Deep Learning

A course about theory & practice

UNIVERSITÀ DI PAVIA

Artificial Neural Networks

Basic ideas, notation and all that

Marco Piastra

Function approximation: linear combination

Function Approximation: linear combination

- Approximating a target function

$$y = f^*(\boldsymbol{x}), \quad \boldsymbol{x} \in \mathbb{R}^d$$

A first approximator: *linear combination*

a.k.a. "*single layer perceptron*"

$$\tilde{y} = \boldsymbol{w} \cdot \boldsymbol{x} + b, \quad \boldsymbol{w} \in \mathbb{R}^d, b \in \mathbb{R}$$

i.e. this is a vector of dimension d

Note that, when the input is scalar, the approximator becomes

$$\tilde{y} = wx + b$$

i.e. a straight line

Function Approximation: linear combination

- Approximating a target function

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A first approximator: *linear combination*

$$\tilde{y} = \boldsymbol{w} \cdot \boldsymbol{x} + b, \quad \boldsymbol{w} \in \mathbb{R}^d, b \in \mathbb{R}$$

dataset

*A set of input and output pairs (**data items**) is what know about the target function*

$$D := \{(\boldsymbol{x}^{(i)}, y^{(i)})\}_{i=1}^{N}, \qquad y^{(i)} = f^*(\boldsymbol{x}^{(i)}), \forall i$$

A set of data items — *Item index*

Function Approximation: linear combination

- Approximating a target function

$$y = f^*(\boldsymbol{x}), \quad \boldsymbol{x} \in \mathbb{R}^d$$

A first approximator: *linear combination*

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Three fundamental aspects:

- ***representation**: which parametric approximator for a given target function?*
- ***evaluation**: how could you tell that some parameter values are better than others?*
- ***optimization**: how can we learn optimal values for the parameters?*

Function Approximation: linear combination

- Example: XOR

$$y = \mathrm{XOR}(\boldsymbol{x}), \quad \boldsymbol{x} \in \{0,1\}^2$$

Approximator: *linear combination*

$$\tilde{y} = \boldsymbol{w} \cdot \boldsymbol{x} + b, \quad \boldsymbol{w} \in \mathbb{R}^d, b \in \mathbb{R}$$

Dataset:

$$D := \{(\boldsymbol{x}^{(i)}, y^{(i)})\}_{i=1}^{N}$$

x_1	x_2	$x_1 \oplus x_2$
0	0	0
0	1	1
1	0	1
1	1	0

this is our *dataset* ($N = 4$)

Function Approximation: linear combination

- Example: XOR

$$y = \mathrm{XOR}(\boldsymbol{x}), \quad \boldsymbol{x} \in \{0,1\}^2$$

x_1	x_2	$x_1 \oplus x_2$
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Approximator: *linear combination*

$$\tilde{y} = \boldsymbol{w} \cdot \boldsymbol{x} + b, \quad \boldsymbol{w} \in \mathbb{R}^d, b \in \mathbb{R}$$

Dataset:

$$D := \{(\boldsymbol{x}^{(i)}, y^{(i)})\}_{i=1}^{N}$$

Loss function (evaluation):

$$L(\boldsymbol{x}^{(i)}, y^{(i)}) := (\tilde{y}(\boldsymbol{x}^{(i)}) - y^{(i)})^2$$

Squared Error (one data item)

$$L(D) := \frac{1}{N} \sum_{(\boldsymbol{x}^{(i)}, y^{(i)}) \in D} L(\boldsymbol{x}^{(i)}, y^{(i)})$$

Mean Squared Error (MSE – entire dataset)

Function Approximation: linear combination

- Example: XOR

$$y = \mathrm{XOR}(\boldsymbol{x}), \quad \boldsymbol{x} \in \{0,1\}^2$$

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this is our *dataset* ($N = 4$)

Approximator: *linear combination*

$$\tilde{y} = \boldsymbol{w} \cdot \boldsymbol{x} + b, \quad \boldsymbol{w} \in \mathbb{R}^d, b \in \mathbb{R}$$

Dataset:

$$D := \{(\boldsymbol{x}^{(i)}, y^{(i)})\}_{i=1}^{N}$$

Optimization problem:

We need to find

parameter values that minimize the loss w.r.t. to the dataset

$$(\boldsymbol{w}, b)^* := \underset{(\boldsymbol{w},b)}{\operatorname{argmin}} \, L(D)$$

Function Approximation: linear combination

- Loss minimization

  Approximator: *linear combination*

$$\tilde{y} = \boldsymbol{w} \cdot \boldsymbol{x} + b, \quad \boldsymbol{w} \in \mathbb{R}^d, b \in \mathbb{R}$$

  Loss function:

$$\begin{aligned} L(D) &= \frac{1}{N} \sum_{i=1}^{N} L(\boldsymbol{x}^{(i)}, y^{(i)}) \\ &= \frac{1}{N} \sum_{i=1}^{N} (\tilde{y}(\boldsymbol{x}^{(i)}) - y^{(i)})^2 \\ &= \frac{1}{N} \sum_{i=1}^{N} ((\boldsymbol{w} \cdot \boldsymbol{x}^{(i)} + b) - y^{(i)})^2 \end{aligned}$$

  Can we express this summation by using linear algebra?

  *Matrix representation leads to a better **parallelization** of computations (more on this, later on)*

Function Approximation: linear combination

- Loss minimization

  Approximator: *linear combination*

$$\tilde{y} = \boldsymbol{w} \cdot \boldsymbol{x} + b, \quad \boldsymbol{w} \in \mathbb{R}^d, b \in \mathbb{R}$$

  Loss function:

$$L(D) = \frac{1}{N} \sum_{i=1}^{N} ((\boldsymbol{w} \cdot \boldsymbol{x}^{(i)} + b) - y^{(i)})^2$$

  define:

$$\boldsymbol{X} := \begin{bmatrix} x_1^{(1)} & \dots & x_d^{(1)} \\ \vdots & \ddots & \vdots \\ x_1^{(N)} & \dots & x_d^{(N)} \end{bmatrix}$$

input data in matrix form (item index first)

Function Approximation: linear combination

- Loss minimization

  Approximator: *linear combination*

$$\tilde{y} = \boldsymbol{w} \cdot \boldsymbol{x} + b, \quad \boldsymbol{w} \in \mathbb{R}^d, b \in \mathbb{R}$$

  Loss function:

$$L(D) = \frac{1}{N} \sum_{i=1}^{N} ((\boldsymbol{w} \cdot \boldsymbol{x}^{(i)} + b) - y^{(i)})^2$$

  define:

$$\hat{\boldsymbol{X}} := \begin{bmatrix} x_1^{(1)} & \dots & x_d^{(1)} & 1 \\ \vdots & \ddots & \vdots & \vdots \\ x_1^{(N)} & \dots & x_d^{(N)} & 1 \end{bmatrix} \quad \boldsymbol{\vartheta} := \begin{bmatrix} w_1 \\ \vdots \\ w_d \\ b \end{bmatrix} \qquad \boldsymbol{y} := \begin{bmatrix} y^{(1)} \\ \vdots \\ y^{(N)} \end{bmatrix}$$

  The loss function becomes:

$$L(D) = \frac{1}{N}(\hat{\boldsymbol{X}}\boldsymbol{\vartheta} - \boldsymbol{y})^2$$

loss function in matrix form

This is a positive-definite quadratic form

Function Approximation: linear combination

- Loss minimization

  Approximator: *linear combination*

$$\tilde{y} = \boldsymbol{w} \cdot \boldsymbol{x} + b, \quad \boldsymbol{w} \in \mathbb{R}^d, b \in \mathbb{R}$$

  Loss function:

$$L(D) = \frac{1}{N} \sum_{i=1}^{N} ((\boldsymbol{w} \cdot \boldsymbol{x}^{(i)} + b) - y^{(i)})^2$$

  define:

$$\hat{\boldsymbol{X}} := \begin{bmatrix} x_1^{(1)} & \dots & x_d^{(1)} & 1 \\ \vdots & \ddots & \vdots & \vdots \\ x_1^{(N)} & \dots & x_d^{(N)} & 1 \end{bmatrix} \quad \boldsymbol{\vartheta} := \begin{bmatrix} w_1 \\ \vdots \\ w_d \\ b \end{bmatrix} \quad \boldsymbol{y} := \begin{bmatrix} y^{(1)} \\ \vdots \\ y^{(N)} \end{bmatrix}$$

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Function Approximation: linear combination

