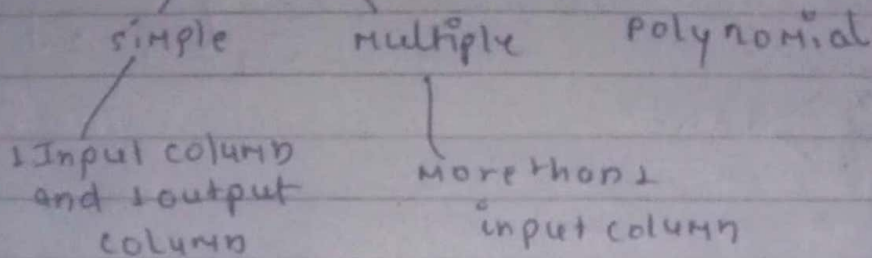


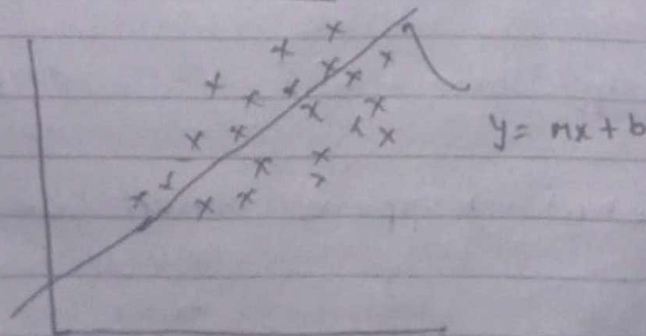
Linear Regression (simple + Multiple)



LR is a supervised learning algo. used to model the relationship between a dependent variable (target) and one or more independent variable (predictors). It assumes a linear relationship b/w the variables and aims to find a best fit line that minimize the error b/w predicted and actual values.

$$y = mx + b$$

How to find m and b



Two method to find m and b

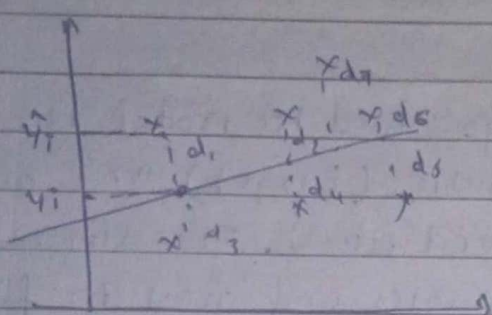
- (1) closed form solution \rightarrow OLS
- (2) Non-closed form \rightarrow Gradient descent

Through OW

$$b = \bar{y} - m\bar{x}$$

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

\bar{x} \bar{y} \rightarrow Mean



$$\text{Error} = d_1 + d_2 + d_3 + \dots + d_n$$

$$E = d_1^2 + d_2^2 + \dots + d_n^2$$

$$E = \sum_{i=1}^n d_i^2$$

← Error function

• Represent with J

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

$$E = \sum_{i=1}^n (y_i - mx_i - b)^2$$

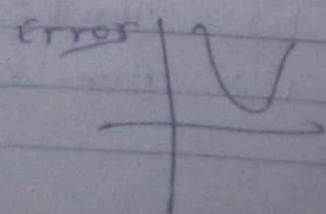
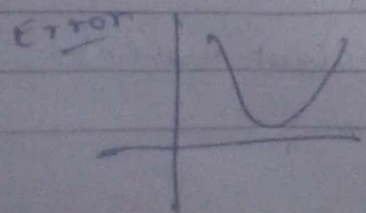
let's see what happens when we change m and b

suppose
 $b = 0$

$m = 1$

$$E(m) = \sum_{i=1}^n (y_i - mx_i)^2$$

$$E(b) = \sum_{i=1}^n (y_i - x_i - b)^2$$



$$\frac{\partial E}{\partial b} = \frac{\partial}{\partial b} \sum_{i=1}^n (y_i - mx_i - b)^2 = 0$$

