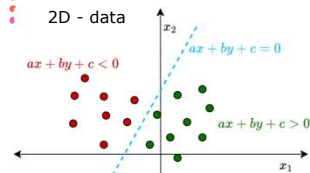


# Dissecting Neural Networks

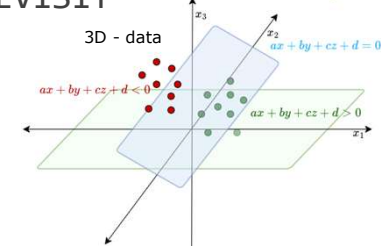
DR. SHAILESH S

## CLASSIFICATION : REVISIT



Predict(p, q)

```
if ap + bq + c < 0:
    return N
else:
    return P
```



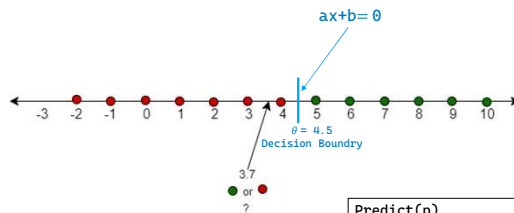
Predict(p, q, r)

```
if ap + bq + cr + d < 0:
    return N
else:
    return P
```

## CLASSIFICATION : REVISIT

Dataset

x	y
-2	N
6	P
7	P
3	N
2	N
-1	N
0	N
5	P
10	P
4	N
9	P
8	P



Binary Classification

Predict(p)

```
if 2p + -9 < 0:
    return N
else:
    return P
```

## CLASSIFICATION : REVISIT

**Generalizing to n-dimensional space**

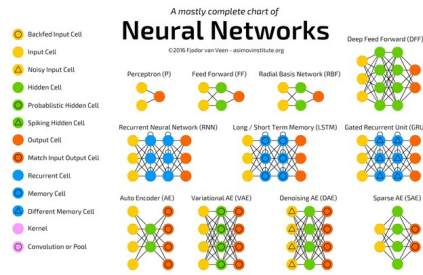
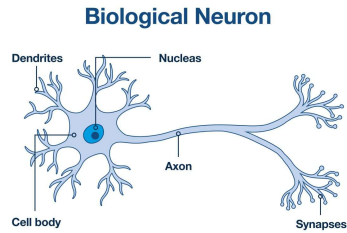
Data :  $X = \langle x_1, x_2, x_3, \dots, x_n \rangle$

Decision Boundary :  $w_1x_1 + w_2x_2 + w_3x_3 + \dots + w_nx_n + \theta = \sum_{i=1}^n w_i x_i + \theta$

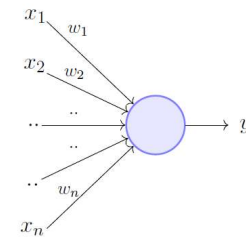
Predict( $p_1, p_2, p_3, \dots, p_n$ )

```
if  $\sum_{i=1}^n w_i p_i + \theta < 0$ :
    return N
else:
    return P
```

## NEURAL NETWORKS



## PERCEPTRON MODEL



$$y = 1 \quad \text{if } \sum_{i=1}^n w_i * x_i \geq \theta$$

$$= 0 \quad \text{if } \sum_{i=1}^n w_i * x_i < \theta$$

Rewriting the above,

$$y = 1 \quad \text{if } \sum_{i=1}^n w_i * x_i - \theta \geq 0$$

$$= 0 \quad \text{if } \sum_{i=1}^n w_i * x_i - \theta < 0$$

## PERCEPTRON

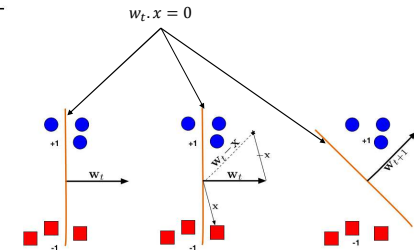
- First neural network learning model in the 1960's
- Simple and limited (single layer models)
- Basic concepts are similar for multi-layer models so this is a good learning tool
- Still used in many current applications (modems, etc.)