- Loss minimization

  Approximator: *linear combination*

$$\tilde{y} = \boldsymbol{w} \cdot \boldsymbol{x} + b, \quad \boldsymbol{w} \in \mathbb{R}^d, b \in \mathbb{R}$$

  Loss function:

$$L(D) = \frac{1}{N}(\hat{\boldsymbol{X}}\boldsymbol{\vartheta} - \boldsymbol{y})^2$$

  For XOR:

$$\hat{\boldsymbol{X}} := \begin{bmatrix} 0 & 0 & 1 \\ 0 & 1 & 1 \\ 1 & 0 & 1 \\ 1 & 1 & 1 \end{bmatrix} \qquad \boldsymbol{\vartheta} := \begin{bmatrix} w_1 \\ w_2 \\ b \end{bmatrix} \qquad \boldsymbol{y} := \begin{bmatrix} 0 \\ 1 \\ 1 \\ 0 \end{bmatrix}$$

XOR

x_1	x_2	$x_1 \oplus x_2$
0	0	0
0	1	1
1	0	1
1	1	0

this is our *dataset* ($N = 4$)

Function Approximation: linear combination

- Loss minimization

  Approximator: *linear combination*

$$\tilde{y} = \boldsymbol{w} \cdot \boldsymbol{x} + b, \quad \boldsymbol{w} \in \mathbb{R}^d, b \in \mathbb{R}$$

representation

  Loss function:

$$L(D) = \frac{1}{N}(\hat{\boldsymbol{X}}\boldsymbol{\vartheta} - \boldsymbol{y})^2$$

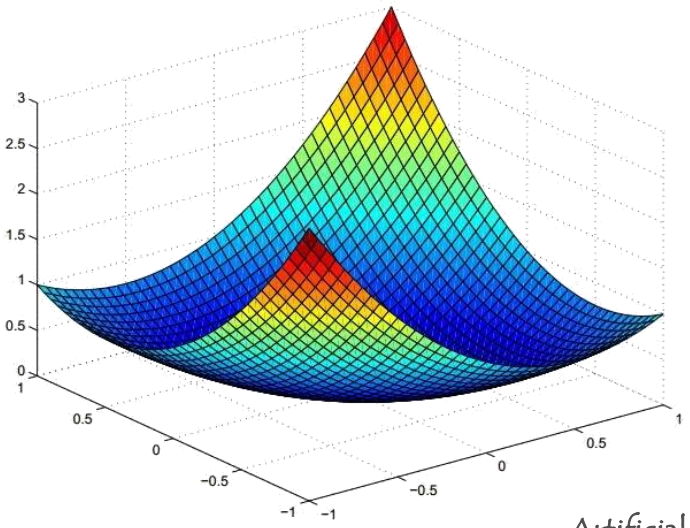
evaluation

  Optimization:

optimization

$$\frac{\partial}{\partial \boldsymbol{\vartheta}} L(D) = 0$$

this loss function is convex:
by solving this equation, we can find $\boldsymbol{\vartheta}^*$
i.e. the optimal parameter values



Function Approximation: linear combination

- Loss minimization

Approximator: *linear combination*

$$\tilde{y} = \boldsymbol{w} \cdot \boldsymbol{x} + b, \quad \boldsymbol{w} \in \mathbb{R}^d, b \in \mathbb{R}$$

Optimization:

$$\begin{aligned}
\frac{\partial}{\partial \boldsymbol{\vartheta}} L(D) &= \frac{1}{N} \frac{\partial}{\partial \boldsymbol{\vartheta}} (\hat{\boldsymbol{X}}\boldsymbol{\vartheta} - \boldsymbol{y})^2 \\
&= \frac{1}{N} \frac{\partial}{\partial \boldsymbol{\vartheta}} (\hat{\boldsymbol{X}}\boldsymbol{\vartheta} - \boldsymbol{y})^T (\hat{\boldsymbol{X}}\boldsymbol{\vartheta} - \boldsymbol{y}) = \frac{1}{N} \frac{\partial}{\partial \boldsymbol{\vartheta}} (\boldsymbol{\vartheta}^T \hat{\boldsymbol{X}}^T - \boldsymbol{y}^T)(\hat{\boldsymbol{X}}\boldsymbol{\vartheta} - \boldsymbol{y}) \\
&= \frac{1}{N} \frac{\partial}{\partial \boldsymbol{\vartheta}} (\boldsymbol{\vartheta}^T \hat{\boldsymbol{X}}^T \hat{\boldsymbol{X}}\boldsymbol{\vartheta} - \boldsymbol{\vartheta}^T \hat{\boldsymbol{X}}^T \boldsymbol{y} - \boldsymbol{y}^T \hat{\boldsymbol{X}}\boldsymbol{\vartheta} + \boldsymbol{y}^T \boldsymbol{y}) \\
&= \frac{1}{N} \frac{\partial}{\partial \boldsymbol{\vartheta}} (\boldsymbol{\vartheta}^T \hat{\boldsymbol{X}}^T \hat{\boldsymbol{X}}\boldsymbol{\vartheta} - 2\boldsymbol{\vartheta}^T \hat{\boldsymbol{X}}^T \boldsymbol{y} + \boldsymbol{y}^T \boldsymbol{y}) \\
&= \frac{1}{N} (2\hat{\boldsymbol{X}}^T \hat{\boldsymbol{X}}\boldsymbol{\vartheta} - 2\hat{\boldsymbol{X}}^T \boldsymbol{y})
\end{aligned}$$

all these terms are scalars

Function Approximation: linear combination

- Loss minimization

  Approximator: *linear combination*

$$\tilde{y} = \boldsymbol{w} \cdot \boldsymbol{x} + b, \quad \boldsymbol{w} \in \mathbb{R}^d, b \in \mathbb{R}$$

  Optimization:

$$\frac{\partial}{\partial \boldsymbol{\vartheta}} L(D) = \frac{1}{N}(2\hat{\boldsymbol{X}}^T \hat{\boldsymbol{X}} \boldsymbol{\vartheta} - 2\hat{\boldsymbol{X}}^T \boldsymbol{y})$$

$$\frac{\partial}{\partial \boldsymbol{\vartheta}} L(D) = 0 \quad \Longrightarrow \quad 2\hat{\boldsymbol{X}}^T \hat{\boldsymbol{X}} \boldsymbol{\vartheta} - 2\hat{\boldsymbol{X}}^T \boldsymbol{y} = 0$$

$$\hat{\boldsymbol{X}}^T \hat{\boldsymbol{X}} \boldsymbol{\vartheta} = \hat{\boldsymbol{X}}^T \boldsymbol{y}$$

$$\boldsymbol{\vartheta} = (\hat{\boldsymbol{X}}^T \hat{\boldsymbol{X}})^{-1} \hat{\boldsymbol{X}}^T \boldsymbol{y}$$

this is what we need

this matrix is SQUARE and SYMMETRIC and, typically, with actual datasets is invertible (i.e. full rank)

Function Approximation: linear combination

- Loss minimization

  Approximator: *linear combination*

$$\tilde{y} = \boldsymbol{w} \cdot \boldsymbol{x} + b, \quad \boldsymbol{w} \in \mathbb{R}^d, b \in \mathbb{R}$$

  For XOR:

$$\boldsymbol{\vartheta} = (\hat{\boldsymbol{X}}^T \hat{\boldsymbol{X}})^{-1} \hat{\boldsymbol{X}}^T \boldsymbol{y}$$

XOR

x_1	x_2	$x_1 \oplus x_2$
0	0	0
0	1	1
1	0	1
1	1	0

$$\hat{\boldsymbol{X}} := \begin{bmatrix} 0 & 0 & 1 \\ 0 & 1 & 1 \\ 1 & 0 & 1 \\ 1 & 1 & 1 \end{bmatrix} \qquad \boldsymbol{\vartheta} := \begin{bmatrix} w_1 \\ w_2 \\ b \end{bmatrix} \qquad \boldsymbol{y} := \begin{bmatrix} 0 \\ 1 \\ 1 \\ 0 \end{bmatrix}$$

$$\hat{\boldsymbol{X}}^T \hat{\boldsymbol{X}} = \begin{bmatrix} 2 & 1 & 2 \\ 1 & 2 & 2 \\ 2 & 2 & 4 \end{bmatrix} \qquad (\hat{\boldsymbol{X}}^T \hat{\boldsymbol{X}})^{-1} = \begin{bmatrix} 1 & 0 & 0.5 \\ 0 & 1 & 0.5 \\ 0.5 & 0.5 & 0.75 \end{bmatrix}$$

$$(\hat{\boldsymbol{X}}^T \hat{\boldsymbol{X}})^{-1} \hat{\boldsymbol{X}}^T \boldsymbol{y} = \begin{bmatrix} 0 \\ 0 \\ 0.5 \end{bmatrix}$$

Function Approximation: linear combination

- Loss minimization

  Approximator: *linear combination*

$$\tilde{y} = \boldsymbol{w} \cdot \boldsymbol{x} + b, \quad \boldsymbol{w} \in \mathbb{R}^d, b \in \mathbb{R}$$

  For XOR:

$$\boldsymbol{\vartheta} := \begin{bmatrix} w_1 \\ w_2 \\ b \end{bmatrix} = \begin{bmatrix} 0 \\ 0 \\ 0.5 \end{bmatrix}$$

  hence the XOR linear approximator becomes:

$$\tilde{y} = 0.5$$

What ???
A constant function?

