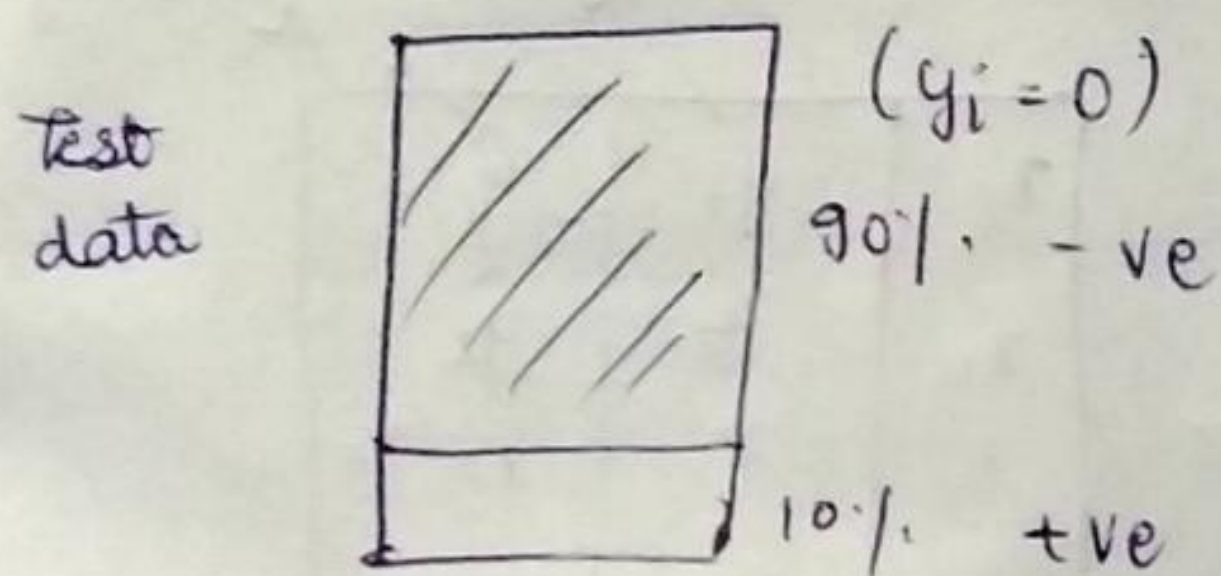


Performance prediction of models: Accuracy  $\rightarrow$  metric  
 $\downarrow$   
 classification, regression (KNN)

Accuracy =  $\frac{\text{no. of correctly classified pts}}{\text{total no. of points in } D_{\text{test}}}$   
 easy to understand measure

Performance is always measured on test data.

21/1  
 ① Imbalanced data  $\rightarrow$



dumb model  $x_q \rightarrow -ve$   
 $\{ \text{return -ve} \}$

$\rightarrow$  model  $\rightarrow$  accuracy = 90% = 0.9  
 high accuracy

So, accuracy is not <sup>at all</sup> a good measure to predict if a model is performing good or not when dataset is imbalanced.

②

	x	y	M <sub>1</sub>	M <sub>2</sub>	$\hat{y}_1$	$\hat{y}_2$
+ve {	$x_1$	1	0.9	0.6	1	1
	$x_2$	1	0.8	0.65	1	1
-ve {	$x_3$	0	0.1	0.45	0	0
	$x_4$	0	0.15	0.48	0	0

