

Neural Nets and Symbolic Reasoning

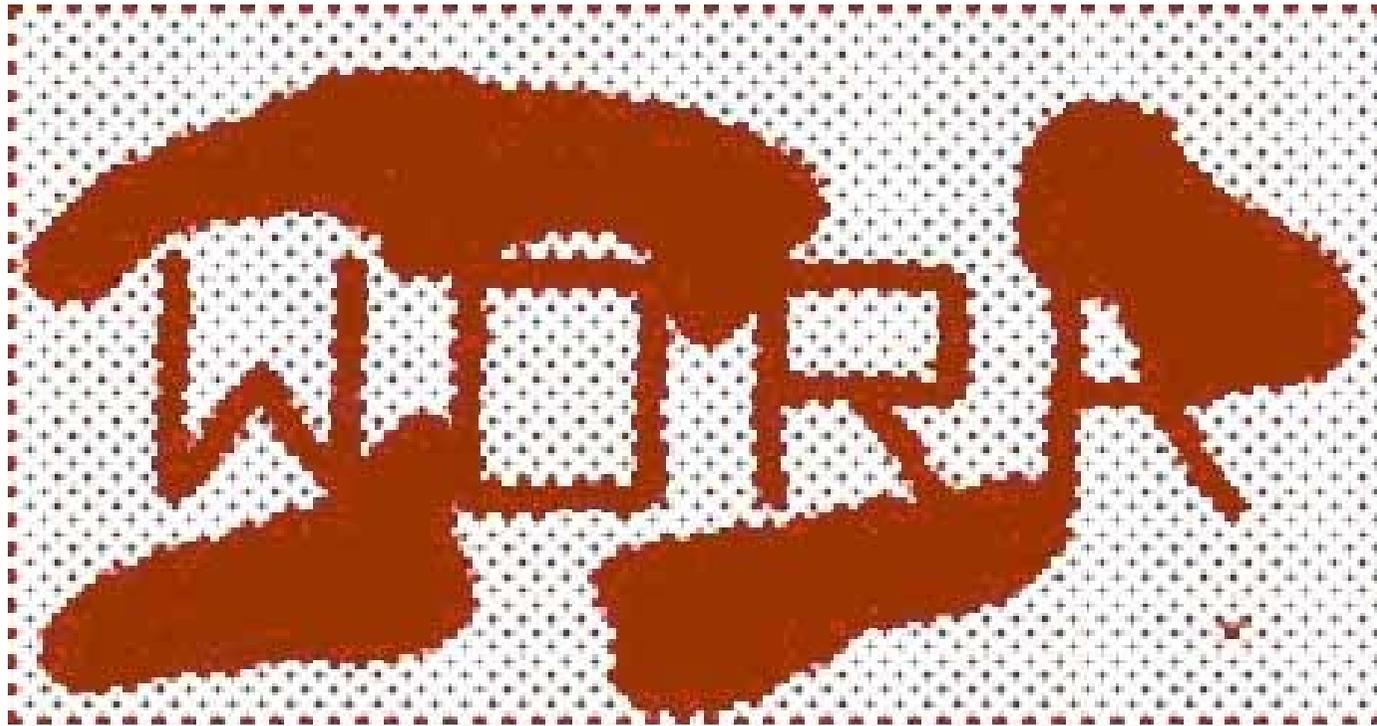
Hopfield Networks

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Outline

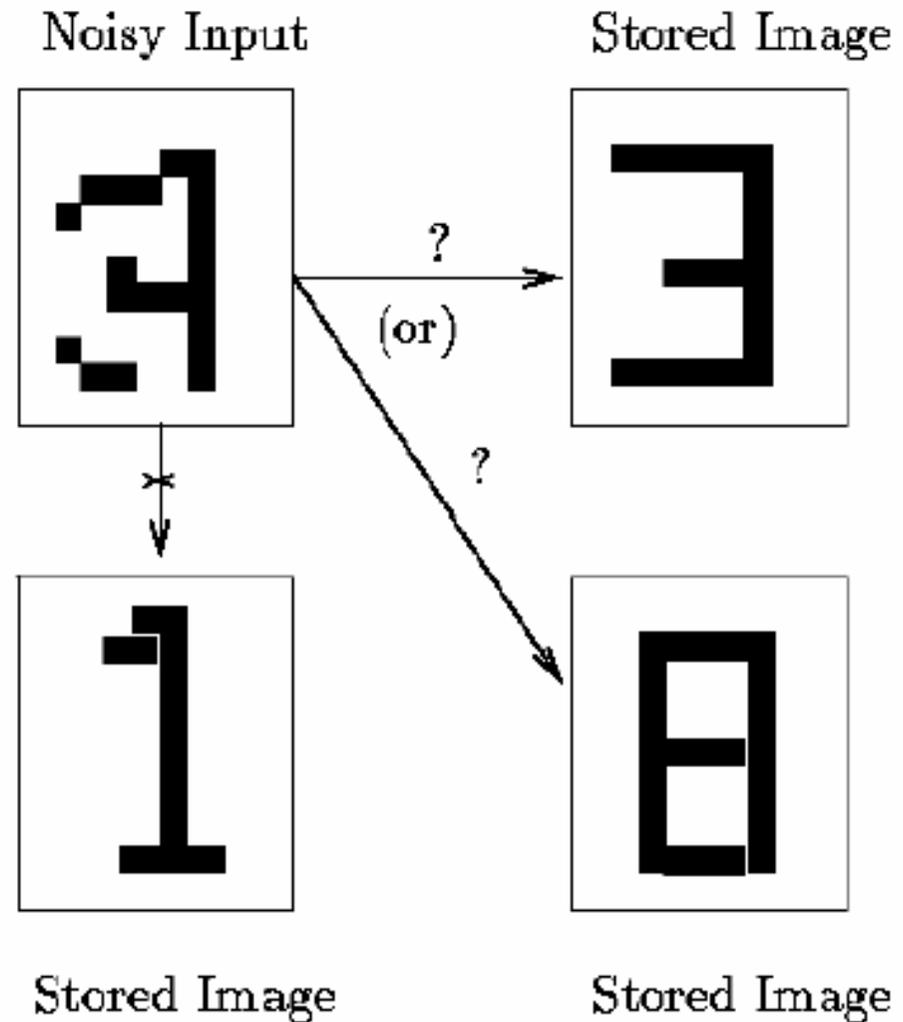
-  The idea of pattern completion
-  The fast dynamics of Hopfield networks
-  Learning with Hopfield networks
-  Emerging properties of Hopfield networks
-  Conclusions

