

CAUSE
EDA.

Line of Code

steps before model training

```
from sklearn.preprocessing import MinMaxScaler
```

```
scaler = MinMaxScaler()
```

```
df = pd.DataFrame(scaler.fit_transform(df), columns=df.columns)
```

All Numerical Columns.

- ① cat → Num
- ② Feature Scaling



Data Notation

sample - m

dimension / # features - d

Predictions - (m, d) matrix X

i th sample - $X^{[i]}$ - one row

j th feature - X_j - one column

True output / target - y^i

Predicted output / target \hat{y}^i

Goal of generalization in ML

m samples - $\{x^i, y^i\}^m$

Historical Data \rightarrow Train an ML Model

$$x^i \xrightarrow[\text{Model}]{\text{ML}} \hat{y}^i \leftarrow \text{Predicted value}$$

Ideally, $y^i \approx \hat{y}^i \rightarrow$ Predicted output
 \hookrightarrow True output

\Rightarrow For all m samples $y^i \approx \hat{y}^i$ \rightarrow GOOD TRAINING

DO WE WANT TO PERFORM PREDICTION ON TRAINING

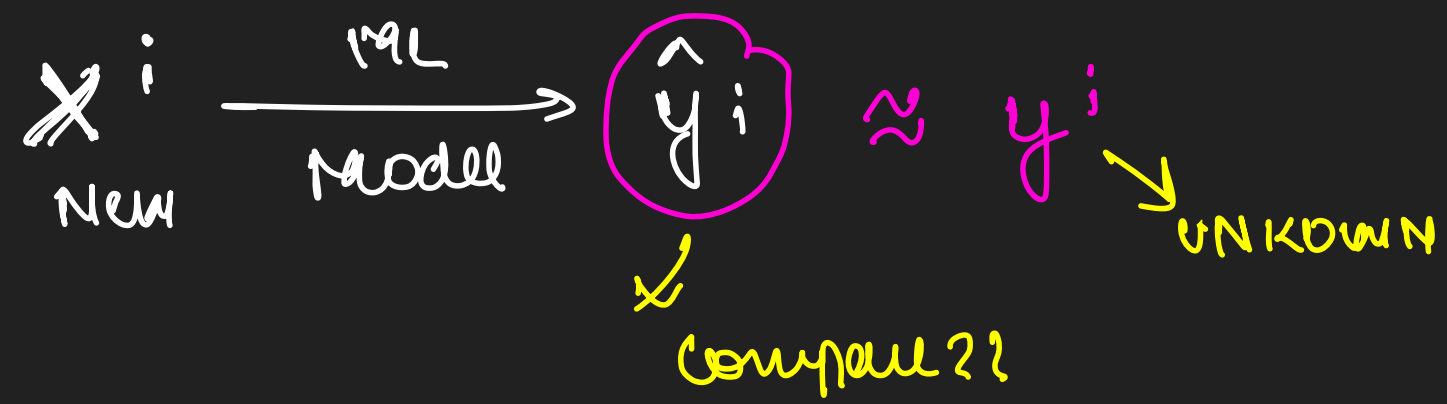
DATA??

for training data y is already present?
Why to even do prediction.

Goal of ML: GENERALISATION.

True
output

Your model should perform well
both on training and NEW data.

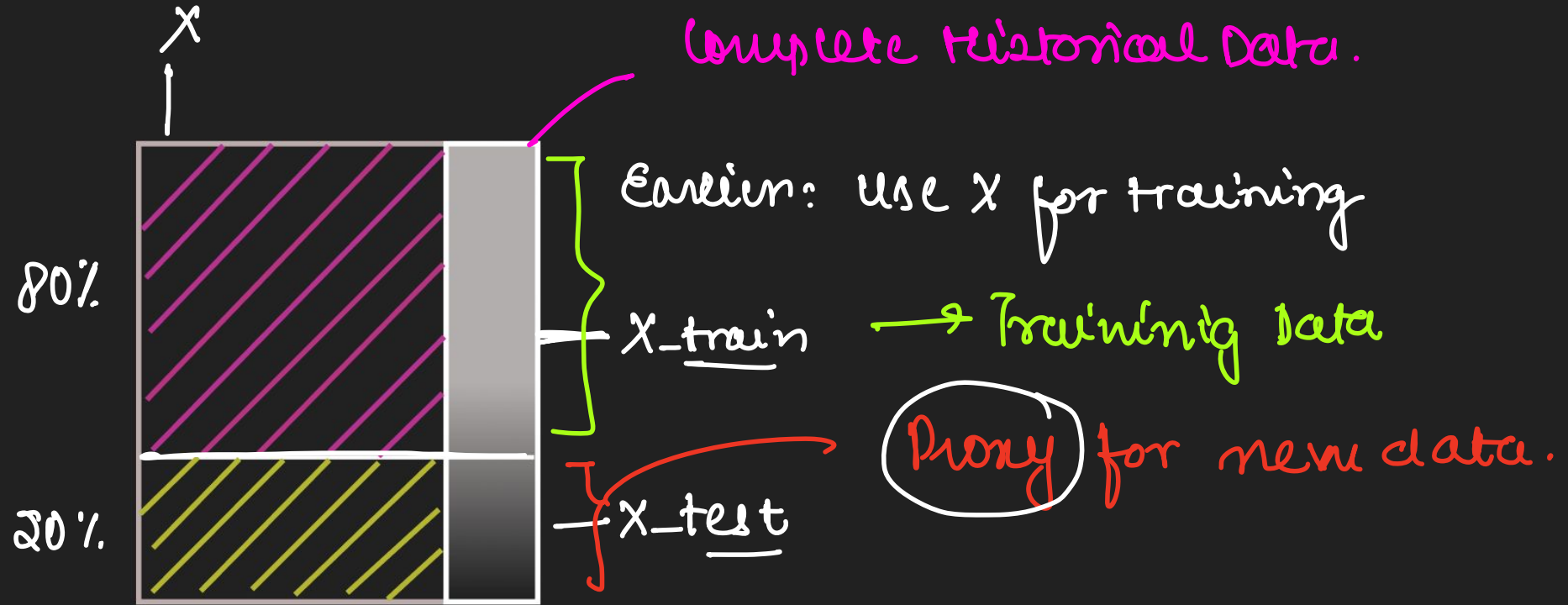


Check if model is GENERALISING??

Problem: No labels for new data.

Solution:

Because we don't know the ground truth (y) for New Sample



Now: Use only X_{train} for training

Phases for Model Development

① Training - use train split for model training
 $y_{\text{train}} \approx \hat{y}_{\text{train}}$ - Model is learning

② Testing/Evaluation - test split

$y_{\text{test}} \approx \hat{y}_{\text{test}}$ - Model is generalising

```
from sklearn.model_selection import train_test_split  
X_train, X_test, y_train, y_test = train_test_split(X, y,  
                                                    test_size=0.3,  
                                                    random_state=100)
```

Random split

80% 20%

0.2
↑
test_size=0.3,
random_state=100
seed.