## PERCEPTRON ALGORITHM

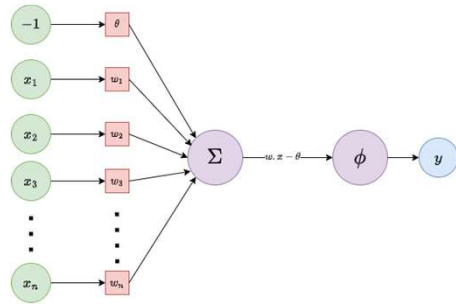
### Algorithm: Perceptron Learning Algorithm

```

P ← inputs with label 1;
N ← inputs with label 0;
Initialize w randomly;
while !convergence do
  Pick random x ∈ P ∪ N;
  if x ∈ P and w.x < 0 then
    w = w + x;
  end
  if x ∈ N and w.x ≥ 0 then
    w = w - x;
  end
end
end
//the algorithm converges when all the
inputs are classified correctly
  
```



## PERCEPTRON MODEL



## LEARNING AND GATE

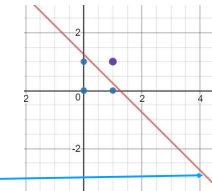
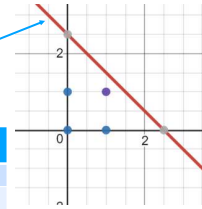
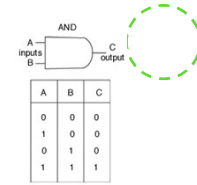
$$z = w_1 \cdot x_1 + w_2 \cdot x_2 - \theta$$

$$w_1=1, w_2=1, \theta=2.5$$

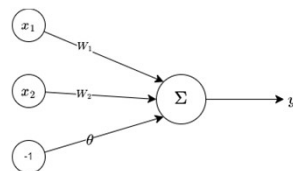
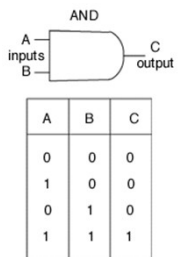
$$1 x_1 + 1 x_2 - 2.5 = 0$$

$w_1$	$w_2$	$- \theta$	$(x_1, x_2)$	$z$
1	1	-2.5	(0,1) -	$1 \times 0 + 1 \times 1 - 2.5 = -1.5$
1	1	-2.5	(1,1) +	$1 \times 1 + 1 \times 1 - 2.5 = -0.5$
2	2	-2.5	(0,0) -	$2 \times 0 + 2 \times 0 - 2.5 = -2.5$
2	2	-2.5	(1,0) -	$2 \times 1 + 2 \times 0 - 2.5 = -0.5$
2	2			

$$2 x_1 + 2 x_2 - 2.5 = 0$$

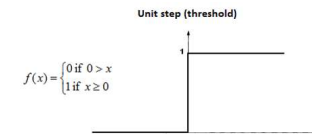


## LEARNING AND GATE



## LEARNING AND GATE

$x_1$	$x_2$	$y = 2x_1 + 2x_2 - 2.5$	$f(y)$
0	0	-2.5	0
0	1	-0.5	0
1	0	-0.5	0
1	1	1.5	1



## IMPLEMENTING AND GATE

```
#importing perceptron model from sklearn
from sklearn.linear_model import Perceptron
```

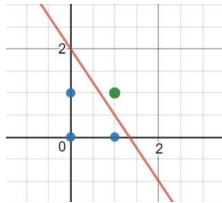
```
#training data for AND
X_train = [[0,0],[0,1],[1,0],[1,1]]
y_train = [0,0,0,1]
```

```
#model creation
clf = Perceptron(tol=1e-3, random_state=0)
clf.fit(X_train, y_train)
```

```
#prediction
y_pred=clf.predict(X_train)
print(y_pred)
```

```
print(clf.coef_,clf.intercept_)
```

```
[0 0 0 1]
[[3. 2.]] [-4.]
```



### Tools

- Python
- sklearn

### Try It For

OR, NAND, NOR, XOR

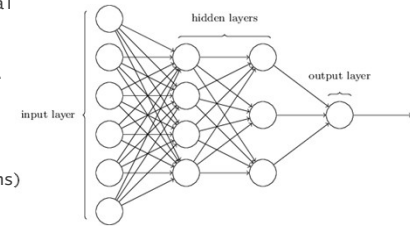
## MULTI LAYER PERCEPTRON(MLP)

■ Feedforward network: The neurons in each layer feed their output forward to the next layer until we get the final output from the neural network.

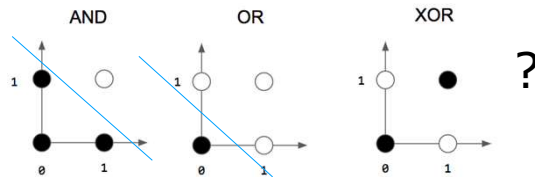
■ There can be any number of hidden layers within a feedforward network.

■ The number of neurons can be completely arbitrary.

