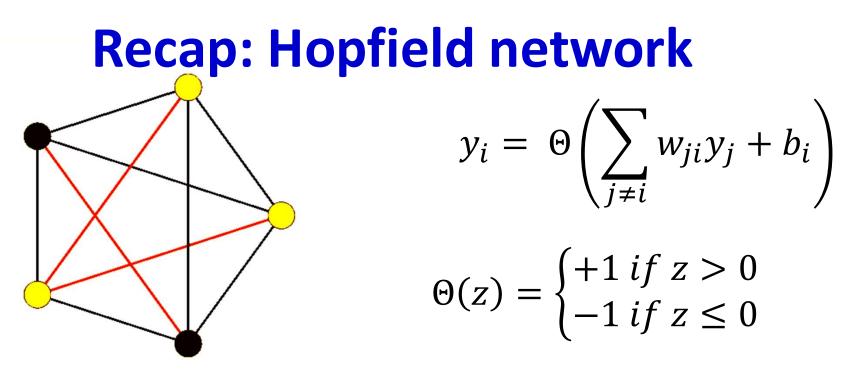
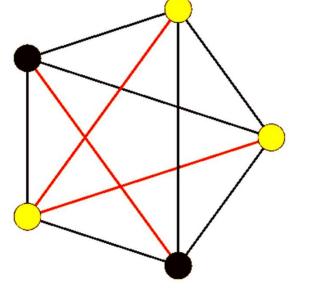
Neural Networks

Hopfield Nets and Boltzmann Machines



- At each time each neuron receives a "field" $\sum_{j \neq i} w_{ji} y_j + b_i$
- If the sign of the field matches its own sign, it does not respond
- If the sign of the field opposes its own sign, it "flips" to match the sign of the field

Recap: Energy of a Hopfield Network



$$y_i = \Theta\left(\sum_{j\neq i} w_{ji} y_j + b_i\right)$$

 $\Theta(z) = \begin{cases} +1 \text{ if } z > 0\\ -1 \text{ if } z \le 0 \end{cases}$

$$E = -\sum_{i,j$$

- The system will evolve until the energy hits a local minimum
- In vector form
 - Bias term may be viewed as an extra input pegged to 1.0

$$E = -\frac{1}{2}\mathbf{y}^T\mathbf{W}\mathbf{y} - \mathbf{b}^T\mathbf{y}$$

Recap: Hopfield net computation

1. Initialize network with initial pattern

$$y_i(0) = x_i, \qquad 0 \le i \le N - 1$$

2. Iterate until convergence $\langle \mathbf{\nabla} \rangle$

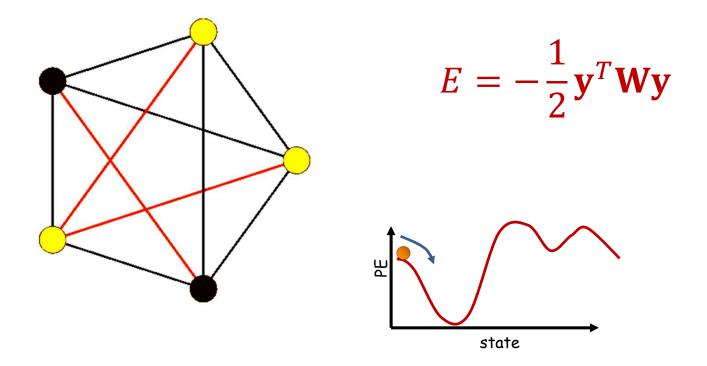
$$y_i(t+1) = \Theta\left(\sum_{j\neq i} w_{ji}y_j\right), \qquad 0 \le i \le N-1$$

- Very simple
- Updates can be done sequentially, or all at once
- Convergence

$$E = -\sum_{i} \sum_{j>i} w_{ji} y_j y_i$$

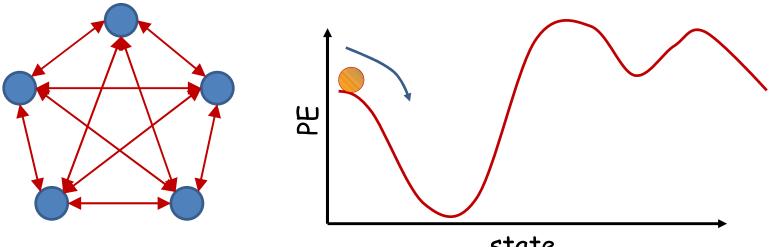
does not change significantly any more

Recap: Evolution



• The network will evolve until it arrives at a local minimum in the energy contour

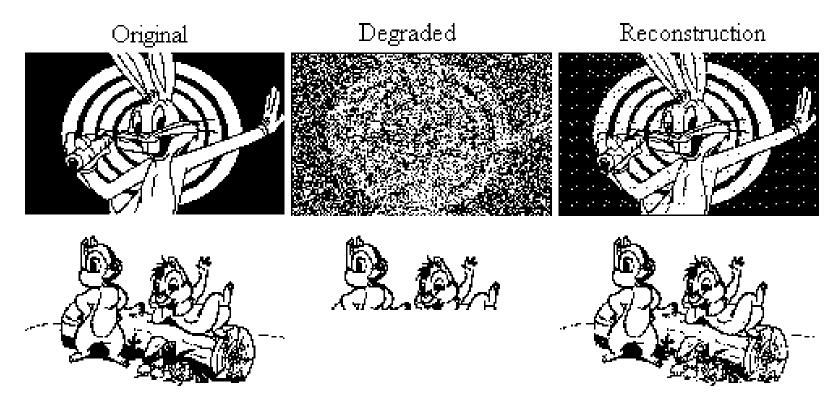
Recap: Content-addressable memory



state

- Each of the minima is a "stored" pattern
 - If the network is initialized close to a stored pattern, it will inevitably evolve to the pattern
- This is a *content addressable memory*
 - Recall memory content from partial or corrupt values
- Also called *associative memory*

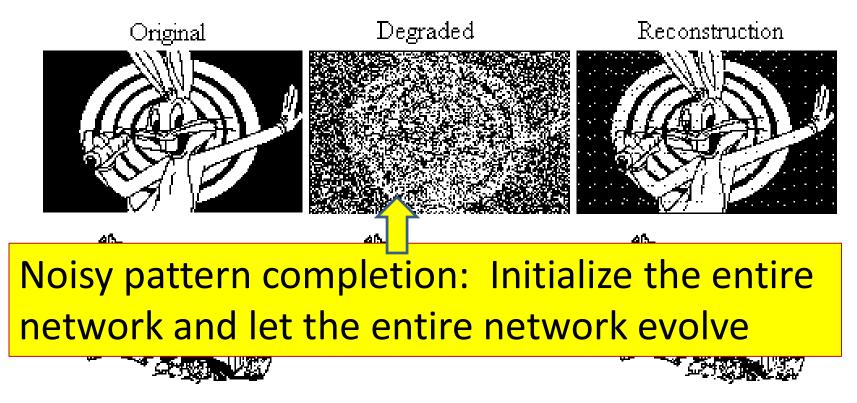
Examples: Content addressable memory



Hopfield network reconstructing degraded images from noisy (top) or partial (bottom) cues.

http://staff.itee.uq.edu.au/janetw/cmc/chapters/Hopfield/ 7

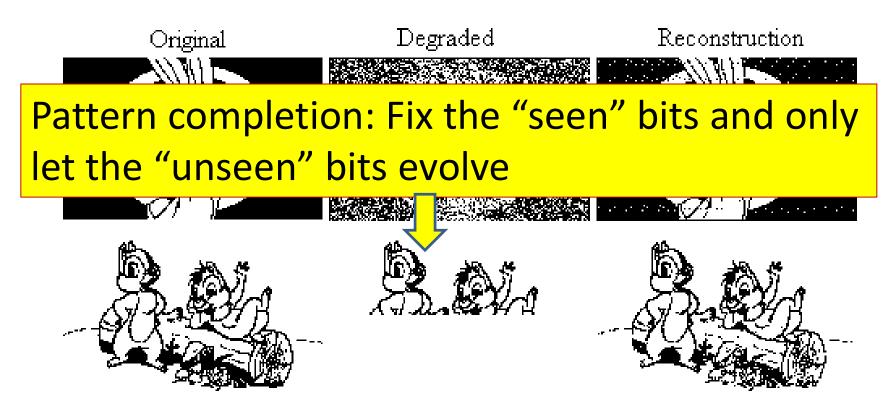
Examples: Content addressable memory



Hopfield network reconstructing degraded images from noisy (top) or partial (bottom) cues.

http://staff.itee.uq.edu.au/janetw/cmc/chapters/Hopfield/ 8

Examples: Content addressable memory



Hopfield network reconstructing degraded images from noisy (top) or partial (bottom) cues.

http://staff.itee.uq.edu.au/janetw/cmc/chapters/Hopfield/ 9

Training a Hopfield Net to "Memorize" target patterns

• The Hopfield network can be *trained* to remember specific "target" patterns

– E.g. the pictures in the previous example

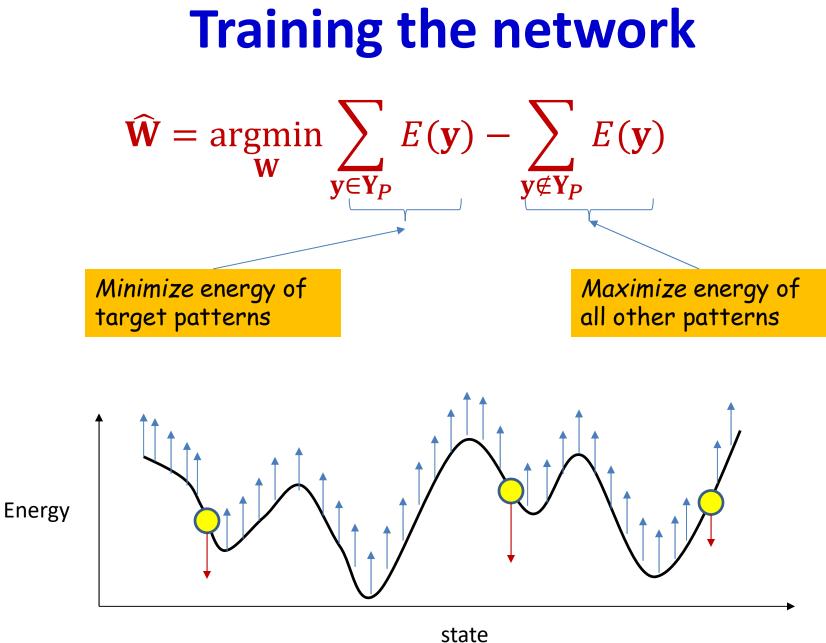
This can be done by setting the weights W appropriately

Random Question: Can you use backprop to train Hopfield nets?

Hint: Think unwrapping...