1 The idea of pattern completion



Example from visual recognition

- Noisy or underspecified input
- Mechanism of pattern completion (using *stored* images)
- Stored patterns are addressable by *content*, not *pointers* (as in traditional computer memories)



Example from semantics: *A fast car*

- a. *a fast car* [one that moves quickly]
- b. *a fast typist* [a person that performs the act of typing quickly]
- c. *a fast book* [one that can be read in a short time]
- d. *a fast driver* [one who drives quickly]

Example from semantics: *A red apple*



A red apple?

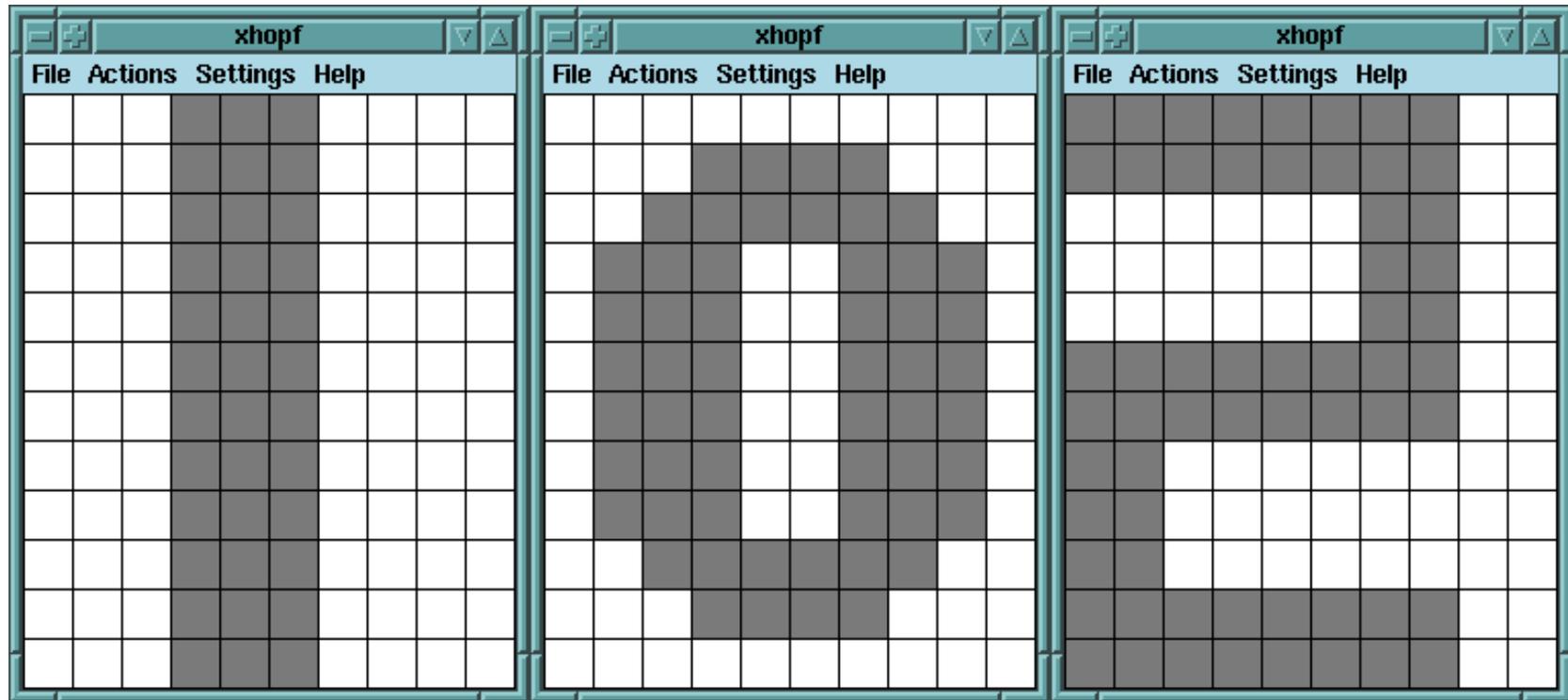
What color is an apple?

Q₁ What color is its peel?

