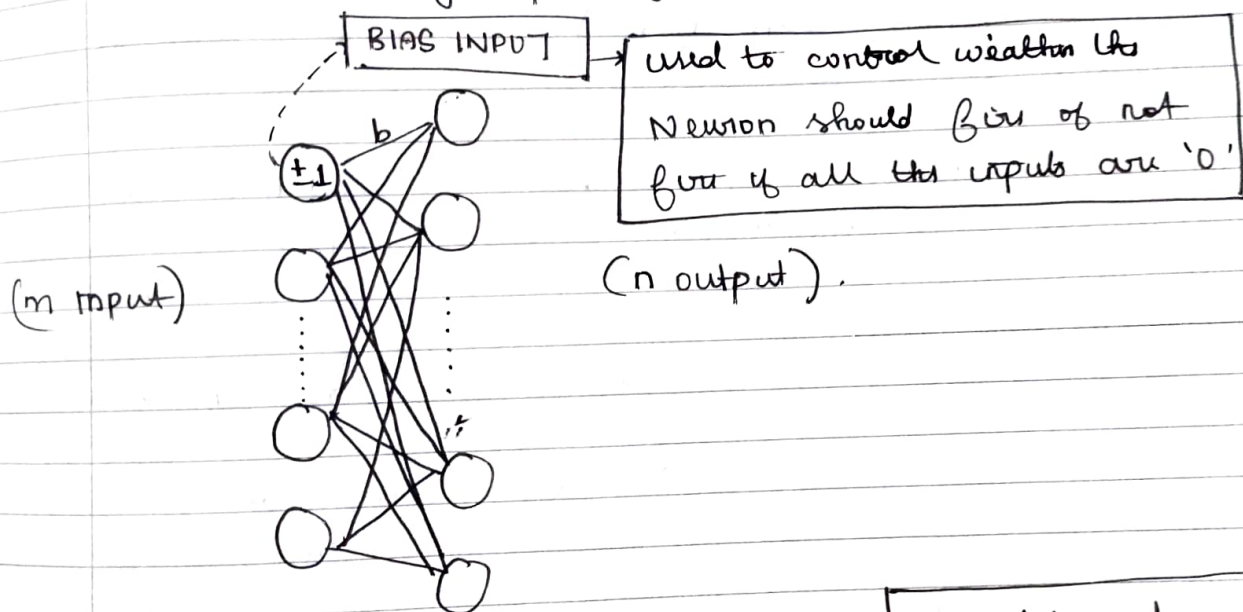


* PERCEPTRON: (SINGLE LAYER PERCEPTRON)

→ A linear model used for binary classification

→ Step Function for activation i.e $f(x) = \begin{cases} 0, & x \leq \alpha \\ 1, & x > \alpha \end{cases}$
 $\alpha \Rightarrow$ THRESHOLD VALUE.

→ Contains only input layer and output layer.



At each output neuron k we have,

$x_0 = +1$ and $w_{0k} = b$ which is weight of bias

$$\hat{y}_k = f\left(\sum_{i=0}^m x_i w_{ik}\right) \Rightarrow \text{Output of } k\text{th neuron.}$$

⇒ The weights of the connections are updated through Perceptron Learning algorithm.

At the k^{th} output neuron the error E_k is Given by

$$E_k = \hat{y}_k - y_k \quad \text{where } \hat{y}_k = \text{predicted value}$$

$y_k = \text{actual value.}$

if $E_k > 0 \Rightarrow$ weights should be reduced

if $E_k < 0 \Rightarrow$ weights should be increased

$$\Rightarrow \Delta w_{ik} = -\eta_i (\hat{y}_k - y_k)$$

Therefore the weight updation is given by.

$$w_{ik} = w_{ik} + \eta \Delta w_{ik} = w_{ik} + \eta \cdot x_i (\hat{y}_k - y_k)$$

The parameter η is called as learning Rate. It determines how fast should the perceptron learn by controlling the amount of change in weight.

if η is too high \Rightarrow network becomes unstable

if η is too low \Rightarrow learning will take large amount of time

$$0.1 < \eta < 0.4 \Rightarrow \text{optimal Range}$$

~~PERCEPTRON~~

* MULTILAYER PERCEPTRON (MULTILAYER FEED FORWARD NETWORK):

