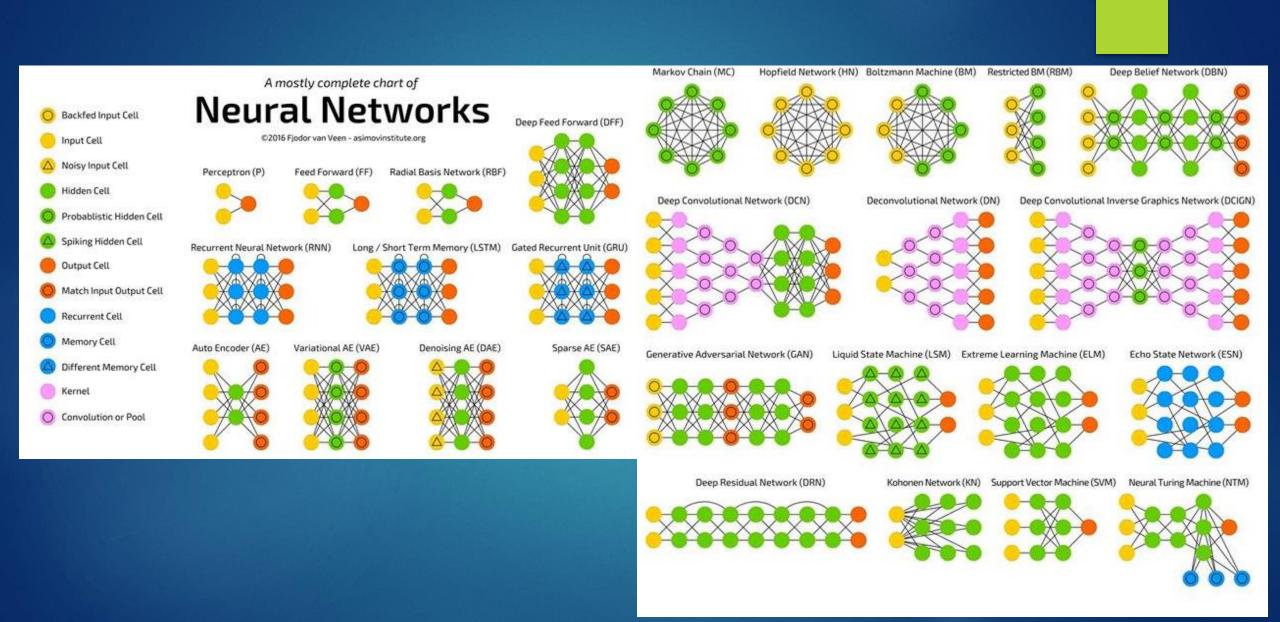
Building Time Series Predictive Model with Hyper-Parameters Tuning using Differential Evolution Algorithm

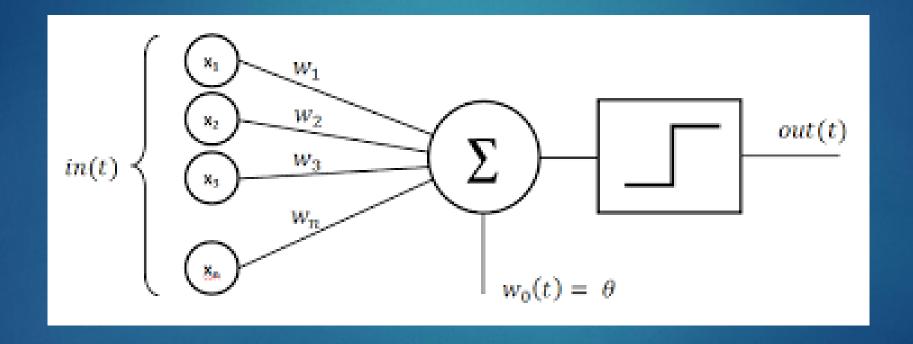
10/27/16

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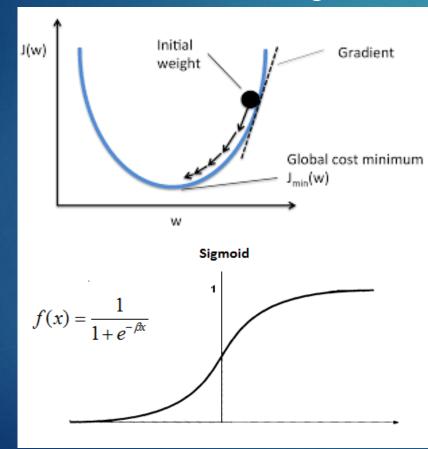