Q₂ What color is its pulp?

- | | | |
|----|-----------------------------|------------------|
| a. | a red apple | [red peel] |
| b. | a sweet apple | [sweet pulp] |
| c. | a reddish grapefruit | [reddish pulp] |
| d. | a white room/ a white house | [inside/outside] |

2 The fast dynamics of Hopfield networks



Abstract description of neural networks

A neural network N can be defined as a quadruple $\langle S, F, W, G \rangle$:

- S : Space of all possible states
- W : Set of possible configurations. $w \in W$ describes for each pair i, j of "neurons" the connection w_{ij} between i and j
- F : Set of activation functions. For a given configuration $w \in W$: the function $f_w \in F$ describes how the neuron activities spread through that network (fast dynamics)
- G : Set of learning functions (slow dynamics)

Discrete dynamic systems

$$\vec{s}(t+1) = g(\vec{s}(t)) \quad \vec{s}(t) \text{ is a vector of the space of states } S$$

Continuous dynamic systems

$$\frac{d}{dt} \vec{s}(t) = g(\vec{s}(t)) \quad \text{the function } g \text{ describes the fast dynamics}$$

Dynamical Systems + Neural Networks = Neurodynamics

- Tools from dynamical systems, statistics and statistical physics can be used. Very rich field.
- The triple $\langle S, F, W \rangle$ corresponds to the *fast* neurodynamics

The importance of recurrent systems

What can recurrent networks do?

- Associative memories
- Pattern completion
- Noise removal
- General networks (can implement everything feedforward networks can do, and even emulate Turing machines)
- Spatio-temporal pattern recognition
- Dynamic reconstruction

J.J.Hopfield (1982), "Neural networks and physical systems with emergent collective computational abilities", *Proceedings of the National Academy of Sciences* 79, 2554-2558.

An autoassociative, fully connected network with binary neurons, asynchronous updates and a Hebbian learning rule. The “classic” recurrent network

Computational properties of use to biological organisms or to the construction of computers can emerge as collective properties of systems having a large number of simple equivalent components (or neurons). The physical meaning of content-addressable memory is described by an appropriate phase space flow of the state of the system. ...

Concise description of the fast dynamics

Let the interval $[-1,+1]$ be the *working range* of each neuron

+1: maximal firing rate

0: resting

-1 : minimal firing rate)

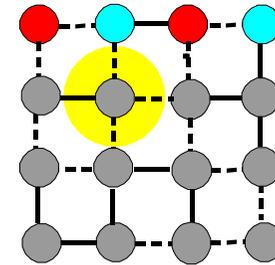
$$S = [-1, 1]^n$$

$$W_{ij} = W_{ji}, W_{ii} = 0$$

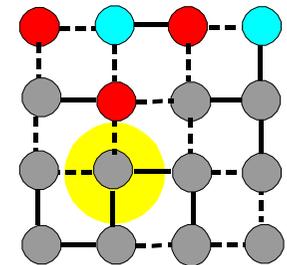
ASYNCHRONOUS UPDATING:

$$s_i(t+1) = \begin{cases} \theta(\sum_j w_{ij} \cdot s_j(t)), & \text{if } i = \text{rand}(1,n) \\ s_i(t), & \text{otherwise} \end{cases}$$

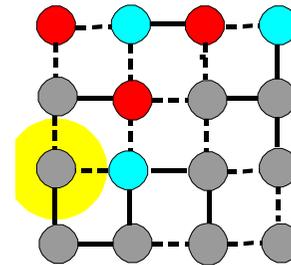
Step 1



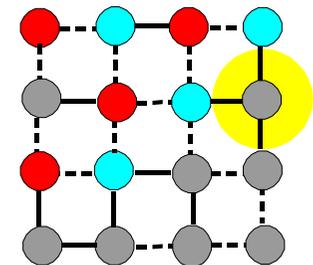
Step 2



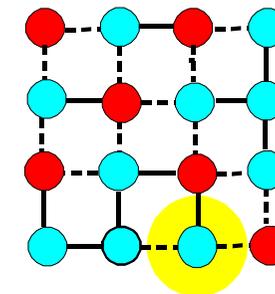
Step 3



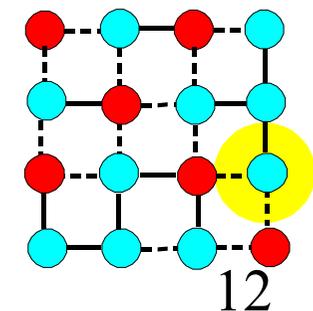
Step 4



Step 98651



Step 98652



Definition of resonances of a dynamical system

A state s in S is called a resonance of a dynamic system $[S, f]$ iff

1. $f(s) = s$ (**equilibrium**)

