

# **AI6102: Machine Learning Methodologies & Applications**

## **L3: Linear Models: Regression**

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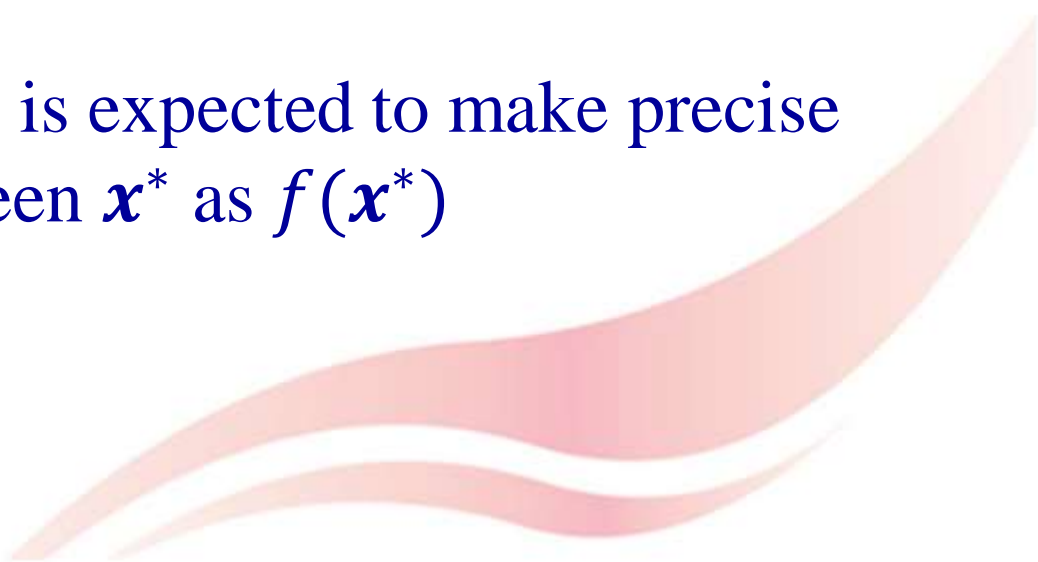
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Homepage: <https://mreallab.github.io/>