XOR

x_1	x_2	$x_1 \oplus x_2$
0	0	0
0	1	1
1	0	1
1	1	0

Function approximation: Feed-Forward Neural Network

Feed-Forward Neural Network

- Approximating a target function

$$y = f^*(\boldsymbol{x}), \quad \boldsymbol{x} \in \mathbb{R}^d$$

Second attempt: ***(shallow) feed-forward neural network***

$$\tilde{y} = \boldsymbol{w} \cdot g(\boldsymbol{W}\boldsymbol{x} + \boldsymbol{b}) + b, \quad \boldsymbol{W} \in \mathbb{R}^{h \times d}, \ \boldsymbol{w}, \boldsymbol{b} \in \mathbb{R}^h, b \in \mathbb{R}$$

this is a matrix of dimensions $h \times d$

this is a non-linear scalar function, applied *elementwise*

Feed-Forward Neural Network

- Approximating a target function

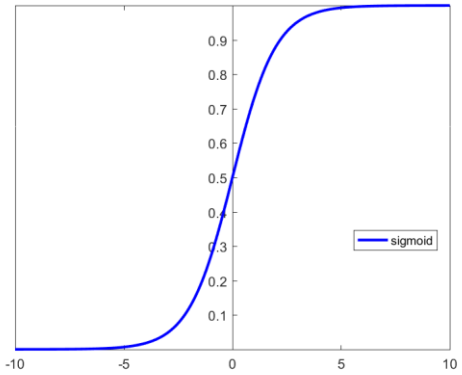
$$y = f^*(\boldsymbol{x}), \quad \boldsymbol{x} \in \mathbb{R}^d$$

Second attempt: ***(shallow) feed-forward neural network***

$$\tilde{y} = \boldsymbol{w} \cdot g(\boldsymbol{W}\boldsymbol{x} + \boldsymbol{b}) + b, \quad \boldsymbol{W} \in \mathbb{R}^{h \times d}, \ \boldsymbol{w}, \boldsymbol{b} \in \mathbb{R}^h, b \in \mathbb{R}$$

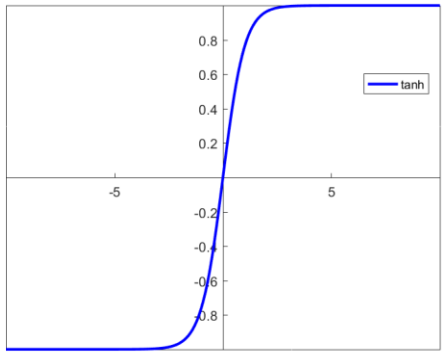
Popular choices for the non-linear function:

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



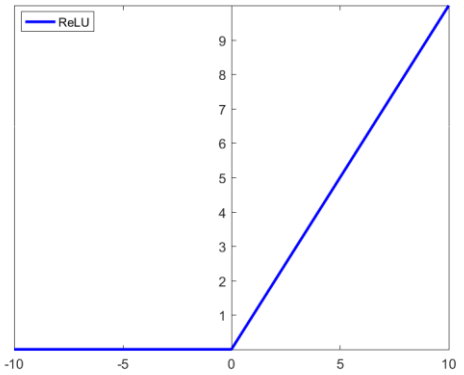
Sigmoid

$$g(x) = \tanh(x)$$



Hyperbolic Tangent

$$g(x) = \max(0, x)$$



Rectified Linear Unit (ReLU)

Feed-Forward Neural Network

- Approximating a target function

$$y = f^*(\boldsymbol{x}), \quad \boldsymbol{x} \in \mathbb{R}^d$$

Second attempt: ***(shallow) feed-forward neural network***

$$\tilde{y} = \boldsymbol{w} \cdot g(\boldsymbol{W}\boldsymbol{x} + \boldsymbol{b}) + b, \quad \boldsymbol{W} \in \mathbb{R}^{h \times d}, \ \boldsymbol{w}, \boldsymbol{b} \in \mathbb{R}^h, b \in \mathbb{R}$$

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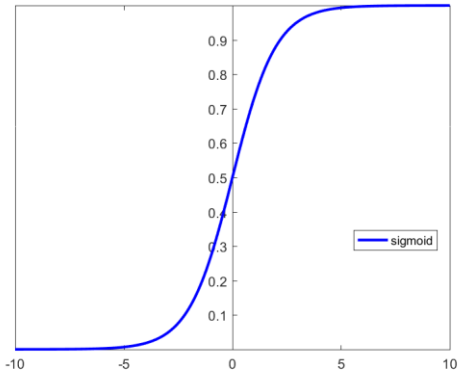

Sigmoid

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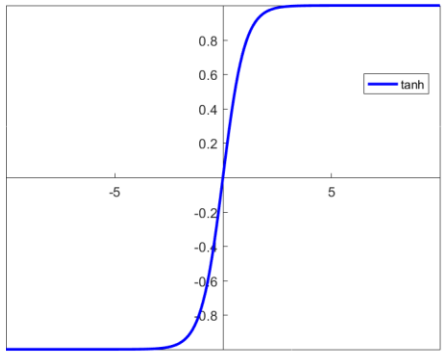
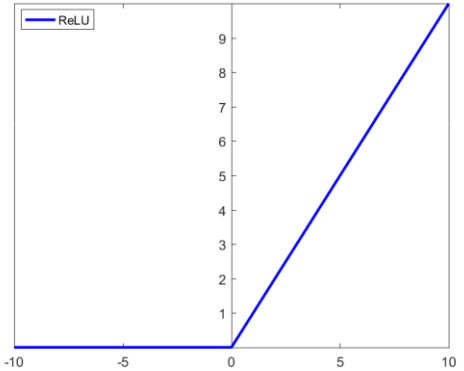


Hyperbolic Tangent

this is somewhat special...

$$g(x) = \max(0, x)$$



Rectified Linear Unit (ReLU)

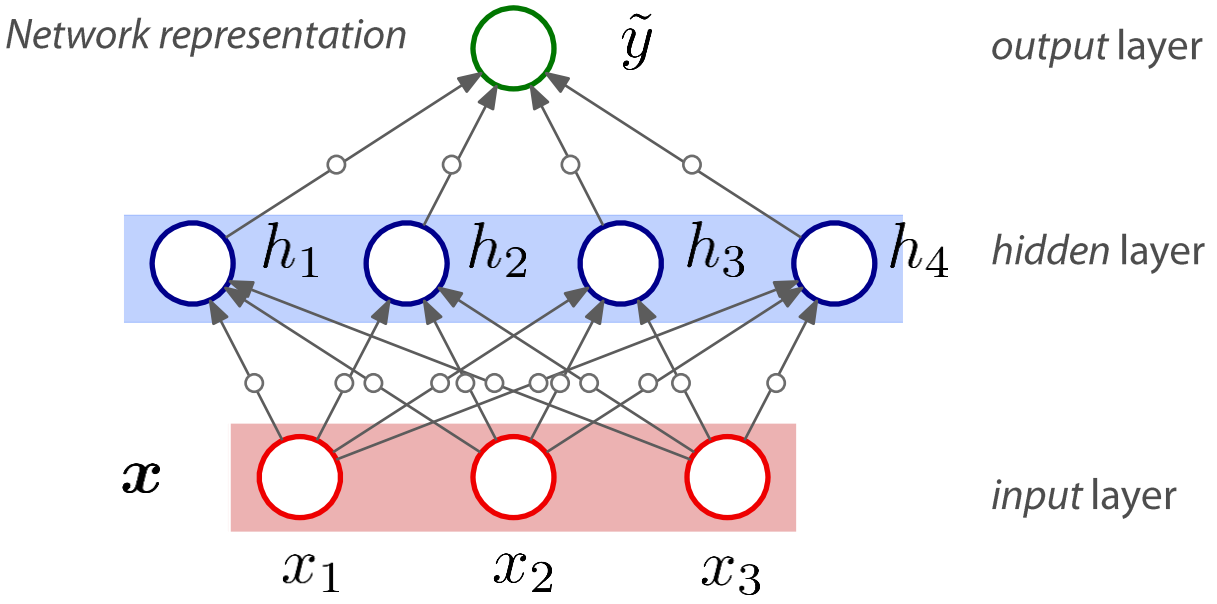
Feed-Forward Neural Network

- Approximating a target function

$$y = f^*(\boldsymbol{x}), \quad \boldsymbol{x} \in \mathbb{R}^d$$

Second attempt: ***(shallow) feed-forward neural network***

$$\tilde{y} = \boldsymbol{w} \cdot g(\boldsymbol{W}\boldsymbol{x} + \boldsymbol{b}) + b, \quad \boldsymbol{W} \in \mathbb{R}^{h \times d}, \ \boldsymbol{w}, \boldsymbol{b} \in \mathbb{R}^h, b \in \mathbb{R}$$