2. For each $\varepsilon > 0$ there exists a

$0 < \delta \leq \varepsilon$ such that for all $n \geq 1$

$|f_n(s') - s| < \varepsilon$ whenever $|s' - s| < \delta$

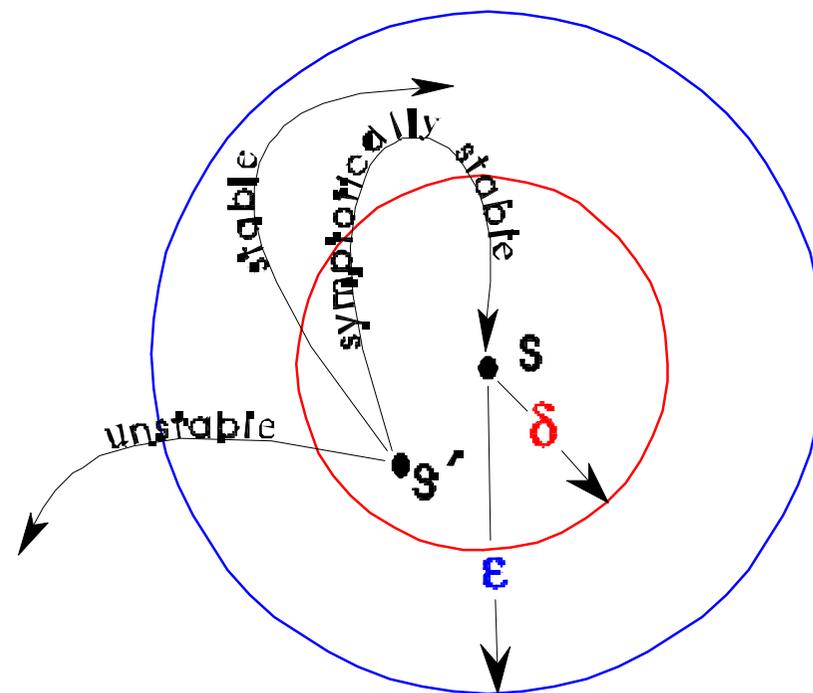
(**stability**)

3. For each $\varepsilon > 0$ there exists a

$0 < \delta \leq \varepsilon$ such that $\lim_{n \rightarrow \infty} f^n(s') = s$

whenever $|s' - s| < \delta$

(**asymptotic stability**)



Energy and convergence

- How can we be sure the dynamics converges to any attractor? Why cannot it enter an endless loop $s^1 \rightarrow s^2 \rightarrow s^3 \rightarrow s^1 \rightarrow \dots$?
- Answer (and the secret of the Hopfield network's popularity): the energy function (also called *Lyapunov* function).
- An energy function (Lyapunov function) always decreases monotonically as we change state and is bounded below. The descent to lower energy levels will have to end eventually at a local minimum.
- Energy landscapes are a popular (and somewhat dangerous) analogy.



Definition A neural network $[S,W,F]$ is called a resonance system iff $\lim_{n \rightarrow \infty} f^n(s)$ exists and is a resonance for each $s \in S$ and $f \in F$.

Theorem 1 (Cohen & Grossberg 1983)

Hopfield networks are resonance systems.

(The same holds for a large class of other systems: The McCulloch-Pitts model (1943), Cohen-Grossberg models (1983), Rumelhart's Interactive Activation model (1986), Smolensky's Harmony networks (1986), etc.)

Lemma (Hopfield 1982)

The function $E(s) = -\frac{1}{2} \sum_{i,j} w_{ij} s_i s_j$ is a *Ljapunov-function* of the system in the case of asynchronous updates.

That means:

- when the activation state of the network changes, E can either decrease or remain the same.

Consequence: The output states $\lim_{n \rightarrow \infty} f^n(s)$ can be characterized as *the local minima* of the Ljapunov-function.

Remark: $E(s) = -\frac{1}{2} \sum_{i,j} w_{ij} s_i s_j = -\sum_{i < j} w_{ij} s_i s_j$ (symmetry!)

For the proof we assume a discrete working space ($s_i = \pm 1$)

Let node 1 be selected for update:

$$s_1(t+1) = \theta (\sum_j w_{1j} \cdot s_j(t)); s_j(t+1) = s_j(t) \text{ for } j \neq 1$$

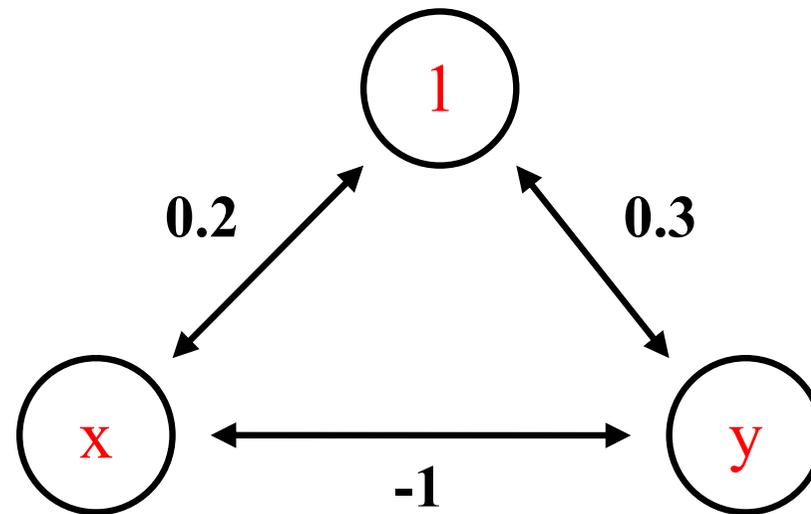
Case 1: $s_1(t+1) = s_1(t)$, then $E(t+1) - E(t) = 0$

Case 2: $s_1(t+1) = -s_1(t)$ [binary threshold!, working space $\{-1, +1\}$]