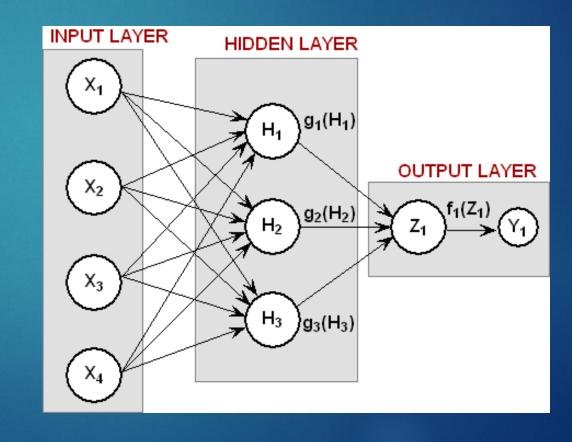
# Perceptron



#### Feed Forward

#### Gradient Descent Algorithm!





#### Momentum

Regular Momentum

$$v_t = \gamma v_{t-1} + \eta \nabla_{\theta} J(\theta).$$
  $heta = \theta - v_t.$ 

- Nesterov Accelerated Gradient (NAG)
  - Look ahead and then correct the position if it makes a mistakes

$$v_t = \gamma v_{t-1} + \eta \nabla_{\theta} J(\theta - \gamma v_{t-1}).$$
  $heta = \theta - v_t.$ 

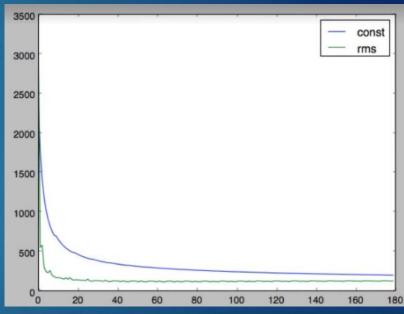
- https://www.youtube.com/watch?v=OWzkRD6MjYl
- And gifs

#### Learning Rate (Hyper-Parameter1)

- Point: Decay the learning rate as we close to the solution
- Step Decay
  - Ex) Reduce every 5, 10 or 20 steps
- Exponential Decay
  - $\triangleright$  Ex) LR = exp(LR)
- 1/t decay
  - $\triangleright$  EX) LR = 1/(1 + LR)

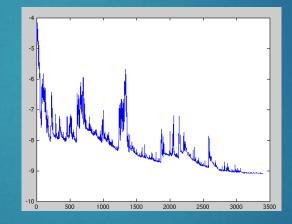
# Learning Rate (Hyper-Parameter1)

- Adagrad It adapts the learning rate to the parameters.
  - ▶ Keep the current changes in to the cache
    - ▶ More changes now and slower changes later
  - w -= LR \* gradient / sqrt(cache) + epsilon)
  - But it's too aggressive
- RMSprop
  - Make cache self decay (leaky cache)
  - cache = decay\_rate \* cache + (1-decay\_rate)\*gradient^2
  - w -= LR \* gradient / sqrt(cache) + epsilon)



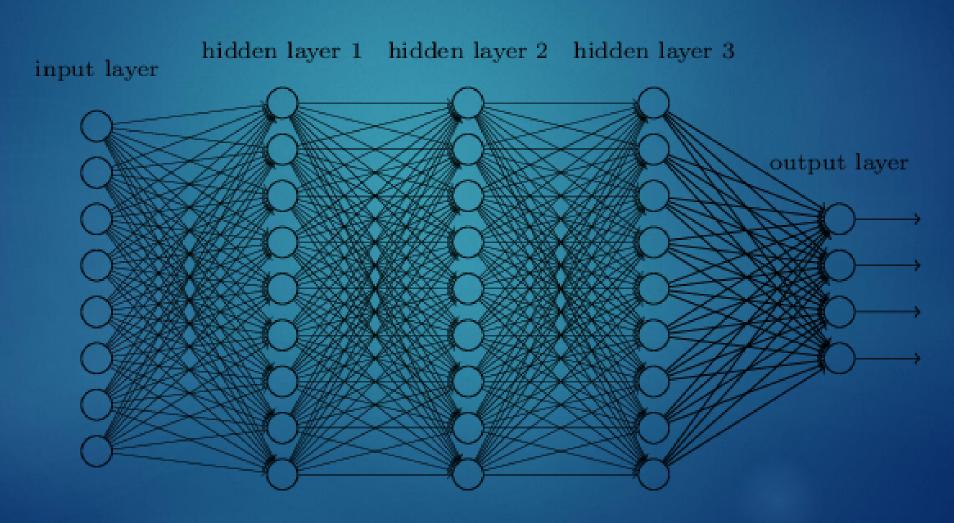
# Full vs Batch vs Stochastic GD (Hyper-Parameter 2)

- Full gradient descent
- ► O(N)
- Stochastic gradient descent



- Batch gradient descent
- O(batch size)  $\theta = \theta \eta \cdot \nabla_{\theta} J(\theta)$ .
- Batch size is another hyper-parameter

# Number of Node and Layers (Hyper-Parameter 3)



## Why RNN?

- Feed Forward
  - Can't take sequences
    - Ex) "I love dogs and cats" vs "Dogs love cats and I"
  - Number of layer grow exponentially as the number of input grows
- Markov Chain
  - longer seq. have prob. that approach 0 because it keeps multiplying values < 1</p>
  - Markov Assumption, is less accurate
    - ► Ex) p(cats) ~= p(cat | and)