{ Test set }

M<sub>1</sub> & M<sub>2</sub> return a prob score  
KNN

$\{ x_q \rightarrow \text{prob}(y_q = 1) \}$   
 $(0 \leq p \leq 1)$

$\hat{y}$  = Predicted value

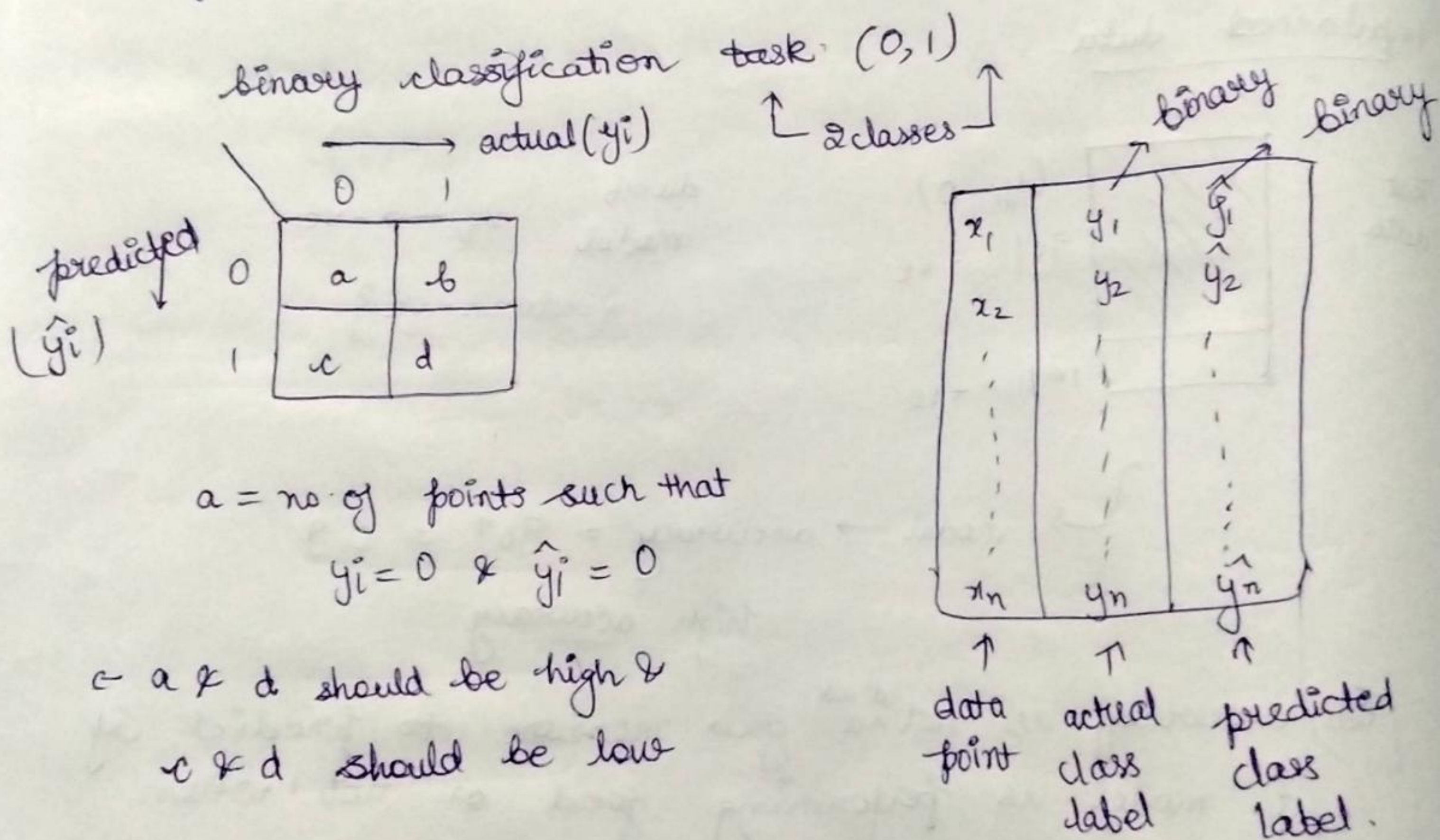


( $\Rightarrow$ ) predicted class labels are exactly the same in  $M_1$  &  $M_2$ .  
 $\Rightarrow M_1$  is better than  $M_2$  by looking at probability scores.

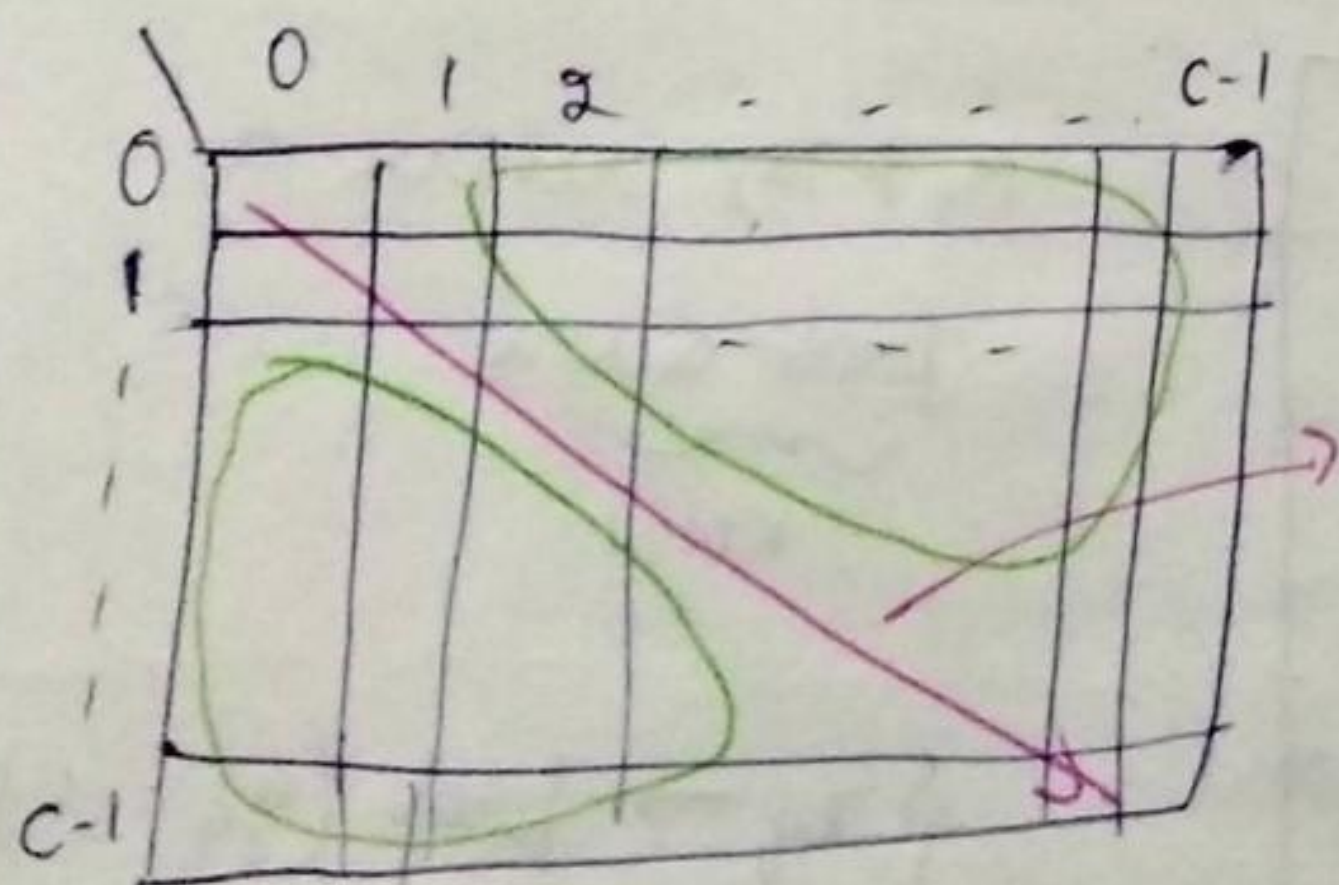
accuracy  $\rightarrow$  cannot use prob-score

$\hat{y}_1 \neq \hat{y}_2 \rightarrow$  by looking at  $\hat{y}_1$  &  $\hat{y}_2$ ,  $M_1$  &  $M_2$  are having same accuracies.