NOTE: *biases* $\boldsymbol{b}$ and b are NOT represented in the graph

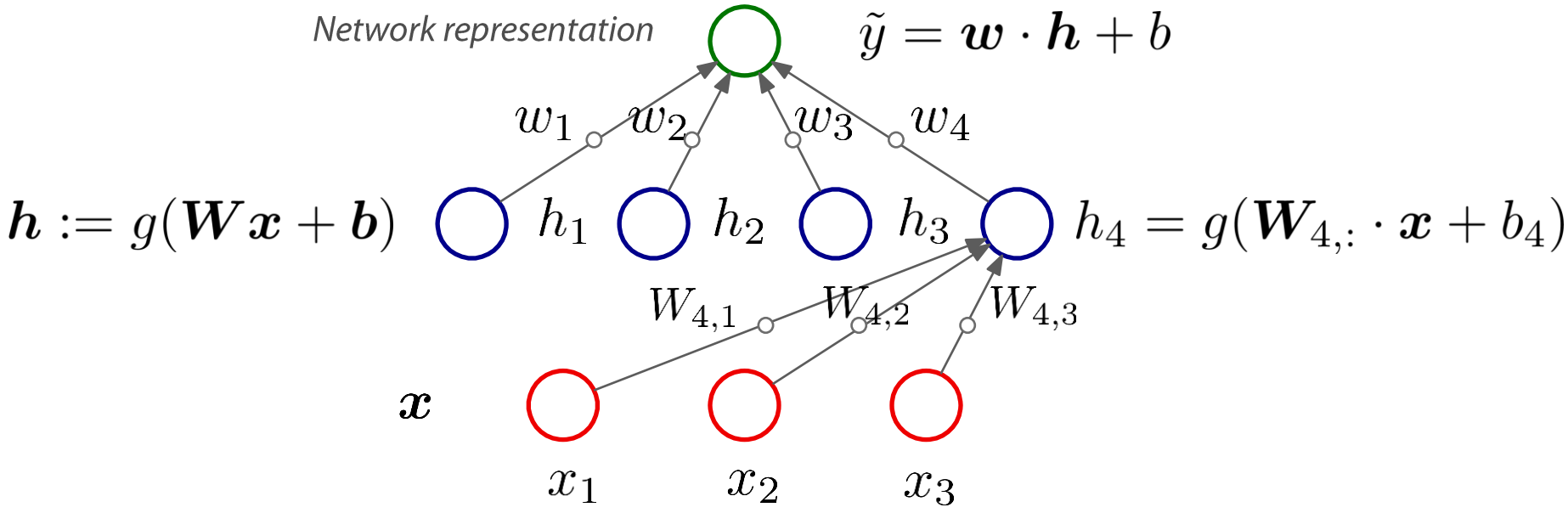
Feed-Forward Neural Network

- Approximating a target function

$$y = f^*(\boldsymbol{x}), \quad \boldsymbol{x} \in \mathbb{R}^d$$

Second attempt: ***(shallow) feed-forward neural network***

$$\tilde{y} = \boldsymbol{w} \cdot g(\boldsymbol{W}\boldsymbol{x} + \boldsymbol{b}) + b, \quad \boldsymbol{W} \in \mathbb{R}^{h \times d}, \ \boldsymbol{w}, \boldsymbol{b} \in \mathbb{R}^h, b \in \mathbb{R}$$



NOTE: *biases* $\boldsymbol{b}$ and b are NOT represented in the graph

Universality of FF Neural Networks

- **Universal approximation theorem** (Cybenko, 1989; Hornik, 1991; Leshno et al. 1991)

  For any target function

  $$y = f^*(\boldsymbol{x}), \quad \boldsymbol{x} \in \mathbb{R}^d$$ *(which is continuous and Borel measurable)*

  and any $\varepsilon > 0$ there exists parameters

  $$h \in \mathbb{Z}^+, \boldsymbol{W} \in \mathbb{R}^{h \times d}, \ \boldsymbol{w}, \boldsymbol{b} \in \mathbb{R}^h, b \in \mathbb{R}$$

  this is the dimension of the hidden layer: it is a parameter in the theorem

  such that the ***(shallow) feed-forward neural network***

  $$\tilde{y} = \boldsymbol{w} \cdot g(\boldsymbol{W}\boldsymbol{x} + \boldsymbol{b}) + b$$

  approximates the target function by less than ε

  $$\sup_{\boldsymbol{x}} \left| f^*(\boldsymbol{x}) - (\boldsymbol{w} \cdot g(\boldsymbol{W}\boldsymbol{x} + \boldsymbol{b}) + b) \right| < \varepsilon$$

  (on any compact subset of $\mathbb{R}^d$)

  This theorem holds with any of the non-linear functions seen before

Universality of FF Neural Networks

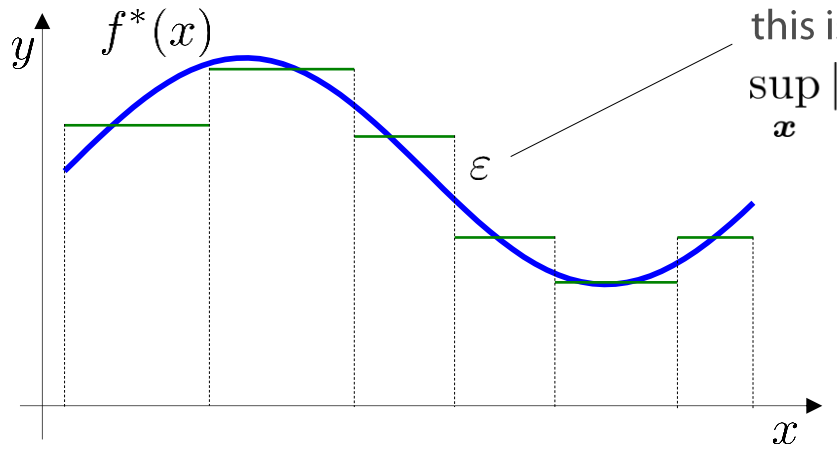
- **Universal approximation theorem** (Cybenko, 1989; Hornik, 1991; Leshno et al. 1991)

  Intuitive rationale

  Any continuous target function

$$y = f^*(x), \quad x \in \mathbb{R}$$

  can be approximated arbitrarily well by a stepwise function



this is the largest difference

$$\sup_{\boldsymbol{x}} | f^*(\boldsymbol{x}) - (\boldsymbol{w} \cdot g(\boldsymbol{W}\boldsymbol{x} + \boldsymbol{b}) + b) |$$

  for simplicity, assume now that x is *scalar* (hence $\boldsymbol{W}$ is *vector*)

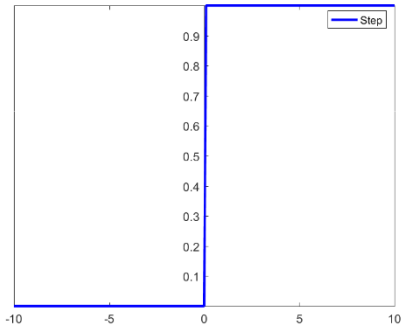
$$\tilde{y} = \boldsymbol{w} \cdot g(\boldsymbol{W}x + \boldsymbol{b}) + b$$

Universality of FF Neural Networks

- **Universal approximation theorem** (Cybenko, 1989; Hornik, 1991; Leshno et al. 1991)