■ MLP used to describe any general feedforward (no recurrent connections) network



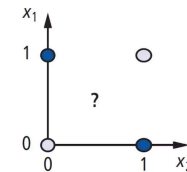
## MORE GATES

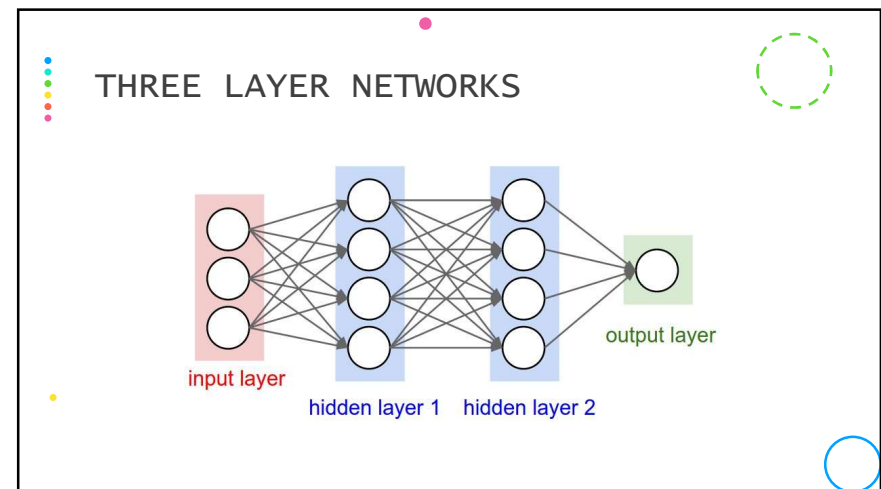
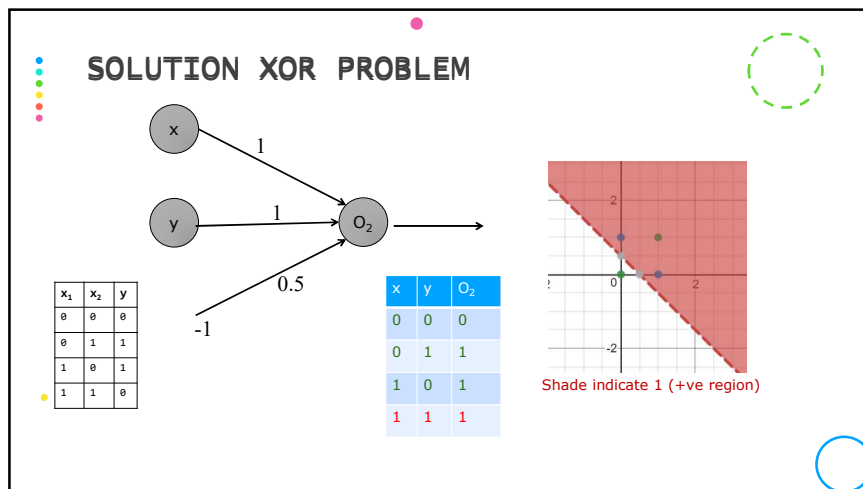
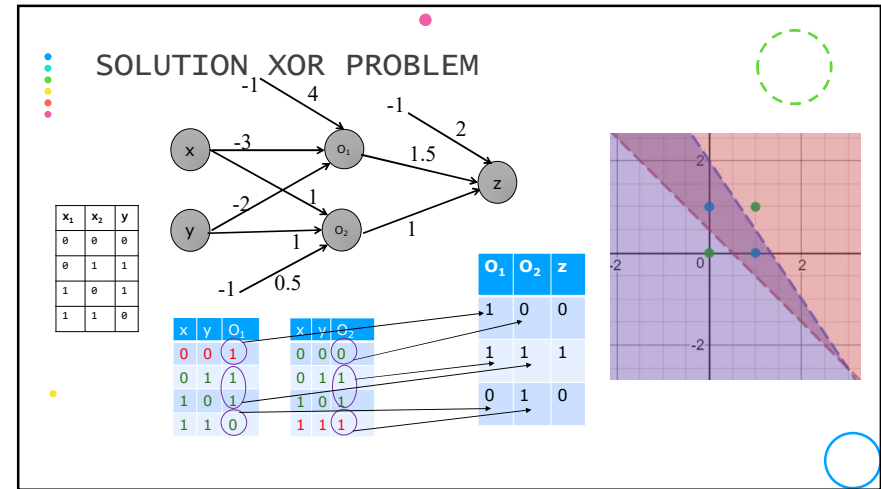
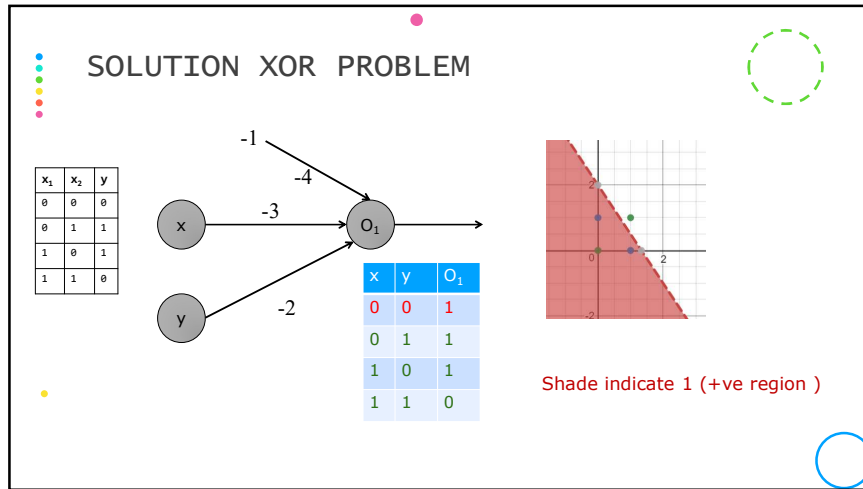