We have $E(\mathbf{s}(t)) = -\sum_{i < j} w_{ij} s_i(t) s_j(t)$. For the difference $E(t+1) - E(t)$ only the terms with index $i=1, 1 < j$ matter. Consequently,

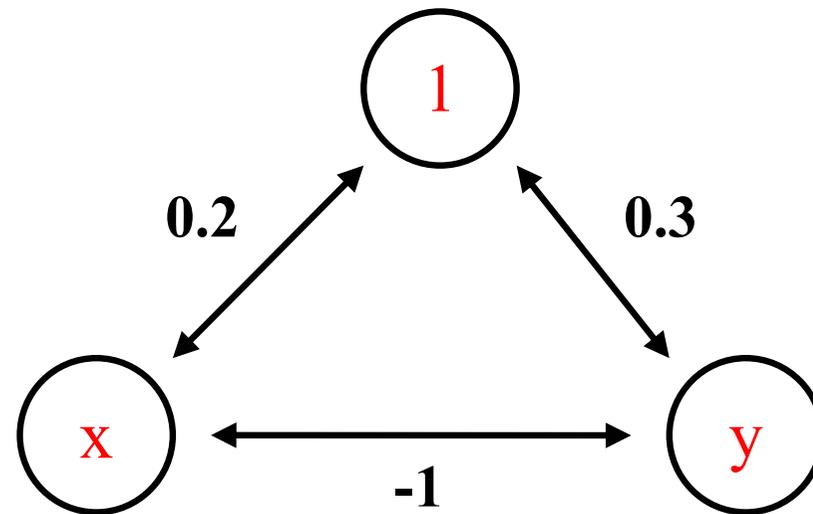
$$\begin{aligned} E(t+1) - E(t) &= -\sum_{j > 1} w_{1j} s_1(t+1) s_j(t) + \sum_{j > 1} w_{1j} s_1(t) s_j(t) = \\ &= -\sum_j w_{1j} s_1(t+1) s_j(t) - \sum_j w_{1j} s_1(t) s_j(t) = -2 s_1(t+1) \cdot \sum_j w_{1j} s_j(t) < 0 \end{aligned}$$

because the two factors have the same sign!



- What is the Ljapunov-function $E(x,y)$ for this system?
- What is the global minimum? (assuming binary activations ± 1)

Example



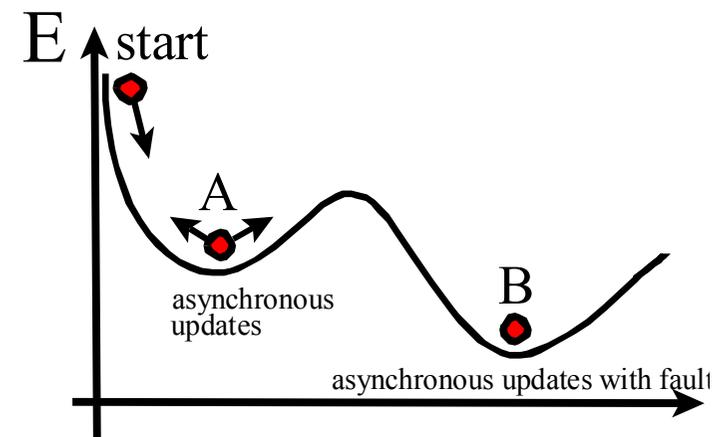
$$E = -\sum_{i < j} w_{ij} s_i s_j$$

$$E(x,y) = -0.2x - 0.3y + xy$$

x	y	E
1	1	.5
1	-1	-.9
-1	1	-1.1
-1	-1	1.5

Theorem 2 (Hopfield 1982)

The output states $\lim_{n \rightarrow \infty} f^n(s)$ can be characterized as *the global minima* of the Lyapunov-function if certain stochastic update functions f are considered ("simulated annealing").



What we need is a probability distribution for the states $P(s)$ for each time and a stochastic update rule which respects that there is some stochastic disturbance during updating the activation vectors.

How to escape local minima?

- One idea: add randomness, so that we can go uphill sometimes. Then we can escape shallow local minima and more likely end up in deep minima. We can use a *probabilistic update rule*.
- Too little randomness: we end up in local minima. Too much, and we jump around instead of converging.
- Solution by analogy from thermodynamics: annealing through slow cooling. Start the network in a high temperature state, and slowly decrease the temperature according to an annealing schedule.

Deriving a probability distribution

What is a plausible probability distribution for activation states?

1. Assume that the probability of an activation state is an function of its *energy*: $P(\mathbf{s}) = f(E(\mathbf{s}))$
2. Assume independent probability distributions for independent subnets
 $E(\mathbf{s} \oplus \mathbf{s}') = E(\mathbf{s}) + E(\mathbf{s}')$; $P(\mathbf{s} \oplus \mathbf{s}') = P(\mathbf{s}) \cdot P(\mathbf{s}')$

$$(*) f(E+E') = f(E) \cdot f(E')$$

Assume f is a continuous function that satisfies f , then it must be an exponential function. Hence,