Training a Hopfield Net to "Memorize" target patterns

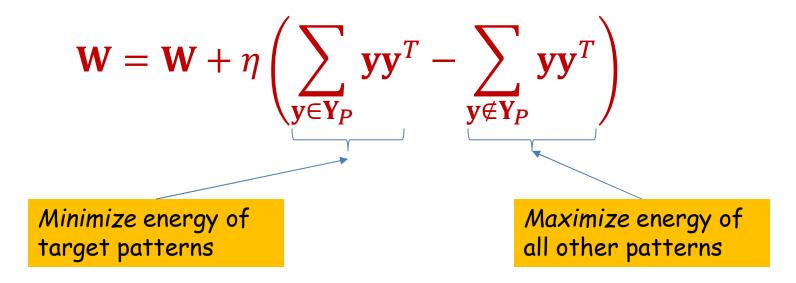
- The Hopfield network can be *trained* to remember specific "target" patterns
 - E.g. the pictures in the previous example
- A Hopfield net with N neurons can designed to store up to N target N-bit memories
 - But can store an exponential number of unwanted "parasitic" memories along with the target patterns
- Training the network: Design weights matrix W such that the energy of ...
 - Target patterns is minimized, so that they are in energy wells
 - Other untargeted potentially parasitic patterns is maximized so that they don't become parasitic

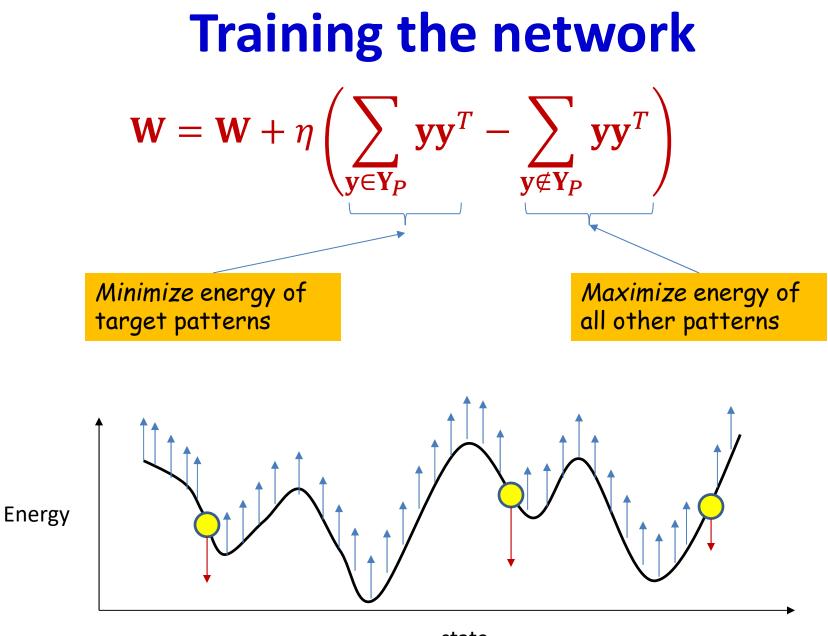


Optimizing W

$$E(\mathbf{y}) = -\frac{1}{2}\mathbf{y}^T \mathbf{W} \mathbf{y} \quad \widehat{\mathbf{W}} = \underset{\mathbf{W}}{\operatorname{argmin}} \sum_{\mathbf{y} \in \mathbf{Y}_P} E(\mathbf{y}) - \sum_{\mathbf{y} \notin \mathbf{Y}_P} E(\mathbf{y})$$

• Simple gradient descent:

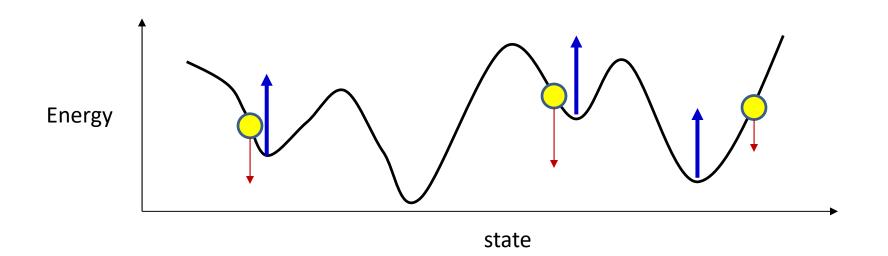




Simpler: Focus on confusing parasites

$$\mathbf{W} = \mathbf{W} + \eta \left(\sum_{\mathbf{y} \in \mathbf{Y}_{P}} \mathbf{y} \mathbf{y}^{T} - \sum_{\mathbf{y} \notin \mathbf{Y}_{P} \& \mathbf{y} = valley} \mathbf{y} \mathbf{y}^{T} \right)$$

- Focus on minimizing parasites that can prevent the net from remembering target patterns
 - Energy valleys in the neighborhood of target patterns

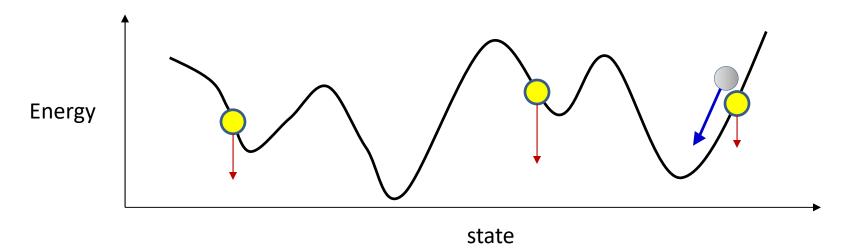


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Simpler: Focus on confusing patterns

$$\mathbf{W} = \mathbf{W} + \eta \left(\sum_{\mathbf{y} \in \mathbf{Y}_{P}} \mathbf{y} \mathbf{y}^{T} - \sum_{\mathbf{y} \notin \mathbf{Y}_{P} \& \mathbf{y} = valley} \mathbf{y} \mathbf{y}^{T} \right)$$

- Lower energy at valid memories
- Initialize the network at valid memories and let it evolve
 - It will settle in a valley. If this is not the target pattern, raise it



Training the Hopfield network $\mathbf{W} = \mathbf{W} + \eta \left(\sum_{\mathbf{y} \in \mathbf{Y}_P} \mathbf{y} \mathbf{y}^T - \sum_{\mathbf{y} \notin \mathbf{Y}_P \& \mathbf{y} = valley} \mathbf{y} \mathbf{y}^T \right)$

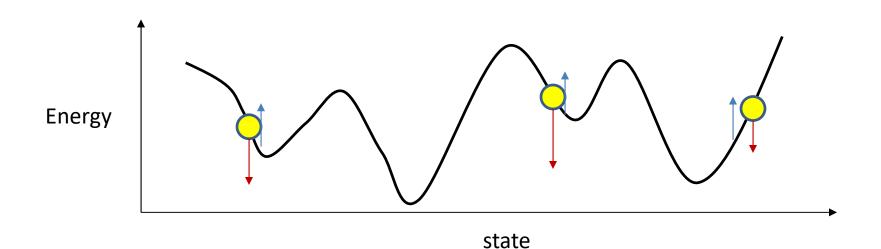
- Initialize W
- Compute the total outer product of all target patterns
 - More important patterns presented more frequently
- Initialize the network with each target pattern and let it evolve
 - And settle at a valley
- Compute the total outer product of valley patterns
- Update weights

Training the Hopfield network: SGD version $\mathbf{W} = \mathbf{W} + \eta \left(\sum_{\mathbf{y} \in \mathbf{Y}_P} \mathbf{y} \mathbf{y}^T - \sum_{\mathbf{y} \notin \mathbf{Y}_P \& \mathbf{y} = valley} \mathbf{y} \mathbf{y}^T \right)$

- Initialize W
- Do until convergence, satisfaction, or death from boredom:
 - Sample a target pattern \mathbf{y}_p
 - Sampling frequency of pattern must reflect importance of pattern
 - Initialize the network at \mathbf{y}_p and let it evolve
 - And settle at a valley $y_{\boldsymbol{\mathcal{V}}}$
 - Update weights
 - $\mathbf{W} = \mathbf{W} + \eta (\mathbf{y}_p \mathbf{y}_p^T \mathbf{y}_v \mathbf{y}_v^T)$

More efficient training

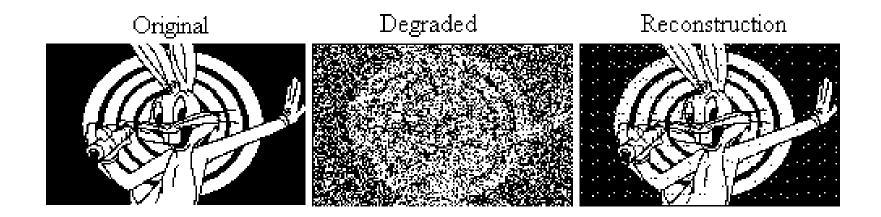
- Really no need to raise the entire surface, or even every valley
- Raise the *neighborhood* of each target memory
 - Sufficient to make the memory a valley
 - The broader the neighborhood considered, the broader the valley



Training the Hopfield network: SGD version $\mathbf{W} = \mathbf{W} + \eta \left(\sum_{\mathbf{y} \in \mathbf{Y}_P} \mathbf{y} \mathbf{y}^T - \sum_{\mathbf{y} \notin \mathbf{Y}_P \& \mathbf{y} = valley} \mathbf{y} \mathbf{y}^T \right)$

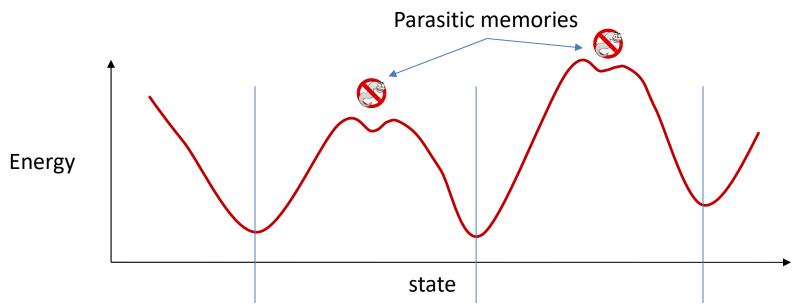
- Initialize W
- Do until convergence, satisfaction, or death from boredom:
 - Sample a target pattern \mathbf{y}_p
 - Sampling frequency of pattern must reflect importance of pattern
 - Initialize the network at \mathbf{y}_p and let it evolve **a** few steps (2-4)
 - And arrive at a down-valley position \mathbf{y}_d
 - Update weights
 - $\mathbf{W} = \mathbf{W} + \eta (\mathbf{y}_p \mathbf{y}_p^T \mathbf{y}_d \mathbf{y}_d^T)$

Problem with Hopfield net



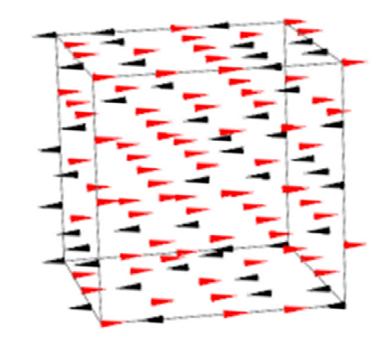
• Why is the recalled pattern not perfect?