## AGAIN TO XOR PROBLEM

■ A Perceptron cannot represent Exclusive XOR since it is not linearly separable.

■ PERCEPTRON -----> MLP

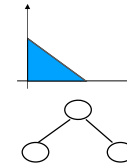




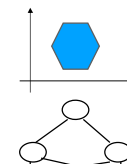
## THREE LAYER NETWORKS

- No connections within a layer
- No direct connections between input and output layers
- Fully connected between layers
- Often more than 3 layers
- Number of output units need not equal number of input units
- Number of hidden units per layer can be more or less than input or output units

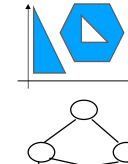
## WHAT DO EACH OF THESE LAYER DO ?



1st layer draws linear boundaries



2nd layer combines the boundaries

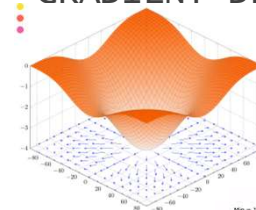


3rd layer can generate arbitrarily complex boundaries

## THREE LAYER NETWORKS

- No connections within a layer
- No direct connections between input and output layers
- Fully connected between layers
- Often more than 3 layers
- Number of output units need not equal number of input units
- Number of hidden units per layer can be more or less than input or output units

## GRADIENT DESCENT METHOD

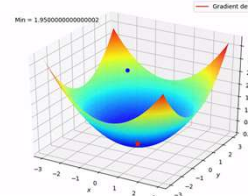


Function:  $f(x,y) = x^2 - y^2$   
 Starting point: (2, 3)  
 Learning rate ( $\alpha$ ): 0.1  
 Maximum iterations: 100

Repeat until convergence

$$\theta_j \leftarrow \theta_j - \alpha \frac{\partial}{\partial \theta_j} J(\theta)$$

}



1. Calculate the gradient:

1.  $\nabla f(x,y) = (2x, -2y)$
2. At (2, 3),  $\nabla f(2,3) = (4, -6)$

2. Update position:

1.  $x_{t+1} = x_t - \alpha * \nabla f_x$
2.  $y_{t+1} = y_t - \alpha * \nabla f_y$
3.  $x_{t+1} = 2 - 0.1 * 4 = 1.6$
4.  $y_{t+1} = 3 - 0.1 * (-6) = 3.6$

3. Repeat for maximum iterations:

1. Iteration 2: (1.6, 3.6)  $\rightarrow$  (1.24, 3.72)
2. Iteration 3: (1.24, 3.72)  $\rightarrow$  (1.008, 3.792)
3. ...
4. Iteration 100: (0.0000, 248453923.5660)

Final minimum point found: (0.0000, 248453923.5660)

## BACKPROPAGATION LEARNING ALGORITHM 'BP'

BP has two phases:

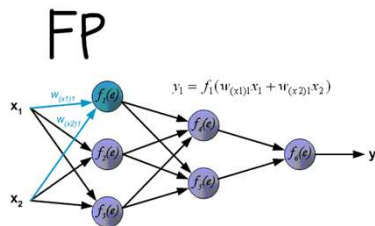
Forward pass phase: computes 'functional signal', feed forward propagation of input pattern signals through network

Backward pass phase: computes 'error signal', propagates the error backwards through network starting at output units (where the error is the difference between actual and desired output values)