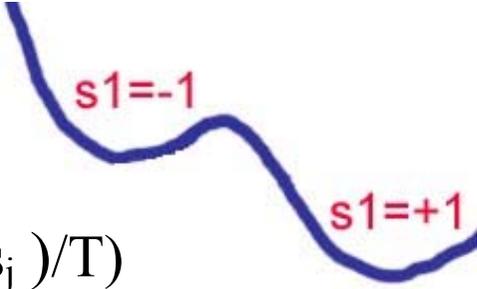
$$P(\mathbf{s}) = \text{const} \cdot e^{k \cdot E(\mathbf{s})}, \text{ or with } k = -1/T:$$

$$P(\mathbf{s}) = \text{const} \cdot e^{-E(\mathbf{s})/T}$$

A probabilistic update rule

Assume (without restricting generality) that unit 1 is selected for updating at time t and has activity -1 at time t . Should it flip from -1 to $+1$ at time $t+1$? The energy $E = -\sum_{i<j} w_{ij} s_i s_j$ is relevant!

t:	$s_1 = -1, s_2, \dots, s_n$	
t+1	$s_1 = -1, s_2, \dots, s_n$	$s_1 = +1, s_2, \dots, s_n$
Energy	E	$E - 2\sum_j w_{1j} s_j$
Prob	$c \cdot \exp(-E/T)$	$c \cdot \exp((-E + 2\sum_j w_{1j} s_j)/T)$



If $\sum_j w_{1j} s_j$ is positive, then flip the activation with a probability that increases exponentially with the energy difference $2\sum_j w_{1j} s_j$:

$P(s_1 = +1) = \sigma(2(\sum_j w_{1j} s_j)/T)$ with the sigmoid function σ .

Classical rule for $T \rightarrow 0$

The Metropolis algorithm

At each step, select a unit and calculate the energy difference ΔE between its current state and its flipped state.

If $\Delta E \geq 0$ don't flip the unit.

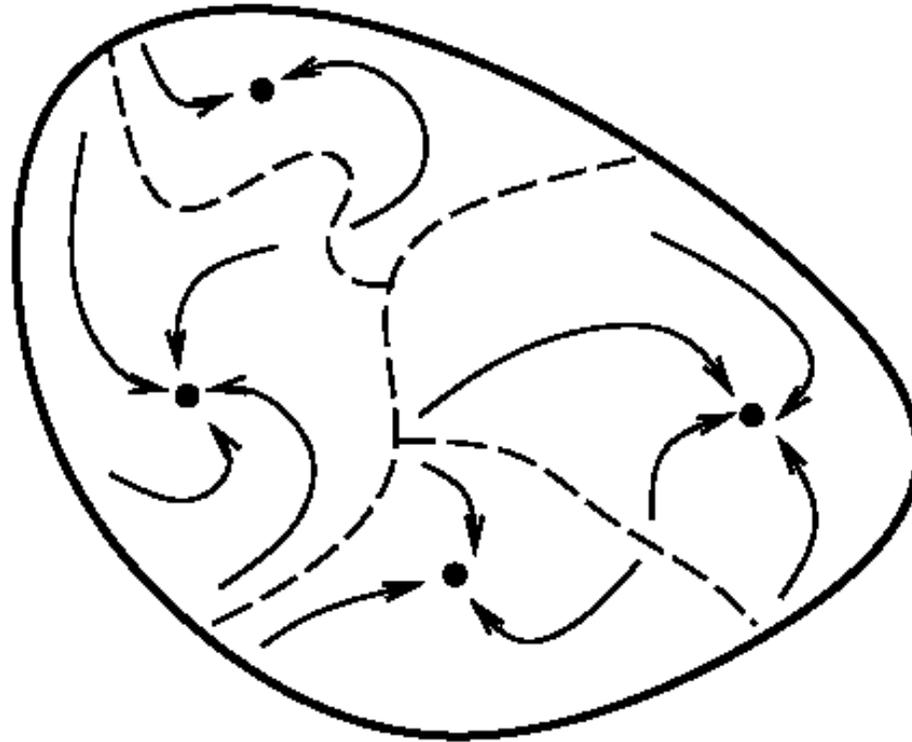
If $\Delta E < 0$, flip the unit with probability $\sigma(-\Delta E/T)$.

After around n such steps, lower the temperature further and repeat the cycle again.

Geman and Geman (1984)

If the temperature in cycle k satisfies $T_k \geq \frac{T_0}{\log(1+k)}$ for every k and T_0 is large enough, then the system will with probability one converge to the minimum energy configuration.

3 Learning with Hopfield networks



Generalized Hebbian rule for a single neuron confronted with a input vector \mathbf{s}^d , working space $S = \{-1, 1\}$

$$\Delta \mathbf{w} = \eta \cdot \mathbf{s}^d \cdot \mathbf{r}^d \quad \mathbf{d} \in D$$

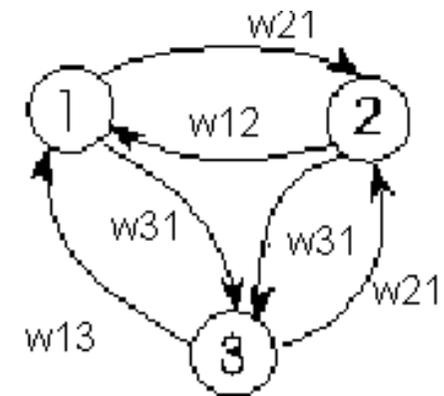
Hopfield used this rule for his networks:

$$\Delta w_{ij} = \eta \cdot s_j^d \cdot r_i^d \quad \text{or equivalently}$$

$$\Delta w_{ij} = \eta \cdot s_j^d \cdot s_i^d \quad \mathbf{d} \in D \quad \text{and } i \neq j!$$

In this case, the resulting connection matrix can be shown to be

$$w_{ij} = 1/N \cdot \sum_{\mathbf{d} \in D} s_j^d \cdot s_i^d \quad \text{for } i \neq j; \text{ zero for } i=j$$



- Consider a network with 3 neurons. Teach the system with the input vector $(1 \ 1 \ 1)$. What is the weight matrix?
- Take the same system, but now teach the system with two input vectors $(1 \ 1 \ 1)$ and $(-1 \ -1 \ -1)$. Is there a behavioural difference that corresponds by adding the second input vector? Take the activation function to be a binary threshold.