\rightarrow Neurons are arranged into groups called as layer

\rightarrow Contains 3 main layers:

- ① Input layer
- ② Hidden layers
- ③ Output layer

- ⇒ In a given layer all the neurons will have same activation function.
- ⇒ For input layer, input will be the actual data used. For other layers, input will be activation of previous layer.
- ⇒ Number of neurons in the input layer will be equal to number of features in the input.
- ⇒ we have two main steps:

① Forward Propagation (Predicting Output)

② Backward Propagation (Reducing Error by tuning weight and bias value)
(Backpropagation of error)

⇒ Each neuron can be given a separate bias value (or) each layer can have a common bias value, this is based on the implementation.

⇒ BACKPROPAGATION: it is the method of calculating and propagating error from output layer to all hidden layers.

→ To reduce the error we use "optimisation algorithm"



error is calculated using "loss function" ⇒ The goal of optimiser is to ~~to~~ minimize the loss function.

linear

Non-linear

DATE

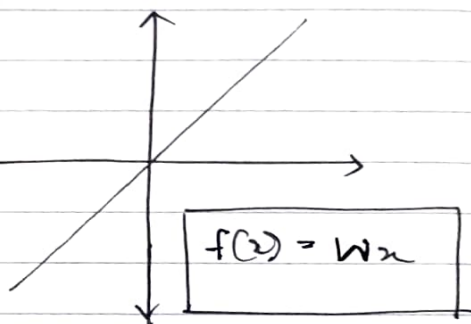
⇒ ACTIVATION FUNCTION: used to introduce non-linearity in the decision boundary ('Because not all problems are linearly separable')

without ~~any~~ activation function the neural network will just perform linear weighted sum (similar to linear Regression).

① LINEAR: $f(x) = wx$.

⇒ identity function

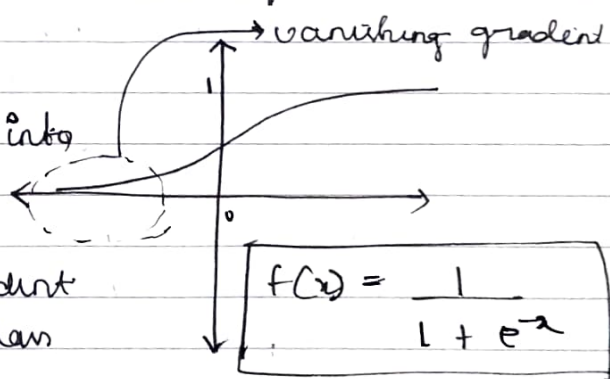
⇒ used in input layer?



② SIGMOID:

⇒ Reduce the input into the Range 0 to 1.

⇒ Outputs independent probability for each class

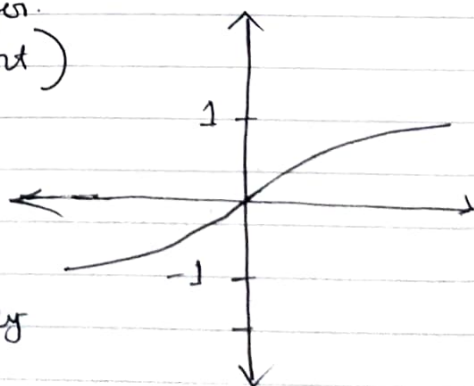


~~used~~: used in binary classification output layer.

③ TANH: (Hyperbolic Tangent)

⇒ Reduce input to the Range -1 to 1

⇒ it can deal more easily with negative values.



$$f(x) = \frac{e^x - e^{-x}}{e^x + e^{-x}}$$

Both Sigmoid and tanh can cause vanishing Gradient problem (For extreme high / low value) \Rightarrow So we use ReLU. But it can cause exploding Gradient \Rightarrow need to scale up.