21.2  
 \* Confusion matrix is does not / cannot process prob-scores.



multiclass - classification  $\rightarrow$  (C-classes)



principal diagonal values should be high and off-diagonal values should be small for a sensible model.



	actual 0      1	
predicted 0      1	True negative	False negative
	False positive	True positive
	N	P

True positive (TP)   
 ↓   
 what is the predicted label?   
 are your predicted label matching with actual? (T/F)

$N$  = sum of false positive and true negative

$P$  = sum of false negative and true positive

$n$  = total no. of points ( $N+P$ )

$$\text{True positive Rate (TPR)} = \frac{\text{True positives}}{P}$$

$$\text{True negative Rate (TNR)} = \frac{\text{True negatives}}{N}$$

$$\text{False positive Rate (FPR)} = \frac{\text{false positives}}{N}$$

$$\text{False negative Rate (FNR)} = \frac{\text{false negatives}}{P}$$

For ex: →

	actual 0      1	
predicted 0      1	850 (TN)	6 (FN)
	50 (FP)	94 (TP)
	$N=900$	$P=100$

Test: → 900 -ve } imbalanced  
 100 +ve }

$$\uparrow \text{TPR} = 94\%$$

$$\uparrow \text{TNR} = \frac{850}{900}$$

$$\text{FPR} = \frac{50}{900} \downarrow$$

$$\text{FNR} = 6\% \downarrow$$



for a sensible model, TPR & TNR should be high &  
FPR & FNR should be low.

dumb model  $\rightarrow$  all -ve

		actual	
		0	1
predicted	0	900 TN	100 FN
	1	0 FP	0 TP
		900 = N	P = 100

900 -ve

100 +ve

$x_q \rightarrow$  -ve ; '0'

terribly low  $\checkmark$  TPR = 0%.

$\checkmark$  TNR = 100%.

FPR = 0%  $\checkmark$

FNR = 100%  $\times$

Goal  $\rightarrow$  To diagnose a cancer patient if he/she has cancer or not.

Domain specific

		act	
		0	1
pred.	0	TN	FN
	1	FP	TP

our objective is not to miss any cancerous patient.

Low FNR should be there as, FNR means to say that patient does not have cancer but it actually has cancer.

If FPR is high  $\rightarrow$  that's okay also, as if a patient doesn't have cancer and it predicted to be cancer patient then he/she will go through more powerful tests to prove him



non-cancerous.

interpretable in english sentences.

21-3

\* Precision, Recall & F<sub>1</sub>-Score :->

		act 0 → 1	
pred ↓ 0 1	0	TN	FN
	1	FP	TP
		N	P

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

of all the points the model declared / predicted to be +ve, what %age of them are actually +ve

Recall :-> same as  $TPR = \frac{TP}{P}$

↓  
of all the actual positive points, what %age of them, are predicted positive by the model.

We want precision to be high & Recall should be high.

Precision ↑  
(0-1)

Recall ↑  
(0-1)

F<sub>1</sub>-Score is combination of Precision & Recall.

$$F_1\text{-Score} = \left( 2 * \frac{\text{precision} * \text{recall}}{\text{precision} + \text{recall}} \right)$$

↓  
harmonic mean  
of precision &  
recall

F<sub>1</sub>-score ↑ if precision ↑ & recall ↑.



31.4 Receiver Operating characteristic curve & AUC  
(ROC) curve.  $\hookrightarrow$  designed by electronics & radio engineers during second world war.

$\hat{y}_{z_2} = 0.92$   $\hat{y}_{z_1} = 0.95$

	$\hat{y}$	$y$	$\hat{y}$
1	1	$x_1$	1
1	0	$x_2$	1
0	0	$x_3$	0
0	0	$x_4$	1
0	0	$x_5$	1

$\rightarrow z_1$   
 $\rightarrow z_2$

model  $\rightarrow 1, 0$

gives score (like prob-score)