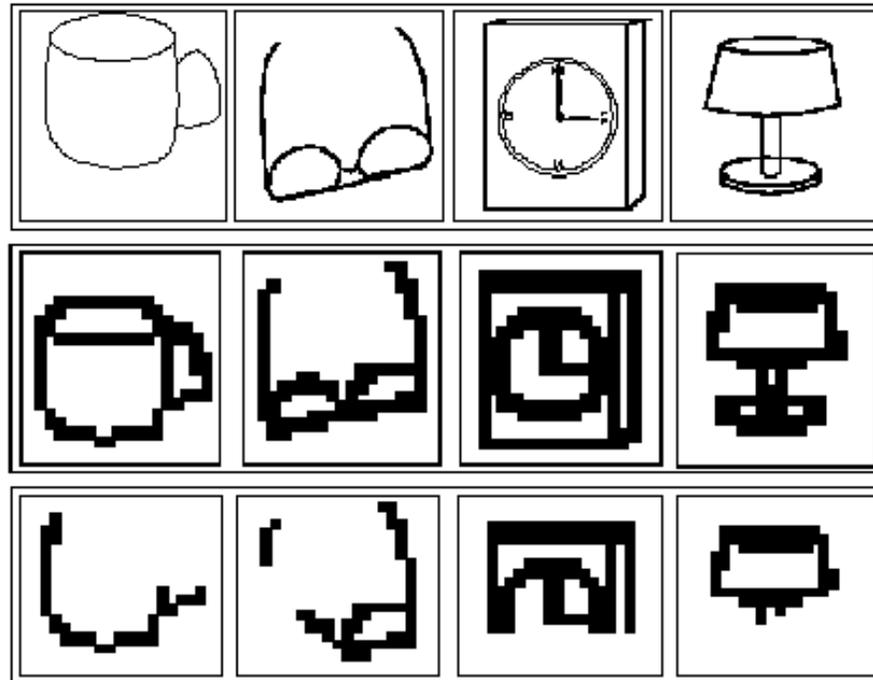
- Consider a network with 3 neurons. Teach the system with the input vector (1 1 1). What is the weight matrix?

$$w = \begin{pmatrix} 0 & 1 & 1 \\ 1 & 0 & 1 \\ 1 & 1 & 0 \end{pmatrix}$$

- Take the same system, but now teach the system with two input vectors (1 1 1) and (-1 -1 -1). Is there a behavioural difference that corresponds by adding the second input vector? Take the activation function to be a binary threshold.

$$w = \begin{pmatrix} 0 & 1 & 1 \\ 1 & 0 & 1 \\ 1 & 1 & 0 \end{pmatrix} \quad \text{No behavioural difference to the case before}$$

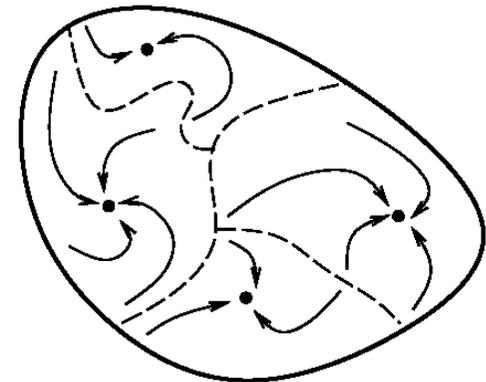
4 Emerging properties of Hopfield networks



Auto-associative memory

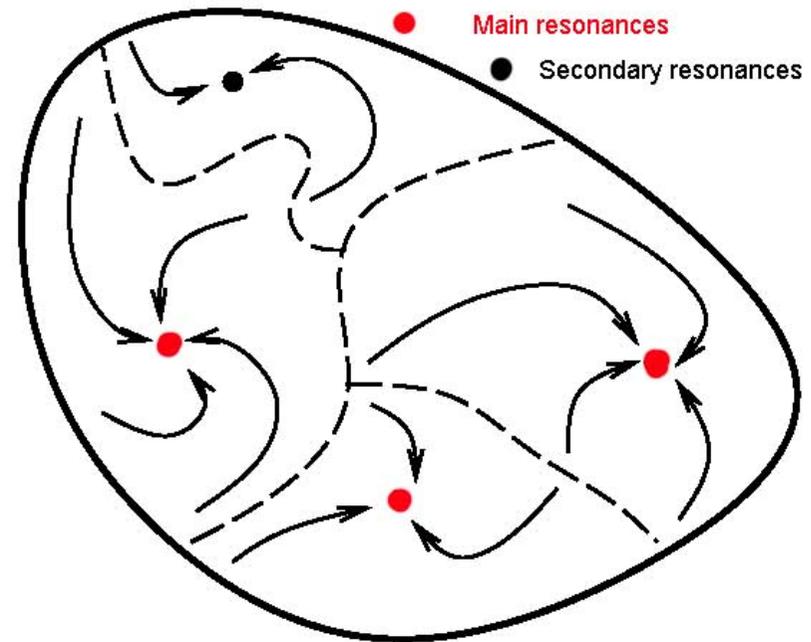
Store a set of patterns $\{ \mathbf{s}^d \}$ in such a way that when presented with a pattern \mathbf{s}^x it will respond with the stored pattern that is most similar to it. Maps patterns to patterns of the same type. (e.g. a noisy or incomplete record maps to a clear record).

