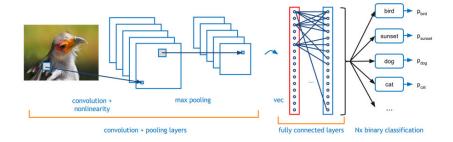
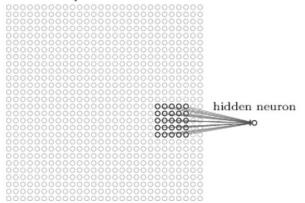
#### **Convolutional network**



# **Convolutional layers**





Every neuron is connected with a (typically small) *receptive field* of neurons in the lower layer.

Neuron is "standard": Computes a weighted sum of its inputs, applies an activation function.

#### **Convolutional layers**

#### input neurons

	000000000000000000000000000000000000000
000000000000000000000000000000000000000	000000000000000000000000000000000000000
00000	
	000000000000000000000000000000000000000

#### input neurons

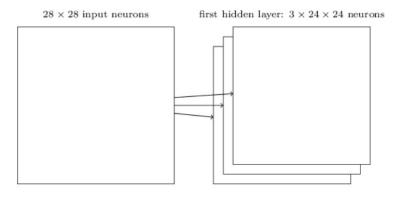
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#### first hidden layer

#### first hidden layer

# Neurons grouped into *feature maps* sharing weights.

# **Convolutional layers**

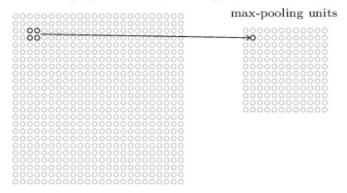


Each feature map represents a property of the input that is supposed to be spatially invariant.

Typically, we consider several feature maps in a single layer.

# **Pooling layers**

#### hidden neurons (output from feature map)



Neurons in the pooling layer compute simple functions of their receptive fields (the fields are typically disjoint):

- Max-pooling : maximum of inputs
- L2-pooling : square root of the sum of squres
- Average-pooling : mean

• • • •

#### **Convolutional networks – architecture**

Neurons organized in layers,  $L_0, L_1, \ldots, L_n$ , connections (typically) only from  $L_m$  to  $L_{m+1}$ .

Several types of layers:

- input layer L<sub>0</sub>
- **dense** layer  $L_m$ : Each neuron of  $L_m$  connected with each neuron of  $L_{m-1}$ .
- convolutional & pooling layer L<sub>m</sub>: Contains two sub-layers:
  - convolutional layer: Neurons organized into disjoint feature maps, all neurons of a given feature map share weights (but have different inputs)
  - pooling layer: Each (convolutional) feature map F has a corresponding pooling map P. Neurons of P
    - ► have inputs only from *F* (typically few of them),
    - compute a simple aggregate function (such as max),
    - have disjoint inputs.

#### **Convolutional networks – architecture**

- Denote
  - X a set of input neurons
  - Y a set of output neurons
  - Z a set of all neurons  $(X, Y \subseteq Z)$
- individual neurons denoted by indices i, j etc.
  - $\xi_j$  is the inner potential of the neuron *j* after the computation stops
  - ► *y<sub>j</sub>* is the output of the neuron *j* after the computation stops

(define  $y_0 = 1$  is the value of the formal unit input)

► *w<sub>ji</sub>* is the weight of the connection **from** *i* **to** *j* 

(in particular,  $w_{j0}$  is the weight of the connection from the formal unit input, i.e.  $w_{j0} = -b_j$  where  $b_j$  is the bias of the neuron *j*)

- *j*← is a set of all *i* such that *j* is adjacent from *i* (i.e. there is an arc **to** *j* from *i*)
- j<sup>→</sup> is a set of all *i* such that *j* is adjacent to *i* (i.e. there is an arc **from** *j* to *i*)
- *j*<sub>share</sub> is a set of neurons sharing weights with *j* i.e. neurons that belong to the same feature map as *j*

#### **Convolutional networks – activity**

- neurons of dense and convolutional layers:
  - inner potential of neuron j:

$$\xi_j = \sum_{i \in j_{\leftarrow}} w_{ji} y_i$$

• activation function  $\sigma_j$  for neuron *j* (arbitrary differentiable):

 $\mathbf{y}_j = \sigma_j(\xi_j)$ 

- Neurons of pooling layers: Apply the "pooling" function:
  - max-pooling:

$$y_j = \max_{i \in j_{\leftarrow}} y_i$$

avg-pooling:

$$y_j = \frac{\sum_{i \in j_{\leftarrow}} y_i}{|j_{\leftarrow}|}$$

A convolutional network is evaluated layer-wise (as MLP), for each  $j \in Y$  we have that  $y_j(\vec{w}, \vec{x})$  is the value of the output neuron j after evaluating the network with weights  $\vec{w}$  and input  $\vec{x}$ .