A Problem with Hopfield Nets



- Many local minima
 - Parasitic memories
- May be escaped by adding some *noise* during evolution
 - Permit changes in state even if energy increases..
 - Particularly if the increase in energy is small

Recap – Analogy: Spin Glasses



Total field at current dipole:

$$f(p_i) = \sum_{j \neq i} J_{ij} x_j + b_i$$

Response of current diplose

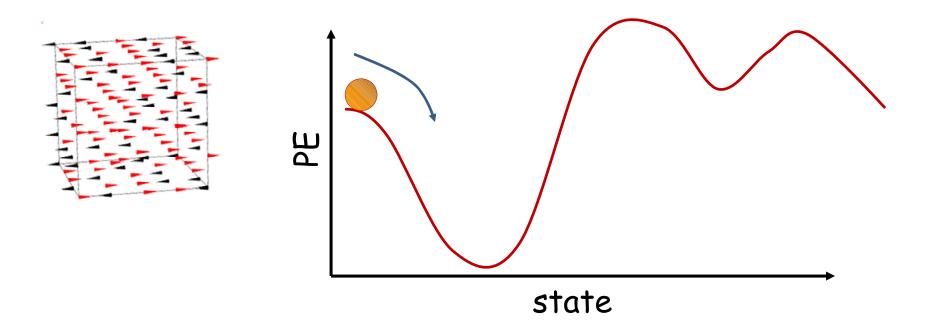
$$x_{i} = \begin{cases} x_{i} \text{ if } sign(x_{i} f(p_{i})) = 1 \\ -x_{i} \text{ otherwise} \end{cases}$$

• The total energy of the system

$$E(s) = C - \frac{1}{2} \sum_{i} x_{i} f(p_{i}) = -\sum_{i} \sum_{j>i} J_{ij} x_{i} x_{j} - \sum_{i} b_{i} x_{j}$$

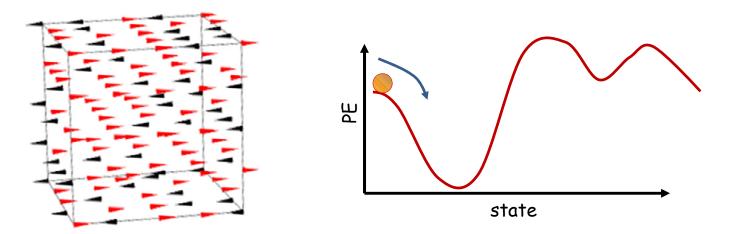
- The system *evolves* to minimize the energy
 - Dipoles stop flipping if flips result in increase of energy

Recap : Spin Glasses



- The system stops at one of its *stable* configurations
 - Where energy is a local minimum

Revisiting Thermodynamic Phenomena



- Is the system actually in a specific state at any time?
- No the state is actually continuously changing
 - Based on the temperature of the system
 - At higher temperatures, state changes more rapidly
- What is actually being characterized is the *probability* of the state at equilibrium
 - The system "prefers" low energy states
 - Evolution of the system favors transitions towards lower-energy states

- A thermodynamic system at temperature *T* can exist in one of many states
 - Potentially infinite states
 - At any time, the probability of finding the system in state sat temperature T is $P_T(s)$
- At each state s it has a potential energy E_s
- The *internal energy* of the system, representing its capacity to do work, is the average:

$$U_T = \sum_{s} P_T(s) E_s$$

• The capacity to do work is counteracted by the internal disorder of the system, i.e. its entropy

$$H_T = -\sum_s P_T(s) \log P_T(s)$$

• The *Helmholtz* free energy of the system measures the *useful* work derivable from it and combines the two terms

$$F_T = U_T + kTH_T$$

$$= \sum_{s} P_T(s) E_s - kT \sum_{s} P_T(s) \log P_T(s)$$

$$F_T = \sum_{s} P_T(s) E_s - kT \sum_{s} P_T(s) \log P_T(s)$$

- A system held at a specific temperature *anneals* by varying the rate at which it visits the various states, to reduce the free energy in the system, until a minimum free-energy state is achieved
- The probability distribution of the states at steady state is known as the *Boltzmann distribution*

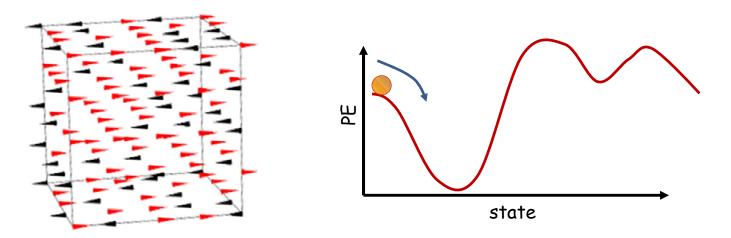
$$F_T = \sum_{s} P_T(s) E_s - kT \sum_{s} P_T(s) \log P_T(s)$$

• Minimizing this w.r.t $P_T(s)$, we get

$$P_T(s) = \frac{1}{Z} exp\left(\frac{-E_s}{kT}\right)$$

- Also known as the *Gibbs* distribution
- -Z is a normalizing constant
- Note the dependence on T
- A T = 0, the system will always remain at the lowestenergy configuration with prob = 1.

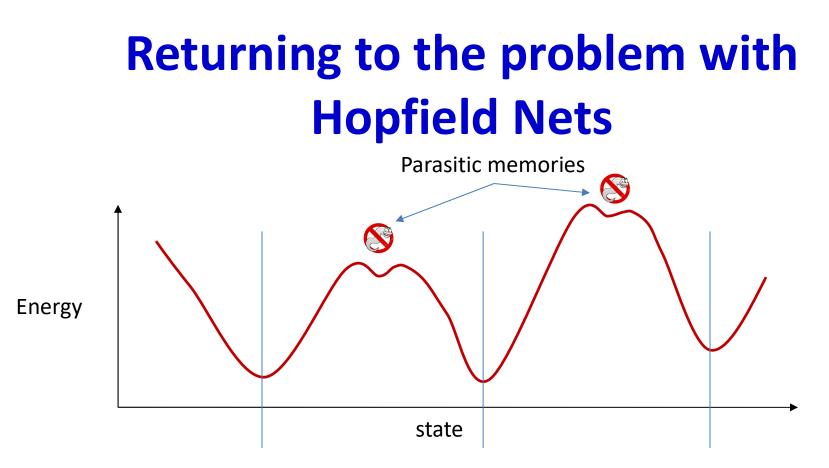
Revisiting Thermodynamic Phenomena



- The evolution of the system is actually *stochastic*
- At equilibrium the system visits various states according to the Boltzmann distribution
 - The probability of any state is inversely related to its energy

• and also temperatures:
$$P(s) \propto exp\left(\frac{-E_s}{kT}\right)$$

• The most likely state is the lowest energy state

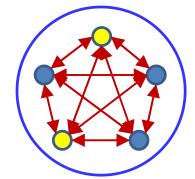


- Many local minima
 - Parasitic memories
- May be escaped by adding some *noise* during evolution
 - Permit changes in state even if energy increases..
 - Particularly if the increase in energy is small

The Hopfield net as a distribution

Visible Neurons

$$E(S) = -\sum_{i < j} w_{ij} s_i s_j - b_i s_i$$

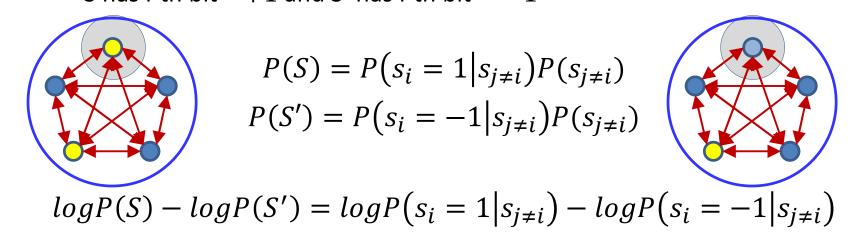


$$P(S) = \frac{exp(-E(S))}{\sum_{S'} exp(-E(S'))}$$

- Mimics the Spin glass system
- The stochastic Hopfield network models a *probability distribution* over states
 - Where a state is a binary string
 - Specifically, it models a *Boltzmann distribution*
 - The parameters of the model are the weights of the network
- The probability that (at equilibrium) the network will be in any state is P(S)
 - It is a *generative* model: generates states according to P(S)

The field at a single node

Let S and S' be otherwise identical states that only differ in the i-th bit
 S has i-th bit = +1 and S' has i-th bit = -1



$$logP(S) - logP(S') = log \frac{P(s_i = 1|s_{j\neq i})}{1 - P(s_i = 1|s_{j\neq i})}$$

The field at a single node

• Let S and S' be the states with the ith bit in the +1 and - 1 states $\log P(S) = -E(S) + C$

$$\log P(S) = -E(S) + C$$

$$E(S) = -\frac{1}{2} \left(E_{not \, i} + \sum_{j \neq i} w_{ij} s_j + b_i \right)$$

$$E(S') = -\frac{1}{2} \left(E_{not \, i} - \sum_{j \neq i} w_{ij} s_j - b_i \right)$$

• $logP(S) - logP(S') = E(S') - E(S) = \sum_{j \neq i} w_{ij}s_j + b_i$

The field at a single node

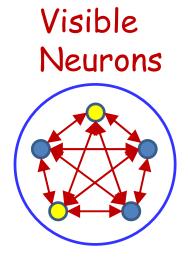
$$log\left(\frac{P(s_{i}=1|s_{j\neq i})}{1-P(s_{i}=1|s_{j\neq i})}\right) = \sum_{j\neq i} w_{ij}s_{j} + b_{i}$$

• Giving us

$$P(s_{i} = 1 | s_{j \neq i}) = \frac{1}{1 + e^{-(\sum_{j \neq i} w_{ij}s_{j} + b_{i})}}$$

• The probability of any node taking value 1 given other node values is a logistic

Redefining the network



$$z_i = \sum_j w_{ij} s_j + b_i$$

$$P(s_i = 1 | s_{j \neq i}) = \frac{1}{1 + e^{-z_i}}$$

- First try: Redefine a regular Hopfield net as a stochastic system
- Each neuron is now a stochastic unit with a binary state s_i, which can take value 0 or 1 with a probability that depends on the local field
 - Note the slight change from Hopfield nets
 - Not actually necessary; only a matter of convenience

The Hopfield net is a distribution

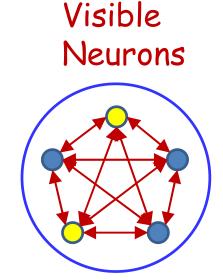
Visible Neurons

$$z_i = \sum_j w_{ij} s_j + b_i$$

$$P(s_i = 1 | s_{j \neq i}) = \frac{1}{1 + e^{-z_i}}$$

- The Hopfield net is a probability distribution over binary sequences
 - The Boltzmann distribution
- The *conditional* distribution of individual bits in the sequence is a logistic

Running the network



$$z_i = \sum_j w_{ij} s_j + b_i$$

$$P(s_i = 1 | s_{j \neq i}) = \frac{1}{1 + e^{-z_i}}$$

- Initialize the neurons
- Cycle through the neurons and randomly set the neuron to 1 or 0 according to the probability given above
 - Gibbs sampling: Fix N-1 variables and sample the remaining variable
 - As opposed to energy-based update (mean field approximation): run the test $z_i > 0$?
- After many many iterations (until "convergence"), *sample* the individual neurons

1. Initialize network with initial pattern

 $y_i(0) = x_i, \qquad 0 \le i \le N - 1$

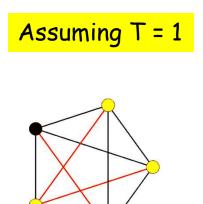
2. Iterate $0 \le i \le N - 1$ $P = \sigma \left(\sum_{j \ne i} w_{ji} y_j \right)$ $y_i(t+1) \sim Binomial(P)$ Assuming T = 1

1. Initialize network with initial pattern

 $y_i(0) = x_i, \qquad 0 \le i \le N - 1$

2. Iterate
$$0 \le i \le N - 1$$

 $P = \sigma\left(\sum_{j \ne i} w_{ji}y_j\right)$
 $y_i(t+1) \sim Binomial(P)$



- When do we stop?
- What is the final state of the system
 - How do we "recall" a memory?