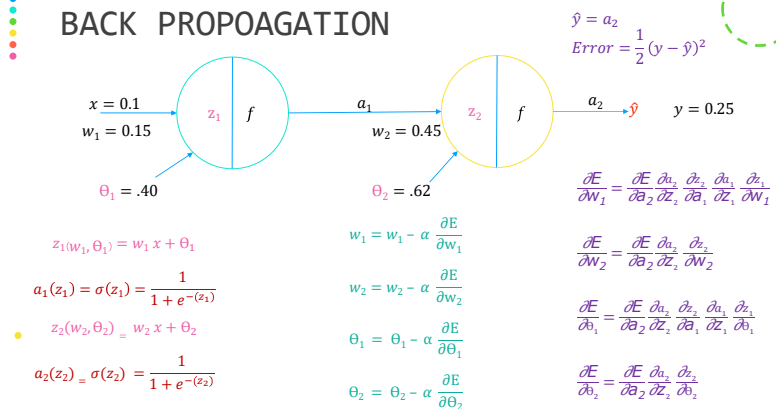
## BACKPROPAGATION LEARNING ALGORITHM 'BP'

- Error gradient along all connection weights were measured by propagating the error from output layer.
- First, a forward pass is performed - output of every neuron in every layer is computed.
- Output error is estimated.
- Then compute how much each neuron in last hidden layer contributed to output error.
- This is repeated backwards until input layer.
- Last step is Gradient Descent on all connection weights using error gradients estimated in previous steps.

## BACKPROPAGATION LEARNING ALGORITHM 'BP'



## BACK PROPOGATION



$z_1(w_1, \theta_1) = w_1 x + \theta_1 = 0.015 * 0.1 + 0.40 = 0.415$   
 $a_1(z_1) = \sigma(z_1) = \frac{1}{1 + e^{-(z_1)}} = \frac{1}{1 + e^{-(0.415)}} = 0.6022$   
 $z_2(w_2, \theta_2) = w_2 x + \theta_2 = 0.45 * 0.6022 + 0.63 = 0.890$   
 $a_2(z_2) = \sigma(z_2) = \frac{1}{1 + e^{-(z_2)}} = \frac{1}{1 + e^{-(0.890)}} = 0.7088$

$w_1 = w_1 - \alpha \frac{\partial E}{\partial w_1} = 0.15 - 1 * 0.00102 = 0.148$   
 $w_2 = w_2 - \alpha \frac{\partial E}{\partial w_2} = 0.45 - 1 * 0.0570 = 0.393$   
 $\theta_1 = \theta_1 - \alpha \frac{\partial E}{\partial \theta_1} = 0.40 - 1 * 0.102 = 0.389$   
 $\theta_2 = \theta_2 - \alpha \frac{\partial E}{\partial \theta_2} = 0.62 - 1 * 0.094 = 0.525$

$\frac{\partial E}{\partial w_1} = \frac{\partial E}{\partial z_2} \frac{\partial z_2}{\partial w_2} \frac{\partial z_2}{\partial a_2} \frac{\partial a_2}{\partial z_1} \frac{\partial z_1}{\partial w_1}$   
 $= (a_2 - y) a_2 (1 - a_2) w_2 a_1 (1 - a_1) x$   
 $= (0.7088 - 0.25) 0.7088 (1 - 0.7088) 0.1$   
 $= 0.00102$

$\frac{\partial E}{\partial w_2} = \frac{\partial E}{\partial z_2} \frac{\partial z_2}{\partial w_2} \frac{\partial z_2}{\partial a_2} \frac{\partial a_2}{\partial z_2}$   
 $= (a_2 - y) \sigma(z_2) (1 - \sigma(z_2)) a_1$   
 $= (a_2 - y) a_2 (1 - a_2) a_1$   
 $= (0.7088 - 0.25) 0.7088 (1 - 0.7088)$   
 $= 0.0570$

$\frac{\partial E}{\partial \theta_1} = \frac{\partial E}{\partial z_2} \frac{\partial z_2}{\partial \theta_1} \frac{\partial z_2}{\partial a_2} \frac{\partial a_2}{\partial z_1} \frac{\partial z_1}{\partial \theta_1}$   
 $= (a_2 - y) a_2 (1 - a_2) w_2 a_1 (1 - a_1) 1$   
 $= (0.7088 - 0.25) 0.7088 (1 - 0.7088) 1$   
 $= 0.0102$