① Sort your data in decreasing order of  $\hat{y}$

② Thresholding  $\rightarrow$  Take any value

( $\tau$ )  
 $\downarrow$   
low

①  $\tau_1 = 0.95$

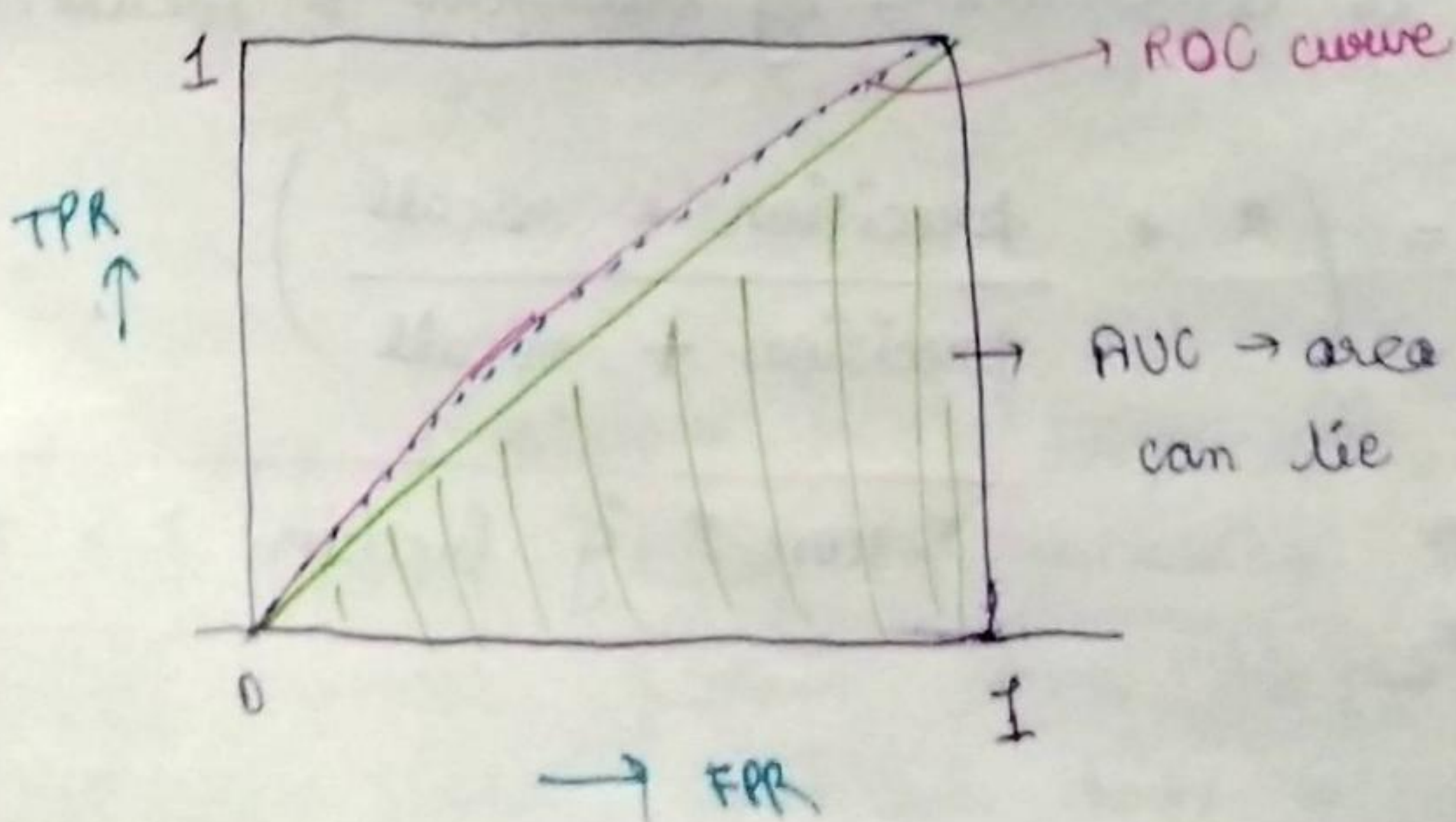
if  $\hat{y} \geq \tau_1$   
①

else ②

For each  $\tau_1, \tau_2, \tau_3, \dots, \tau_n$

$\downarrow \downarrow$   $FPR_2, TPR_2$  and so on.  
we can easily get

$FPR_1, TPR_1$



AUC  $\rightarrow$  area under curve  
can lie b/w ① to ①

$\downarrow \downarrow$   
terrible  $\rightarrow$  good



AUC  $\rightarrow$

① Imbalanced data  $\rightarrow$  AUC can be high with a dumb model too / Simple model.

② AUC is not dependent on the  $\hat{y}$  scores but it depends on the decreasing order.

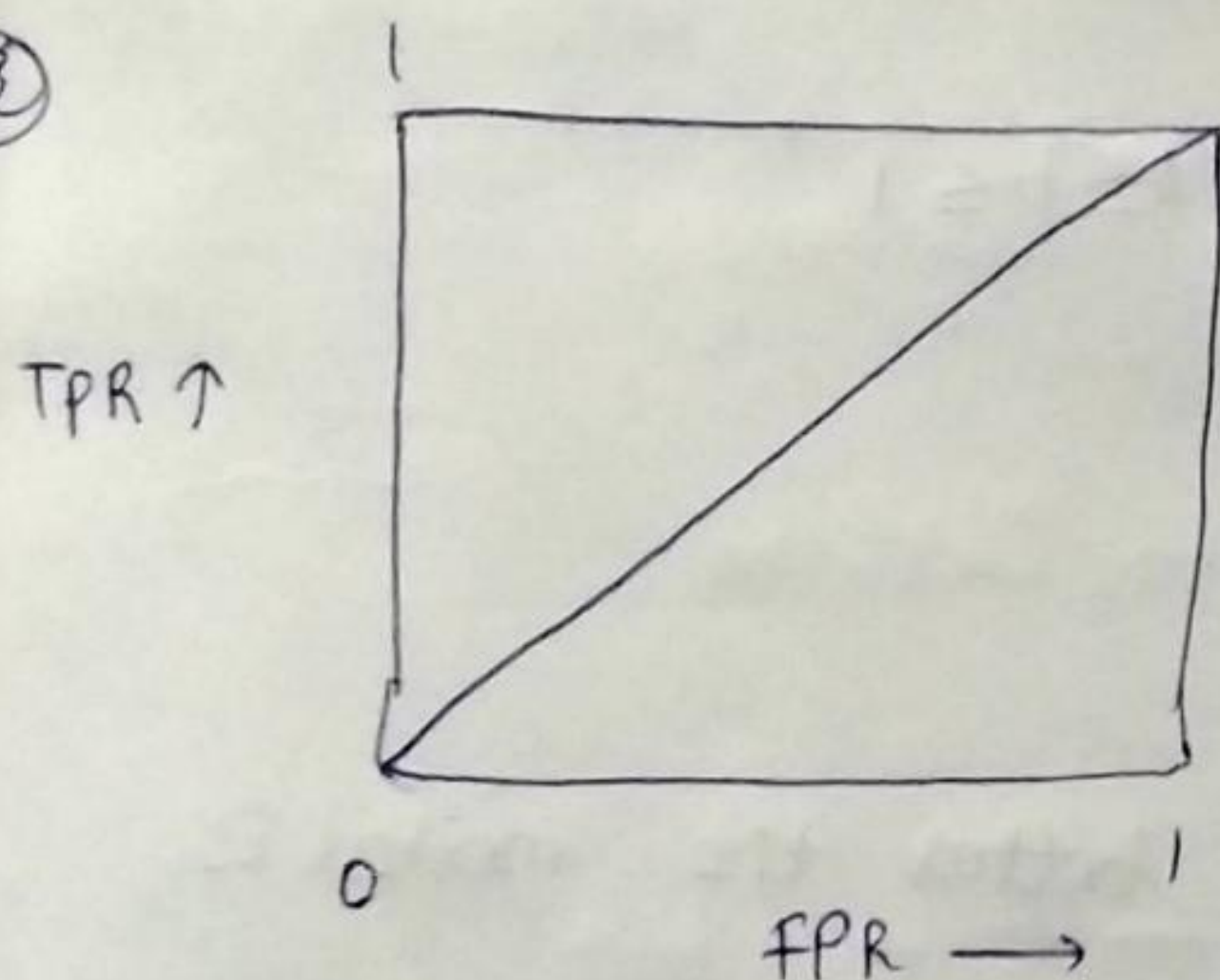
		$M_1$	$M_2$
$x_1$	1	0.95	0.2
$x_2$	1	0.92	0.1
$x_3$	0	0.80	0.08
$x_4$	1	0.76	0.07
$x_5$	1	0.71	0.06