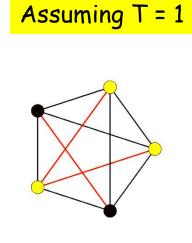
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$$y_i(t+1) \sim Binomial(P)$$



- When do we stop?
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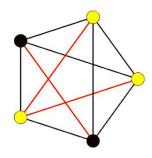
1. Initialize network with initial pattern

$$y_i(0) = x_i, \qquad 0 \le i \le N - 1$$

Iterate
$$0 \le i \le N - 1$$

 $P = \sigma\left(\sum_{j \ne i} w_{ji}y_j\right)$
 $y_i(t+1) \sim Binomial(P)$

Assuming T = 1



• Let the system evolve to "equilibrium"

2.

- Let $\mathbf{y}_0, \mathbf{y}_1, \mathbf{y}_2, \dots, \mathbf{y}_L$ be the sequence of values (L large)
- Final predicted configuration: from the average of the final few iterations

$$\mathbf{y} = \left(\frac{1}{M} \sum_{t=L-M+1}^{L} \mathbf{y}_t\right) > 0?$$

- Estimates the probability that the bit is 1.0.
- If it is greater than 0.5, sets it to 1.0

Evolution of the stochastic network

1. Initialize network with initial pattern

$$y_i(0) = x_i, \qquad 0 \le i \le N - 1$$

2. For $T = T_0$ down to T_{min}

Noisy pattern completion: Initialize the entire network and let the entire network evolve

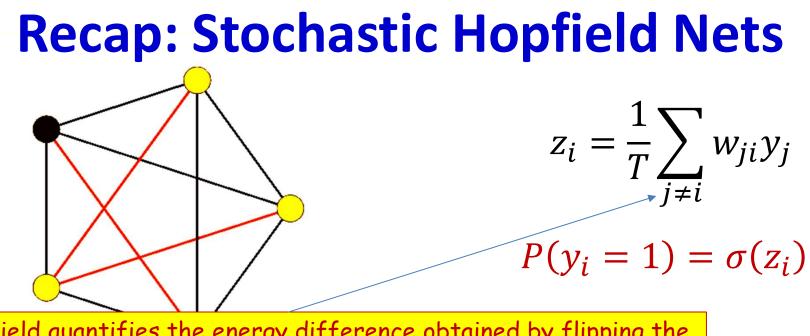
Pattern completion: Fix the "seen" bits and only let the "unseen" bits evolve

- Let the system evolve to "equilibrium"
- Let $\mathbf{y}_0, \mathbf{y}_1, \mathbf{y}_2, \dots, \mathbf{y}_L$ be the sequence of values (L large)
- Final predicted configuration: from the average of the final few iterations

$$\mathbf{y} = \left(\frac{1}{M} \sum_{t=L-M+1}^{L} \mathbf{y}_t\right) > 0?$$

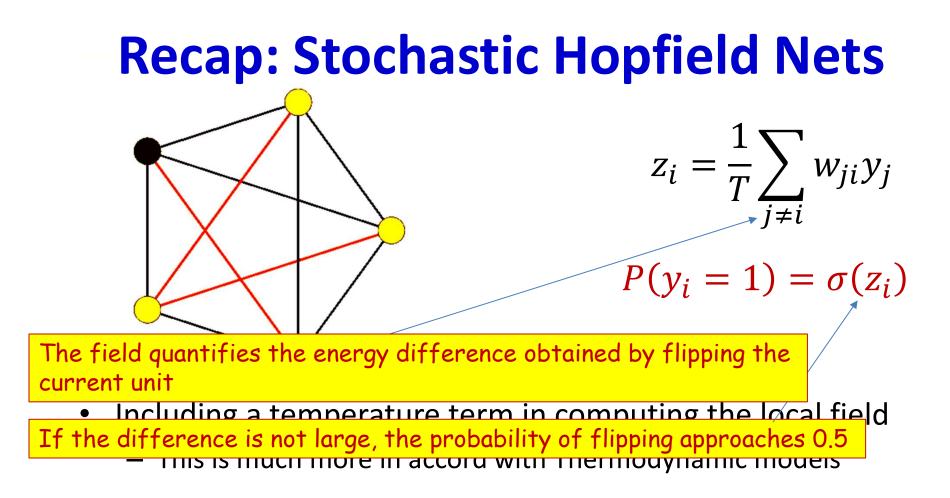
Including a "Temperature" term $z_{i} = \frac{1}{T} \sum_{j \neq i} w_{ij} y_{j}$ $P(y_{i} = 1) = \sigma(z_{i})$ $P(y_{i} = 0) = 1 - \sigma(z_{i})$

- Including a temperature term in computing the local field
 - This is much more in accord with Thermodynamic models
- At $T = \infty$ the energy "surface" will be flat. At T = 1 the surface will be the usual energy surface
 - This can be used to improve the likelihood of finding good (or optimal) minimum-energy states



The field quantifies the energy difference obtained by flipping the current unit

- Including a temperature term in computing the local field
 - This is much more in accord with Thermodynamic models
- At $T = \infty$ the energy "surface" will be flat. At T = 1 the surface will be the usual energy surface
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- At $T = \infty$ the energy "surface" will be flat. At T = 1 the surface will be the usual energy surface
 - This can be used to improve the likelihood of finding good (or optimal) minimum-energy states

Recap: Stochastic Hopfield Nets

The field quantifies the energy difference obtained by flipping the current unit

 Including a temperature term in computing the local field If the difference is not large, the probability of flipping approaches 0.5 — This is machiner maccord with mermodynamic models
 T is a "temperature" parameter: increasing it moves the probability of the bits towards 0.5
 At T=1.0 we get the traditional definition of field and energy
 At T = 0, we get deterministic Hopfield behavior

This can be used to improve the likelihood of finding good (or optimal) minimum-energy states

 $z_i = \frac{1}{T} \sum w_{ji} y_j$

 $(y_i = 1) = \sigma(z_i)$

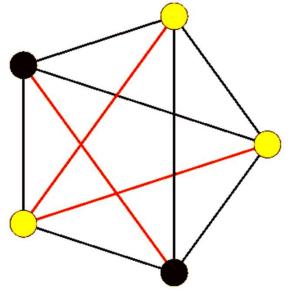
Annealing

- 1. Initialize network with initial pattern $y_i(0) = x_i, \quad 0 \le i \le N - 1$ 2. For $T = T_0$ down to T_{min} i. For iter 1..L a) For $0 \le i \le N - 1$ $P = \sigma \left(\frac{1}{T} \sum_{j \ne i} w_{ji} y_j\right)$ $y_i(t+1) \sim Binomial(P)$
- Let the system evolve to "equilibrium"
- Let $\mathbf{y}_0, \mathbf{y}_1, \mathbf{y}_2, \dots, \mathbf{y}_L$ be the sequence of values (L large)
- Final predicted configuration: from the average of the final few iterations

$$\mathbf{y} = \left(\frac{1}{M} \sum_{t=L-M+1}^{L} \mathbf{y}_t\right) > 0?$$

- 1. Initialize network with initial pattern $y_i(0) = x_i, \quad 0 \le i \le N - 1$ 2. For $T = T_0$ down to T_{min} i. For iter 1..L a) For $0 \le i \le N - 1$ $P = \sigma \left(\frac{1}{T} \sum_{j \ne i} w_{ji} y_j\right)$ $y_i(t+1) \sim Binomial(P)$
- When do we stop?
- What is the final state of the system
 - How do we "recall" a memory?

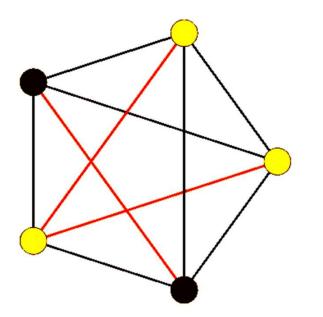
Recap: Stochastic Hopfield Nets



 $z_{i} = \frac{1}{T} \sum_{j \neq i} w_{ji} y_{j}$ $P(y_{i} = 1 | y_{j \neq i}) = \sigma(z_{i})$

- The probability of each neuron is given by a *conditional* distribution
- What is the overall probability of *the entire set of neurons* taking any configuration **y**

The overall probability



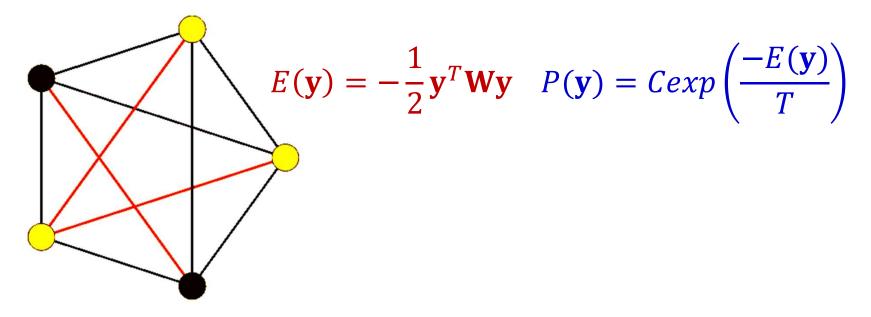
$$z_{i} = \frac{1}{T} \sum_{j \neq i} w_{ji} y_{j}$$
$$P(y_{i} = 1 | y_{j \neq i}) = \sigma(z_{i})$$

 The probability of any state y can be shown to be given by the *Boltzmann distribution*

$$E(\mathbf{y}) = -\frac{1}{2}\mathbf{y}^T \mathbf{W} \mathbf{y} \qquad P(\mathbf{y}) = Cexp\left(\frac{-E(\mathbf{y})}{T}\right)$$

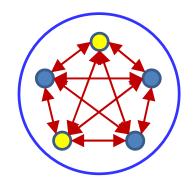
- Minimizing energy maximizes log likelihood

The overall probability



- Stop when the running average of the log probability of patterns stops increasing
 - I.e. when the (running average) of the energy of the patterns stops decreasing

The Hopfield net is a distribution



$$z_i = \frac{1}{T} \sum_j w_{ji} s_j$$

$$P(s_i = 1 | s_{j \neq i}) = \frac{1}{1 + e^{-z_i}}$$

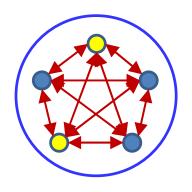
- The Hopfield net is a probability distribution over binary sequences
 - The Boltzmann distribution

$$E(\mathbf{y}) = -\frac{1}{2}\mathbf{y}^{T}\mathbf{W}\mathbf{y}$$
$$P(\mathbf{y}) = Cexp\left(-\frac{E(\mathbf{y})}{T}\right)$$

The parameter of the distribution is the weights matrix W

- The conditional distribution of individual bits in the sequence is a logistic
- We will call this a Boltzmann machine

The Boltzmann Machine



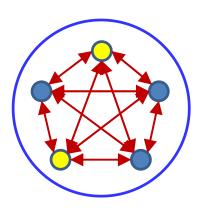
$$z_i = \frac{1}{T} \sum_j w_{ji} s_j$$

$$P(s_i = 1 | s_{j \neq i}) = \frac{1}{1 + e^{-z_i}}$$

- The entire model can be viewed as a *generative model*
- Has a probability of producing any binary vector **y**:

$$E(\mathbf{y}) = -\frac{1}{2}\mathbf{y}^{T}\mathbf{W}\mathbf{y}$$
$$P(\mathbf{y}) = Cexp\left(-\frac{E(\mathbf{y})}{T}\right)$$