#### **Convolutional networks – learning**

#### Learning:

• Given a training set  ${\mathcal T}$  of the form

$$\left\{ \left( \vec{x}_k, \vec{d}_k \right) \mid k = 1, \dots, p \right\}$$

Here, every  $\vec{x}_k \in \mathbb{R}^{|X|}$  is an *input vector* end every  $\vec{d}_k \in \mathbb{R}^{|Y|}$  is the desired network output. For every  $j \in Y$ , denote by  $d_{kj}$  the desired output of the neuron j for a given network input  $\vec{x}_k$  (the vector  $\vec{d}_k$  can be written as  $(d_{kj})_{j \in Y}$ ).

Error function – mean square error (for example):

$$E(\vec{w}) = \frac{1}{p} \sum_{k=1}^{p} E_k(\vec{w})$$

where

$$E_k(\vec{w}) = \frac{1}{2} \sum_{j \in Y} (y_j(\vec{w}, \vec{x}_k) - d_{kj})^2$$

#### **Convolutional networks – SGD**

The algorithm computes a sequence of weight vectors  $\vec{w}^{(0)}, \vec{w}^{(1)}, \vec{w}^{(2)}, \dots$ 

- weights in  $\vec{w}^{(0)}$  are randomly initialized to values close to 0
- In the step t + 1 (here t = 0, 1, 2...), weights w<sup>(t+1)</sup> are computed as follows:
  - Choose (randomly) a set of training examples  $T \subseteq \{1, ..., p\}$
  - Compute

$$\vec{w}^{(t+1)} = \vec{w}^{(t)} + \Delta \vec{w}^{(t)}$$

where

$$\Delta \vec{w}^{(t)} = -\varepsilon(t) \cdot \frac{1}{|T|} \sum_{k \in T} \nabla E_k(\vec{w}^{(t)})$$

Here *T* is a *minibatch* (of a fixed size),

•  $0 < \varepsilon(t) \le 1$  is a *learning rate* in step t + 1

►  $\nabla E_k(\vec{w}^{(t)})$  is the gradient of the error of the example *k* Note that the random choice of the minibatch is typically implemented by randomly shuffling all data and then choosing minibatches sequentially. **Epoch** consists of one round through all data.

#### Backprop

Recall that  $\nabla E_k(\vec{w}^{(t)})$  is a vector of all partial derivatives of the form  $\frac{\partial E_k}{\partial w_i}$ .

How to compute  $\frac{\partial E_k}{\partial w_{ji}}$  ?

First, switch from derivatives w.r.t.  $w_{ji}$  to derivatives w.r.t.  $y_j$ :

Recall that for every w<sub>ji</sub> where j is in a dense layer, i.e. does not share weights:

$$\frac{\partial E_k}{\partial w_{ji}} = \frac{\partial E_k}{\partial y_j} \cdot \sigma'_j(\xi_j) \cdot y_i$$

Now for every w<sub>ji</sub> where j is in a convolutional layer, that is shares w<sub>ji</sub> with neurons of j<sub>share</sub>:

$$\frac{\partial E_k}{\partial \mathbf{w}_{ji}} = \sum_{r \in j_{share}} \frac{\partial E_k}{\partial y_r} \cdot \sigma_r'(\xi_r) \cdot y_r$$

Neurons of pooling layers do not have weights.

#### Backprop

Now compute derivatives w.r.t. y<sub>j</sub>:

• for every  $j \in Y$ :

$$\frac{\partial E_k}{\partial y_j} = y_j - d_{kj}$$

This holds for the mean-square error, for other error functions the derivative w.r.t. outputs will be different.

For every j ∈ Z \ Y such that j<sup>→</sup> is either a dense layer, or a convolutional layer:

$$\frac{\partial E_k}{\partial y_j} = \sum_{r \in j^{\rightarrow}} \frac{\partial E_k}{\partial y_r} \cdot \sigma'_r(\xi_r) \cdot \mathbf{w}_{rj}$$

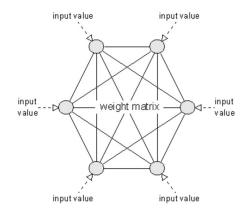
for every j ∈ Z \ Y such that j<sup>→</sup> is max-pooling: Then j<sup>→</sup> = {i} for a single "max" neuron and we have

$$\frac{\partial E_k}{\partial y_j} = \begin{cases} \frac{\partial E_k}{\partial y_i} & \text{if } j = arg \ max_{r \in i_{\leftarrow}} y_r \\ 0 & \text{otherwise} \end{cases}$$

I.e. gradient can be propagated from the output layer downwards as in MLP.

