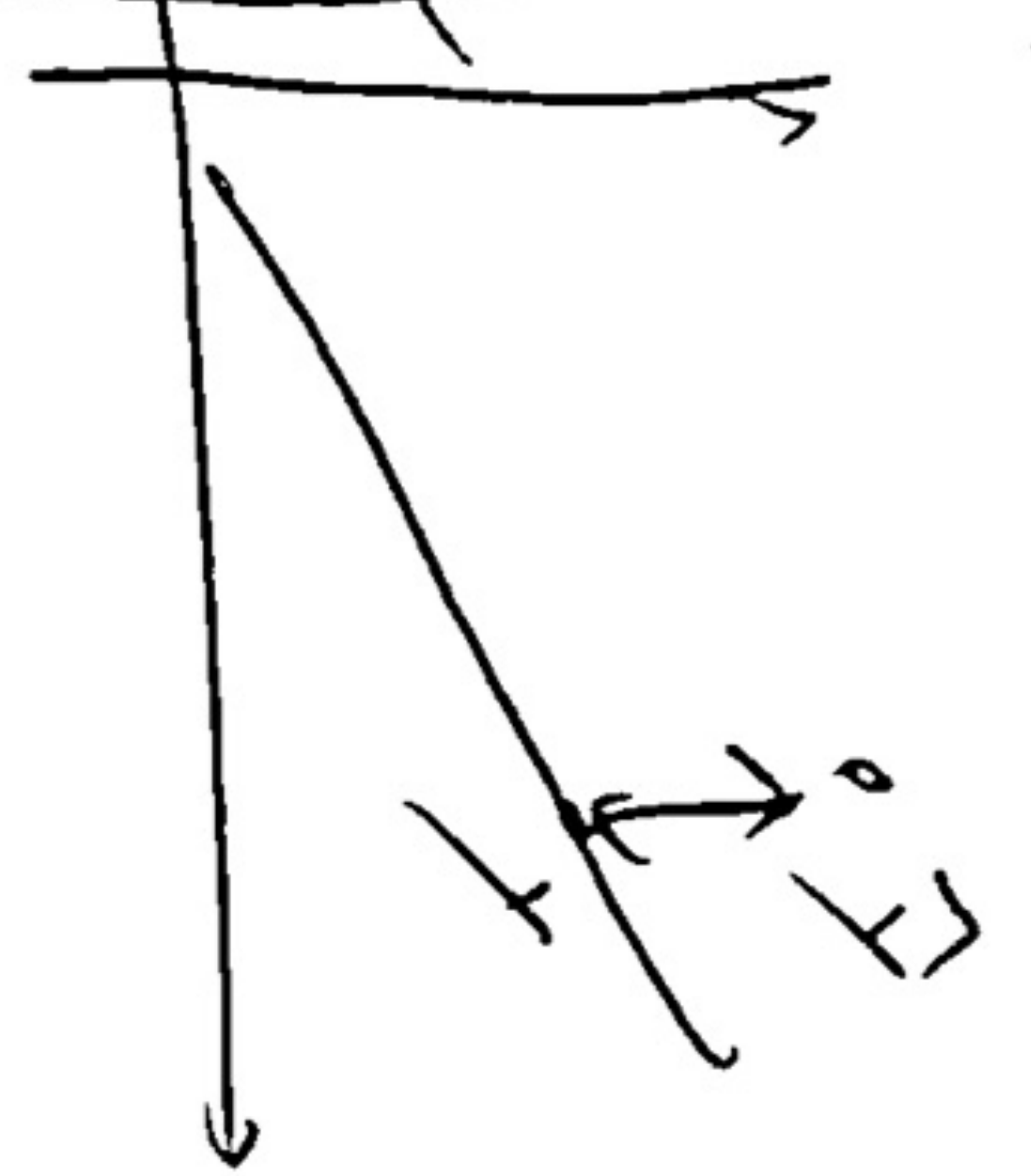
→ Clustering



→ calculate the Mean Squared Error

$$MSE = (y - \hat{y})^2$$

$$MSE = \frac{1}{n} \sum_{i=1}^n (y_i - \hat{y}_i)^2$$



→ finding the best parameters:

$$\hat{y} = \theta_0 + \theta_1 x$$

$$\theta_1 = \frac{\sum_{i=1}^n (x_i - \bar{x})(y_i - \bar{y})}{\sum_{i=1}^n (x_i - \bar{x})^2}$$