④ SOFTMAX:

- \rightarrow provides the probability distribution of each class.
- \rightarrow used in output layer for multi-class classification

~~$f(x) = \frac{e^x}{\sum_l e^{x_l}}$~~ \Rightarrow if output layer contains M neurons and x be the output value/activation

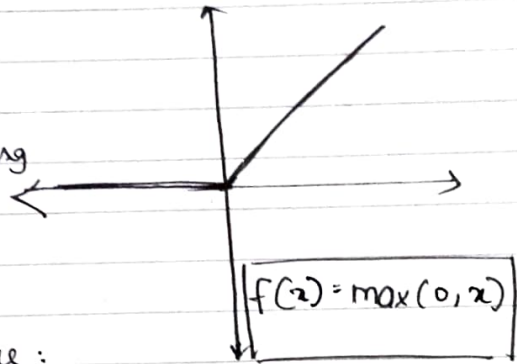
$$\Rightarrow f(x_i) = \frac{e^{x_i}}{\sum_l e^{x_l}}$$

(Sum of all probability will be 1)

(class with maximum probability will be the Result).

⑤ RECTIFIED LINEAR UNIT (ReLU):

- \Rightarrow overcomes vanishing / exploding gradient problem.
- \Rightarrow ReLU is highly used in hidden layer
- \Rightarrow Some variants of ReLU are:



(i) Leaky ReLU = $f(x) = \max(0.01x, x)$.

(ii) Parametric ReLU = $f(x) = \max(\alpha x, x)$.

(iii) Exponential ReLU = $f(x) = \begin{cases} x, & \text{if } x \geq 0 \\ \alpha(e^x - 1), & \text{if } x < 0 \end{cases}$
 \Downarrow
 $\max(\alpha(e^x - 1), x)$.

LOSS FUNCTION: It quantifies how close the predicted Value is to the Real Value.

For Given weights W and Bias B we have $L(W, B)$ as loss function

$\hat{Y}_i = h_{W, B}(X) \Rightarrow$ predicted value for i th Sample
 $Y_i \Rightarrow$ Real value for i th Sample

LOSS FUNCTION FOR REGRESSION

① Mean Squared Error

$$L(W, B) = \frac{1}{N} \sum_{i=1}^N (\hat{Y}_i - Y_i)^2$$

if more than one feature is der in output

$$\Rightarrow \frac{1}{N} \sum_{i=1}^N \left(\sqrt{\sum_{j=1}^M (\hat{Y}_{ij} - Y_{ij})^2} \right)^2$$

$$\Rightarrow \frac{1}{N} \sum_{i=1}^N \sum_{j=1}^M (\hat{Y}_{ij} - Y_{ij})^2 //$$

② Mean Absolute Error

$$L(W, B) = \frac{1}{N} \sum_{i=1}^N |\hat{Y}_i - Y_i|$$

$$\Rightarrow \frac{1}{N} \sum_{i=1}^N \sum_{j=1}^M |\hat{Y}_{ij} - Y_{ij}|$$

③ Root Mean Squared

$$L(W, B) = \sqrt{MSE}$$

LOSS FUNCTION FOR CLASSIFICATION

Cross entropy loss.
(OR)

Negative Log Likelihood

Binary
Categorical.

$$BCE = -\frac{1}{N} \sum_{i=1}^N Y_i \log \hat{Y}_i + (1 - Y_i) \log (1 - \hat{Y}_i)$$

$$CCE = -\frac{1}{N} \sum_{i=1}^N \sum_{j=1}^M Y_{ij} \log (\hat{Y}_{ij})$$

Hing loss:

\hookrightarrow designed for SVM

Result should be $-1/1$.

$$L(W, B) = \frac{1}{N} \sum_{i=1}^N \max(0, 1 - Y_i \hat{Y}_i)$$

~~$$\frac{1}{N} \sum_{i=1}^N \max(0, 1 - Y_i \hat{Y}_i)$$~~

$$\sum_{i=1}^N \sum_{j=1}^M \max(0, y_{ij} \times \hat{y}_{ij}) \Rightarrow \text{only used for binary classification}$$

MSE and MAE are prone to outliers due to which we can use **HUBER LOSS** \rightarrow For Regression

$$L(W, B) = \begin{cases} \frac{1}{2} (y_i - \hat{y}_i)^2, & |y_i - \hat{y}_i| \leq \delta \\ \delta |y_i - \hat{y}_i| - \frac{1}{2} \delta^2, & \text{else.} \end{cases}$$

Here δ is a hyperparameter which depicts the tolerance to outliers.