# Recall: Supervised Learning

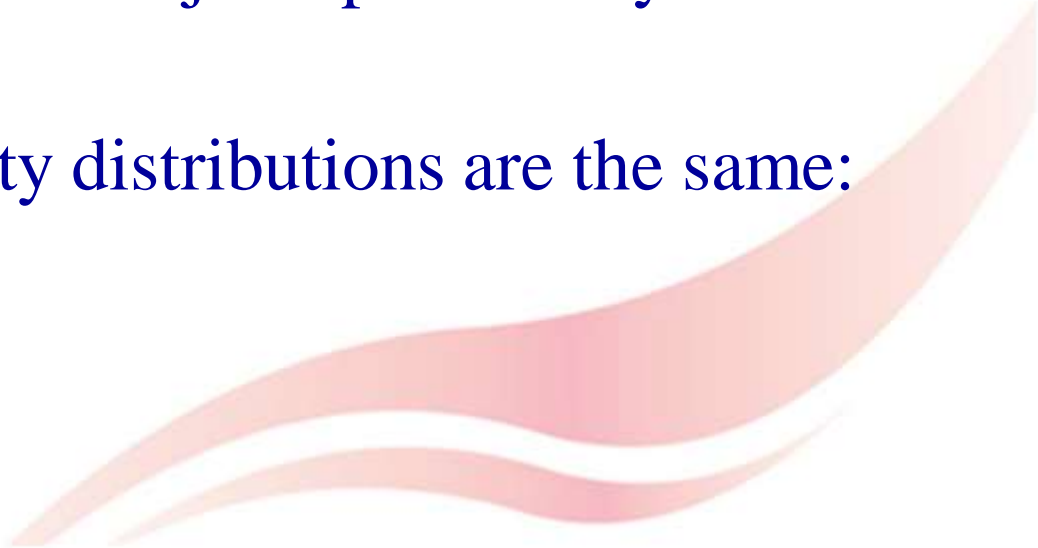
In mathematics

- Given: a set of  $N$  labeled data  $\{(\mathbf{x}_1, y_1), \dots, (\mathbf{x}_N, y_N)\}$ , where  $\mathbf{x}_i$  is  $m$ -dimensional vector of numerical values, and  $y_i$  is a scalar
  - We aim to learn a mapping  $f: \mathbf{x} \rightarrow y$  by requiring  $f(\mathbf{x}_i) = y_i$
  - The learned mapping  $f$  is expected to make precise predictions on any unseen  $\mathbf{x}^*$  as  $f(\mathbf{x}^*)$
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# Hypothesis

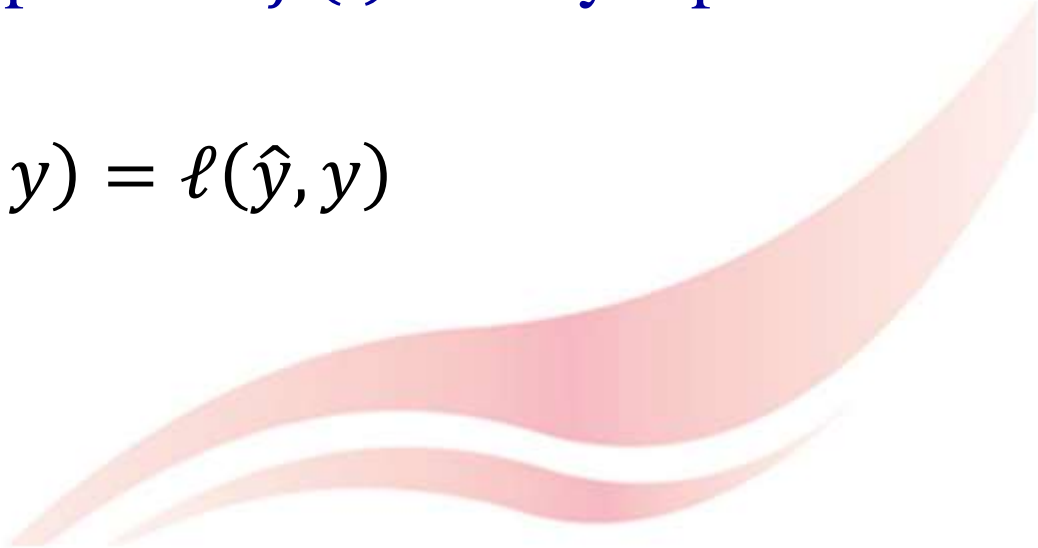
- A mapping or function  $f: \mathbf{x} \rightarrow y$  can be considered as an element of some space of possible functions  $\mathcal{H}: \mathbb{R}^m \rightarrow \mathbb{R}$ , often called hypothesis space
  - Supervised learning aims to find a hypothesis  $f \in \mathcal{H}$  from training data  $\{(\mathbf{x}_1, y_1), \dots, (\mathbf{x}_N, y_N)\}$ , s.t. for any test  $\mathbf{x}^*$ ,  $f(\mathbf{x}^*) = y^*$
- 

# Assumption

- The training data instances  $(\mathbf{x}_1, y_1), \dots, (\mathbf{x}_N, y_N)$ , are independent and identically distributed (i.i.d.), and drawn from an unknown joint probability distribution  $P_{tr}(\mathbf{x}, y)$   
Independent -> probability definition  
Identical ->
  - Unseen test data instances  $\{(\mathbf{x}^*, y^*)\}$  are also i.i.d, and drawn from an unknown joint probability distribution  $P_{ts}(\mathbf{x}, y)$
  - The two joint probability distributions are the same:  
 $P_{tr}(\mathbf{x}, y) = P_{ts}(\mathbf{x}, y)$
- 