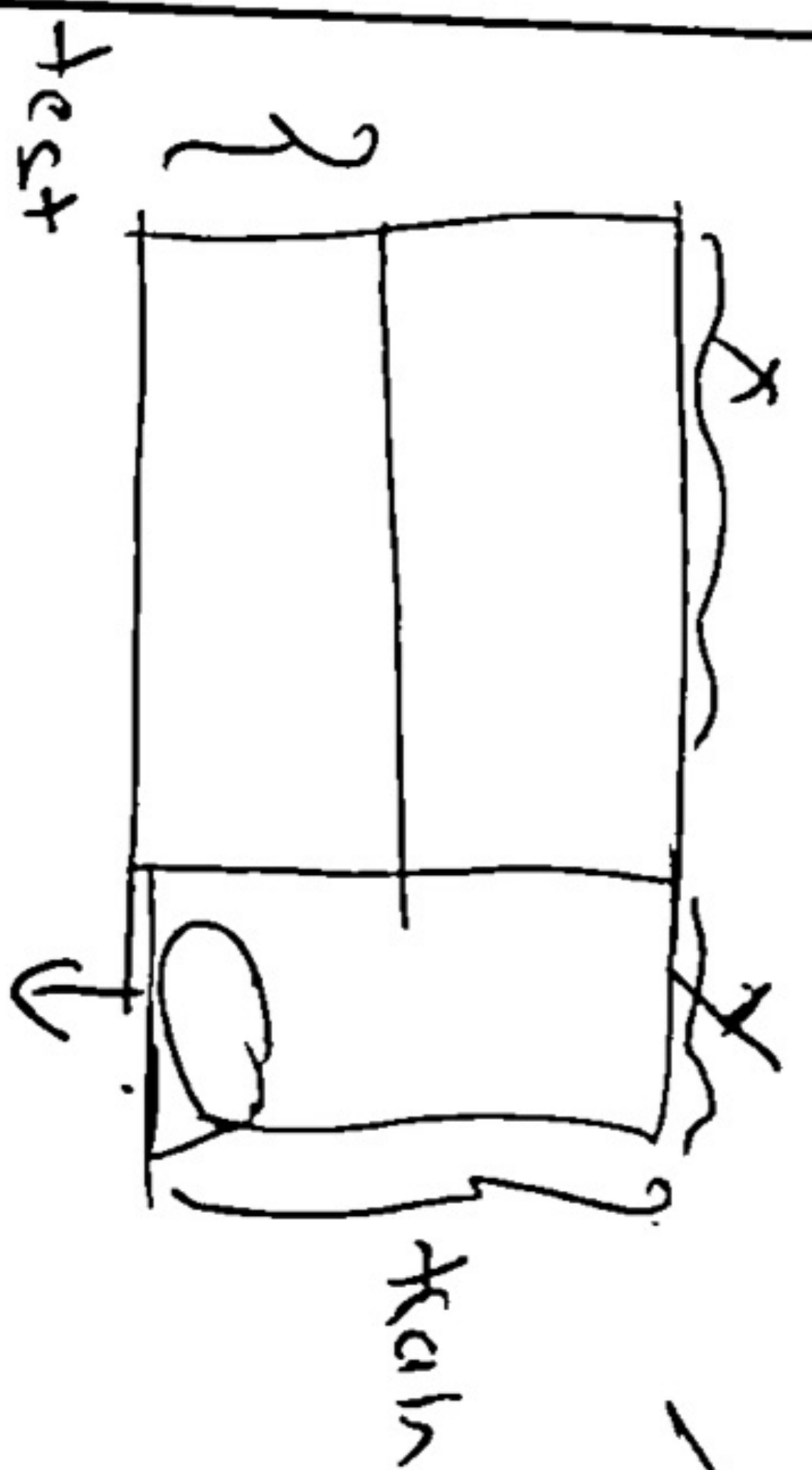
→ Pros of linear regression

- very fast
- no parameter tuning
- Easy to understand
- High accuracy

Pre-

→ Model Evaluation