- Mechanism of pattern completion: The stored patterns are attractors. If the system starts outside any of the attractors it will begin to move towards one of them.
- Stored patterns are addressable by *content*, not *pointers* (as in traditional computer memories)

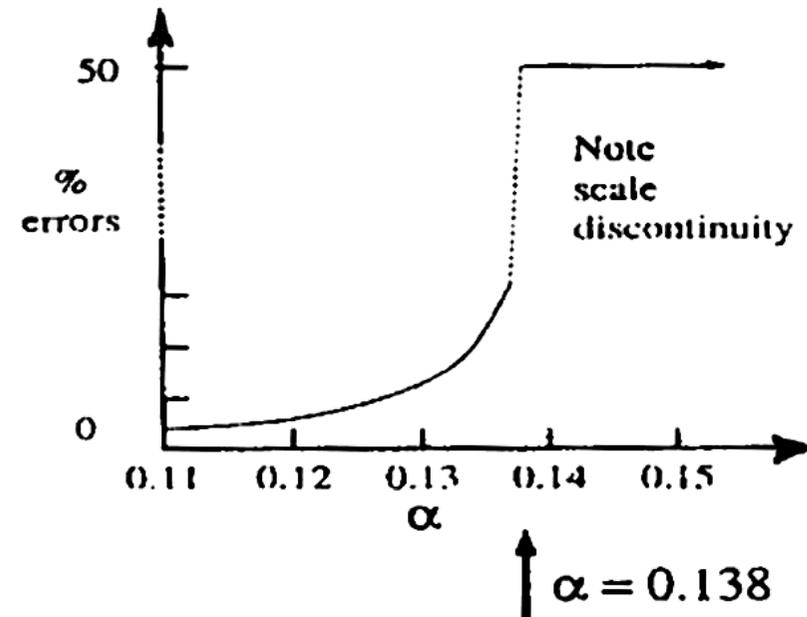


See the link [Hopfield network as associative memory](#) on the website

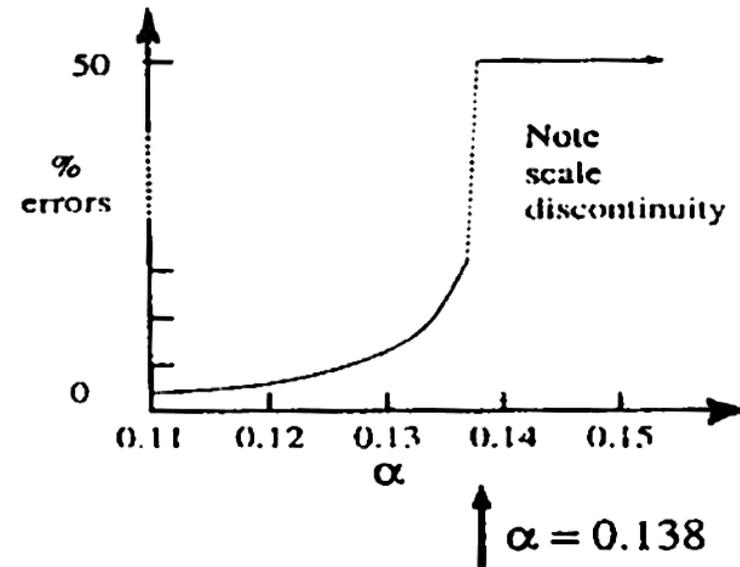
- The inputted patterns of activation are resonances. However, they are not the only resonances of the system
- The state $(-1, -1, \dots, -1)$ is always a resonance
- If \mathbf{s} is a resonance, so is $-\mathbf{s}$



- How many patterns can a n unit network store?
- The more patterns added, the more crosstalk and spurious states (second. resonances). The larger the network, the greater the capacity.
- It turns out that the capacity is roughly proportional to n :
 $M = \alpha \cdot n$, where M is the number of inputted pattern that can be correctly reproduced (with an error probability of p)



- If we want $p_{\text{error}} = 0.01$, then $M = 0.105 \cdot n$. This is an upper bound.
- At $M = 0.138 \cdot n$ patterns the network suddenly breaks down and cannot recall anything useful (“catastrophic forgetting”).
- The behaviour around $\alpha = 0.138$ corresponds to a phase transition in physics (solid – liquid).
- Physical analogy: spin glasses. Unit states correspond to spin states in a solid. Each spin is affected by the spins of the others plus thermal noise, and can flip between two states (1 and -1).



5 Conclusions

- Hopfield nets are "the harmonic oscillator" in modern neurodynamics
- the idea of resonance systems
- the idea of simulated annealing
- the idea of content addressable memory
- very simple learning theory based on the generalized Hebbian rule.