#### **SIMULATING RNNs on GPUs**

**WORKING TITLE** 

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#### **MULT-LAYER PERCEPTRON**

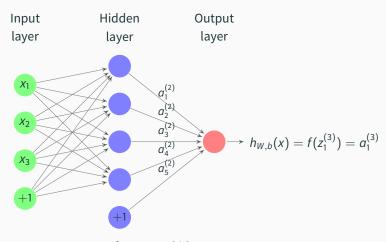
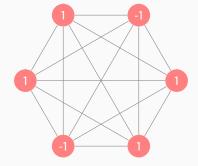
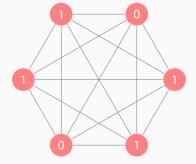


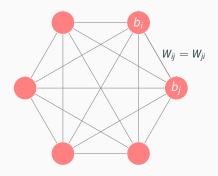
Figure 1: Multi-layer perceptron

**HOPFIELD NETWORKS AND** 

**BOLTZMANN MACHINES** 

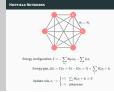




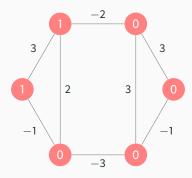


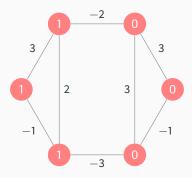
Energy configuration, 
$$E = -\sum_{i < j} W_{ij} x_i x_j - \sum_i b_i x_i$$
  
Energy gap,  $\Delta E_i = E(x_i = 0) - E(x_i = 1) = \sum_j W_{ij} x_j + b_i$   
Update rule,  $x_i := \begin{cases} +1 & \sum_j W_{ij} x_j + b_i \ge 0 \\ -1 & \text{otherwise} \end{cases}$ 

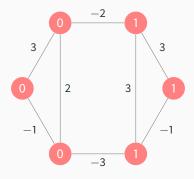
#### Hopfield Networks



- · composed of primitive computing elements called units
- units has two states, on or off, represented by  $\{1, -1\}$  or  $\{1, 0\}$
- connected to each other by bi-directional links
- adopts these states as a function of the states of its neighbouring units and weights of its links to them, it is a probabilistic function for a Boltzmann machine.
- · weights can take on any real value
- a unit being on or off is taken to mean that the system currently accepts or rejects some elemental hypothesis of the domain
- weight on a link represents a weak pairwise constrain between two hypothesis
- positive (negative) weights indicate that two hypothesis support (contradict) one another with other things being equal
- · link weights are symmetric, having the same strength in both directions







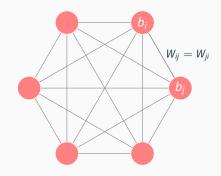
#### Simulating RNNs on GPUs

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#### └─ Hopfield Networks

- Updating of Hopfield networks is done sequentially usually in a randomized order. Parallel
  updating might increase the energy instead.
- Hopfield networks always make decisions to reduce the energy and makes it impossible to escape from local minima.
- random noise can help us escape from poor minima, by starting with lots of noise so its easy to
  cross energy barriers and gradually decrease the noise so the system ends in a deep minimum. This
  is called simulated annealing.

#### **BOLTZMANN MACHINES**



$$E = -\sum_{i < j} W_{ij} x_i x_j - \sum_i b_i x_i$$

$$\Delta E_i = E(x_i = 0) - E(x_i = 1) = \sum_j W_{ij} x_j + b_i$$

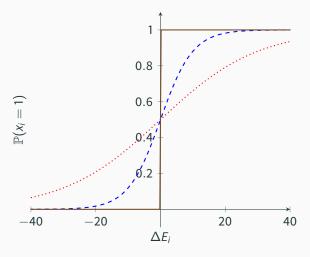
$$\mathbb{P}(x_i = 1) = \frac{1}{1 + e^{-\Delta E_i / \tau}}$$

#### **Boltzmann Machines**

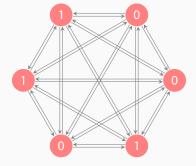


- replace the binary threshold units by binary stochastic units that make biased random decisions
- temperature variable controls the amount of noise
- when au o 0 we get back the Hopfield network
- for  $\tau_1 > \tau_2$ , we are less likely to go to a lower energy state compared to in  $\tau_1$  compared to  $\tau_2$ , i.e. more likely to go to a higher energy state when the temperature is higher. This allows us to escape from local minimum and arrive at the global minimum