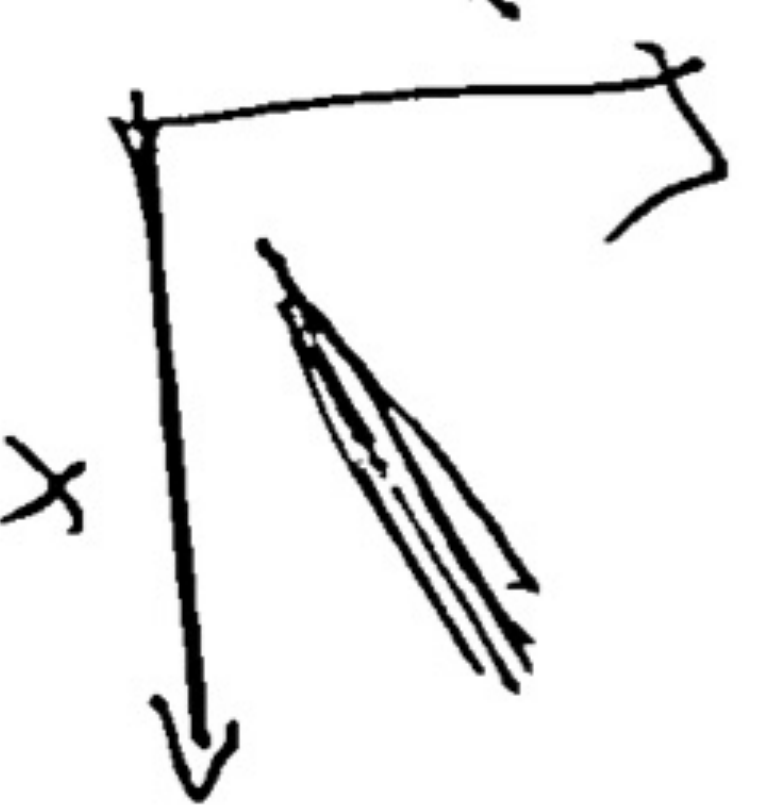
- Train/Test Data



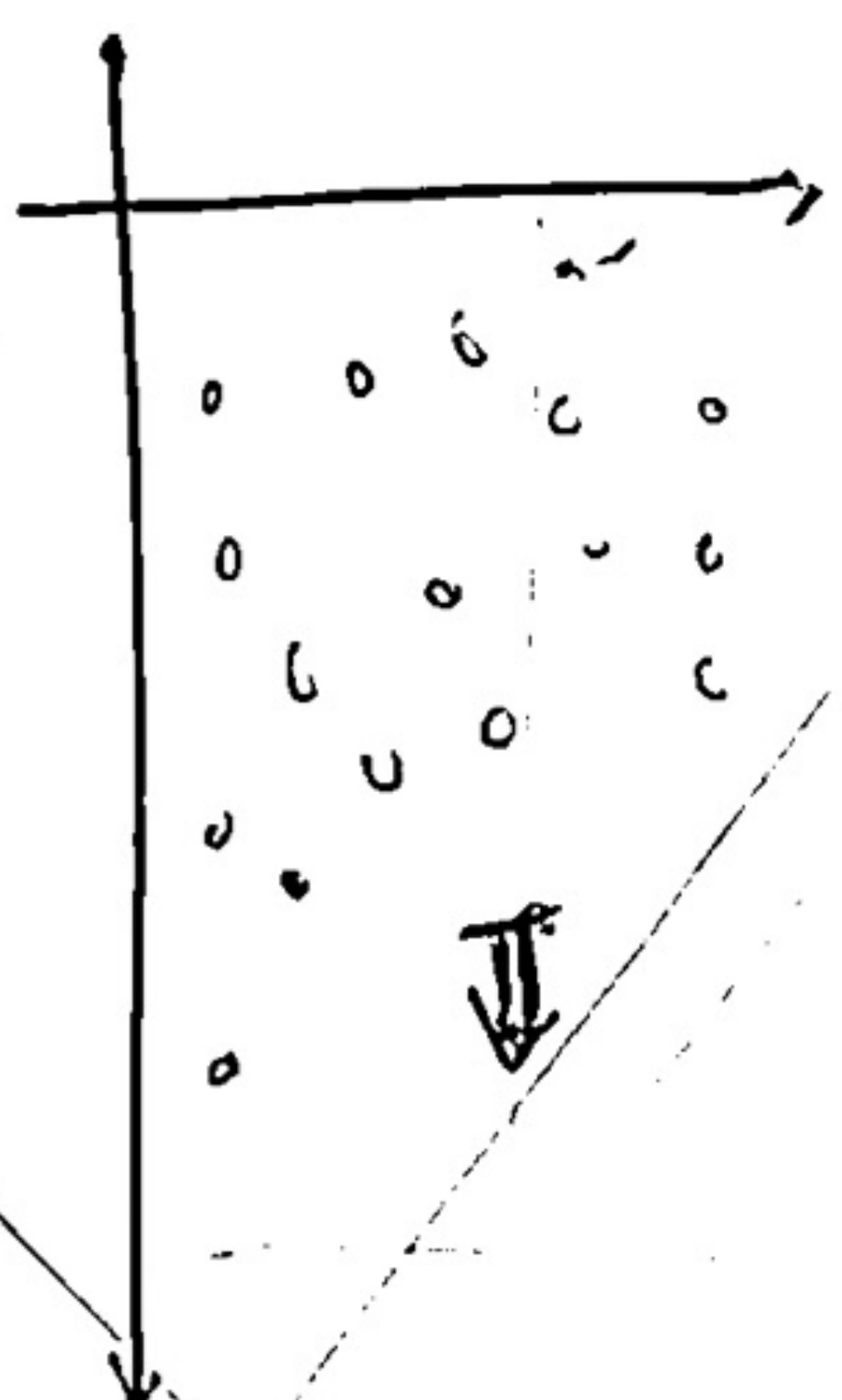
→ calculate the Error

$$Error = \frac{1}{n} \sum_{i=1}^n |y_i - \hat{y}_i|$$