Training the network



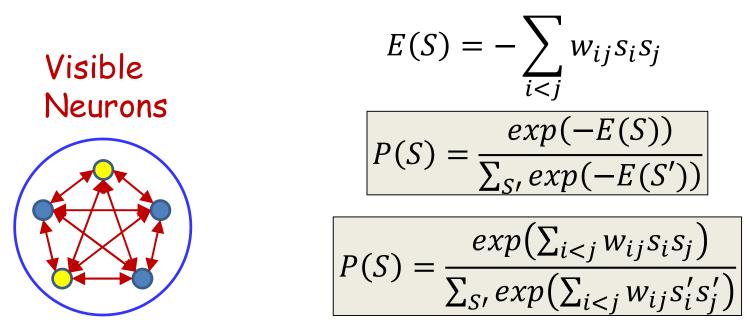
$$E(S) = -\sum_{i < j} w_{ij} s_i s_j$$

$$P(S) = \frac{exp(-E(S))}{\sum_{s, exp(-E(S'))}}$$

$$P(S) = \frac{exp(\sum_{i < j} w_{ij} s_i s_j)}{\sum_{s, exp(\sum_{i < j} w_{ij} s_i' s_j')}}$$

- Training a Hopfield net: Must learn weights to "remember" target states and "dislike" other states
 - "State" == binary pattern of all the neurons
- Training Boltzmann machine: Must learn weights to assign a desired probability distribution to states
 - (vectors y, which we will now calls S because I'm too lazy to normalize the notation)
 - This should assign more probability to patterns we "like" (or try to memorize) and less to other patterns

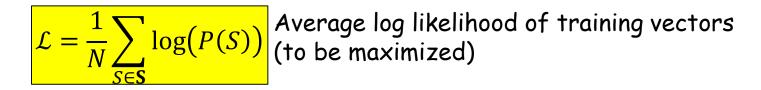
Training the network



- Must train the network to assign a desired probability distribution to states
- Given a set of "training" inputs S_1, \ldots, S_N
 - Assign higher probability to patterns seen more frequently
 - Assign lower probability to patterns that are not seen at all
- Alternately viewed: *maximize likelihood of stored states*

Maximum Likelihood Training

$$\log(P(S)) = \left(\sum_{i < j} w_{ij} s_i s_j\right) - \log\left(\sum_{S'} exp\left(\sum_{i < j} w_{ij} s'_i s'_j\right)\right)$$



$$= \frac{1}{N} \sum_{S} \left(\sum_{i < j} w_{ij} s_i s_j \right) - \log \left(\sum_{S'} exp\left(\sum_{i < j} w_{ij} s'_i s'_j \right) \right)$$

- Maximize the average log likelihood of all "training" vectors S = {S₁, S₂, ..., SN}
 - In the first summation, s_i and s_j are bits of S
 - In the second, s_i' and s_j' are bits of S'

Maximum Likelihood Training

$$\mathcal{L} = \frac{1}{N} \sum_{S} \left(\sum_{i < j} w_{ij} s_i s_j \right) - \log \left(\sum_{S'} exp\left(\sum_{i < j} w_{ij} s'_i s'_j \right) \right)$$

$$\frac{d\mathcal{L}}{dw_{ij}} = \frac{1}{N} \sum_{S} s_i s_j - ???$$

- We will use gradient ascent, but we run into a problem..
- The first term is just the average s_is_j over all training patterns
- But the second term is summed over *all* states
 - Of which there can be an exponential number!

The second term

$$\frac{d\log(\sum_{S'} exp(\sum_{i < j} w_{ij}s'_is'_j))}{dw_{ij}} = \frac{1}{\sum_{S''} exp(\sum_{i < j} w_{ij}s''_is''_j)} \frac{d\log\sum_{S'} exp(\sum_{i < j} w_{ij}s'_is'_j)}{dw_{ij}}$$

$$= \frac{1}{\sum_{S''} exp\left(\sum_{i < j} w_{ij} s_i^{"} s_j^{"}\right)} \sum_{S'} exp\left(\sum_{i < j} w_{ij} s_i^{'} s_j^{'}\right) s_i^{'} s_j^{'}}$$

$$\frac{d\log(\sum_{s'} exp(\sum_{i < j} w_{ij}s'_is'_j))}{dw_{ij}} = \sum_{s'} \frac{exp(\sum_{i < j} w_{ij}s'_is'_j)}{\sum_{s''} exp(\sum_{i < j} w_{ij}s''_is''_j)} s'_is'_j$$

The second term

$$\frac{d\log(\sum_{S'} exp(\sum_{i < j} w_{ij}s'_is'_j))}{dw_{ij}} = \frac{1}{\sum_{S''} exp(\sum_{i < j} w_{ij}s''_is''_j)} \frac{d\log\sum_{S'} exp(\sum_{i < j} w_{ij}s'_is'_j)}{dw_{ij}}$$

$$\frac{1}{\sum_{S''} exp(\sum_{i < j} w_{ij}s_i^{"}s_j^{"})} \sum_{S'} exp\left(\sum_{i < j} w_{ij}s_i's_j'\right) s_i's_j'}{\frac{d\log(\sum_{S'} exp(\sum_{i < j} w_{ij}s_i's_j'))}{dw_{ij}}} = \sum_{S'} \frac{exp(\sum_{i < j} w_{ij}s_i's_j')}{\sum_{S''} exp(\sum_{i < j} w_{ij}s_i^{"}s_j^{"})} s_i's_j'}$$

The second term

$$\frac{d\log(\sum_{S'} exp(\sum_{i < j} w_{ij}s'_is'_j))}{dw_{ij}} = \frac{1}{\sum_{S''} exp(\sum_{i < j} w_{ij}s''_is''_j)} \frac{d\log\sum_{S'} exp(\sum_{i < j} w_{ij}s'_is'_j)}{dw_{ij}}$$

$$= \frac{1}{\sum_{S''} exp\left(\sum_{i < j} w_{ij} s_i^{"} s_j^{"}\right)} \sum_{S'} exp\left(\sum_{i < j} w_{ij} s_i^{'} s_j^{'}\right) s_i^{'} s_j^{'}}$$

$$\frac{d\log(\sum_{s,exp}(\sum_{i$$

$$\frac{d\log(\sum_{S'} exp(\sum_{i < j} w_{ij}s'_is'_j))}{dw_{ij}} = \sum_{S'} P(S')s'_is'_j$$

$$\frac{d\log(\sum_{s'} exp(\sum_{i < j} w_{ij}s'_{i}s'_{j}))}{dw_{ij}} = \sum_{s'} P(S')s'_{i}s'_{j}$$

- The second term is simply the *expected value* of s_is_j, over all possible values of the state
- We cannot compute it exhaustively, but we can compute it by sampling!

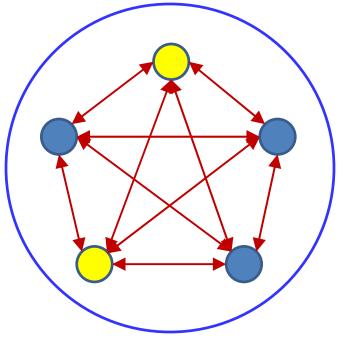
Estimating the second term

$$\frac{d\log(\sum_{S'} exp(\sum_{i < j} w_{ij}s'_is'_j))}{dw_{ij}} = \sum_{S'} P(S')s'_is'_j$$

$$\sum_{S'} P(S') s'_i s'_j \approx \frac{1}{M} \sum_{S' \in \mathbf{S}_{samples}} s'_i s'_j$$

- The expectation can be estimated as the average of samples drawn from the distribution
- Question: How do we draw samples from the Boltzmann distribution?
 - How do we draw samples from the network?

The simulation solution



- Initialize the network randomly and let it "evolve"
 - By probabilistically selecting state values according to our model
- After many many epochs, take a snapshot of the state
- Repeat this many many times
- Let the collection of states be

$$\mathbf{S}_{simul} = \{S_{simul,1}, S_{simul,1=2}, \dots, S_{simul,M}\}$$

The simulation solution for the second term

$$\frac{d\log(\sum_{S'} exp(\sum_{i < j} w_{ij}s'_is'_j))}{dw_{ij}} = \sum_{S'} P(S')s'_is'_j$$

$$\sum_{S'} P(S') s'_i s'_j \approx \frac{1}{M} \sum_{S' \in \mathbf{S}_{simul}} s'_i s'_j$$

 The second term in the derivative is computed as the average of sampled states when the network is running "freely"

Maximum Likelihood Training

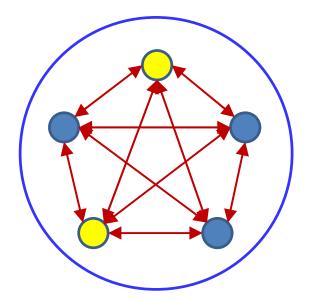
Sampled estimate

$$\frac{d\langle \log(P(\mathbf{S}))\rangle}{dw_{ij}} = \frac{1}{N} \sum_{S} s_i s_j - \frac{1}{M} \sum_{S' \in \mathbf{S}_{simul}} s'_i s'_j$$

$$w_{ij} = w_{ij} + \eta \frac{d\langle \log(P(\mathbf{S})) \rangle}{dw_{ij}}$$

• The overall gradient ascent rule

Overall Training

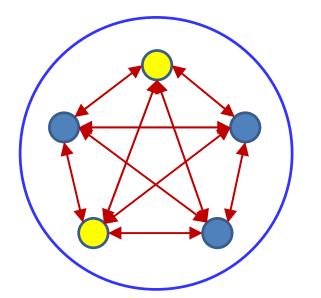


$$\frac{d\langle \log(P(\mathbf{S})) \rangle}{dw_{ij}} = \frac{1}{N} \sum_{S} s_i s_j - \frac{1}{M} \sum_{S' \in \mathbf{S}_{simul}} s'_i s'_j$$

$$w_{ij} = w_{ij} + \eta \frac{d\langle \log(P(\mathbf{S})) \rangle}{dw_{ij}}$$

- Initialize weights
- Let the network run to obtain simulated state samples
- Compute gradient and update weights
- Iterate