## Why RNN?

- Predict a label over entire sequence
  - Ex) "hello world" -> male, "bob hi" -> female
- Predict a label for every step of input sequence
  - ▶ No need to wait till the end of the node

## Simple RNN

$$h(t) = f(W_h^T h(t-1) + W_x^T x(t) + b_h)$$

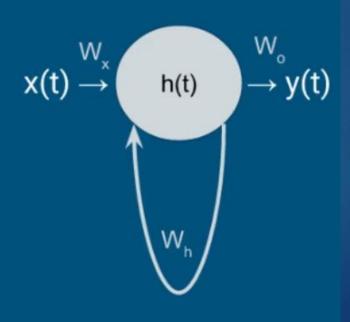
$$y(t) = softmax(W_o^Th(t) + b_o)$$

f = sigmoid, tanh, relu

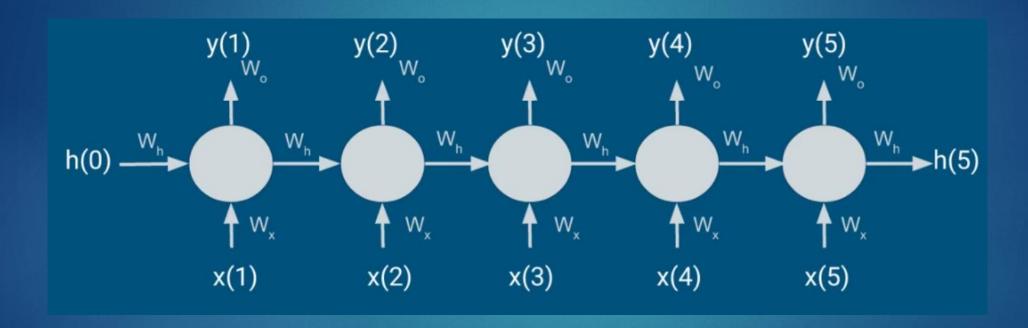
#### Regular feedforward net:

#### Recurrent net:



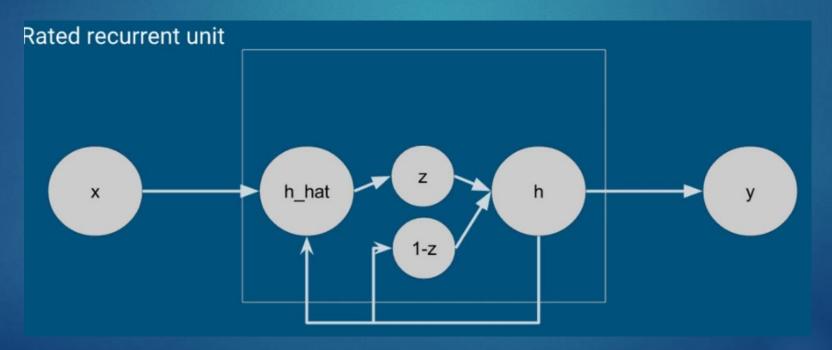


## Unfold Simple RNN



### Rated Recurrent Unit (RRU)

- Simple RNN has lack of variation, since it's multiplying the previous output over and over.
- RRU solves the problem by adding more variation



### Gate Recurrent Unit (GRU) - 2014

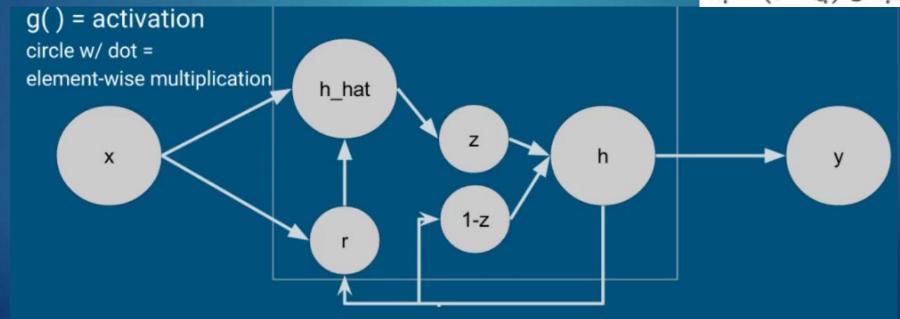
- Regular RNN has to take entire previous history
  - lots of unnecessary data -> wastes of memory
  - Need to forget some of the data (r = reset gate)

$$r_{t} = \sigma(x_{t}W_{xr} + h_{t-1}W_{hr} + b_{r})$$

$$z_{t} = \sigma(x_{t}W_{xz} + h_{t-1}W_{hz} + b_{z})$$

$$\hat{h}_{t} = g(x_{t}W_{xh} + (r_{t} \odot h_{t-1})W_{hh} + b_{h})$$

$$h_{t} = (1 - z_{t}) \odot h_{t-1} + z_{t} \odot \hat{h}_{t}.$$



#### Conclusion

- Why make RNN more complex?
  - Adding more parameters makes the model more expressive and able to solve more complex problem.