Prediction

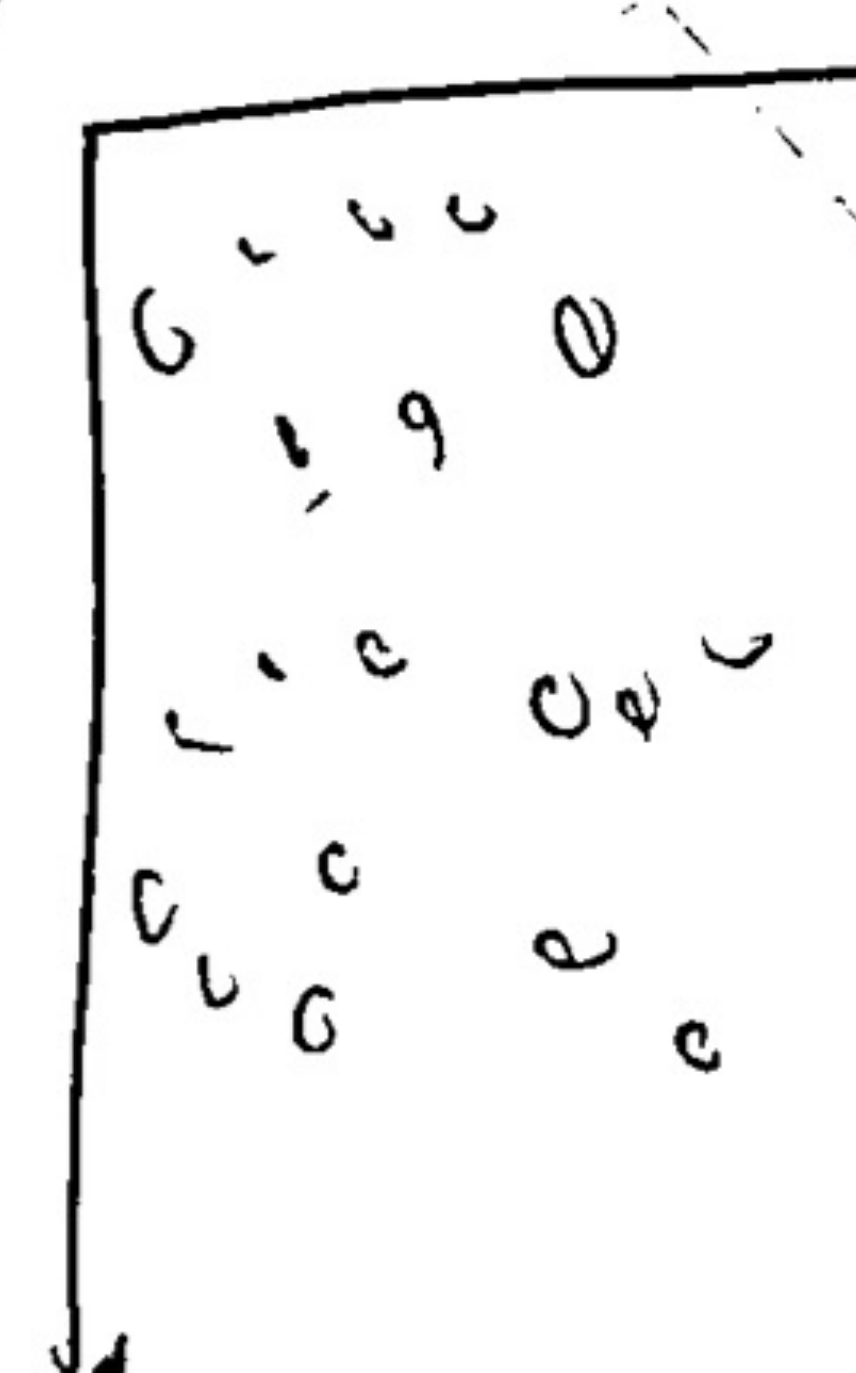


→ All DATA



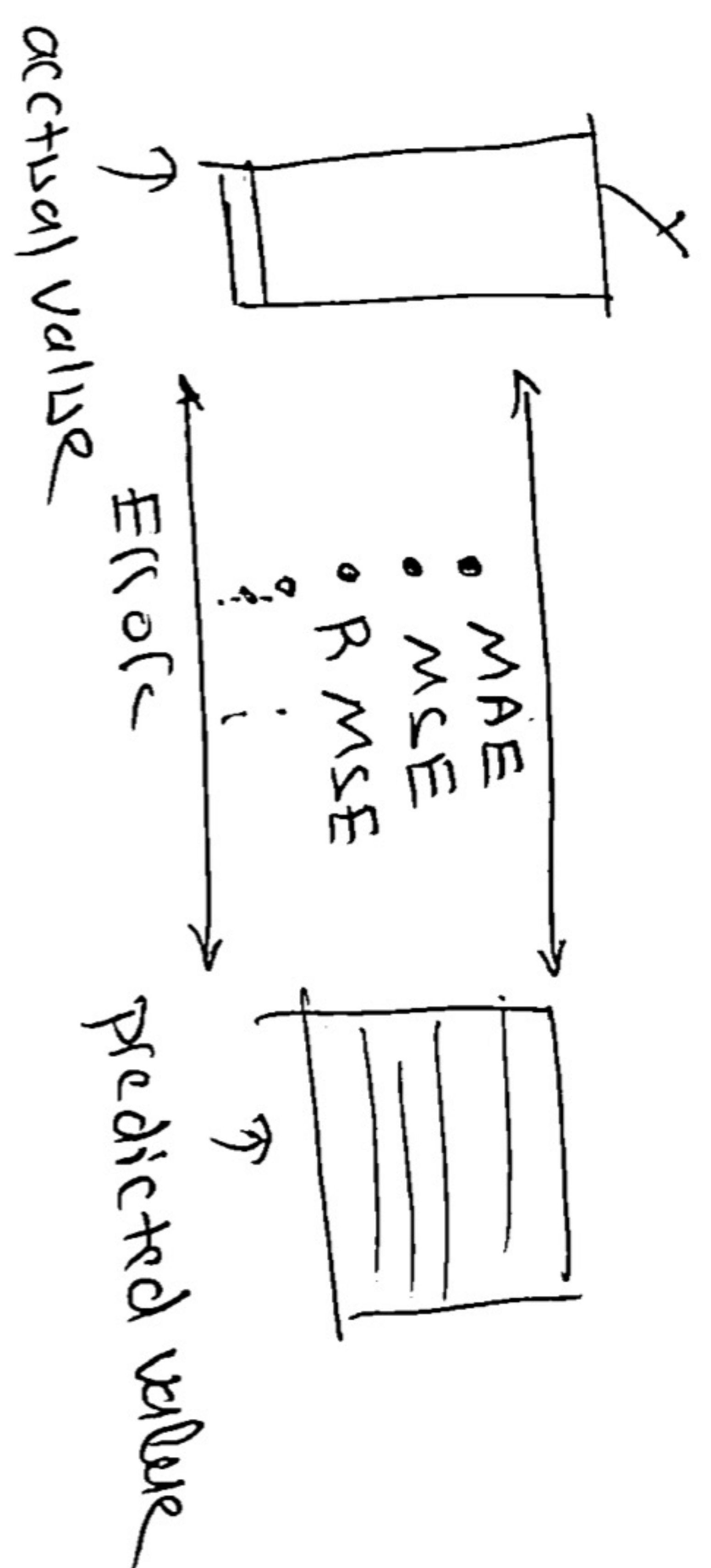
Testing set

Training set



- High Error
- Low accuracy
- overfit