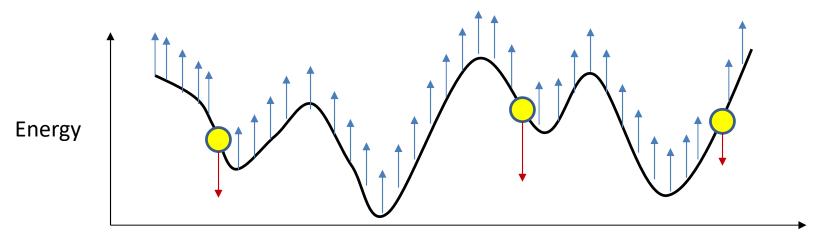
Overall Training



$$\frac{d\langle \log(P(\mathbf{S}))\rangle}{dw_{ij}} = \frac{1}{N} \sum_{S} s_i s_j - \frac{1}{M} \sum_{S' \in \mathbf{S}_{simul}} s'_i s'_j$$

$$w_{ij} = w_{ij} + \eta \frac{d\langle \log(P(\mathbf{S})) \rangle}{dw_{ij}}$$

Note the similarity to the update rule for the Hopfield network

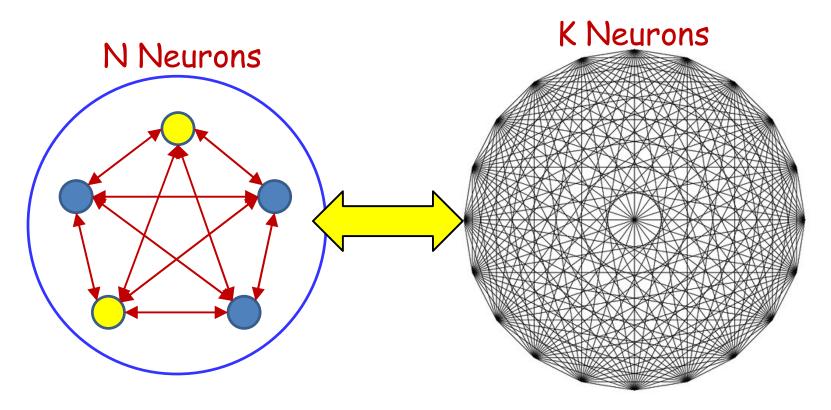




Adding Capacity to the Hopfield Network / Boltzmann Machine

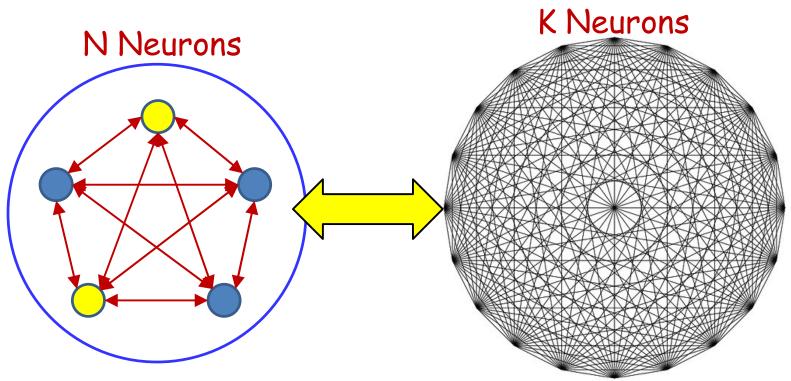
- The network can store up to *N N*-bit patterns
- How do we increase the capacity

Expanding the network

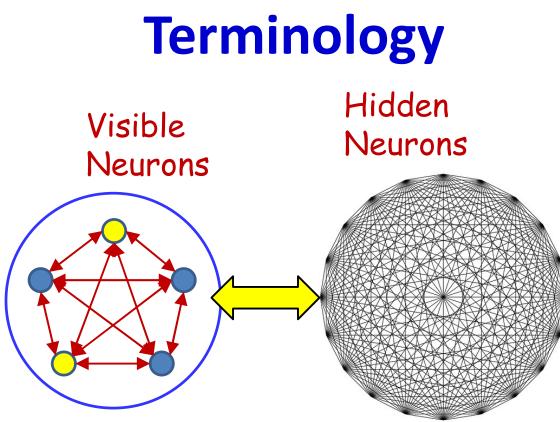


 Add a large number of neurons whose actual values you don't care about!

Expanded Network

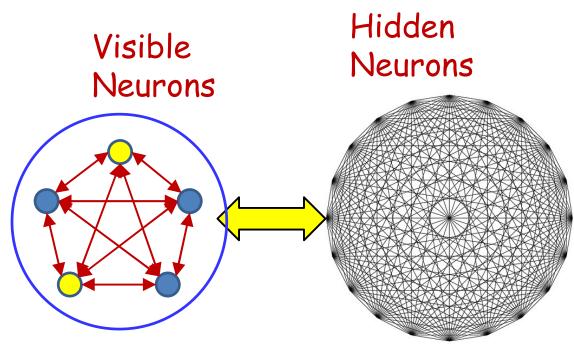


- New capacity: $\sim (N + K)$ patterns
 - Although we only care about the pattern of the first N neurons
 - We're interested in *N-bit* patterns



- Terminology:
 - The neurons that store the actual patterns of interest: Visible neurons
 - The neurons that only serve to increase the capacity but whose actual values are not important: *Hidden neurons*
 - These can be set to anything in order to store a visible pattern

Training the network

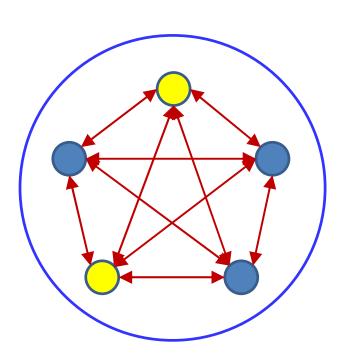


- For a given pattern of *visible* neurons, there are any number of *hidden* patterns (2^K)
- Which of these do we choose?
 - Ideally choose the one that results in the lowest energy
 - But that's an exponential search space!

The patterns

- In fact we could have *multiple* hidden patterns coupled with any visible pattern
 - These would be multiple stored patterns that all give the same visible output
 - How many do we permit
- Do we need to specify one or more particular hidden patterns?
 - How about *all* of them
 - What do I mean by this bizarre statement?

Boltzmann machine without hidden

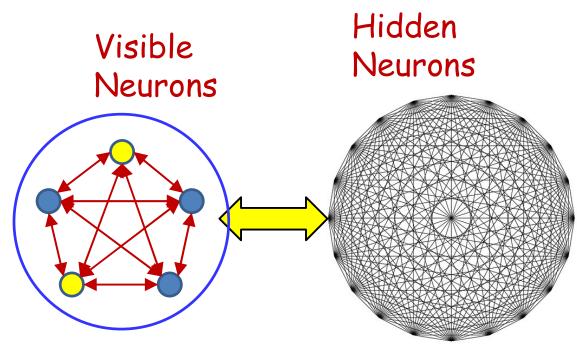


$$\frac{d\langle \log(P(\mathbf{S})) \rangle}{dw_{ij}} = \frac{1}{N} \sum_{S} s_i s_j - \frac{1}{M} \sum_{S' \in \mathbf{S}_{simul}} s'_i s'_j$$

$$w_{ij} = w_{ij} + \eta \frac{d\langle \log(P(\mathbf{S})) \rangle}{dw_{ij}}$$

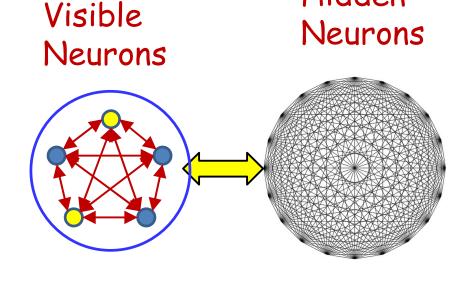
- This basic framework has no hidden units
- Extended to have hidden units

With hidden neurons



- Now, with hidden neurons the complete state pattern for even the *training* patterns is unknown
 - Since they are only defined over visible neurons

With hidden neurons Hidden



$$P(S) = \frac{exp(-E(S))}{\sum_{S'} exp(-E(S'))}$$

$$P(S) = P(V, H)$$

$$P(V) = \sum_{H} P(S)$$

- We are interested in the *marginal* probabilities over *visible* bits
 - We want to learn to represent the visible bits
 - The hidden bits are the "latent" representation learned by the network
- S = (V, H)
 - V = visible bits
 - H = hidden bits

With hidden neurons Hidden Neurons

We are interested in the marginal probabilities over visible bits We want to learn to represent the visible bits We want to learn to represent the visible bits The hidden bits are the "latent" representation learned by the network S = (V, H) V = visible bits H = hidden bits

$$P(S) = \frac{exp(-E(S))}{\sum_{S'} exp(-E(S'))}$$

$$P(S) = P(V, H)$$

$$P(V) = \sum_{H} P(S)$$

- We are interested in the *marginal* probabilities over *visible* bits
 - We want to learn to represent the visible bits
 - The hidden bits are the "latent" representation learned by the network
- S = (V, H)

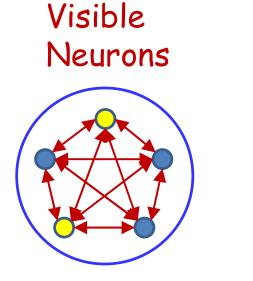
Visible

Neurons

- V = visible bits
- H = hidden bits

Must train to maximize probability of desired patterns of *visible* bits

Training the network



$$E(S) = -\sum_{i < j} w_{ij} s_i s_j$$

$$P(S) = \frac{exp(\sum_{i < j} w_{ij} s_i s_j)}{\sum_{S'} exp(\sum_{i < j} w_{ij} s'_i s'_j)}$$

$$P(V) = \sum_{H} \frac{exp(\sum_{i < j} w_{ij} s_i s_j)}{\sum_{S'} exp(\sum_{i < j} w_{ij} s'_i s'_j)}$$

- Must train the network to assign a desired probability distribution to visible states
- Probability of visible state sums over all hidden states

Maximum Likelihood Training

$$\log(P(V)) = \log\left(\sum_{H} exp\left(\sum_{i < j} w_{ij}s_is_j\right)\right) - \log\left(\sum_{S'} exp\left(\sum_{i < j} w_{ij}s'_is'_j\right)\right)$$

 $\mathcal{L} = \frac{1}{N} \sum_{V \in \mathbf{V}} \log(P(V))$ (to be maximized)

$$= \frac{1}{N} \sum_{V \in \mathbf{V}} \log \left(\sum_{H} exp\left(\sum_{i < j} w_{ij} s_i s_j \right) \right) - \log \left(\sum_{S'} exp\left(\sum_{i < j} w_{ij} s'_i s'_j \right) \right)$$

- Maximize the average log likelihood of all visible bits of "training" vectors V = {V₁, V₂, ..., V_N}
 - The first term also has the same format as the second term
 - Log of a sum
 - Derivatives of the first term will have the same form as for the second term

Maximum Likelihood Training

$$\mathcal{L} = \frac{1}{N} \sum_{V \in \mathbf{V}} \log \left(\sum_{H} exp\left(\sum_{i < j} w_{ij} s_i s_j \right) \right) - \log \left(\sum_{S'} exp\left(\sum_{i < j} w_{ij} s_i' s_j' \right) \right)$$

$$\frac{d\mathcal{L}}{dw_{ij}} = \frac{1}{N} \sum_{V \in \mathbf{V}} \sum_{H} \frac{exp(\sum_{k < l} w_{kl} s_k s_l)}{\sum_{H'} exp(\sum_{k < l} w_{kl} s_k^{"} s_l^{"})} s_i s_j - \sum_{S'} \frac{exp(\sum_{k < l} w_{kl} s_k^{'} s_l^{'})}{\sum_{S''} exp(\sum_{k < l} w_{ij} s_k^{"} s_l^{"})} s_i^{'} s_j^{'}$$

$$\frac{d\mathcal{L}}{dw_{ij}} = \frac{1}{N} \sum_{V \in \mathbf{V}} \sum_{H} P(S|V) s_i s_j - \sum_{S'} P(S') s'_i s'_j$$

- We've derived this math earlier
- But now *both* terms require summing over an exponential number of states
 - The first term fixes visible bits, and sums over all configurations of hidden states for each visible configuration in our training set
 - But the second term is summed over all states

The simulation solution

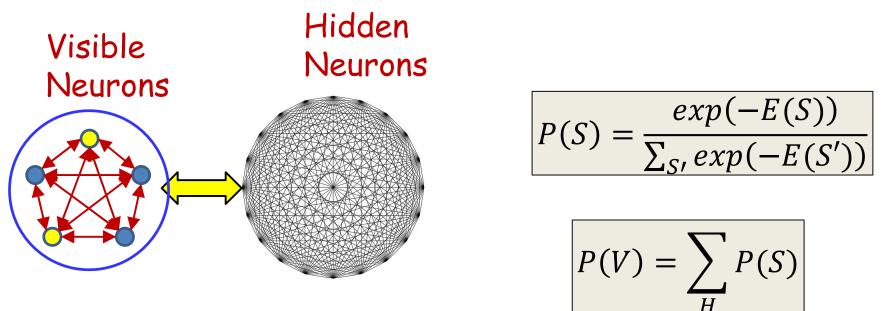
$$\frac{d\mathcal{L}}{dw_{ij}} = \frac{1}{N} \sum_{V \in \mathbf{V}} \sum_{H} P(S|V) s_i s_j - \sum_{S'} P(S') s'_i s'_j$$

$$\sum_{H} P(S|V) s_i s_j \approx \frac{1}{K} \sum_{H \in \mathbf{H}_{simul}} s_i s_j$$