  Intuitive rationale

  $g(x) = \text{step}(x)$

  Consider the *step function* as the non-linearity;

  $$\tilde{y} = \boldsymbol{w} \cdot \text{step}(\boldsymbol{W}x + \boldsymbol{b}) + b$$

  then, by expanding the scalar product:

  $$\tilde{y} = w_1 \,\text{step}(W_1 x + b_1) + \cdots + w_h \,\text{step}(W_h x + b_h) + b$$

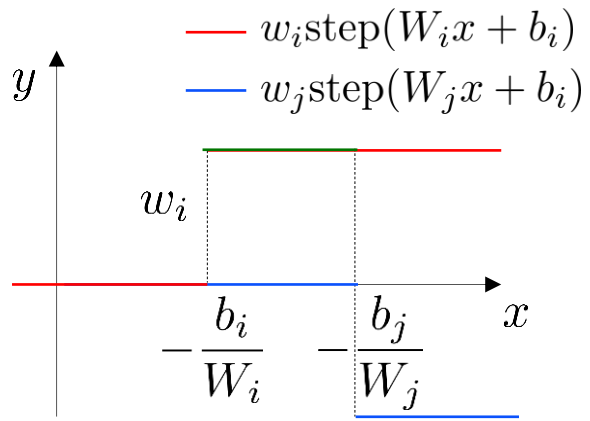
  where each step occurs at

  $$W_i \cdot x + b_i = 0 \quad \Longrightarrow \quad W_i \cdot x = -b_i \quad \Longrightarrow \quad x = -\frac{b_i}{W_i}$$

  Consider *pairs* of steps i and j and impose

  $$-\frac{b_i}{W_i} < -\frac{b_j}{W_j}, \quad W_i, W_j > 0, \quad w_i = -w_j$$

  in this way we can construct $\frac{h}{2}$ *such function steps*

Learning with Feed-Forward Neural Networks

Learning with FF Neural Networks

- Approximating a target function

$$y = f^*(\boldsymbol{x}), \quad \boldsymbol{x} \in \mathbb{R}^d$$

Second attempt: ***(shallow) feed-forward neural network***

$$\tilde{y} = \boldsymbol{w} \cdot g(\boldsymbol{W}\boldsymbol{x} + \boldsymbol{b}) + b, \quad \boldsymbol{W} \in \mathbb{R}^{h \times d}, \ \boldsymbol{w}, \boldsymbol{b} \in \mathbb{R}^h, b \in \mathbb{R}$$

Optimization problem (learning)

Given a dataset $D := \{(\boldsymbol{x}^{(i)}, y^{(i)})\}_{i=1}^N, \quad y^{(i)} = f^*(\boldsymbol{x}^{(i)}), \ \forall i$

the dimension of the hidden layer is pre-defined

we want to find parameter values $\boldsymbol{W} \in \mathbb{R}^{h \times d}, \ \boldsymbol{w}, \boldsymbol{b} \in \mathbb{R}^h, b \in \mathbb{R}$

that *minimize* the loss function $L(D) := \frac{1}{N} \sum_D (\tilde{y}^{(i)} - y^{(i)})^2$

where: $\tilde{y}^{(i)} := \boldsymbol{w} \cdot g(\boldsymbol{W}\boldsymbol{x}^{(i)} + \boldsymbol{b}) + b$

Learning with FF Neural Networks

- Approximating a target function

$$y = f^*(\boldsymbol{x}), \quad \boldsymbol{x} \in \mathbb{R}^d$$

Second attempt: ***(shallow) feed-forward neural network***

$$\tilde{y} = \boldsymbol{w} \cdot g(\boldsymbol{W}\boldsymbol{x} + \boldsymbol{b}) + b, \quad \boldsymbol{W} \in \mathbb{R}^{h \times d}, \ \boldsymbol{w}, \boldsymbol{b} \in \mathbb{R}^h, b \in \mathbb{R}$$

Difficulty

In general, *minimizing* the loss function

$$L(D) = \frac{1}{N} \sum_{D} ((\boldsymbol{w} \cdot g(\boldsymbol{W}\boldsymbol{x}^{(i)} + \boldsymbol{b}) + b) - y^{(i)})^2$$

this loss function is <u>not</u> convex, in general

cannot be done directly, since

$$\frac{\partial}{\partial \boldsymbol{\vartheta}} L(D) = 0$$

cannot be solved analytically

We need to find another way…

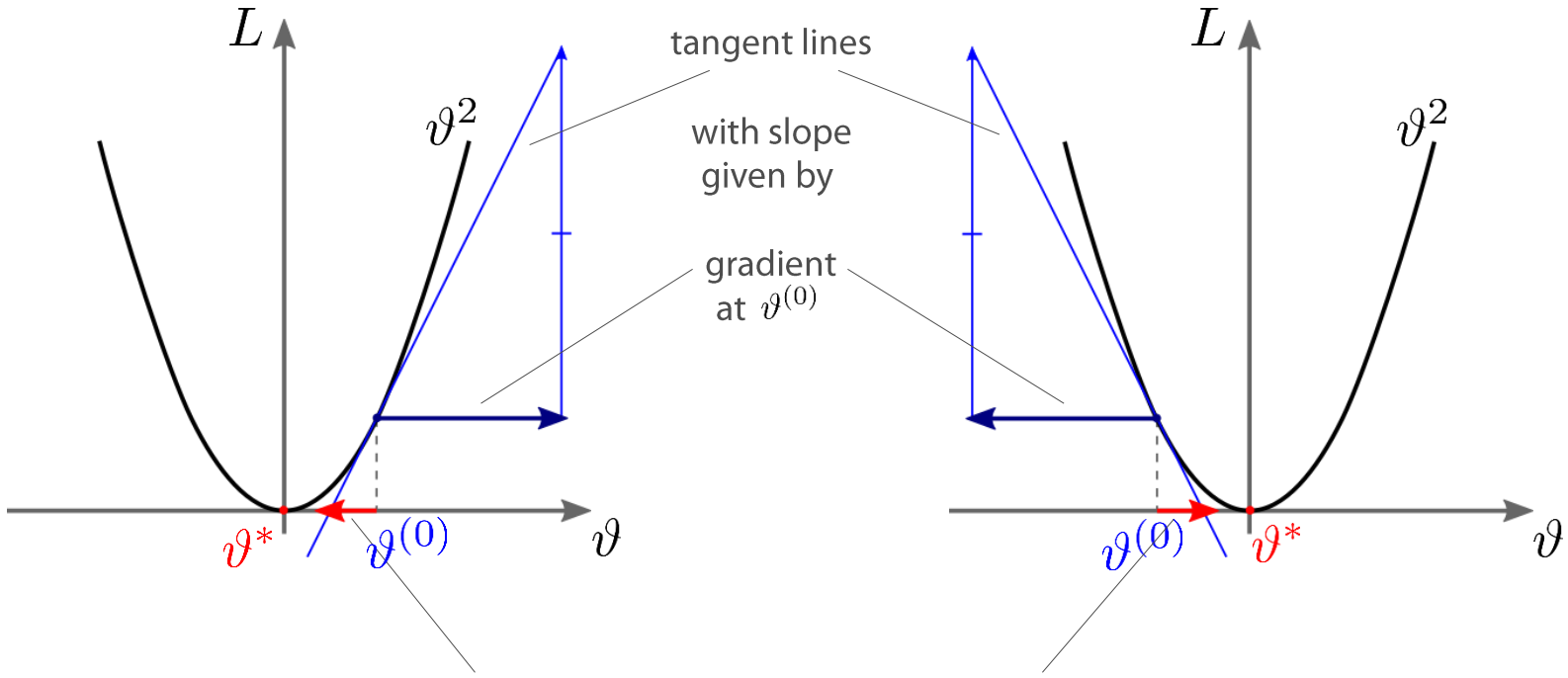
Gradient Descent (GD): intuition

- *Optimization problem*

$$\boldsymbol{\vartheta}^* := \text{argmin}_{\boldsymbol{\vartheta}} \; L(D, \boldsymbol{\vartheta})$$

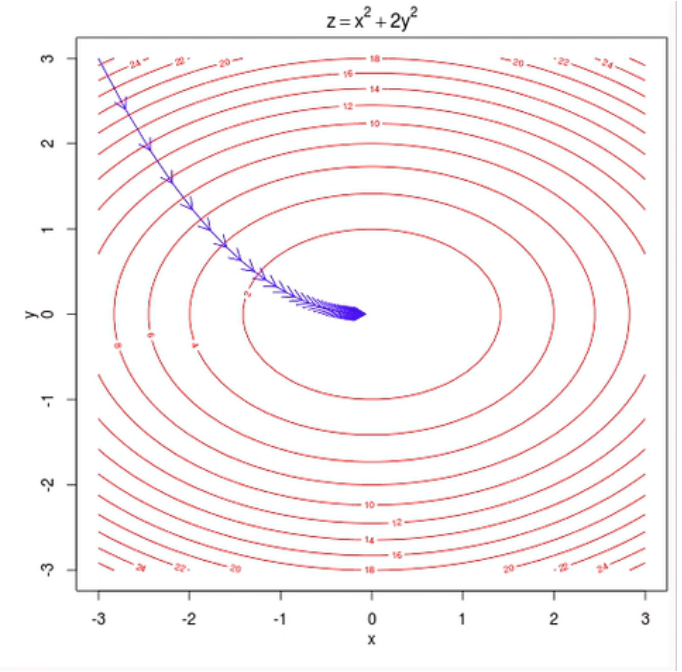
Just making the dependence explicit

- *Minimizing a generic function*



Follow the opposite of the gradient!