$$\Rightarrow \sum \frac{\partial}{\partial b} (y_i - mx_i - b)^2 = 0$$

$$\Rightarrow \sum -2 (y_i - mx_i - b) = 0$$

$$\Rightarrow \sum y_i - \sum mx_i - \sum b = 0$$

$$\frac{\sum y_i}{n} - \frac{\sum mx_i}{n} - \frac{\sum b}{n} = 0$$

$$\bar{y} - m\bar{x} - \frac{1}{n}\sum b = 0$$

$$\bar{y} - m\bar{x} - b = 0$$

$$\boxed{b = \bar{y} - m\bar{x}}$$

$$E = \sum (y_i - mx_i - \bar{y} + m\bar{x})^2$$

$$\frac{\partial E}{\partial m} = \sum \frac{\partial}{\partial m} (y_i - mx_i - \bar{y} + m\bar{x})^2 = 0$$

$$\Rightarrow \sum 2 (y_i - mx_i - \bar{y} + m\bar{x}) (-x_i + \bar{x}) = 0$$

$$\Rightarrow \sum -2 (y_i - mx_i - \bar{y} + m\bar{x}) (x_i - \bar{x}) = 0$$

$$\Rightarrow \sum (y_i - mx_i - \bar{y} + m\bar{x}) (x_i - \bar{x}) = 0$$

$$\Rightarrow \sum [(y_i - \bar{y}) - m(x_i - \bar{x})] (x_i - \bar{x}) = 0$$

$$\Rightarrow \sum [(y_i - \bar{y})(x_i - \bar{x}) - m(x_i - \bar{x})^2] = 0$$

$$\Rightarrow \sum (y_i - \bar{y})(x_i - \bar{x}) = m \sum (x_i - \bar{x})^2$$

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

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

Multiple Linear Regression

More than one column in input

$$x_1 | x_2 | x_3 | y$$

$$\begin{aligned}
 & \left. \begin{array}{l} 3D \rightarrow \text{plane} \\ 4D \rightarrow \text{hyperplane} \\ nD \rightarrow \text{---} \end{array} \right\} \begin{aligned} & y = mx_1 + nx_2 + b \\ & y = \beta_0 + \beta_1 x_1 + \beta_2 x_2 \\ & y = \beta_0 + \beta_1 x_1 + \beta_2 x_2 + \beta_3 x_3 \\ & y = \beta_0 + \sum_{i=1}^n \beta_i x_i \end{aligned}
 \end{aligned}$$

Gradient Descent

Gradient descent is a first order iterative optimization algorithm for finding a local minimum of a differentiable fxn.

Types

Batch GD

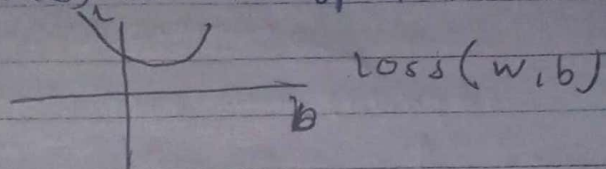
SGD

MBGD

SGD Regression

Loss fcn $J(\theta)$

opposite direction of the gradient of the objective function $\nabla J(\theta)$ with respect to the parameters



Learning rate η

$$\eta \frac{\partial L}{\partial w}$$

determine the size of the steps we take to reach a (local) minimum.

Back propagation Algo

epochs = 5

for i in range(epochs):

for j in range(x.shape[0]):

→ select row (random)

→ predict (using forward prop)

→ calculate loss (using loss function → mse)

→ update weights and bias using GD

$$w_n = w_0 - \eta \frac{\partial L}{\partial w}$$

→ calculate avg loss for the epoch

★ $\left\{ \frac{\partial L}{\partial w} \right\}$ derivative ^{How much data we use to compute} 3 types

accuracy → time — tradeoff

Batch 40 (vanilla 40)

↳ we take the entire dataset and then we make update

Exo

current weigh
50 points → predict →
↑
dot product
 $\hat{y} = \text{np.dot}(x, w) + b$
↳ so predict

$y = 50$ actual outcome

- using this y and \hat{y} we calculate loss $\sum_{i=1}^{50}$
- then basis of that loss we update single time weigh and bias