$\frac{\partial E}{\partial \theta_2} = \frac{\partial E}{\partial z_2} \frac{\partial z_2}{\partial \theta_2} \frac{\partial z_2}{\partial a_2} \frac{\partial a_2}{\partial z_2}$   
 $= (a_2 - y) \sigma(z_2) (1 - \sigma(z_2)) 1$   
 $= (a_2 - 0.25) a_2 (1 - a_2)$   
 $= 0.0946$

$\sigma(x) = \frac{1}{1 + e^{-x}}$   
 $\sigma'(x) = \frac{1 + e^{-x} - 1}{(1 + e^{-x})(1 + e^{-x})} = \frac{1}{(1 + e^{-x})(1 + e^{-x})}$   
 $\sigma'(x) = \frac{1}{(1 + e^{-x})} - \frac{1}{(1 + e^{-x})(1 + e^{-x})}$   
 $\sigma'(x) = \frac{1}{(1 + e^{-x})} \left( 1 - \frac{1}{(1 + e^{-x})} \right)$   
 $\sigma'(x) = \sigma(x) \cdot (1 - \sigma(x))$

## STOCHASTIC GRADIENT DESCENT

Stochastic means randomness on which the algorithm is based upon

A variant of gradient descent that involves updating the parameters based on a small, randomly-selected subset of the data rather than the full dataset.

$$\theta_{t+1} = \theta_t - \alpha \nabla_{\theta} L_i(\theta_t)$$

$\theta_t$  represents the parameter vector at iteration  $t$ .

$\nabla_{\theta} L_i(\theta_t)$  is the gradient of the loss function for a randomly chosen training example at the current parameter vector  $\theta_t$ .

$\alpha$  is the learning rate, determining the step size of the parameter updates.

## MORE OPTIMIZERS

- **Optimizers**
- Algorithms or methods used to minimize an error function (Loss function) or to maximize the efficiency of production.
- Mathematical functions which are dependent on model's learnable parameters i.e. Weights & Biases.
- **Gradient Descent**
- **Stochastic Gradient Descent**
- **Stochastic Gradient Descent with Momentum**
- **Mini-Batch Gradient Descent**
- **AdaGrad (Adaptive Gradient Descent)**
- **RMS-Prop (Root Mean Square Propagation)**
- **AdaDelta**
- **Adam (Adaptive Moment Estimation)**

## MINI-BATCH GRADIENT DESCENT

Best among all the variations of gradient descent algorithms

Mini-batch gradient descent is similar to SGD, but instead of using a single sample to compute the gradient, it uses a small, fixed-size "mini-batch" of samples

$$\theta_{t+1} = \theta_t - \alpha \frac{1}{|B|} \nabla_{\theta} \sum_{i \in B} L_i(\theta_t)$$

$\theta_t$  represents the parameter vector at iteration  $t$ .

$\frac{1}{|B|} \nabla_{\theta} \sum_{i \in B} L_i(\theta_t)$  is the average gradient of the loss function for a randomly chosen training example at the current parameter vector  $\theta_t$ .

$\alpha$  is the learning rate, determining the step size of the parameter updates.



## SGD WITH MOMENTUM

### Momentum

Momentum was invented for reducing high variance in SGD and softens the convergence.

It accelerates the convergence towards the relevant direction and reduces the fluctuation to the irrelevant direction.

$\theta_t$  represents the parameter vector at iteration  $t$ .

$$\begin{aligned}\theta_{t+1} &= \theta_t - v_{t+1} \\ v_{t+1} &= \eta v_t + \alpha \nabla_{\theta} L(\theta_t)\end{aligned}$$

$\nabla_{\theta} L(\theta_t)$  is the gradient of the loss function for a randomly chosen training example at the current parameter vector  $\theta_t$ .

$\alpha$  - is the learning rate determining the step size of parameter updates.

$v$  - is the momentum term at iteration  $t$ .

$\eta$  - is the momentum coefficient (typically between 0 and 1), determining how much of the previous momentum to retain.

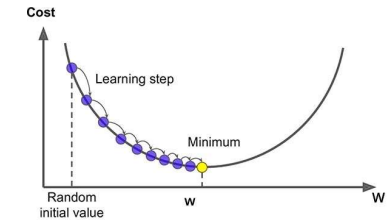
## ADAGRAD (ADAPTIVE GRADIENT DESCENT)

•Adagrad is an optimization algorithm that uses an adaptive learning rate per parameter.

•The learning rate is updated based on the historical gradient information so that parameters that receive many updates have a lower learning rate, and parameters that receive fewer updates have a larger learning rate.