Gradient Descent (GD): intuition



- *Optimization problem*

$$\boldsymbol{\vartheta}^* := \text{argmin}_{\boldsymbol{\vartheta}} \; L(D, \boldsymbol{\vartheta})$$

Just making the dependence explicit

- *Iterative method*

Step in the method

1. Initialize $\boldsymbol{\vartheta}^{(0)}$ at random
2. Update $\boldsymbol{\vartheta}^{(t)} = \boldsymbol{\vartheta}^{(t-1)} - \eta \; \dfrac{\partial}{\partial \boldsymbol{\vartheta}} L(D, \boldsymbol{\vartheta}^{(t-1)})$
3. Unless some termination criterion has been met, go back to step 2.

where

$$\frac{\partial}{\partial \boldsymbol{\vartheta}} L(D, \boldsymbol{\vartheta}) := \frac{1}{N} \sum_{D} \frac{\partial}{\partial \boldsymbol{\vartheta}} L(\hat{y}^{(i)}, y^{(i)}, \boldsymbol{\vartheta})$$

The gradient of the loss over the dataset D is the average of gradients over each data item

$$\eta \ll 1$$

A *learning rate*, it is arbitrary (i.e., an *hyperparameter*)

Gradient Descent (GD): convergence

- *Convergence*

  When $L(D, \boldsymbol{\vartheta})$ is *convex*, *derivable,* and its gradient is *Lipschitz continuous*, that is

$$\left\| \frac{\partial}{\partial \boldsymbol{\vartheta}} L(D, \boldsymbol{\vartheta}_1) - \frac{\partial}{\partial \boldsymbol{\vartheta}} L(D, \boldsymbol{\vartheta}_2) \right\| \leq C \, \|\boldsymbol{\vartheta}_1 - \boldsymbol{\vartheta}_2\|, \quad C > 0$$

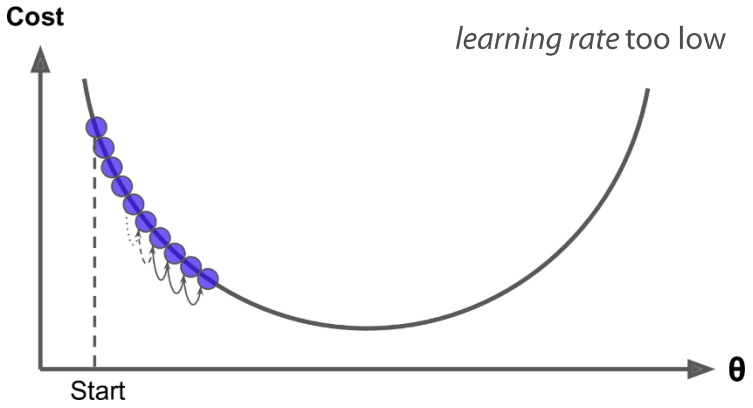
  the gradient descent method converges to the optimal $\boldsymbol{\vartheta}^*$ for $t \to \infty$
  provided that $\eta \leq 1/C$

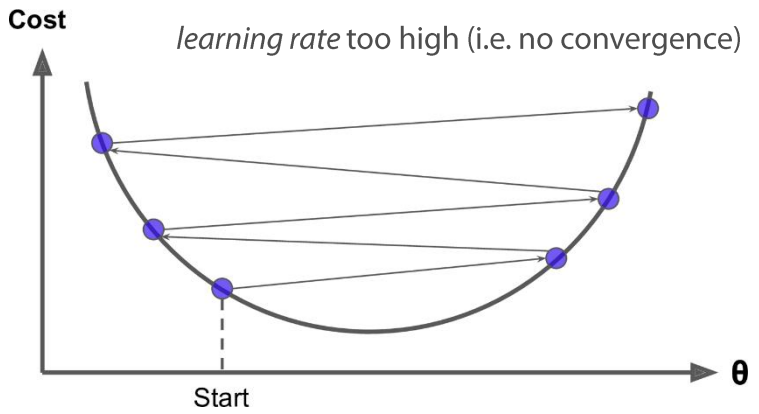
  When $L(D, \boldsymbol{\vartheta})$ is *derivable* but <u>not</u> *convex,* and its gradient is *Lipschitz continuous,*
  the gradient descent method converges to a <u>local minimum</u> of $L(D, \boldsymbol{\vartheta})$
  under the same conditions

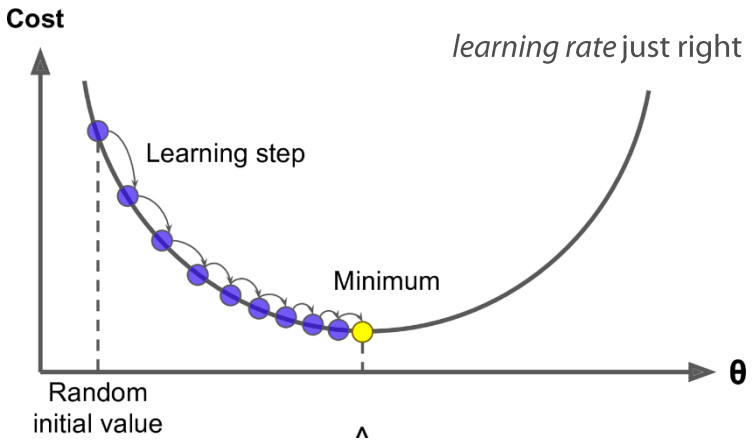
Gradient Descent (GD): practicalities

- *Convergence in practice*

  The choice of the *learning rate* η is crucial





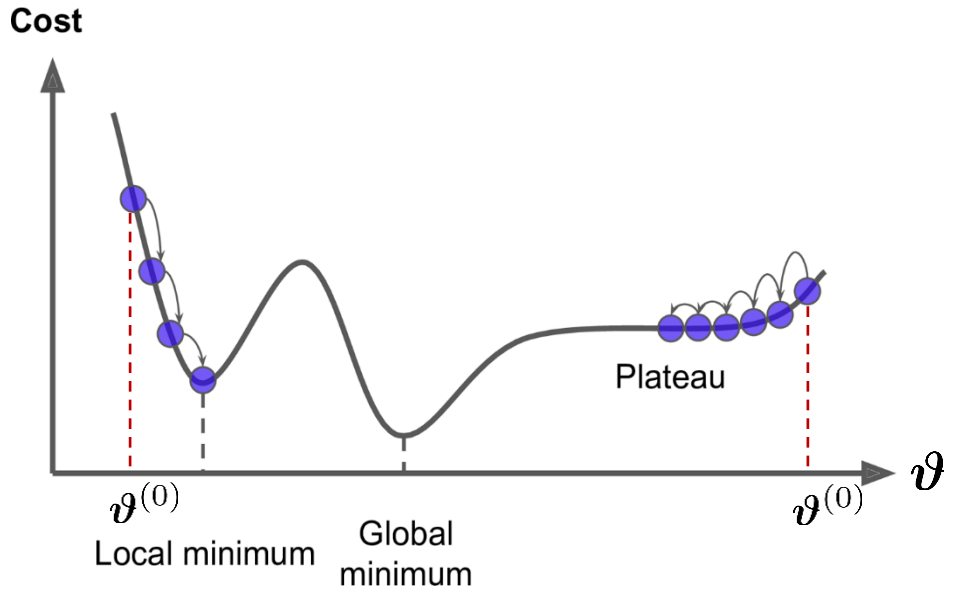


Images from https://www.safaribooksonline.com/library/view/hands-on-machine-learning/9781491962282/ch04.html

Gradient Descent (GD): practicalities

- *Convergence in practice*

  When $L(D, \boldsymbol{\vartheta})$ is <u>not</u> convex, the **initial estimate** $\boldsymbol{\vartheta}^{(0)}$ is crucial



The outcome of the method will depend on which $\boldsymbol{\vartheta}^{(0)}$ *is picked*

Image from https://www.safaribooksonline.com/library/view/hands-on-machine-learning/9781491962282/ch04.html

Learning Feed-Forward Neural Networks (contd.)