# Loss Function

- Denote by  $\hat{y} = f(\mathbf{x})$  the prediction of the function  $f(\cdot)$  on a data instance  $\mathbf{x}$ , and  $y$  is the ground-truth output of  $\mathbf{x}$
- Let  $\ell: \mathbb{R} \times \mathbb{R} \rightarrow \mathbb{R}_+ \geq 0$  be a loss function to measure the difference between the ground-truth output  $y$  and the prediction  $\hat{y}$  of a hypothesis  $f(\cdot)$  on any input data instance  $\mathbf{x}$

$$\ell(f(\mathbf{x}), y) = \ell(\hat{y}, y)$$


# Risk Minimization

- The risk associated with a hypothesis  $f(\cdot)$  is defined as the expectation of the loss function over all possible input-output pairs drawn from a joint probabilistic distribution  $P(\mathbf{x}, y)$ :

$$R(f) = \mathbb{E}_{(\mathbf{x}, y) \sim P} [\ell(f(\mathbf{x}), y)]$$

- Recall: in supervised learning, the learned hypothesis  $f(\cdot)$  is expected to make precise predictions on any test data instance  $\mathbf{x}^*$ , i.e.,  $f(\mathbf{x}^*) = y^*$

$$f^* = \arg \min_f R_{ts}(f) = \arg \min_f \mathbb{E}_{(\mathbf{x}, y) \sim P_{ts}} [\ell(f(\mathbf{x}), y)]$$



Test data is unseen in training! And even in the test phase,  $y$  is not observed!

# Risk Minimization (cont.)

$$f^* = \arg \min_f R_{ts}(f)$$

$$= \arg \min_f \mathbb{E}_{(\mathbf{x}, y) \sim P_{ts}} [\ell(f(\mathbf{x}), y)]$$

$$= \arg \min_f \mathbb{E}_{(\mathbf{x}, y) \sim P_{ts}} \left[ \frac{P_{tr}(\mathbf{x}, y)}{P_{tr}(\mathbf{x}, y)} \ell(f(\mathbf{x}), y) \right]$$


$$= \arg \min_f \int_y \int_{\mathbf{x}} P_{ts}(\mathbf{x}, y) \left( \frac{P_{tr}(\mathbf{x}, y)}{P_{tr}(\mathbf{x}, y)} \ell(f(\mathbf{x}), y) \right) d\mathbf{x} dy$$

Definition of expectation

$$\mathbb{E}_{\mathbf{x} \sim P}[g(\mathbf{x})] = \int_{\mathbf{x}} P(\mathbf{x}) g(\mathbf{x}) d\mathbf{x} = \int_{x_1} \dots \int_{x_m} P(\mathbf{x}) g(\mathbf{x}) dx_1 \dots dx_m$$

# Risk Minimization (cont.)

$$f^* = \arg \min_f R_{ts}(f)$$

$$= \arg \min_f \int_y \int_x \boxed{P_{ts}(\mathbf{x}, y)} \left( \boxed{\frac{P_{tr}(\mathbf{x}, y)}{P_{tr}(\mathbf{x}, y)}} \ell(f(\mathbf{x}), y) \right) d\mathbf{x} dy$$


$$= \arg \min_f \int_y \int_x P_{tr}(\mathbf{x}, y) \left( \boxed{\frac{P_{ts}(\mathbf{x}, y)}{P_{tr}(\mathbf{x}, y)}} \ell(f(\mathbf{x}), y) \right) d\mathbf{x} dy$$

$= 1$  Assumption  
 $P_{tr}(\mathbf{x}, y) = P_{ts}(\mathbf{x}, y)$

$$= \arg \min_f \int_y \int_x P_{tr}(\mathbf{x}, y) \ell(f(\mathbf{x}), y) d\mathbf{x} dy$$

Definition of expectation

$$= \arg \min_f \mathbb{E}_{(\mathbf{x}, y) \sim P_{tr}} [\ell(f(\mathbf{x}), y)]$$



# Empirical Risk Minimization

$$f^* = \arg \min_f \mathbb{E}_{(\mathbf{x}, y) \sim P_{tr}} [\ell(f(\mathbf{x}), y)] = \arg \min_f R_{tr}(f)$$