$$AUC(M_1) = AUC(M_2)$$

as AUC only depends on order.

and not at all depend on actual value

③



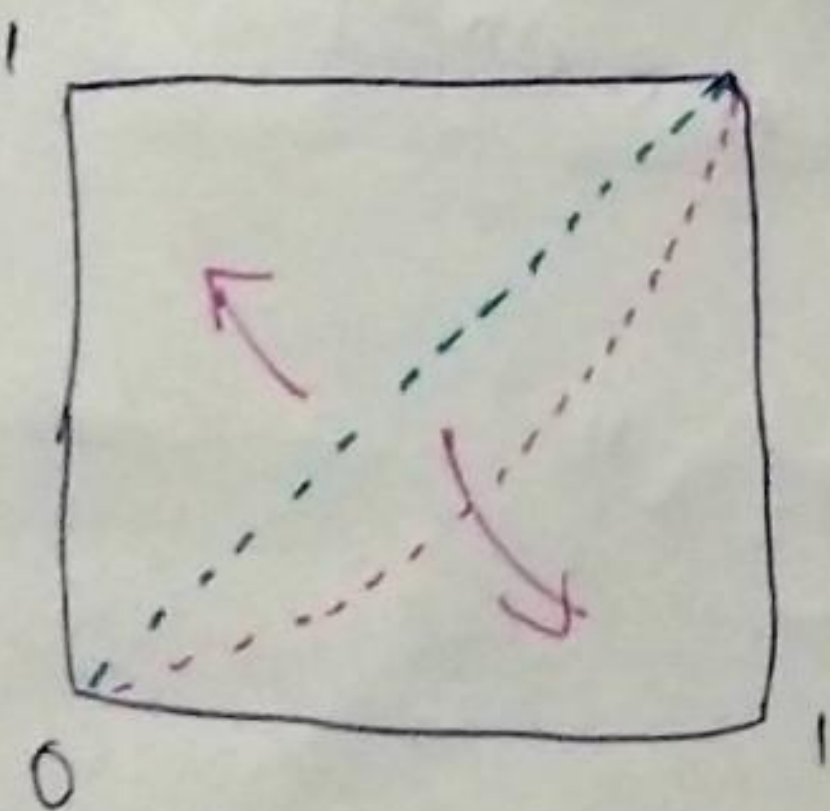
random model

$\hookrightarrow x_q \rightarrow 1 \text{ or } 0$

$$\left\{ AUC(\text{random-model}) \right\} = 0.5$$

④

Model  $M \rightarrow$



$$\text{Let } AUC(M) = 0.2$$

$\rightarrow$  means you did something wrong in modelling

$\hookrightarrow$  worse than random model.

AUC  $\rightarrow$

0.5 to 1  $\rightarrow \simeq$

0.5  $\rightarrow$  random model

0.0 to 0.5  $\rightarrow$  worst

$\hookrightarrow$  swap class labels,  $AUC = 1 - 0.2$

$\hat{y}_i = 0 \rightarrow 1$   
 $\hat{y}_i = 1 \rightarrow 0$  } swap -pping



21.5

→ (0 to 1)