- Conv. nets. are nowadays the most used networks in image processing (and also in other areas where input has some local, "spatially" invariant properties)
- Typically trained using backpropagation.
- Due to the weight sharing allow (very) deep architectures.
- Typically extended with more adjustments and tricks in their topologies.

#### **Recurrent networks – Hopfield network**



Auto-associative network: Given an input, the network outputs a training example (encoded in its weights) "similar" to the given input.

#### Architecture:

- complete topology, i.e. output of each neuron is input to all neurons
- all neurons are both input and output
- denote by ξ<sub>1</sub>,..., ξ<sub>n</sub> inner potentials and by y<sub>1</sub>,..., y<sub>n</sub> outputs (states) of individual neurons
- ► denote by  $w_{ji}$  the weight of connection from a neuron  $i \in \{1, ..., n\}$  to a neuron  $j \in \{1, ..., n\}$
- assume  $w_{jj} = 0$  for every j = 1, ..., n
- For now: no neuron has a bias

#### **Hopfield network**

#### Learning: Training set

 $\mathcal{T} = \{\vec{x}_k \mid \vec{x}_k = (x_{k1}, \dots, x_{kn}) \in \{-1, 1\}^n, k = 1, \dots, p\}$ 

The goal is to "store" the training examples of  $\mathcal{T}$  so that the network is able to *associate* similar examples.

**Hebb's learning rule:** If the inputs to a system cause the same pattern of activity to occur repeatedly, the set of active elements constituting that pattern will become increasingly strongly interassociated. That is, each element will tend to turn on every other element and (with negative weights) to turn off the elements that do not form part of the pattern. To put it another way, the pattern as a whole will become "auto-associated".

#### Mathematically speaking:

$$w_{ji} = \sum_{k=1}^{p} x_{kj} x_{ki} \qquad 1 \le j \ne i \le n$$

Intuition: "Neurons that fire together, wire together".

# **Hopfield network**

Learning: Training set

$$\mathcal{T} = \{\vec{x}_k \mid \vec{x}_k = (x_{k1}, \dots, x_{kn}) \in \{-1, 1\}^n, k = 1, \dots, p\}$$

Hebb's rule:

$$w_{ji} = \sum_{k=1}^{p} x_{kj} x_{ki} \qquad 1 \le j \ne i \le n$$

Note that  $w_{ii} = w_{ij}$ , i.e. the weight matrix is symmetric.

Learning can be seen as poll about equality of inputs:

- If x<sub>kj</sub> = x<sub>ki</sub>, then the training example votes for "*i* equals *j*" by adding one to w<sub>ji</sub>.
- If x<sub>kj</sub> ≠ x<sub>ki</sub>, then the training example votes for "*i* does not equal *j*" by subtracting one from w<sub>ji</sub>.

#### **Hopfield network**

**Activity:** Initially, neurons set to the network input  $\vec{x} = (x_1, ..., x_n)$ , thus  $y_j^{(0)} = x_j$  for every j = 1, ..., n.

Cyclically update states of neurons, i.e. in step t + 1 compute the value of a neuron *j* such that  $j = (t \mod p) + 1$ , as follows: Compute the inner potential:

$$\xi_j^{(t)} = \sum_{i=1}^n w_{ji} y_i^{(t)}$$

then

$$y_j^{(t+1)} = \begin{cases} 1 & \xi_j^{(t)} > 0 \\ y_j^{(t)} & \xi_j^{(t)} = 0 \\ -1 & \xi_j^{(t)} < 0 \end{cases}$$

#### Hopfield network – activity

The computation stops in a step  $t^*$  if the network is for the first time in a *stable* state, i.e.

$$y_j^{(t^*+n)} = y_j^{(t^*)}$$
  $(j = 1, ..., n)$ 

#### Theorem

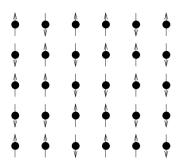
Assuming symmetric weights, computation of a Hopfiled network always stops for every input.

This implies that a given Hopfiled network computes a function from  $\{-1, 1\}^n$  to  $\{-1, 1\}^n$  (determined by its weights).