- The distribution  $P_{tr}(\mathbf{x}, y)$  is unknown, thus we are not able to sample (infinite) input-output pairs  $\{(\mathbf{x}, y)\}$  to learn a hypothesis  $f(\cdot)$  !
- In practice, only a finite number of training pairs are available,  $(\mathbf{x}_1, y_1), \dots, (\mathbf{x}_N, y_N)$
- Approximate the expected risk by empirical risk:


$$\mathbb{E}_{(\mathbf{x}, y) \sim P_{tr}} [\ell(f(\mathbf{x}), y)] \approx \frac{1}{N} \sum_{i=1}^N \ell(f(\mathbf{x}_i), y_i) = \hat{R}_{tr}(f)$$

# Empirical Risk Minimization (cont.)

- In practice, the hypothesis  $f$  can be learned by minimizing the empirical risk

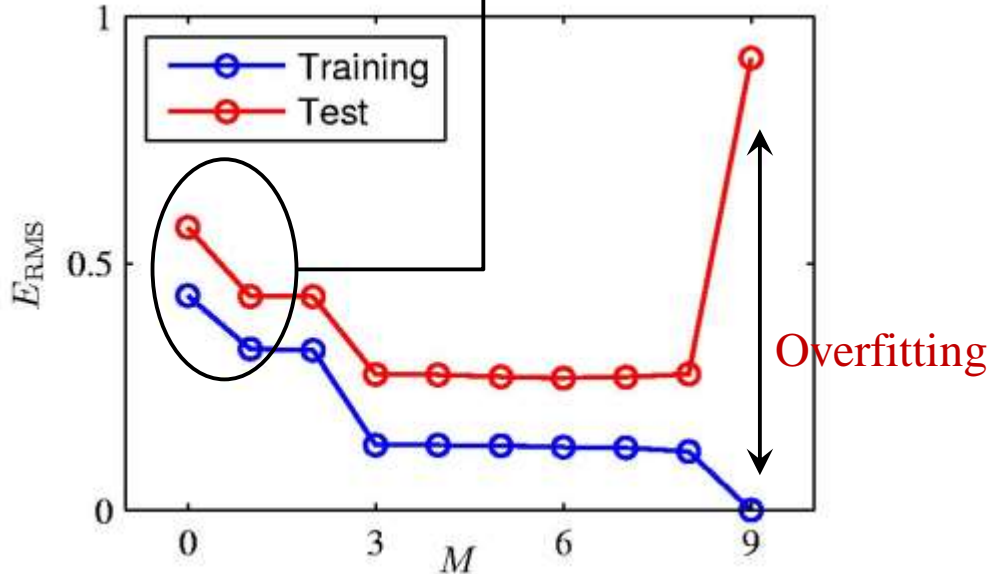
$$\hat{f} = \arg \min_f \hat{R}_{tr}(f) = \arg \min_f \frac{1}{N} \sum_{i=1}^N \ell(f(\mathbf{x}_i), y_i)$$

- Given a training data set,  $N$  is a constant, thus for convenience in presentation,  $\frac{1}{N}$  is dropped

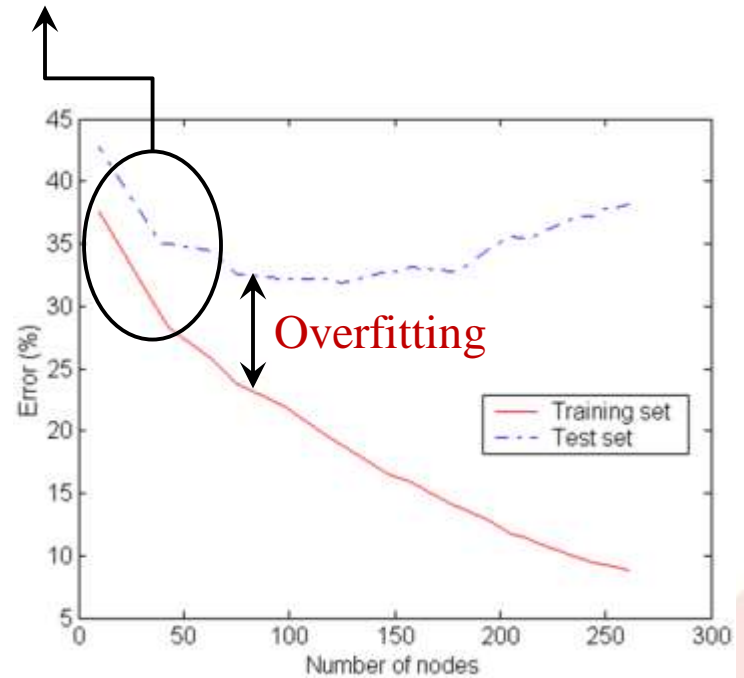
$$\hat{f} = \arg \min_f \sum_{i=1}^N \ell(f(\mathbf{x}_i), y_i)$$