log-loss → uses prob-scores.  
as small as possible

Binary classification →

Test set of  $n$ -points

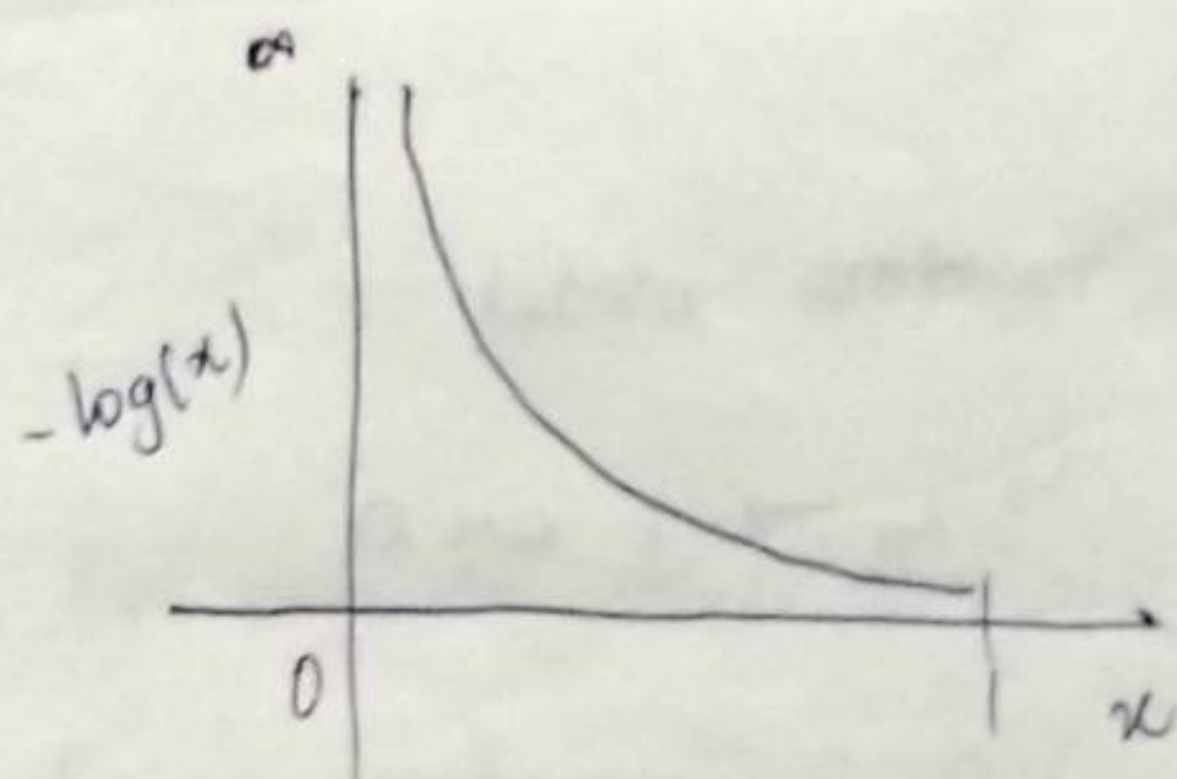
$$-\log(0.6) = 0.51$$

$$-\log(0.6) = 0.51$$

$x$	$y$	$\hat{y} = p$	
$x_1$	1	0.9	→ $-\log(0.9) = 0.1054$
$x_2$	1	0.6	
$x_3$	0	0.1	→ $-\log(0.9) = 0.1054$
$x_4$	0	0.4	

$$\text{log-loss} = -\frac{1}{n} \sum_{i=1}^n \left\{ \left( \log(p_i) * y_i \right) + \left( (1-y_i) * \log(1-p_i) \right) \right\}$$

log-loss → average of negative log prob of correct class label.



$$0 \leq p \leq 1$$

smaller the log loss, better the model is.

Multiclass Log loss →

$$-\frac{1}{n} \sum_{i=1}^n \sum_{j=1}^K y_{ij} \log(p_{ij}) \rightarrow \text{probability that } x_i \in \text{class } j$$

↓

$$= 1 \text{ if } x_i \in \text{class } j$$

otherwise 0

Best case of log loss is 0, but other values are not easy to make sense.



21.6

$R^2$  (or) coefficient of determination  $\rightarrow$   
 for regression

Test  $\rightarrow$ 

$$y_i \in \mathbb{R}$$

 $i = 1 \text{ to } n$ 

$$x_i, y_i, \hat{y}_i$$

$$e_i \Rightarrow y_i - \hat{y}_i$$

error

actual

actual

output

model

predicted

output

$$SS_{\text{total}} = \sum_{i=1}^n (y_i - \bar{y})^2$$

Sum of squares

actual

mean of actual class labels

$$\text{here } \bar{y} = \frac{1}{n} \sum_{i=1}^n y_i$$

$\bar{y} \rightarrow$  avg. value of  $y_i$ 's  
in test data.

for regression,

the simplest model you can build is  
called the average model  $\rightarrow$  simple mean model

$$x_q \rightarrow \text{mean}(y_i) = \bar{y} = y_q$$

$SS_{\text{total}} \rightarrow$  sum of squared errors using simple-mean model.

residuals

$$SS_{\text{res}} = \sum_{i=1}^n (y_i - \hat{y}_i)^2 = \sum_{i=1}^n e_i^2$$

$$\text{residual} = e_i = y_i - \hat{y}_i$$

actual

predicted



$$R^2 = \left( 1 - \frac{SS_{res}}{SS_{tot}} \right)$$

$$\text{Case 1: } \rightarrow SS_{res} = 0 \leftarrow (e_i = 0) \rightarrow R^2 = 1$$

phenomenal model

Best value

$$\text{Case 2: } \rightarrow \text{if } SS_{res} < SS_{tot} ; R^2 = 0 \text{ to } 1$$

$$\text{Case 3: } \rightarrow \text{if } SS_{res} = SS_{tot} ; R^2 = 1 - 1 = 0 \rightarrow \text{model is same as simple mean model.}$$

$$\text{Case 4: } \rightarrow \text{if } SS_{res} > SS_{tot} ; R^2 = 1 - (\text{greater } > 1) = -ve$$

model is worse than simple mean model.

21.7

Median absolute deviation of errors:

$$SS^2_{res} = \sum_{i=1}^n e_i^2 \quad \text{if one } e_i \text{ is very large}$$

$R^2$  is not very robust to outliers.