Metrics in Regression Model



$$MAE = \frac{1}{n} \sum_{i=1}^n |y_i - \hat{y}_i|$$

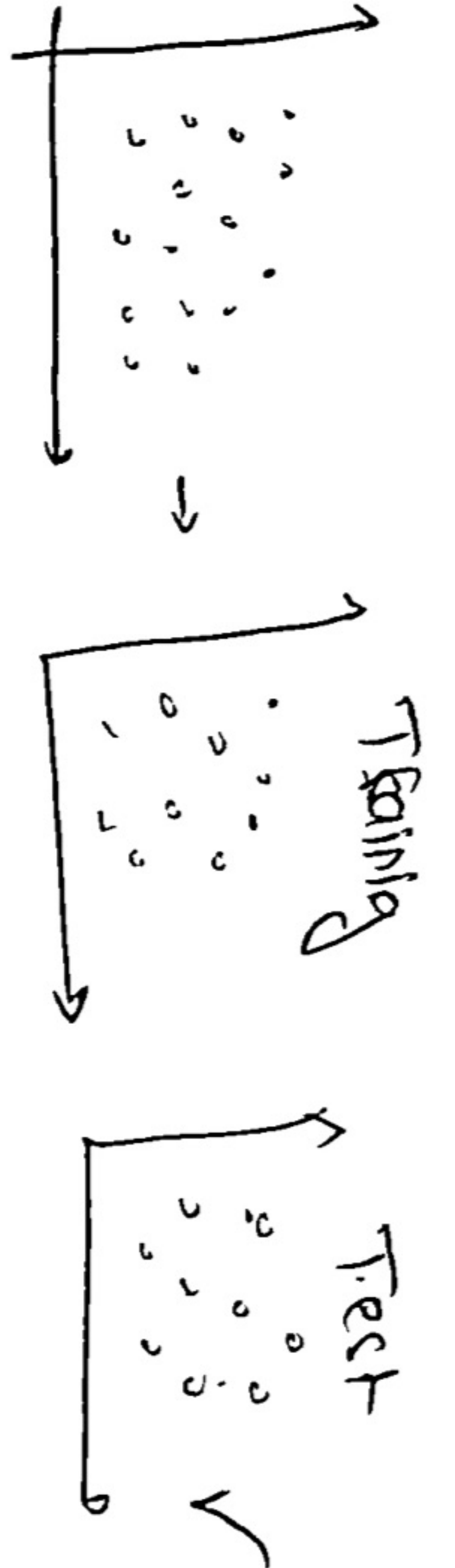
$$MSE = \frac{1}{n} \sum_{i=1}^n (y_i - \hat{y}_i)^2$$