* GRADIENT DESCENT:

Huber loss is like a combination of MSE and MAE

\rightarrow An optimisation algorithm for ~~minimising~~ minimising the cost / loss function

- (i) Batch Gradient Descent
- (ii) Stochastic Gradient Descent
- (iii) Mini Batch Gradient Descent

- \Rightarrow In (i) The complete training set is used for every iteration / epoch.
- \Rightarrow In (ii) Only single sample is used for one iteration (i.e. Number of iterations / epochs = Number of samples in training set \rightarrow minimum).
- \Rightarrow In (iii) A Batch of samples used for each iteration from the training dataset.
- \Rightarrow Some other optimisers are: RMSProp, ADADelta, ADAGRad, Adam.

* HYPERPARAMETERS:

⇒ Configuration options that can be varied or altered to improve the performance of model.

- ① No. of Hidden Layers
- ② No. of Neurons in each layer
- ③ Activation Functions
- ④ Loss Functions
- ⑤ Optimiser Functions
- ⑥ No. of epochs/iterations
- ⑦ Type of Regularisation
- ⑧ Weight initialisation
- ⑨ Learning Rate & Momentum.

* BIAS AND VARIANCE: (Trade off)

⇒ when a model is trained it captures the patterns in the training data and ~~tries~~ to generalise it.

⇒ BIAS is the measure of our model's ability to learn patterns from data set.

⇒ when model has HIGH BIAS ⇒ did not train well on training set

when model has LOW BIAS ⇒ it has identified some patterns from training set properly

⇒ VARIANCE is the measure of model's fluctuation in performance when there is a change in data set.

⇒ UNDERFITTING: when a model does not perform well in training data.
(HIGH BIAS).

⇒ OVERFITTING: when a model performs well in train data but it lacks performance in testing data.
(LOW BIAS, HIGH VARIANCE)

⇒ A Good model should have LOW BIAS and LOW VARIANCE.

MACHINE LEARNING: A Algorithm is said to be learning from experience E with respect to some task T and performance measure P if the performer P increases with increase in experience E .

TYPES OF MACHINE LEARNING:

① SUPERVISED (→ CLASSIFICATION
→ REGRESSION)

② UNSUPERVISED (→ CLUSTERING).

③ SEMI-SUPERVISED

④ REINFORCEMENT

LINEAR SEPERABILITY:

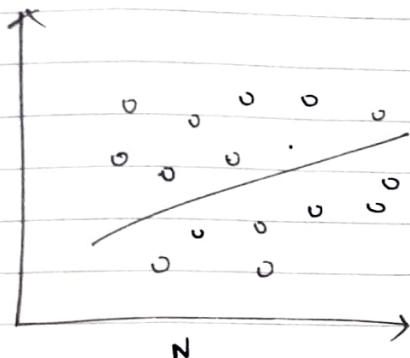
A data is said to be linearly Seperable if only a single line / Plane / Hyperplane can partition the data into 2 classes.

Eg: Perceptron can only work on a data set that is linearly separable.

→ One of the main reason for MLP improvement

LINEAR REGRESSION:

consider the following equation $y = \theta_0 x + \theta_1$



Linear regression
tries to fit a line of
the form on the
Given dataset.

$$E = \frac{1}{N} \sum_{i=1}^N (\hat{y}_i - y_i)^2 \Rightarrow \text{where } \hat{y}_i \text{ is the predicted value and } y \text{ is the actual value}$$

The error E is called as MEAN SQUARED ERROR.
we have to Reduce this error this is called
LEAST SQUARE REGRESSION.

$$E = \frac{1}{N} \sum_{i=1}^N (\hat{y}_i - y_i)^2 = \frac{1}{N} \sum_{i=1}^N (\theta_0 x_i + \theta_1 - y_i)^2$$

$$\frac{\partial E}{\partial \theta_0} = \frac{\partial}{\partial \theta_0} \left(\frac{1}{N} \sum_{i=1}^N (\theta_0 x_i + \theta_1 - y_i)^2 \right) = 0$$

$$= \theta_0 \bar{x}^2 + \theta_1 \bar{x} - \bar{x} \bar{y} = 0 \quad \text{--- (1)}$$

$$\frac{\partial E}{\partial \theta_1} = \frac{\partial}{\partial \theta_1} \left(\frac{1}{N} \sum_{i=1}^N (\theta_0 x_i + \theta_1 - y_i)^2 \right) = 0$$

$$= \theta_0 \bar{X} + \theta_1 - \bar{Y} = 0$$

$$\Rightarrow \theta_1 = \bar{Y} - \theta_0 \bar{X} \quad \text{--- (2)}$$

From (1) and (2) we have.