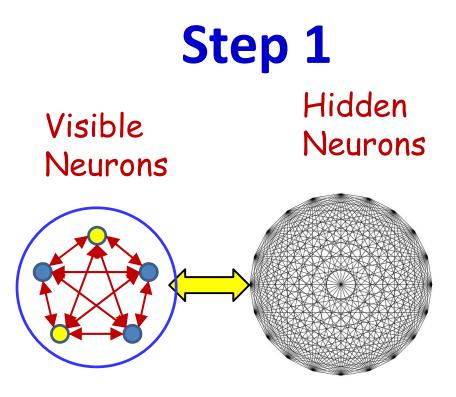
$$\sum_{S'} P(S') s'_i s'_j \approx \frac{1}{M} \sum_{S' \in \mathbf{S}_{simul}} s'_i s'_j$$

- The first term is computed as the average sampled *hidden* state with the visible bits fixed
- The second term in the derivative is computed as the average of sampled states when the network is running "freely"

More simulations



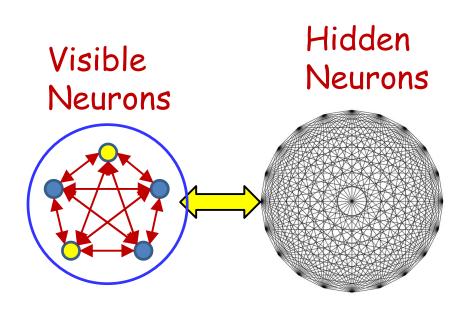
- Maximizing the marginal probability of V requires summing over all values of H
 - An exponential state space
 - So we will use simulations again



- For each training pattern V_i
 - Fix the visible units to V_i
 - Let the hidden neurons evolve from a random initial point to generate H_i
 - Generate $S_i = [V_i, H_i]$
- Repeat K times to generate synthetic training

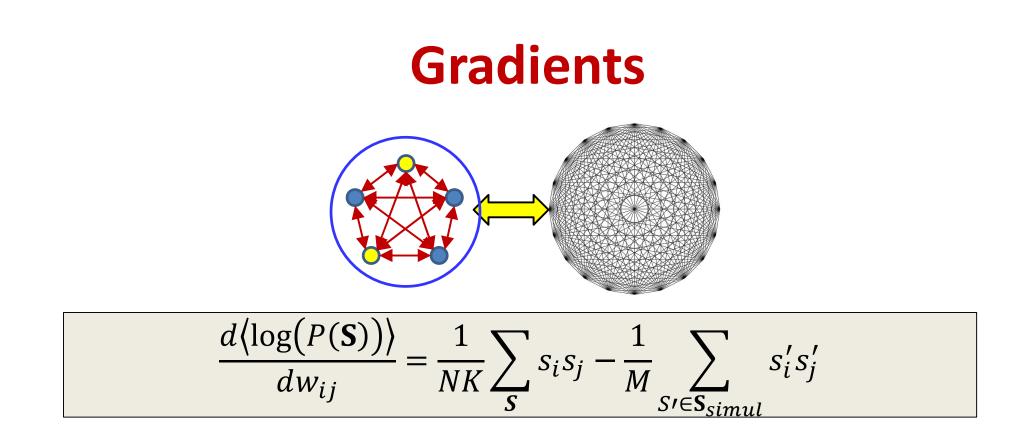
$$\mathbf{S} = \{S_{1,1}, S_{1,2}, \dots, S_{1K}, S_{2,1}, \dots, S_{N,K}\}$$

Step 2



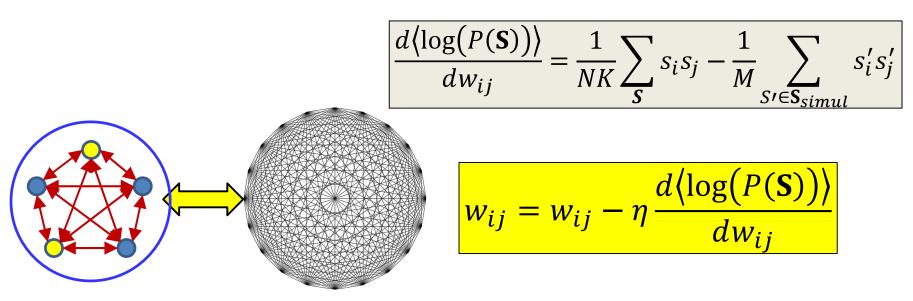
 Now unclamp the visible units and let the entire network evolve several times to generate

$$\mathbf{S}_{simul} = \{S_{simul,1}, S_{simul,1=2}, \dots, S_{simul,M}\}$$



 Gradients are computed as before, except that the first term is now computed over the *expanded* training data

Overall Training



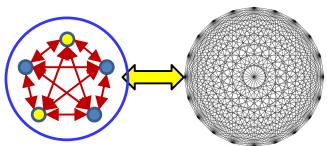
- Initialize weights
- Run simulations to get clamped and unclamped training samples
- Compute gradient and update weights
- Iterate

Boltzmann machines

- Stochastic extension of Hopfield nets
- Enables storage of many more patterns than Hopfield nets
- But also enables computation of probabilities of patterns, and completion of pattern

Boltzmann machines: Overall

$$z_i = \sum_j w_{ji} s_i + b_i$$
$$P(s_i = 1) = \frac{1}{1 + e^{-Z_i}}$$



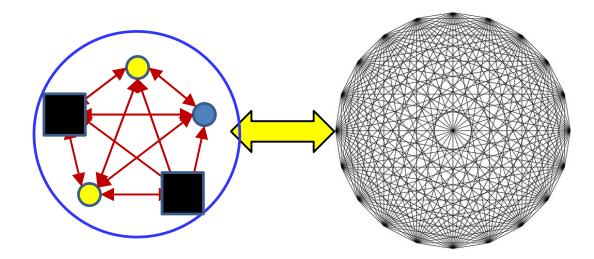
$$\frac{\langle \log(P(\mathbf{S})) \rangle}{dw_{ij}} = \frac{1}{NK} \sum_{\mathbf{S}} s_i s_j - \frac{1}{M} \sum_{S' \in \mathbf{S}_{simul}} s'_i s'_j$$
$$w_{ij} = w_{ij} - \eta \frac{d\langle \log(P(\mathbf{S})) \rangle}{dw_{ij}}$$

- Training: Given a set of training patterns
 - Which could be repeated to represent relative probabilities

d

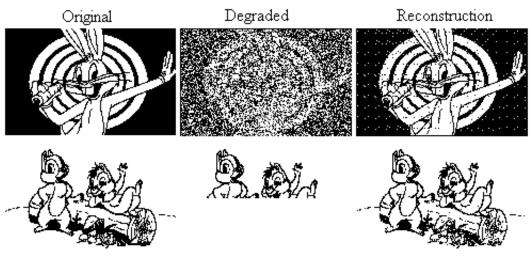
- Initialize weights
- Run simulations to get clamped and unclamped training samples
- Compute gradient and update weights
- Iterate

Boltzmann machines: Overall



- Running: Pattern completion
 - "Anchor" the *known* visible units
 - Let the network evolve
 - Sample the unknown visible units
 - Choose the most probable value

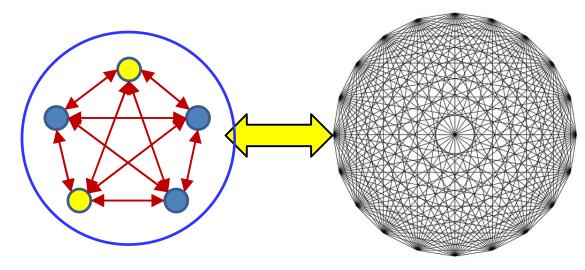




Hopfield network reconstructing degraded images from noisy (top) or partial (bottom) cues.

- Filling out patterns
- Denoising patterns
- Computing conditional probabilities of patterns
- Classification!!
 - How?

Boltzmann machines for classification

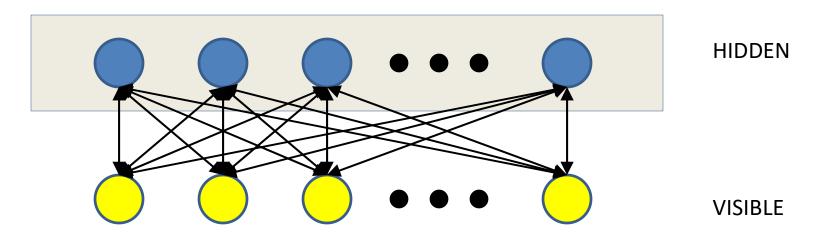


- Training patterns:
 - $[f_1, f_2, f_3,, class]$
 - Features can have binarized or continuous valued representations
 - Classes have "one hot" representation
- Classification:
 - Given features, anchor features, estimate a posteriori probability distribution over classes
 - Or choose most likely class

Boltzmann machines: Issues

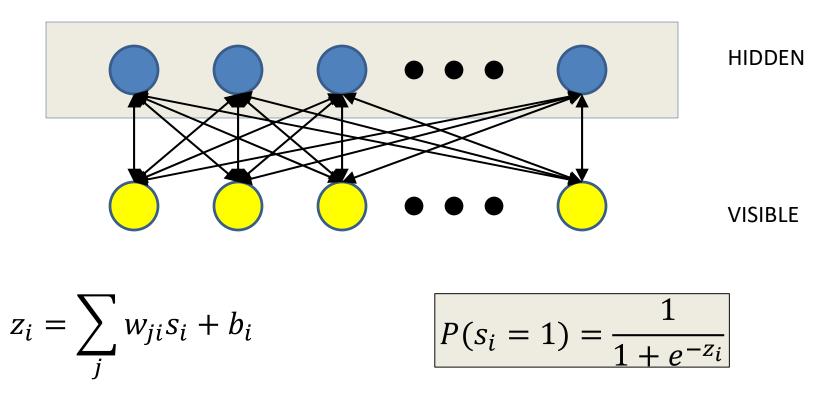
- Training takes for ever
- Doesn't really work for large problems
 - A small number of training instances over a small number of bits

Solution: *Restricted* Boltzmann Machines



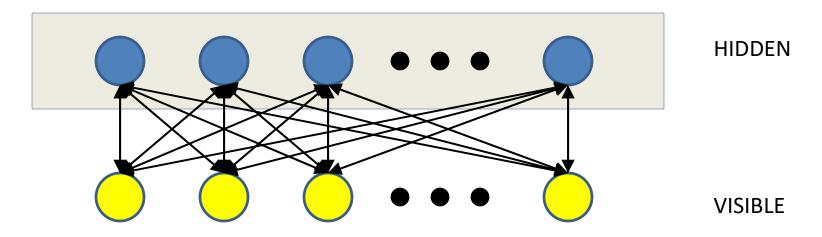
- Partition visible and hidden units
 - Visible units ONLY talk to hidden units
 - Hidden units ONLY talk to visible units
- Restricted Boltzmann machine..
 - Originally proposed as "Harmonium Models" by Paul Smolensky

Solution: *Restricted* Boltzmann Machines



- Still obeys the same rules as a regular Boltzmann machine
- But the modified structure adds a big benefit..