$$RMSE = \sqrt{\frac{1}{n} \sum_{i=1}^n (y_i - \hat{y}_i)^2}$$

$$RAE = \frac{\sum_{i=1}^n |y_i - \hat{y}_i|}{\sum_{i=1}^n |y_i - \bar{y}|}$$

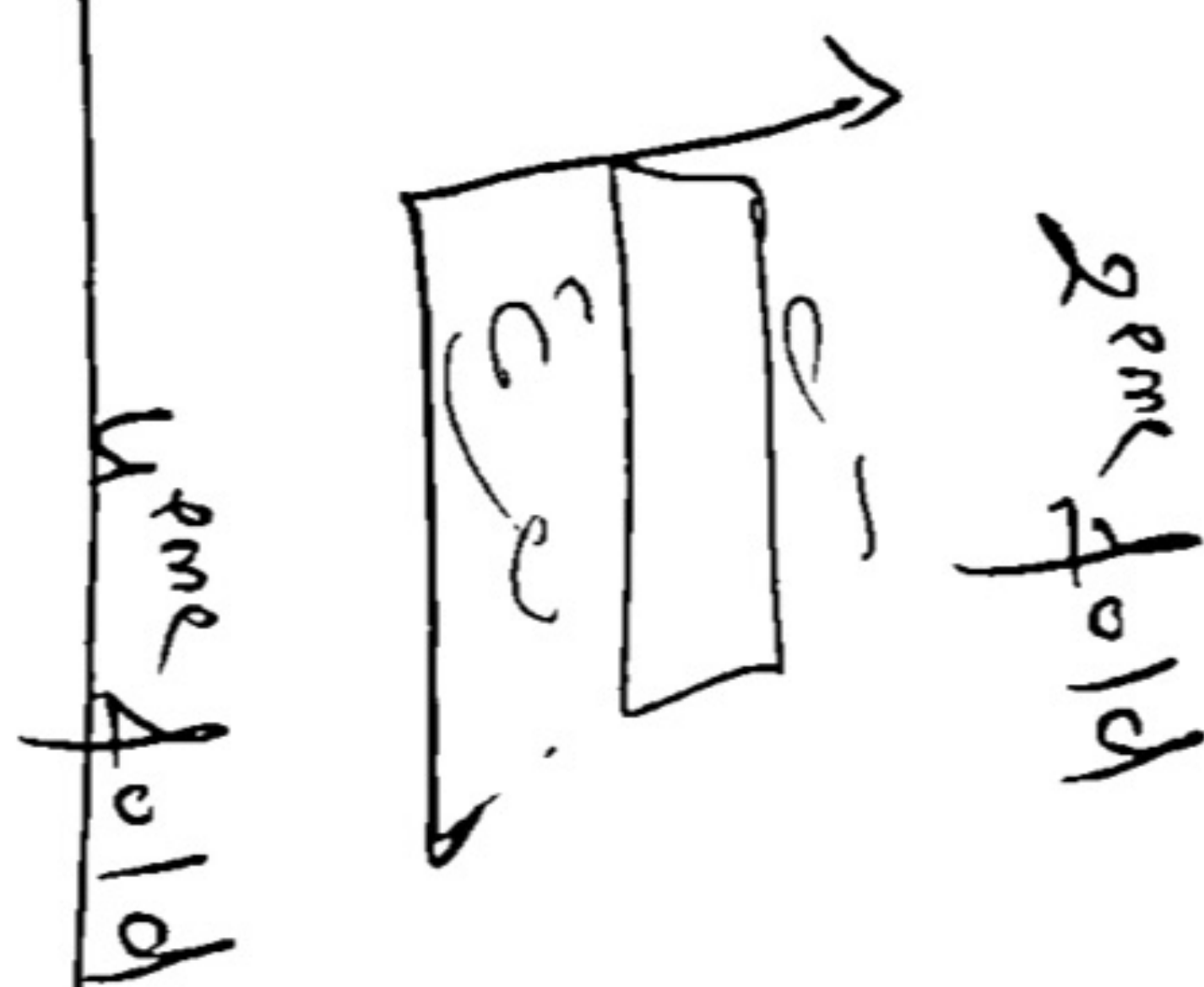
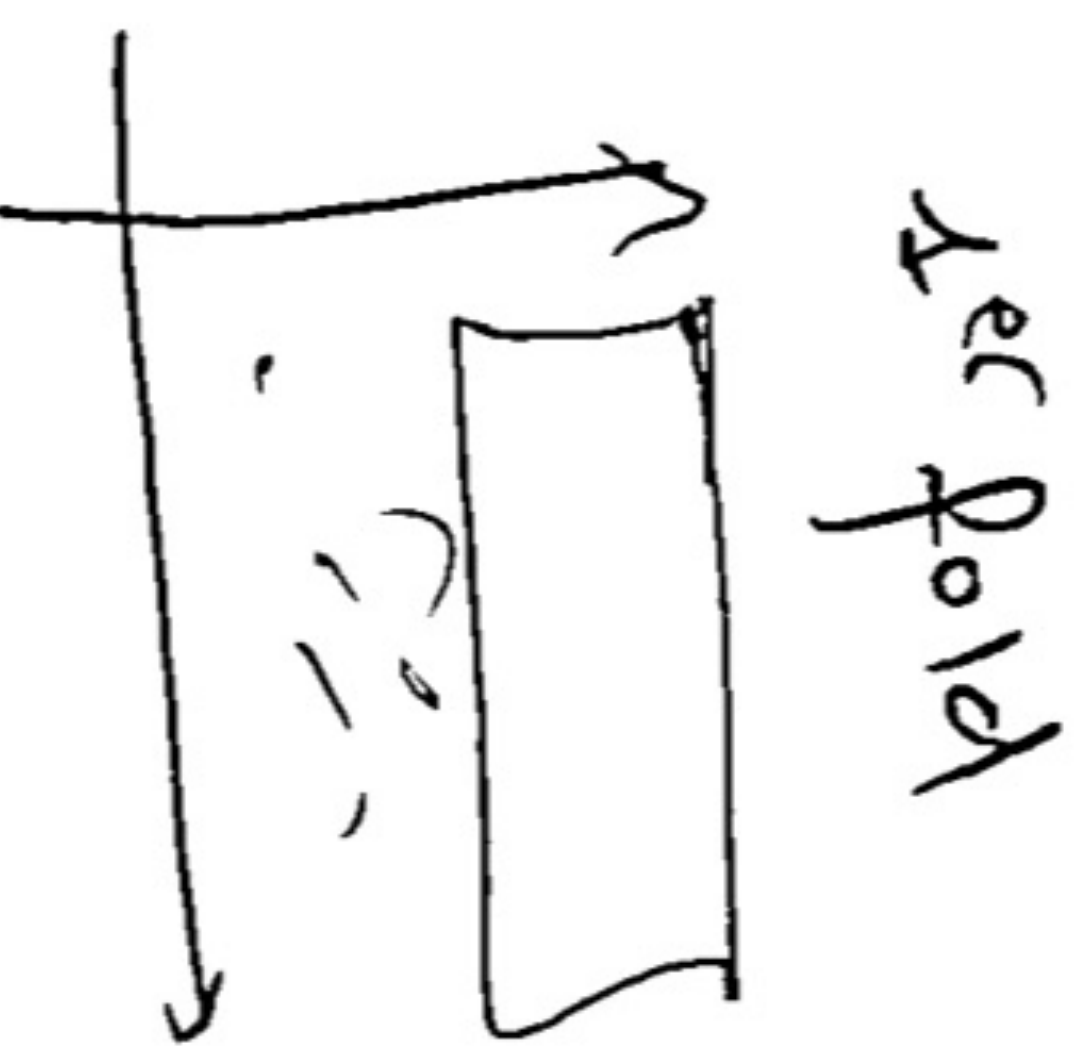
$$RSE = \frac{\sum_{i=1}^n (y_i - \hat{y}_i)^2}{\sum_{i=1}^n (y_i - \bar{y})^2}$$