$$x_i \rightarrow y_i, \hat{y}_i, e_i$$

$|e_i|s \rightarrow 0 \rightarrow \text{great}$   
 $|e_i|s \rightarrow \text{large} \rightarrow \text{not so good}$

if  $e_i \rightarrow$  random variable,

mean  $\leftarrow \text{median}(e_i) = \text{central value of errors} \rightarrow \text{small}$

acting like these

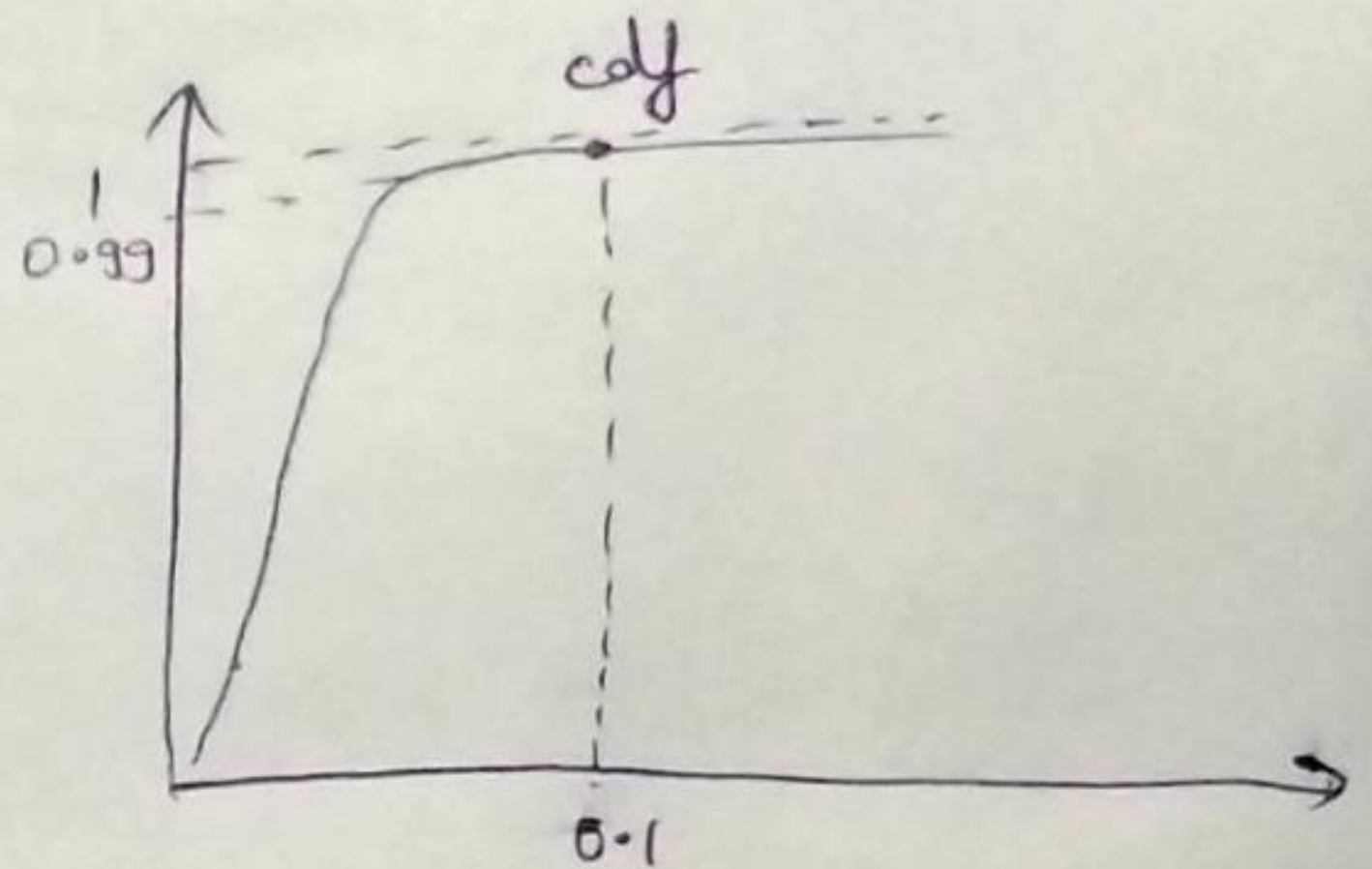
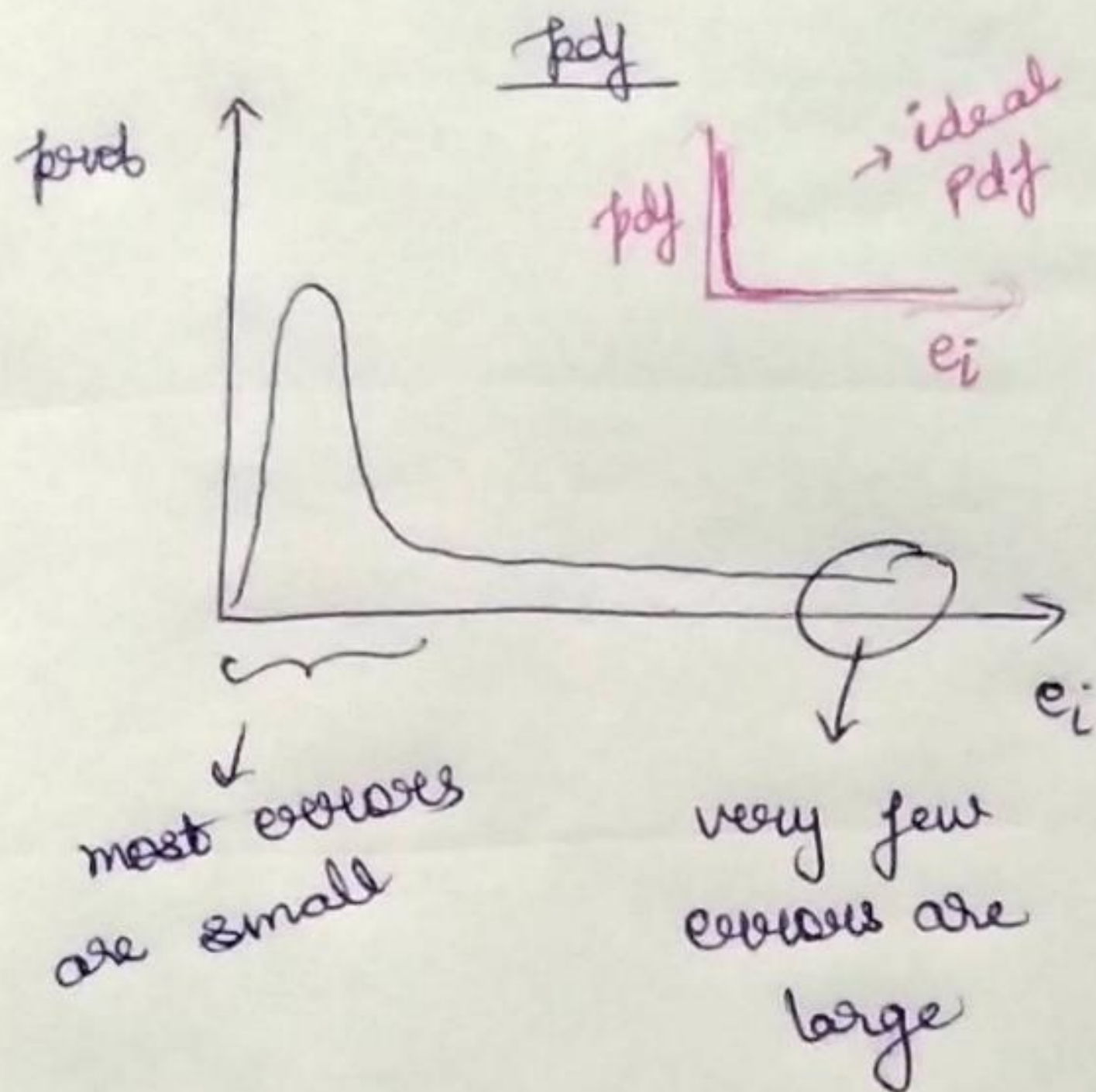
std. dev.  $\leftarrow \text{MAD}(e_i) = \text{median} \left( \underbrace{|e_i - \text{median}(e_i)|}_{\text{deviation}} \right) \rightarrow \text{small}$