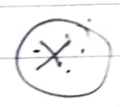
$$\theta_0 \bar{X}^2 + (\bar{Y} - \theta_0 \bar{X}) \cdot \bar{X} - \bar{X} \cdot \bar{Y} = 0$$

$$\theta_0 \bar{X}^2 + \bar{X} \cdot \bar{Y} - \theta_0 \bar{X}^2 - \bar{X} \cdot \bar{Y} = 0$$

$$\theta_0 = \frac{\bar{X} \cdot \bar{Y} - \bar{X} \cdot \bar{Y}}{\bar{X}^2 - \bar{X}^2}$$

\Rightarrow

$$\theta_0 = \frac{\bar{X} \cdot \bar{Y} - \bar{X} \cdot \bar{Y}}{\bar{X}^2 - \bar{X}^2}$$



$$\theta_1 = \bar{Y} - \theta_0 \bar{X}$$

\Rightarrow Why Huber Loss for Regression?

MSE is Highly Sensitive to outliers because we square the over all error values. MAE just take the absolute value of error it give same weight to all errors due to which outliers can be completely neglected. Huber Loss provide Balance (S).

$$\text{Huber Loss} = 0$$

* BAYESIAN STATISTICS:

↳ (methods that use Bayes Theorem)

⇒ BAYES THEOREM: Finds the probability of an event occurring, given the probability of another event that has already occurred

↳ (called as conditional probability)

$$P(A|B) = \frac{P(B|A) \cdot P(A)}{P(B)}$$

⇒ Here, $P(B)$ is called as evidence.

→ $P(A)$ is called as prior probability (i.e., probability occurrence of event A without considering event B)

→ $P(A|B)$ is called as posterior Probability (i.e., probability of occurrence of event A with considering event B)

⇒ we can $P(Y|X)$ using above method.

$Y \Rightarrow$ Target Class and $X =$ Feature Vector.

$$\Rightarrow P(Y_j | x_1, x_2, \dots, x_M) = \cancel{P(x_1) \cdot P(x_2) \cdot \dots \cdot P(x_M)}$$

$$= \frac{P(Y_j | x_1) \cdot P(Y_j | x_2) \cdot \dots \cdot P(Y_j | x_M) \cdot P(Y_j)}{P(x_1) \cdot P(x_2) \cdot \dots \cdot P(x_M)}$$

Since ~~prior~~ evidence is same.

for all target we can ignore this

⇒ Result will be class j having max $P(Y_j | x)$.

⇒ There are various variation of Naive Bayes algorithms like

(i) Gaussian / Normal NB

(ii) Multinomial NB

(iii) Bernoulli NB

;

⇒ Important assumption for Naive Bayes algorithm is all the features are independent.

$$\text{i.e } P(x_i | x_j) = P(x_i)$$

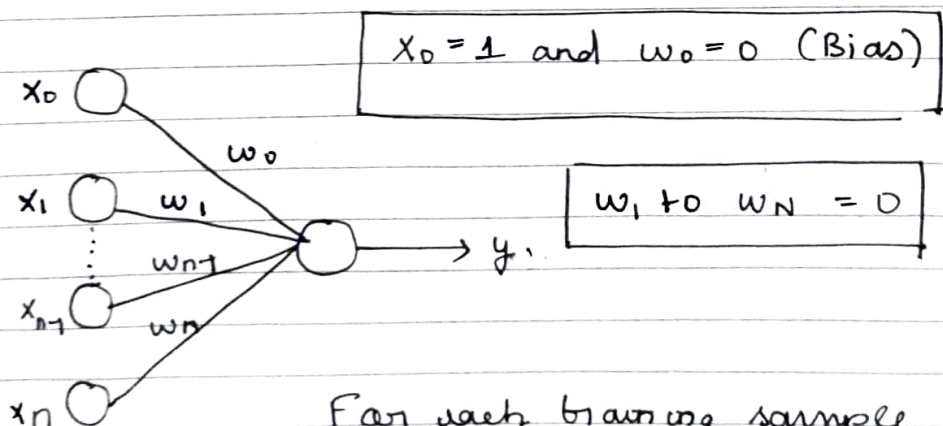
(X)