Gradient Descent for FF Neural Networks

Recall that the *item-wise* loss for a specific data item in the dataset is

$$L(\tilde{y}^{(i)}, y^{(i)}) := (\tilde{y}^{(i)} - y^{(i)})^2$$

then

$$L(D) = \frac{1}{N} \sum_{D} L(\tilde{y}^{(i)}, y^{(i)})$$

and the gradient of the loss function is

$$\frac{\partial}{\partial \boldsymbol{\vartheta}} L(D) = \frac{\partial}{\partial \boldsymbol{\vartheta}} \frac{1}{N} \sum_{D} L(\tilde{y}^{(i)}, y^{(i)})$$

$$= \frac{1}{N} \sum_{D} \frac{\partial}{\partial \boldsymbol{\vartheta}} L(\tilde{y}^{(i)}, y^{(i)})$$

Moral: we need to compute the gradient on <u>each data item</u>

Gradient Descent for FF Neural Networks

Suppose we can compute the four *item-wise gradients*, w.r.t. to the parameters:

$$\frac{\partial}{\partial \boldsymbol{W}} L(\tilde{y}^{(i)}, y^{(i)}) \qquad \frac{\partial}{\partial \boldsymbol{b}} L(\tilde{y}^{(i)}, y^{(i)}) \qquad \frac{\partial}{\partial \boldsymbol{w}} L(\tilde{y}^{(i)}, y^{(i)}) \qquad \frac{\partial}{\partial b} L(\tilde{y}^{(i)}, y^{(i)})$$

then we can apply a *gradient descent* method

- ***Gradient Descent***
  1. Assign initial values to the four parameters $\boldsymbol{W}^{(0)},\ \boldsymbol{b}^{(0)},\ \boldsymbol{w}^{(0)},\ b^{(0)}$
  2. Update the four parameters by adding

$$\Delta \boldsymbol{W} = -\eta \frac{1}{N} \sum_{D} \frac{\partial}{\partial \boldsymbol{W}} L(\tilde{y}^{(i)}, y^{(i)}) \qquad \Delta \boldsymbol{b} = -\eta \frac{1}{N} \sum_{D} \frac{\partial}{\partial \boldsymbol{b}} L(\tilde{y}^{(i)}, y^{(i)})$$

$$\Delta \boldsymbol{w} = -\eta \frac{1}{N} \sum_{D} \frac{\partial}{\partial \boldsymbol{w}} L(\tilde{y}^{(i)}, y^{(i)}) \qquad \Delta b = -\eta \frac{1}{N} \sum_{D} \frac{\partial}{\partial b} L(\tilde{y}^{(i)}, y^{(i)})$$

  3. Unless complete, return to step 2.

Computing Gradients

All we need to apply the descent method is computing the item-wise gradients

For instance:

$$\frac{\partial}{\partial \boldsymbol{W}} L(\tilde{y}^{(i)}, y^{(i)}) = \frac{\partial}{\partial \boldsymbol{W}} (\tilde{y}^{(i)} - y^{(i)})^2$$

$$= \frac{\partial}{\partial \boldsymbol{W}} ((\boldsymbol{w} \cdot g(\boldsymbol{W}\boldsymbol{x}^{(i)} + \boldsymbol{b}) + b) - y^{(i)})^2$$

(similar expressions hold for the other three gradients)

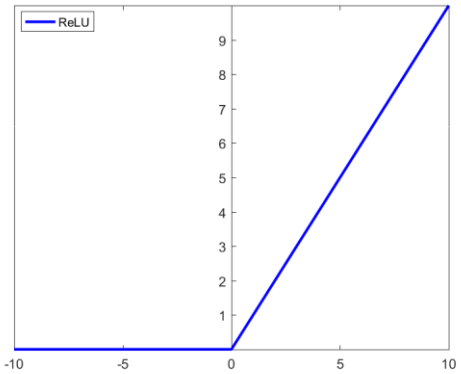
Assume

$$g(x) = \mathrm{ReLU}(x) := \max(0, x)$$

i.e., the non-linearity is ReLU

Easy, huh?

$$g(x) = \max(0, x)$$



Function Approximation: FF Neural Networks

- Loss minimization

  Approximator:
  (shallow) feed-forward neural network

$$\tilde{y} = \boldsymbol{w} \cdot \mathrm{ReLU}(\boldsymbol{W}\boldsymbol{x} + \boldsymbol{b}) + b$$

  Optimal values for XOR and $h = 2$

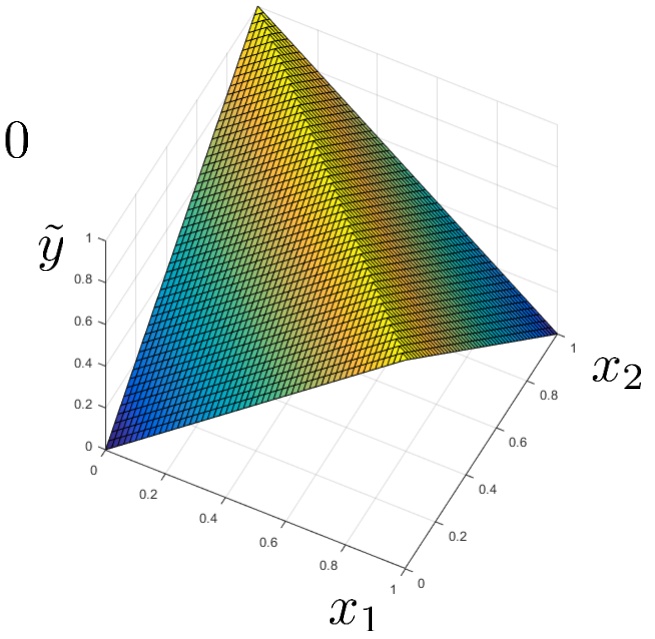
  dimension of the hidden layer

$$\boldsymbol{W} = \begin{bmatrix} 1 & 1 \\ 1 & 1 \end{bmatrix} \quad \boldsymbol{b} = \begin{bmatrix} 0 \\ -1 \end{bmatrix} \quad \boldsymbol{w} = \begin{bmatrix} 1 \\ -2 \end{bmatrix}$$

XOR

x_1	x_2	$x_1 \oplus x_2$
0	0	0
0	1	1
1	0	1
1	1	0

$b = 0$



Stochastic and Mini-Batch Gradient Descent

Function Approximation: FF Neural Networks

- Loss minimization

  Approximator:
  (shallow) feed-forward neural network

$$\tilde{y} = \boldsymbol{w} \cdot \mathrm{ReLU}(\boldsymbol{W}\boldsymbol{x} + \boldsymbol{b}) + b$$

  In this case our dataset was tiny… ($N = 4$)

  What if the dataset was <u>very</u> large?

