

# Deep Learning & Applied AI

Multi-layer perceptron and  
back-propagation

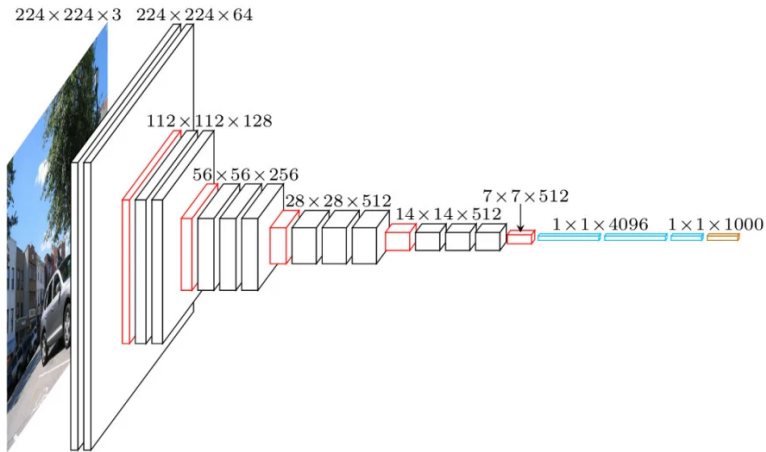
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SAPIENZA  
UNIVERSITÀ DI ROMA

# A glimpse into neural networks

In deep learning, we deal with highly parametrized models called deep neural networks:



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More in general, consider other **activation functions**:

$$\sigma(x) = \frac{1}{1 + e^{-x}}$$

continuous

$$\sigma(x) = \max\{0, x\}$$

discontinuous  
gradient

# Multi-layer perceptron

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Each layer outputs an intermediate **hidden representation**:

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The hidden representations are sometimes called the **activations** at layer  $\ell + 1$ .

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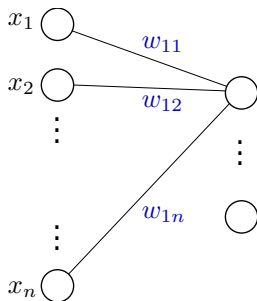
- ① Each layer is a vector-to-vector function  $\mathbb{R}^p \rightarrow \mathbb{R}^q$ .
- ② Each layer has  $q$  neurons acting **in parallel**.  
Each neuron acts as a scalar function  $\mathbb{R}^p \rightarrow \mathbb{R}$ .

# Single layer illustration

$$\sigma(\mathbf{W}\mathbf{x}) = \sigma \circ \begin{pmatrix} w_{11} & w_{12} & \cdots & w_{1n} \\ w_{21} & w_{22} & \cdots & w_{2n} \\ \vdots & \cdots & \ddots & \vdots \\ w_{m1} & w_{m2} & \cdots & w_{mn} \end{pmatrix} \begin{pmatrix} x_1 \\ x_2 \\ \vdots \\ x_n \end{pmatrix} = \sigma \circ \begin{pmatrix} y_1 \\ y_2 \\ \vdots \\ y_m \end{pmatrix}$$

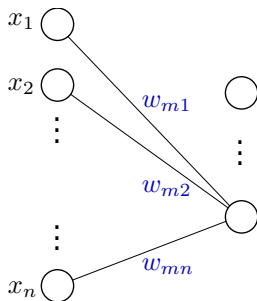
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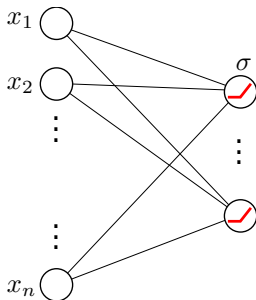
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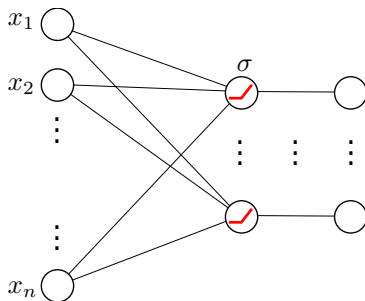
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It is common to have a **linear** layer at the output:

$$\mathbf{y} = f \circ (\sigma \circ f) \circ \cdots \circ (\sigma \circ f)(\mathbf{x})$$

which maps:

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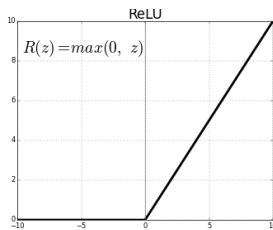
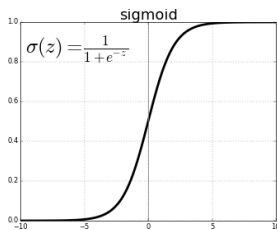
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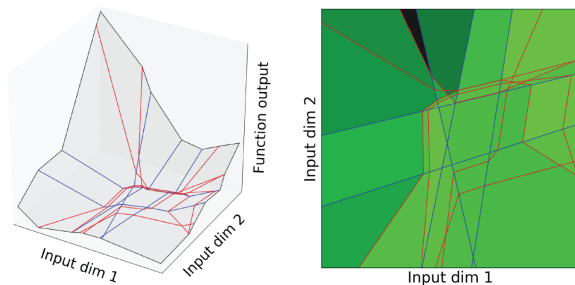
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The blue and red edges are produced by the **first** and **second** layer.

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If  $\sigma$  is sigmoidal, we have the following:

**Universal Approximation Theorem** For any compact set  $\Omega \subset \mathbb{R}^p$ , the space spanned by the functions  $\phi(\mathbf{x}) = \sigma(\mathbf{W}\mathbf{x} + \mathbf{b})$  is dense in  $\mathcal{C}(\Omega)$  for the uniform convergence.



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For large enough  $q$ , the training error can be made **arbitrarily small**.

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- One layer, sigmoid activation, cross-entropy loss ( $\Rightarrow$  logistic regression).



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We want to automatize this **computational step** efficiently.

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$x$   
●



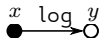
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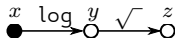
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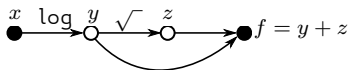
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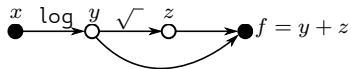
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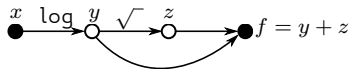
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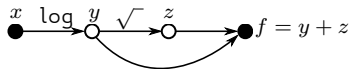
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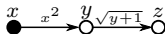
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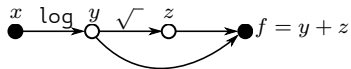
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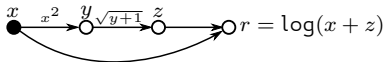
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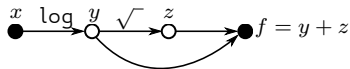
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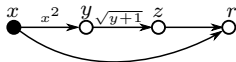
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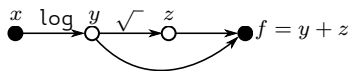
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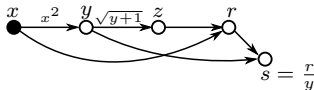
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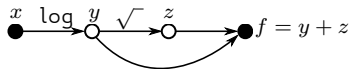
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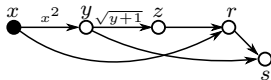
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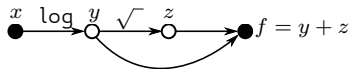
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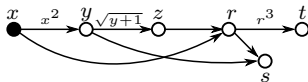
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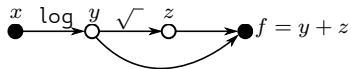
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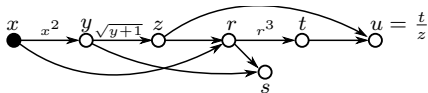
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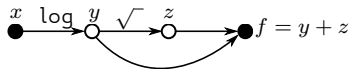
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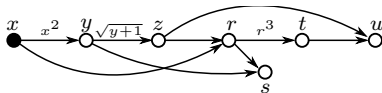
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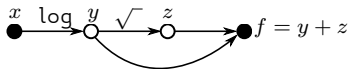
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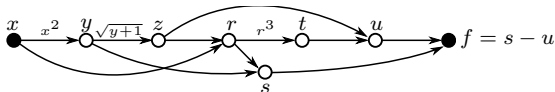
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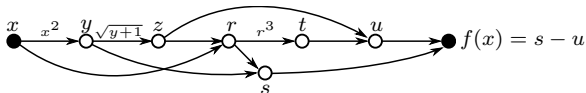
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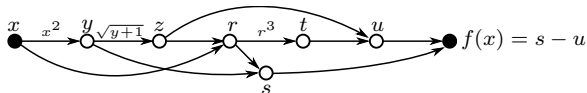
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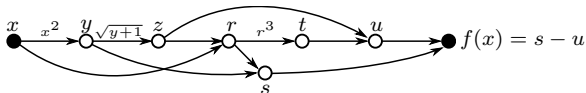
The graph is constructed programmatically, for example:

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z = sqrt(sum(square(x), 1));
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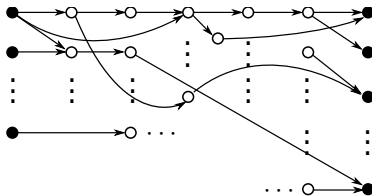
The evaluation of  $f(x)$  corresponds to a **forward traversal** of the graph:



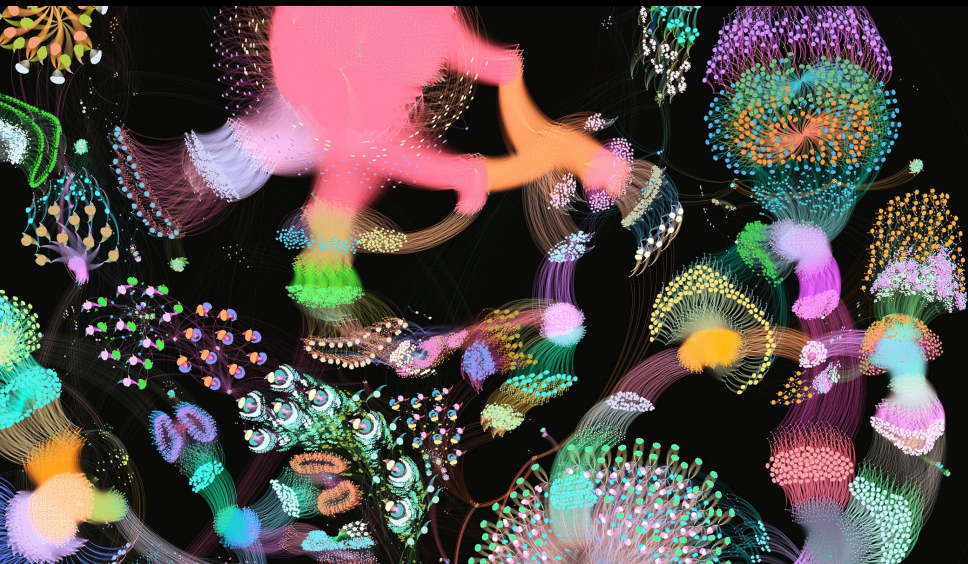
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```
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```

For **high-dimensional** input/output, the graph may be more complex:



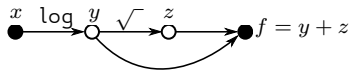
The computational graph gets big quickly.



Poplar visualization, see <https://www.graphcore.ai/products/poplar>

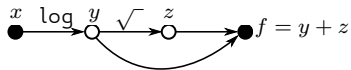
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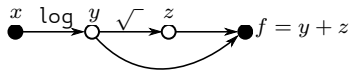
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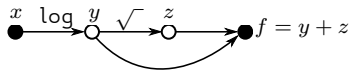


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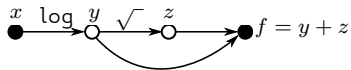


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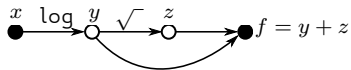