# Overfitting Revisit

Underfitting: when model is too simple, both training and test error are large



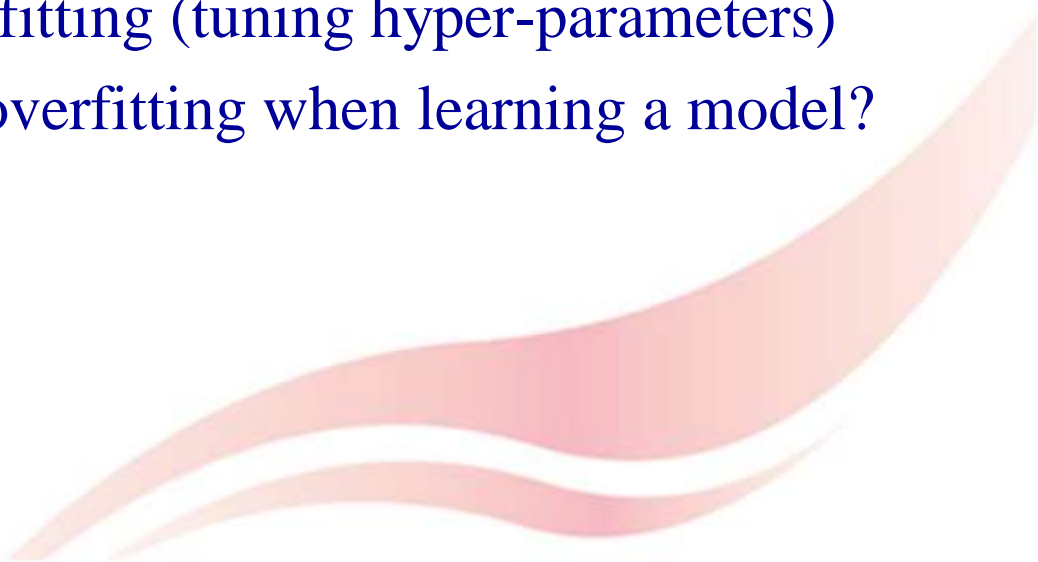
Polynomial curve fitting



Decision tree classification

Overfitting: when test error begins to increase even though training error continues to decrease

# Overfitting v.s. Model Complexity

- Observations of the two examples:
    - Increasing model complexity could make training error or training loss to keep being decreased
    - When model complexity keeps increasing, test error or test loss will increase after some point
  - After a model is learned, we can use validation set to evaluate it to reduce the risk of overfitting (tuning hyper-parameters)
  - Can we reduce the risk of overfitting when learning a model?
- 

# Occam's Razor Principle

- Given two models of similar performance, we should prefer the simpler model over the more complex model
- For complex models, there is a greater chance that it is fitted accidentally by noise in data
  - Overfitting results in models that are more complex than necessary
- Therefore, we should include model complexity when learning a model




# Structural Risk Minimization (cont.)

- Empirical Risk Minimization

$$\hat{f} = \arg \min_f \sum_{i=1}^N \ell(f(\mathbf{x}_i), y_i)$$

- Structural Risk Minimization

$$\hat{f} = \arg \min_f \sum_{i=1}^N \ell(f(\mathbf{x}_i), y_i) + \boxed{\lambda \Omega(f)}$$


- $\Omega(f)$  is known as a penalty or regularization term to control the model complexity of  $f$
- $\lambda > 0$  is a trade-off hyper-parameter

# Structural Risk Minimization (cont.)

$$\hat{f} = \arg \min_f \sum_{i=1}^N \ell(f(\mathbf{x}_i), y_i) + \lambda \Omega(f)$$

- How to learn  $f$ ?