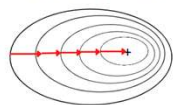
$$v_t = v_{t-1} + \left[ \frac{\partial L}{\partial w_t} \right]^2$$

$$w_{t+1} = w_t - \frac{\alpha}{\sqrt{v_t + \epsilon}} \cdot \frac{\partial L}{\partial w_t}$$

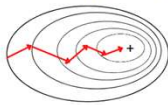


## SGD VARIANTS

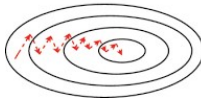
Batch Gradient Descent



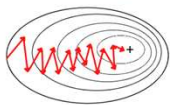
Mini-Batch Gradient Descent



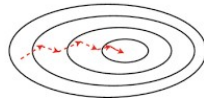
SGD without Momentum



Stochastic Gradient Descent



SGD with Momentum

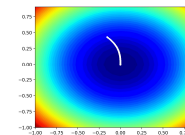


## ADA DELTA

•AdaDelta is an optimization algorithm similar to RMSProp but does not require a hyperparameter learning rate.

•It uses an exponentially decaying average of the gradients and the squares of the gradients to determine the updated scale.

$$\theta_t = \theta_{t-1} - \eta \cdot (\sqrt{G_t + \epsilon})^{-1/2} \cdot g_t$$



• $\theta_t$ : Parameter vector at iteration  $t$

• $\theta_{t-1}$ : Parameter vector at iteration  $t-1$

• $\eta$ : Learning rate

• $G_t$ : Diagonal matrix of accumulated squared gradients up to iteration  $t$

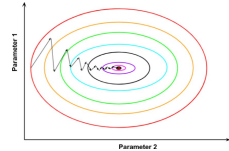
• $\epsilon$ : Small positive value to prevent division by zero

• $g_t$ : Gradient of the loss function with respect to  $\theta_t$  Visualization:

## RMSPROP(ROOT MEAN SQUARE PROBABILITY)

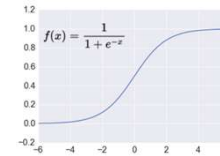
- RMSProp is an optimization algorithm similar to Adagrad, but it uses an exponentially decaying average of the squares of the gradients rather than the sum.
- Helps to reduce the monotonic learning rate decay of Adagrad and improve convergence.

$$\begin{aligned} \mathbf{g}_t &= \nabla_{\theta} L(\theta_t) \\ \mathbf{v}_t &= \gamma \mathbf{v}_{t-1} + (1 - \gamma) \mathbf{g}_t^2 \\ \theta_{t+1} &= \theta_t - \frac{\eta}{\sqrt{\mathbf{v}_t} + \epsilon} \mathbf{g}_t \end{aligned}$$



$\epsilon$  is a small positive constant for numerical stability.  $\eta$  is the learning rate.  $\gamma$  is the decay rate for the RMSProp.

## Activation: Sigmoid



$$R^n \rightarrow [0, 1]$$

Takes a real-valued number and "squashes" it into range between 0 and 1.

+ Nice interpretation as the **firing rate** of a neuron

- $0$  = not firing at all
- $1$  = fully firing
- Sigmoid neurons **saturate** and **kill gradients**, thus NN will barely learn
  - when the neuron's activation are 0 or 1 (saturate)

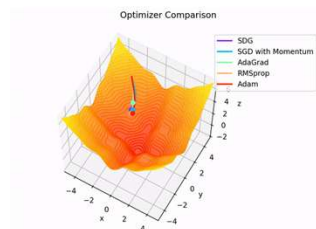
- 🙄 gradient at these regions almost zero
- 🙄 almost no signal will flow to its weights
- 🙄 if initial weights are too large then most neurons would saturate

## ADAM(ADAPTIVE MOMENT ESTIMATION)

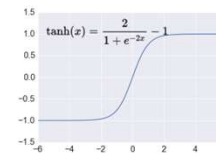
- Adam is an optimization algorithm that combines the ideas of SGD with momentum and RMSProp.
- It uses an exponentially decaying average of the gradients and the squares of the gradients to determine the updated scale, similar to RMSProp

$$\begin{aligned} \nu_t &= \beta_1 * \nu_{t-1} + (1 - \beta_1) * g_t \\ s_t &= \beta_2 * s_{t-1} + (1 - \beta_2) * g_t^2 \\ \Delta \omega_t &= -\eta \frac{\nu_t}{\sqrt{s_t} + \epsilon} * g_t \\ \omega_{t+1} &= \omega_t + \Delta \omega_t \end{aligned}$$

$\eta$ : Initial Learning rate  
 $g_t$ : Gradient at time  $t$  along  $\omega_j$   
 $\nu_t$ : Exponential Average of gradients along  $\omega_j$   
 $s_t$ : Exponential Average of squares of gradients along  $\omega_j$   
 $\beta_1, \beta_2$ : Hyperparameters



## Activation: Tanh

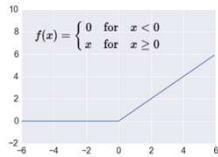


$$R^n \rightarrow [-1, 1]$$

Takes a real-valued number and "squashes" it into range between -1 and 1.