MAE Precision \in

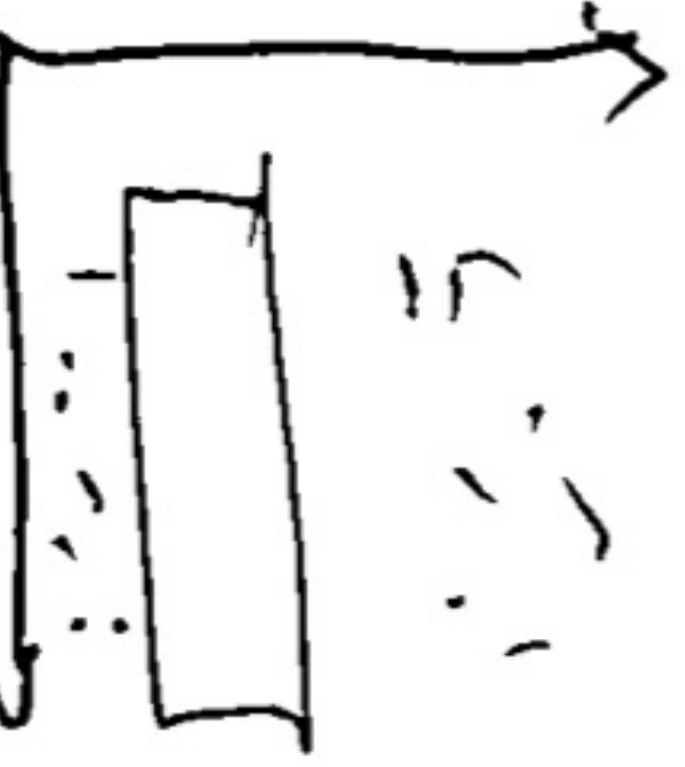


• Highest accuracy

• All points same k -fold cross-validation



3rd fold



• \square : testing
• \square : training

• Highest accuracy (80% \rightarrow 90%)