Exo

epoch = 10
for i in range(10):
 $\hat{y} = \text{np.dot}(x, w) + b$

total
10
times
repeat

so values
 $y = 50$
 $\hat{y}, y \rightarrow$ loss
w, b update $w_n = w_0 - \eta \frac{\partial L}{\partial w}$
→ loss

weight update frequency high

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Stochastic GD

designed to handle large datasets efficiently by introducing randomness in the gradient computation

epoch $\rightarrow 10$ (50 rows)

for i in range(10): ~~shuffle~~
~~for i in range(x.shape[0])~~
~~\hookrightarrow shuffle~~
~~\hookrightarrow point~~

~~\hookrightarrow select random point~~

~~\hookrightarrow y-hat \rightarrow forward~~

~~\hookrightarrow loss~~

~~\hookrightarrow w, b update $\rightarrow w_n = w_0 - \eta \frac{\partial L}{\partial w}$~~

avg loss print \rightarrow for the epoch

★ 50 rows then 50 times ^{weight} update. (500)
 no. of epoch times update (500)

difference between both

	BGD	SGD
update if 10 epochs	10 time	10 \times no. of rows
which is the faster to converge (given same # epochs)		SGD

Mini Batch GD

combines the advantages of both SGD and BGD by dividing training dataset into smaller called mini-batches.

```
for i in epochs ->
    for j in num of batch
        batch
            ↳ y-pred (vector)
            ↳ loss
            ↳ update
```

update of epoch fast slow
speed $\text{bgd} > \text{mbgd} > \text{sgd}$

convergence slow fast
 $\text{bgd} < \text{mbgd} < \text{sgd}$

→ why batch size is provided in multiple of (2)?

actually RAM ka effective use krne ke liye
diya jata hai design to handle binary value

→ what if batch-size doesn't divide # rows properly

e.g. # of rows $n = 400$

batch-size = 150

of batch $= \frac{400}{150} = 2.66$

150, 150, left 100

Batch 1, b2, b3 ← left include in b3

Regression Metrics

(1) Mean Absolute error

used to evaluate the accuracy of prediction
measures the avg absolute diff. b/w the actual values
and the predicted values.

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

total no. of observation. actual predicted

Advantage ① same unit

② Robust to outliers

Disad. modulus fxn graph is not diff. at zero.

(2) Mean squared error

avg squared diff. b/w predicted values and the
actual values in a dataset

A lower MSE indicates better model accuracy,
higher suggests poor performance.

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

Advantage ① use as a loss fxn becoz differentiable

Dis

① Not Robust to outliers.

RMSE

Root of MSE

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

Benefit

RMSE output comes in same unit

Disad - Not robust to outliers.

R² score

Goodness of fit

value of R^2 ranges b/w 0 and 1,

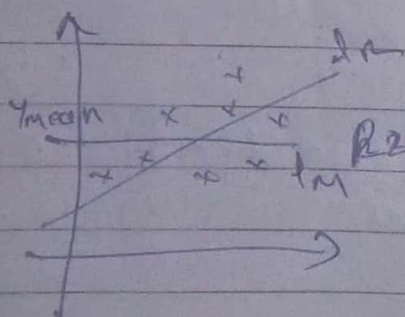
→ $R^2 = 1$ indicates a perfect fit, meaning the model explains all the variability in the dependent variable

→ $R^2 = 0$ means the model does not explain any variability and fails to capture the relationship b/w the variables.

$$R^2 = 1 - (SS_{res} / SS_{tot})$$

residual sum of squares

total sum of squares



$$R^2 = SSR / SS_{tot}$$

sum of squares of Regression

$$R^2 = 1 - \frac{\left[\sum_{i=1}^n (y_i - \hat{y}_i)^2 \right]_{res.}}{\left[\sum_{i=1}^n (y_i - \bar{y})^2 \right]_{M.}}$$

Adjusted R^2 score

To address the limitations of R^2 , Adjusted R^2 is used.

if we add more input R^2 score \uparrow

$$R^2_{adj} = 1 - \left[\frac{(1 - R^2)(n-1)}{(n-1-k)} \right]$$

n - no. of rows
 k - independent
 values.