XOR

x_1	x_2	$x_1 \oplus x_2$
0	0	0
0	1	1
1	0	1
1	1	0

this is our *dataset*

Stochastic Gradient Descent (SGD): intuition

- *Objective*

$$\boldsymbol{\vartheta}^* := \text{argmin}_{\boldsymbol{\vartheta}} \; L(D, \boldsymbol{\vartheta})$$

- *Iterative method*

1. Initialize $\boldsymbol{\vartheta}^{(0)}$ at random
2. Pick a data item $(\boldsymbol{x}^{(i)}, y^{(i)}) \in D$ with uniform probability
3. Update $\boldsymbol{\vartheta}^{(t)} = \boldsymbol{\vartheta}^{(t-1)} - \eta^{(t)} \; \dfrac{\partial}{\partial \boldsymbol{\vartheta}} L(\tilde{y}^{(i)}, y^{(i)}, \boldsymbol{\vartheta}^{(t-1)})$
4. Unless some termination criterion has been met, go back to step 2.

$$\eta^{(t)} \ll 1$$

Note that the *learning rate* may *vary* across iterations…

Stochastic Gradient Descent for FF Neural Networks

With very large datasets, the sum in:

$$\Delta\boldsymbol{\vartheta} = -\eta \frac{1}{N} \sum_{D} \frac{\partial}{\partial \boldsymbol{\vartheta}} L(\tilde{y}^{(i)}, y^{(i)})$$

may take very long to compute (and this must be repeated at each iteration)

- ***Stochastic* Gradient Descent (SGD)** (i.e. *"you don't actually need to sum up them all"*)
  1. Assign initial values to the four parameters $\boldsymbol{W}^{(0)},\ \boldsymbol{b}^{(0)},\ \boldsymbol{w}^{(0)},\ b^{(0)}$
  2. Pick up a data item $(\boldsymbol{x}^{(i)}, y^{(i)})$ from D with uniform probability and update the four parameters (with $\eta \ll 1.0, \quad \eta \to 0$ as iterations progress)

$$\Delta\boldsymbol{W} = -\eta \frac{\partial}{\partial \boldsymbol{W}} L(\tilde{y}^{(i)}, y^{(i)}) \qquad \Delta\boldsymbol{b} = -\eta \frac{\partial}{\partial \boldsymbol{b}} L(\tilde{y}^{(i)}, y^{(i)})$$

$$\Delta\boldsymbol{w} = -\eta \frac{\partial}{\partial \boldsymbol{w}} L(\tilde{y}^{(i)}, y^{(i)}) \qquad \Delta b = -\eta \frac{\partial}{\partial b} L(\tilde{y}^{(i)}, y^{(i)})$$

  3. Unless complete, return to step 2.

Stochastic Gradient Descent (SGD): convergence

- *Convergence*

  When $L(D, \boldsymbol{\vartheta})$ is *convex*, *derivable,* and its gradient is *Lipschitz continuous*, that is

$$\left\| \frac{\partial}{\partial \boldsymbol{\vartheta}} L(D, \boldsymbol{\vartheta}_1) - \frac{\partial}{\partial \boldsymbol{\vartheta}} L(D, \boldsymbol{\vartheta}_2) \right\| \leq C \, \|\boldsymbol{\vartheta}_1 - \boldsymbol{\vartheta}_2\|, \quad C > 0$$

  the stochastic gradient descent method converges to the optimal $\boldsymbol{\vartheta}^*$ for $t \to \infty$ provided that

$$\eta^{(t)} \leq \frac{1}{Ct}$$

  Note that $\eta^{(t)} \to 0$ for $t \to \infty$

  When $L(D, \boldsymbol{\vartheta})$ is *derivable,* and its gradient is *Lipschitz continuous* but not *convex* the stochastic gradient descent method converges to a local minimum of $L(D, \boldsymbol{\vartheta})$ under the same conditions

Speed of Convergence

*Perhaps surprisingly, **stochastic gradient descent** shares the same properties and could be faster than GD …*

Consider a generic loss function $L(\boldsymbol{\vartheta})$ which is *convex* in the parameter $\boldsymbol{\vartheta}$

Define *accuracy* as an upper bound:

optimal value

current parameter estimate

$$\left|L(\boldsymbol{\vartheta}^*) - L(\tilde{\boldsymbol{\vartheta}})\right| < \rho$$

[*Bottou & Bousquet, 2008*]

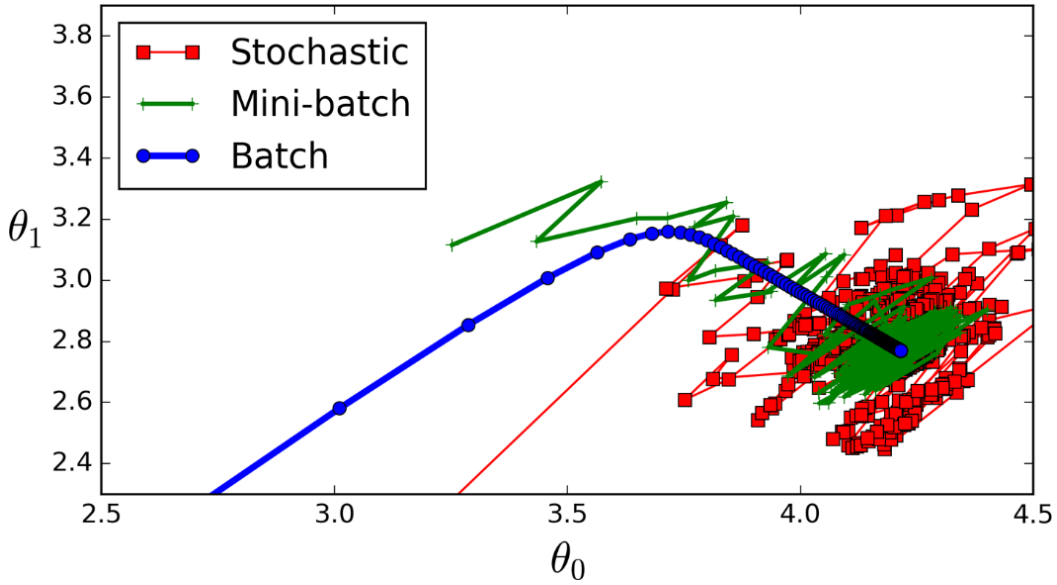
N size of the dataset

q number of (scalar) parameters in $\boldsymbol{\vartheta}$

Algorithm	***Cost per iteration***	***Iterations to reach accuracy ρ***	***Time to reach accuracy ρ***
Gradient descent (GD)	$\mathcal{O}(N\,q)$	$\mathcal{O}\left(\log\frac{1}{\rho}\right)$	$\mathcal{O}\left(N\,q\log\frac{1}{\rho}\right)$
Stochastic gradient descent (SGD)	$\mathcal{O}(q)$	$\mathcal{O}\left(\frac{1}{\rho}\right)$	$\mathcal{O}\left(q\frac{1}{\rho}\right)$

Qualitative comparison of GD methods

Typical traces
of the three methods
(batch = GD)



In general:

- GD is more regular but slower (with large datasets)
- SGD is faster (with large datasets) but noisy
- MBGD is often the right compromise in practice…

Image from https://www.safaribooksonline.com/library/view/hands-on-machine-learning/9781491962282/ch04.html

Mini-batch Gradient Descent (MBGD): intuition

- *Objective*

$$\boldsymbol{\vartheta}^* := \text{argmin}_{\boldsymbol{\vartheta}} \; L(D, \boldsymbol{\vartheta})$$

- *Iterative method*

1. Initialize $\theta^{(0)}$ at random
2. Pick a mini batch $B \subseteq D$ with uniform probability
3. Update $\boldsymbol{\vartheta}^{(t)} = \boldsymbol{\vartheta}^{(t-1)} - \eta^{(t)} \frac{\partial}{\partial \boldsymbol{\vartheta}} L(B, \boldsymbol{\vartheta}^{(t-1)})$
4. Unless some termination criterion has been met, go back to step 2.

where:

$$\frac{\partial}{\partial \boldsymbol{\vartheta}} L(B, \boldsymbol{\vartheta}) := \frac{1}{|B|} \sum_{B} \frac{\partial}{\partial \boldsymbol{\vartheta}} L(\hat{y}^{(i)}, y^{(i)}, \boldsymbol{\vartheta})$$

This method has the same convergence properties of SGD

Mini-batch Gradient Descent for FF Neural Networks

- ***Mini-batch* Gradient Descent (MBGD)**
  1. Assign initial values to the four parameters $\boldsymbol{W}^{(0)},\ \boldsymbol{b}^{(0)},\ \boldsymbol{w}^{(0)},\ b^{(0)}$
  2. Pick a *mini-batch* $B \subseteq D$ with uniform probability
     and update the four parameters (with $\eta \ll 1.0, \quad \eta \to 0$ as iterations progress)

$$\Delta \boldsymbol{W} = -\eta \frac{1}{|B|} \sum_{B} \frac{\partial}{\partial \boldsymbol{W}} L(\tilde{y}^{(i)}, y^{(i)}) \qquad \Delta \boldsymbol{b} = -\eta \frac{1}{|B|} \sum_{B} \frac{\partial}{\partial \boldsymbol{b}} L(\tilde{y}^{(i)}, y^{(i)})$$

$$\Delta \boldsymbol{w} = -\eta \frac{1}{|B|} \sum_{B} \frac{\partial}{\partial \boldsymbol{w}} L(\tilde{y}^{(i)}, y^{(i)}) \qquad \Delta b = -\eta \frac{1}{|B|} \sum_{B} \frac{\partial}{\partial b} L(\tilde{y}^{(i)}, y^{(i)})$$

  3. Unless complete, return to step 2.

This method has the same convergence properties of SGD