Solution: *Restricted* Boltzmann Machines



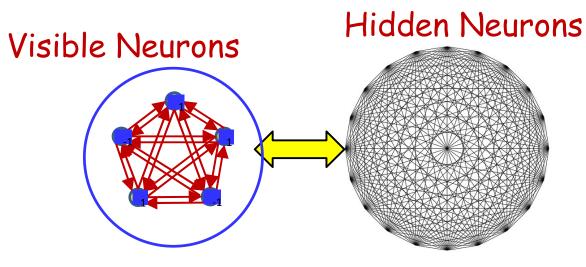
$$z_i = \sum_j w_{ji} v_i + b_i \qquad P(h_i = 1) = \frac{1}{1 + e^{-z_i}}$$

VISIBLE

HIDDEN

$$y_i = \sum_j w_{ji}h_i + b_i$$
 $P(v_i = 1) = \frac{1}{1 + e^{-y_i}}$

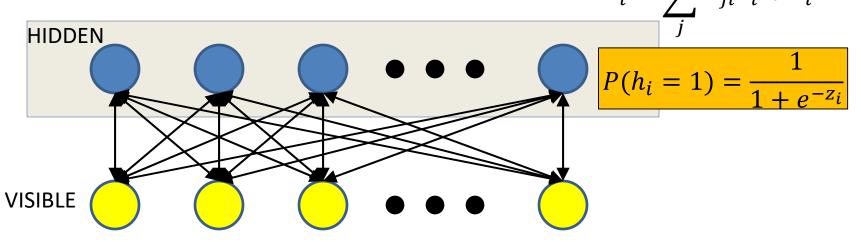
Recap: Training full Boltzmann machines: Step 1



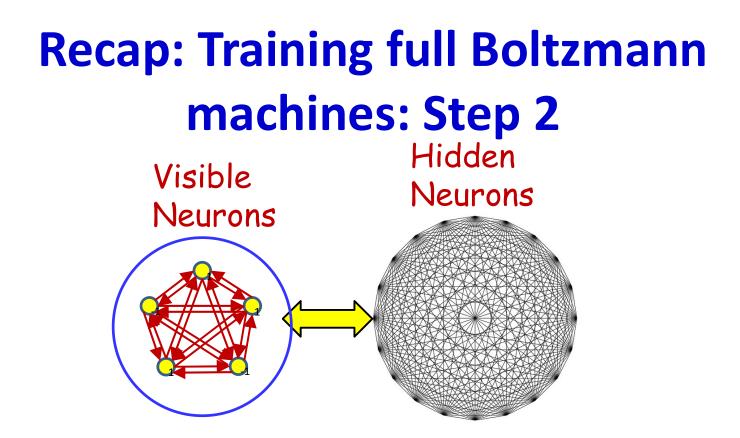
- For each training pattern V_i
 - Fix the visible units to V_i
 - Let the hidden neurons evolve from a random initial point to generate H_i
 - Generate $S_i = [V_i, H_i]$
- Repeat K times to generate synthetic training

$$\mathbf{S} = \{S_{1,1}, S_{1,2}, \dots, S_{1K}, S_{2,1}, \dots, S_{N,K}\}$$

Sampling: Restricted Boltzmann machine $z_i = \sum w_{ji}v_i + b_i$



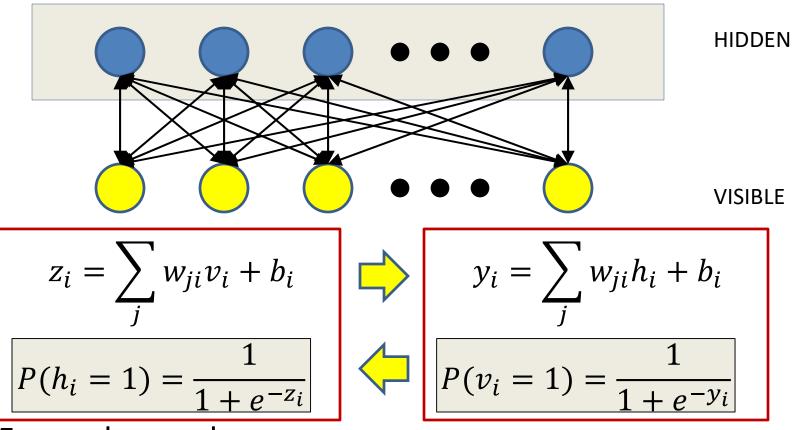
- For each sample:
 - Anchor visible units
 - Sample from hidden units
 - No looping!!



 Now unclamp the visible units and let the entire network evolve several times to generate

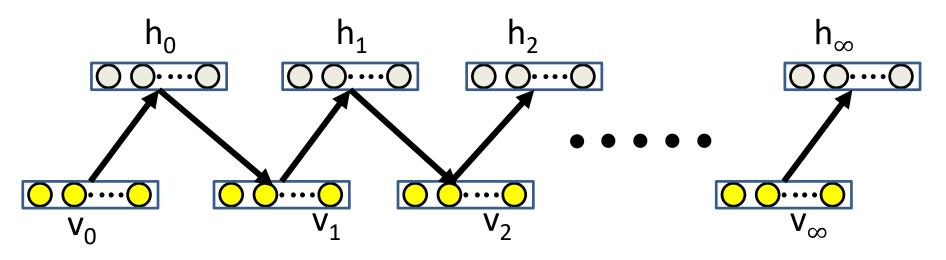
$$\mathbf{S}_{simul} = \{S_{simul,1}, S_{simul,1=2}, \dots, S_{simul,M}\}$$

Sampling: Restricted Boltzmann machine



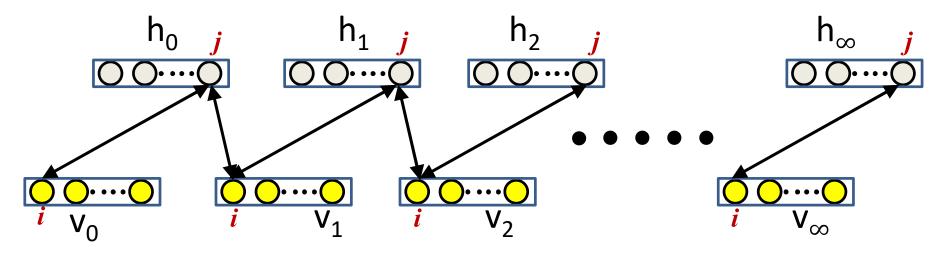
- For each sample:
 - Iteratively sample hidden and visible units for a long time
 - Draw final sample of both hidden and visible units

Pictorial representation of RBM training



- For each sample:
 - Initialize V_0 (visible) to training instance value
 - Iteratively generate hidden and visible units
 - For a very long time

Pictorial representation of RBM training



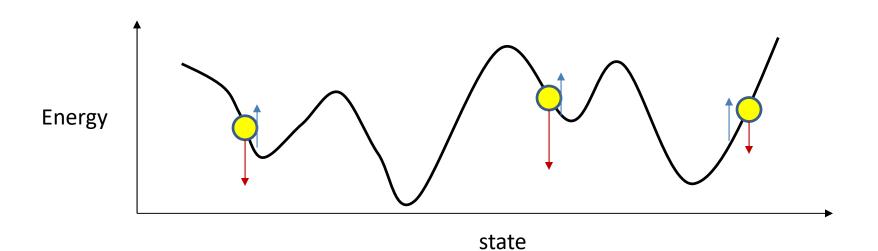
Gradient (showing only one edge from visible node *i* to hidden node *j*)

$$\frac{\partial \log p(v)}{\partial w_{ij}} = \langle v_i h_j \rangle^0 - \langle v_i h_j \rangle^\infty$$

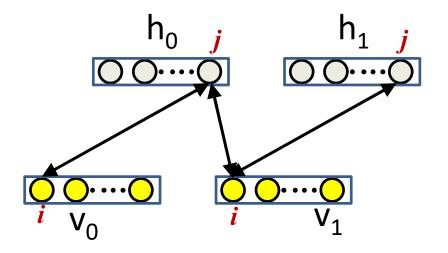
<*v_i*, *h_j*> represents average over many generated training samples

Recall: Hopfield Networks

- Really no need to raise the entire surface, or even every valley
- Raise the *neighborhood* of each target memory
 - Sufficient to make the memory a valley
 - The broader the neighborhood considered, the broader the valley



A Shortcut: Contrastive Divergence

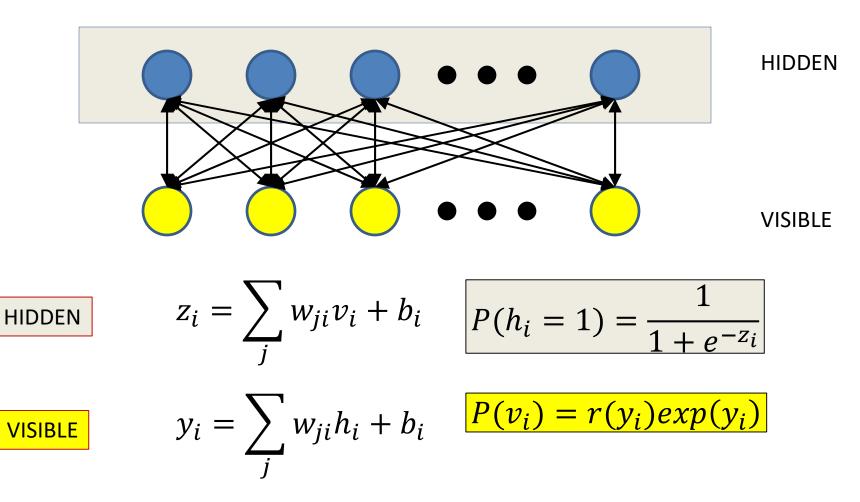


- Sufficient to run one iteration! $\frac{\partial \log p(v)}{\partial w_{ii}} = \langle v_i h_j \rangle^0 - \langle v_i h_j \rangle^1$
- This is sufficient to give you a good estimate of the gradient

Restricted Boltzmann Machines

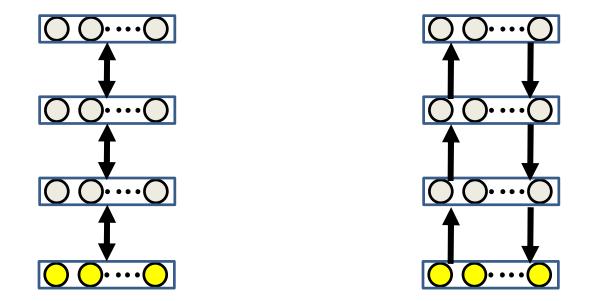
- Excellent generative models for binary (or binarized) data
- Can also be extended to continuous-valued data
 - "Exponential Family Harmoniums with an Application to Information Retrieval", Welling et al., 2004
- Useful for classification and regression
 - How?
 - More commonly used to *pretrain* models

Continuous-values RBMs



Hidden units may also be continuous values

Other variants



- Left: "Deep" Boltzmann machines
- Right: Helmholtz machine
 - Trained by the "wake-sleep" algorithm

Topics missed..

- Other algorithms for Learning and Inference over RBMs
 - Mean field approximations
- RBMs as feature extractors

Pre training

- RBMs as generative models
- More structured DBMs
- ..