— x —

* HEBBIAN LEARNING RULE (HEBB NETWORK)

→ inputs can only be 1 and -1

→ initialise all weights and bias to 0.
in Starting



For each training sample
we do the following

$$\Rightarrow \Delta w_i = x_i y \text{ and } w_i = w_i + \Delta w_i$$

— x —

* FORMULAS

① PERCEPTRON WEIGHT UPDATION:

$$W_{ij} = W_{ij} + \eta (\hat{y}_j - y_j) (x_i)$$

② LOSS FUNCTIONS:

(i) REGRESSION:

⇒ MEAN SQUARED ERROR:

$$E = \frac{\sum_{i=1}^N \sum_{j=1}^M (\hat{y}_{ij} - y_{ij})^2}{N} \quad (\text{OR}) \quad E = \frac{\sum_{i=1}^N (\hat{y}_i - y_i)^2}{N}$$

⇒ MEAN ABSOLUTE ERROR:

$$E = \frac{\sum_{i=1}^N \sum_{j=1}^M |\hat{y}_{ij} - y_{ij}|}{N} \quad (\text{OR}) \quad E = \frac{\sum_{i=1}^N |\hat{y}_i - y_i|}{N}$$


⇒ ROOT MEAN SQUARED ERROR

$$E = \sqrt{\frac{\sum_{i=1}^N \sum_{j=1}^M (\hat{y}_{ij} - y_{ij})^2}{N}} \quad (\text{OR}) \quad E = \sqrt{\frac{\sum_{i=1}^N (\hat{y}_i - y_i)^2}{N}}$$

$$\Rightarrow \boxed{ERMSE = \sqrt{EMSE}}$$

⇒ HUBER LOSS (A Balancer between MSE and MAE).

$$E = \begin{cases} \frac{1}{2} (y_i - \hat{y}_i)^2, & \text{if } |y_i - \hat{y}_i| \leq \delta \\ \delta |y_i - \hat{y}_i| - \frac{1}{2} \delta^2, & \text{otherwise} \end{cases}$$

classmate  δ value controls the behaviour of huber loss (hyperparameter).

(ii) CLASSIFICATION :

⇒ HINGE LOSS :

⇒ CROSS ENTROPY LOSS :

$$E = - \sum_{i=1}^N \sum_{j=1}^M y_{ij} \log \hat{y}_{ij} \quad (\text{Multiple class} \Rightarrow \text{Categorical CE}).$$

$$E = - \sum_{i=1}^N y_i \log \hat{y}_i + (1 - y_i) \log (1 - \hat{y}_i).$$

(Binary Class \Rightarrow Binary CE).

③ SOME ACTIVATION FUNCTION :

(i) Sigmoid :

$$g(x) = \frac{1}{1 + e^{-x}}$$

(ii) TanH :

$$g(x) = \frac{e^x - e^{-x}}{e^x + e^{-x}}$$

(iii) ReLU :

→ Parametric : $g(x) = \max(ax, 1)$

$g(x) = \max(0, x)$ → Leaky : $g(x) = \max(0.01x, x)$

→ Exponential : $g(x) = \max(x(e^x - 1), x)$

(iv) Softmax:

$$g(x_i) = \frac{e^{x_i}}{\sum_{j=1}^M e^{x_j}}$$

(v) LINEAR:

$$g(x) = \alpha x.$$

(4) NAIVE BAYES:

For M target classes ~~from~~ we have to find max of

~~$$P(Y_j | x_1, x_2, x_3, \dots, x_k)$$~~

$$P(Y_j | x_1, x_2, x_3, \dots, x_k) = P(x_1 | Y_j) \cdot P(x_2 | Y_j) \cdots P(x_k | Y_j) \cdot P(Y_j)$$

$$\underbrace{P(x_1) \cdot P(x_2) \cdots P(x_k)}$$

Same for all. So can be ignored.

(5) MULTILAYER PERCEPTRON:

~~$$E_k = O_k (1 - O_k) (O_k - T_k)$$~~

$$E_k = O_k (1 - O_k) (O_k - T_k). \quad (\text{Output layer})$$

$$E_j = O_j (1 - O_j) \left(\sum_{k=1}^N W_{jk} E_k \right). \quad (\text{Hidden layer})$$

$$W_{ij} = W_{ij} + \eta O_i E_j$$