- Like sigmoid, tanh neurons **saturate**
- Unlike sigmoid, output is **zero-centered**
- Tanh is a **scaled sigmoid**:  $\tanh(x) = 2\sigma(2x) - 1$

## Activation: ReLU



$$f(x) = \max(0, x)$$

$$\mathbb{R}^n \rightarrow \mathbb{R}_+^n$$

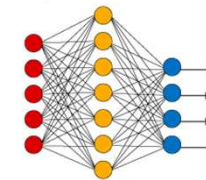
- Takes a real-valued number and thresholds it at zero

Most Deep Networks use ReLU nowadays

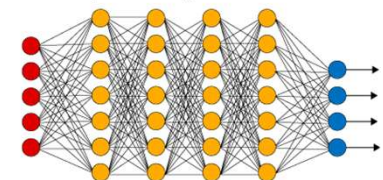
- Trains much **faster**
  - accelerates the convergence of SGD
  - due to linear, non-saturating form
- Less expensive operations
  - compared to sigmoid/tanh (exponentials etc.)
  - implemented by simply thresholding a matrix at zero
- More **expressive**
- Prevents the **gradient vanishing problem**

## DEEP NEURAL NETWORKS

Simple Neural Network



Deep Learning Neural Network



● Input Layer ● Hidden Layer ● Output Layer

## Loss Functions

Regression	Binary Classification	Multinomial Classification
Mean Squared Error (MSE)	Likelihood Loss (LL)	Categorical Cross Entropy (CCE)
Mean Absolute Error (MAE)	Binary Cross Entropy (BCE)	Kullback-Leibler Divergence (KLD)
Root Mean Squared Error (RMSE)	Hinge Loss and Squared Hinge Loss (H and SH)	
Mean Bias Error (MBE)		
Huber Loss (HL)		

True Label	Prediction	Binary Cross-Entropy
1	0.9	0.10536052
0	0.1	0.10536052
1	0.8	0.22314355
0	0.2	0.22314355
1	0.7	0.35667494
0	0.3	0.35667494
1	0.6	0.51082562
0	0.4	0.51082562
1	0.5	0.69314718
0	0.5	0.69314718

Binary Cross Entropy

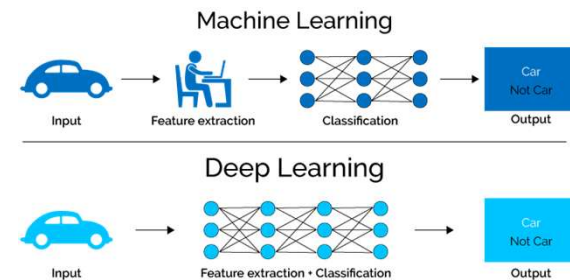
$$Loss = -\frac{1}{n} \sum_{i=1}^n y_i \log \hat{y}_i + (1 - y_i) \log (1 - \hat{y}_i)$$

Categorical Cross Entropy and Sparse Categorical Cross Entropy

$$Loss = -\sum_{i=1}^n y_i \log \hat{y}_i$$

- True label = [1, 0, 0], predicted probabilities = [0.8, 0.1, 0.1]  
 $CCE = -(1 * \log(0.8) + 0 * \log(0.1) + 0 * \log(0.1)) = 0.2231$
- True label = [0, 1, 0], predicted probabilities = [0.4, 0.5, 0.1]  
 $CCE = -(0 * \log(0.4) + 1 * \log(0.5) + 0 * \log(0.1)) = 0.6931$
- True label = [0, 0, 1], predicted probabilities = [0.2, 0.3, 0.5]  
 $CCE = -(0 * \log(0.2) + 0 * \log(0.3) + 1 * \log(0.5)) = 0.6931$

## MACHINE VS DEEP LEARNING





## TAKE AWAYS

- Mathematical Foundations for Optimization
- Neural Networks
- Perceptron
- Multilayer Perceptron
- Deep Neural Networks
- Gradient Descent
- Back Propagation
- Optimizers
- ➤ Activation Functions
- Loss Functions



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