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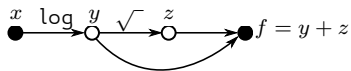
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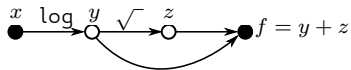
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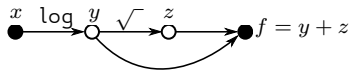
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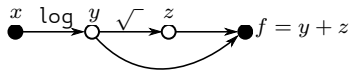
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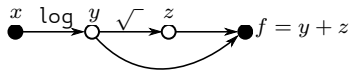
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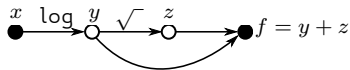
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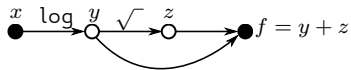
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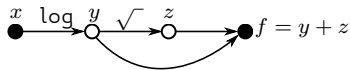
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Assumption: Each partial derivative is a "primitive" accessible in **closed form** and can be computed on the fly.

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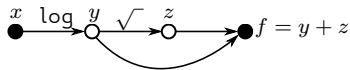
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However, if the input is high-dimensional, i.e.  $f: \mathbb{R}^p \rightarrow \mathbb{R}$ :

$$\text{cost of computing } \nabla f(\mathbf{x}) = p \times \text{cost of computing } f(\mathbf{x})$$

since partial derivatives must be computed w.r.t. each input dimension.

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Computes all the partial derivatives  $\frac{\partial y}{\partial x}, \frac{\partial z}{\partial x}, \dots$  with respect to the **input**  $x$ .

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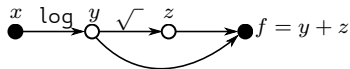
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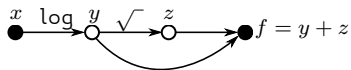
We accumulate values during code execution, to get numerical **evaluations** rather than **expressions** for the derivative.



**Reverse mode:** compute all the partial derivatives  $\frac{\partial f}{\partial z}, \dots, \frac{\partial f}{\partial x}$  with respect to the **inner nodes**.

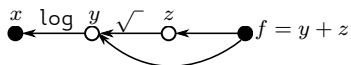
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# Automatic differentiation: Reverse mode

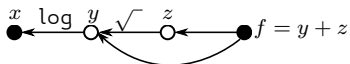
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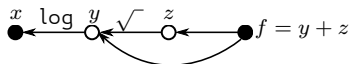
$$\frac{\partial f}{\partial z} =$$

$$\frac{\partial f}{\partial y} =$$

$$\frac{\partial f}{\partial x} =$$

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$$\frac{\partial f}{\partial f} = 1$$

$$\frac{\partial f}{\partial z} = \frac{\partial f}{\partial f} \frac{\partial f}{\partial z}$$

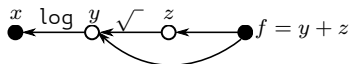
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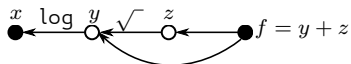
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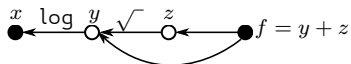
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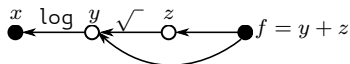
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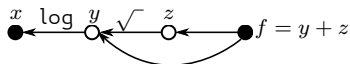
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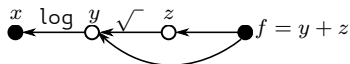
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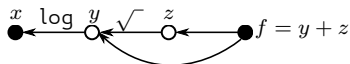
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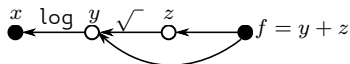
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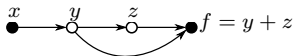
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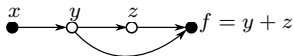
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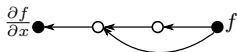
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When training NNs, we compute the gradient of a loss

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$\epsilon$  computes the actual scalar error for the loss.

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Backprop through computational graph of the loss

$\approx$

Backprop “through the network”

## Suggested reading

Nice, accessible survey on automatic differentiation:  
<https://arxiv.org/abs/1502.05767>