Denote by  $\vec{y}(W, \vec{x}) = (y_1^{(t^*)}, \dots, y_n^{(t^*)})$  the value of the network for a given input  $\vec{x}$  and a weight matrix W. Denote by  $y_j(W, \vec{x}) = y_j^{(t^*)}$  the component of the value of the network corresponding to the neuron *j*.

If *W* is clear from the context, we write only  $y(\vec{x})$  a  $y_j(\vec{x})$ .

Simple models of magnetic materials resemble Hopfield network.



- atomic magnets organized into square-lattice
- each magnet may have only one of two possible orientations (in the Hopfield network +1 a -1)
- orientation of each magnet is influenced by an external magnetic field (input of the network) as well as orientation of the other magnets
- weights in the Hopfiled net model determine interaction among magnets

# **Energy function**

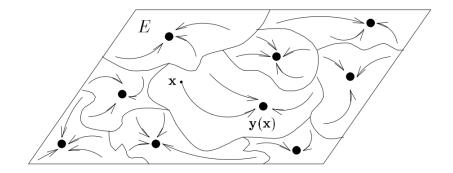
Energy function *E* assigns to every state  $\vec{y} \in \{-1, 1\}^n$  a (potential) energy:

$$E(\vec{y}) = -\frac{1}{2} \sum_{j=1}^{n} \sum_{i=1}^{n} w_{ji} y_j y_i$$

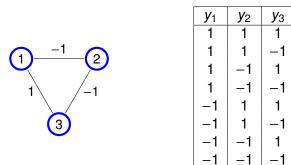
- states with low energy are stable (few neurons "want to" change their states), states with high energy are not stable
- i.e. large (positive) w<sub>ji</sub>y<sub>j</sub>y<sub>i</sub> is stable and small (negative) w<sub>ji</sub>y<sub>j</sub>y<sub>i</sub> is not stable

The energy does not increase during computation:  $E(\vec{y}^{(t)}) \ge E(\vec{y}^{(t+1)})$ , stable states  $\vec{y}^{(t^*)}$  correspond to local minima of *E*.

# **Energy landscape**



# Hopfield – example



- Hopfield network with three neurons
- ► trained on a single training example (1, -1, 1) using Hebb's rule

Е

-3

-3

(note that (-1, 1, -1) has also been "stored" into the network)

Observe that

- ► the energy does not increase during computation:  $E(\vec{y}^{(t)}) ≥ E(\vec{y}^{(t+1)})$
- ► if the state is updated in a step t + 1, then  $E(\vec{y}^{(t)}) > E(\vec{y}^{(t+1)})$
- there are only finitely many states, and thus, eventually, a local minimum of E is reached.

This proves that computation of a Hopfield network always stops.

The energy function E may have local minima that do not correspond to training examples (so called phantoms).

Phantoms can be "unlearned" e.g. using the following rule: Given a phantom  $(x_1, ..., x_n) \in \{-1, 1\}^n$  and weights  $w_{ji}$ , then new weights  $w'_{ii}$  are computed by

$$w_{ji}' = w_{ji} - x_i x_j$$

(i.e. similar to Hebb's rule but with the opposite sign)

#### **Reproduction – statistical analysis**

**Capacity** of Hopfield network is defined as the ratio p/n of number of training examples the net is able to learn over the number of neurons.

Assume that training examples are chosen randomly: each component of  $\vec{x}_k$  is set to 1 with probability 1/2 and to -1 with probability 1/2.

Consider a configuration *W* obtained by learning using the Hebb's rule.

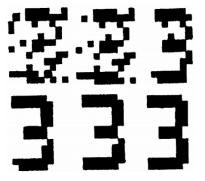
Denote

$$\beta = \mathbf{P} \left[ \vec{x}_k = \vec{y}(W, \vec{x}_k) \text{ pro } k = 1, \dots, p \right]$$

Then for  $n \to \infty$  and  $p \le n/(4 \log n)$  we have  $\beta \to 1$ .

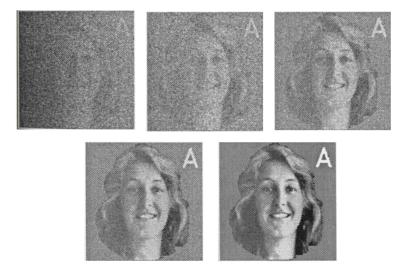
I.e. the maximum number of examples that can be effectively stored in Hopfield net is proportional to  $n/(4 \log n)$ .

# Hopfield network – example



- figures 12 × 10
  (120 neurons, -1 is white and 1 is black)
- learned 8 figures
- input generated with 25% noise
- image shows the activity of the Hopfield network

# Hopfield network – example



# Hopfield network – example