#### **BOLTZMANN MACHINES**



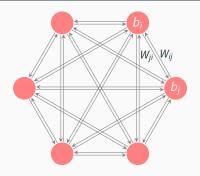
**Figure 2:** au= 0 (solid), au= 5 (dashed), au= 15 (dotted)



#### Simulating RNNs on GPUs



- state 1 is the refractory state, the neuron just fired and is unable to fire till it recovers
- state 0 is the armed state, the neuron just recovered and is waiting to fire
- here we model the units with the Nossenson-Messer neuron model, which explains biological firing rates in response to external stimuli



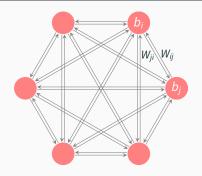
Transition Energy, 
$$E(y, x|\theta) = -\sum_{ji \in E} W_{ji}y_jx_i - \sum_{j \in V} b_js_j - \sum_{i \in V} b_is_i$$

$$\Gamma_{yx} = \exp\left(-\frac{1}{2\tau}E(y, x|\theta) + \frac{1}{2\tau}E(x, x|\theta)\right)$$

#### Simulating RNNs on GPUs

# Transition francy, $\xi(y,s) = -\sum_{j} u_j y_j - \sum_{j} b_j x_j - \sum_{j} b_j x_j - \sum_{j} b_k x_j -$

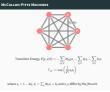
- allow transitions where y and x differ by only one bit
- here the W matrix need not be symmetrical with zero diagonals like what we had in the Hopfield network and Boltzmann machine models
- for each  $y \neq x$ , start a Poisson process with rate  $\Gamma_{yx}$



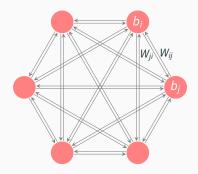
Transition Energy, 
$$E(y, x | \theta) = -\sum_{ji \in E} W_{ji} y_j x_i - \sum_{j \in V} b_j s_j - \sum_{i \in V} b_i s_i$$

$$\Gamma_{yx} := \exp\left(\frac{1}{2\tau} s_j z_j\right)$$

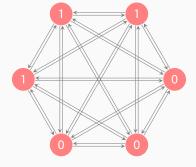
where  $s_j = 1 - 2x_j$ ,  $z_j = \sum_j W_{ji}x_i + b_j$  and x, y differ by the jth unit.

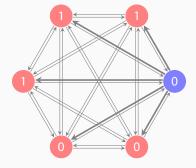


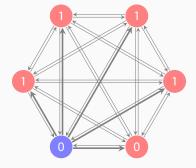
• when doing the updates we can just update the linear responses  $z_j$  and apply softmax on the  $\lambda_j$ 's to get the probability distribution of the transitions.

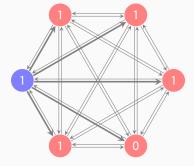


Transition probability from 
$$x$$
 to  $y$ ,  $p_{yx} = \frac{\lambda_j}{\sum_{j'} \lambda_{j'}}$ 









#### Simulating RNNs on GPUs



- digraph, G=(V,E), weights  $W:V\to\mathbb{R}$ , biases  $b:E\to\mathbb{R}$ , binary states  $\mathbb{B}=\{0,1\}$  with an initial distribution  $\mathbb{B}^{|V|}\to\delta$  and a temperature  $\tau$
- allow transitions where y and x differ by only one bit
- for each  $y \neq x$ , start a Poisson process with rate  $\Gamma_{yx}$
- if the Poisson process for state y' is the first to fire, or the interarrival timing from x to y' is the shortest,



$$E(x^{(n+1)}, x^{(n)}|\theta) = -\sum_{ji \in E} W_{ji} x_j^{(n+1)} x_i^{(n)} - \sum_{j \in V} b_j s_j - \sum_{i \in V} b_i s_i$$
$$\Gamma_{x^{(n+1)} x^{(n)}} = \exp\left(-\frac{1}{2\tau} E(x^{(n+1)}, x^{(n)}|\theta) + \frac{1}{2\tau} E(x^{(n)}, x^{(n)}|\theta)\right)$$



#### REFERENCES



J. Macor.

#### A Brief Introduction to Type Theory and the Univalence Axiom

http://math.uchicago.edu/ may/REU2015/REUPapers/Macor.pdf



The Univalent Foundations Program

Homotopy Type Theory: Univalent Foundations of Mathematics.

https://homotopytypetheory.org/book



The n-Category Café

#### From Set Theory to Type Theory

https://golem.ph.utexas.edu/category/2013/01/from\_set\_theory\_to\_type\_theory.html



The nLab

#### Function Type

https://ncatlab.org/nlab/show/function+type