abs



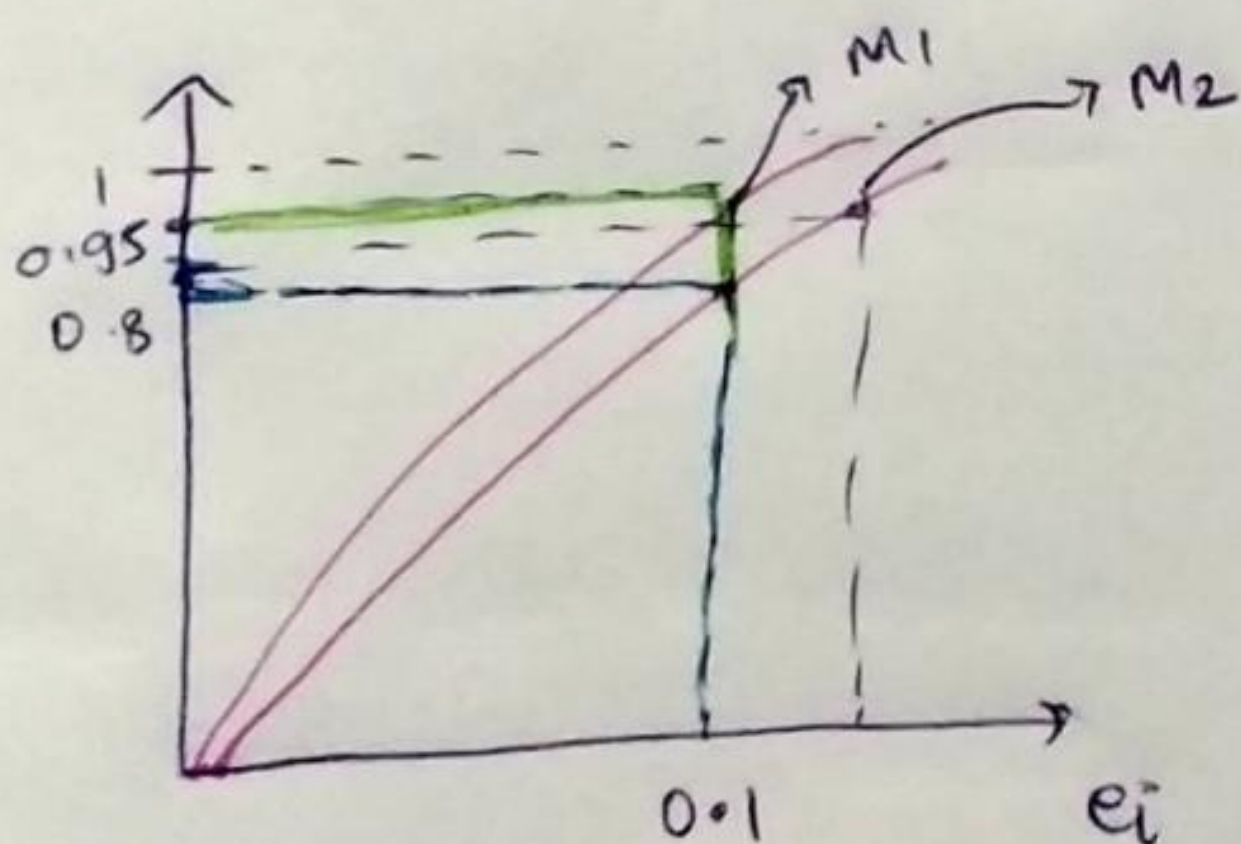
mean or median of  $e_i$ s  $\rightarrow$  used to understand if the errors are small or large.  
 std-dev or MAD  
 robust to outliers

## 21.8 Distribution of errors $\Rightarrow$



99% of errors are  $< 0.1$   
 1% errors are  $\geq 0.1$

Models  $M_1$  &  $M_2$



$M_2$  cdf is below  $M_1$

$M_1 \Rightarrow$  95% errors are below 0.1

$M_2 \Rightarrow$  80% errors are below 0.1

So,  $M_1$  is a better model for regression than  $M_2$