- Design a specific form of  $f$  in terms of some parameters, denoted by a vector  $\boldsymbol{\theta} \in \mathbb{R}^{t \times 1}$ , i.e.,  $f(\mathbf{x}; \boldsymbol{\theta})$
- The parameterized  $f(\mathbf{x}; \boldsymbol{\theta})$  defines a family of functions with different values of  $\boldsymbol{\theta}$
- Learning  $f$  is equivalent to learning the values of  $\boldsymbol{\theta}$

$$\hat{\boldsymbol{\theta}} = \arg \min_{\boldsymbol{\theta}} \sum_{i=1}^N \ell(f(\mathbf{x}_i; \boldsymbol{\theta}), y_i) + \lambda \Omega(\boldsymbol{\theta})$$

# Structural Risk Minimization (cont.)

$$\hat{\boldsymbol{\theta}} = \arg \min_{\boldsymbol{\theta}} \sum_{i=1}^N \ell(f(\mathbf{x}_i; \boldsymbol{\theta}), y_i) + \lambda \Omega(\boldsymbol{\theta})$$

- Popular regularization terms include
  - the squared L2 norm:  $\|\boldsymbol{\theta}\|_2^2$ 
    - $\|\boldsymbol{\theta}\|_2^2 = \sum_{i=1}^t \theta_i^2$
    - Tends to prefer a model with a smaller value for each parameter  $\theta_i$
  - the L1 norm:  $\|\boldsymbol{\theta}\|_1$ 
    - $\|\boldsymbol{\theta}\|_1 = \sum_{i=1}^t |\theta_i|$
    - Tends to prefer a model with a smaller value for each parameter  $\theta_i$ , and fewer parameters with non-zero values
    - Induce sparsity, i.e., some  $\theta_i$ 's tend to be zeros



# Linear Models: Regression

$$\hat{\theta} = \arg \min_{\theta} \sum_{i=1}^N \ell(f(\mathbf{x}_i; \theta), y_i) + \lambda \Omega(\theta)$$

- In general, for regression,  $f(\mathbf{x}_i; \theta)$  is defined as

$$f(\mathbf{x}; \theta) = \mathbf{w} \cdot \mathbf{x} + b$$

$\theta$  is a concatenation of  $\mathbf{w}$  and  $b$

- Given  $\mathbf{x}_i$ , the prediction of  $f(\mathbf{x}; \theta)$  is the linear combination of its  $m$  feature values with weights  $\mathbf{w}$  plus a bias term  $b$

$$\mathbf{x}_i = \begin{bmatrix} x_{1i} \\ x_{2i} \\ \dots \\ x_{mi} \end{bmatrix}$$

$$\mathbf{w} = \begin{bmatrix} w_1 \\ w_2 \\ \dots \\ w_m \end{bmatrix}$$

$$\hat{y} = f(\mathbf{x}; \theta) = \mathbf{w} \cdot \mathbf{x} + b = \sum_{i=1}^m x_i w_i + b$$

# Linear Models: Classification

$$\hat{\boldsymbol{\theta}} = \arg \min_{\boldsymbol{\theta}} \sum_{i=1}^N \ell(f(\mathbf{x}_i; \boldsymbol{\theta}), y_i) + \lambda \Omega(\boldsymbol{\theta})$$

- In general, for classification,  $f(\mathbf{x}_i; \mathbf{w})$  is defined as

$$f(\mathbf{x}; \boldsymbol{\theta}) = h(\mathbf{w} \cdot \mathbf{x} + b)$$

where  $h(z)$  is function to map continuous values to discrete values (denoting different categories)

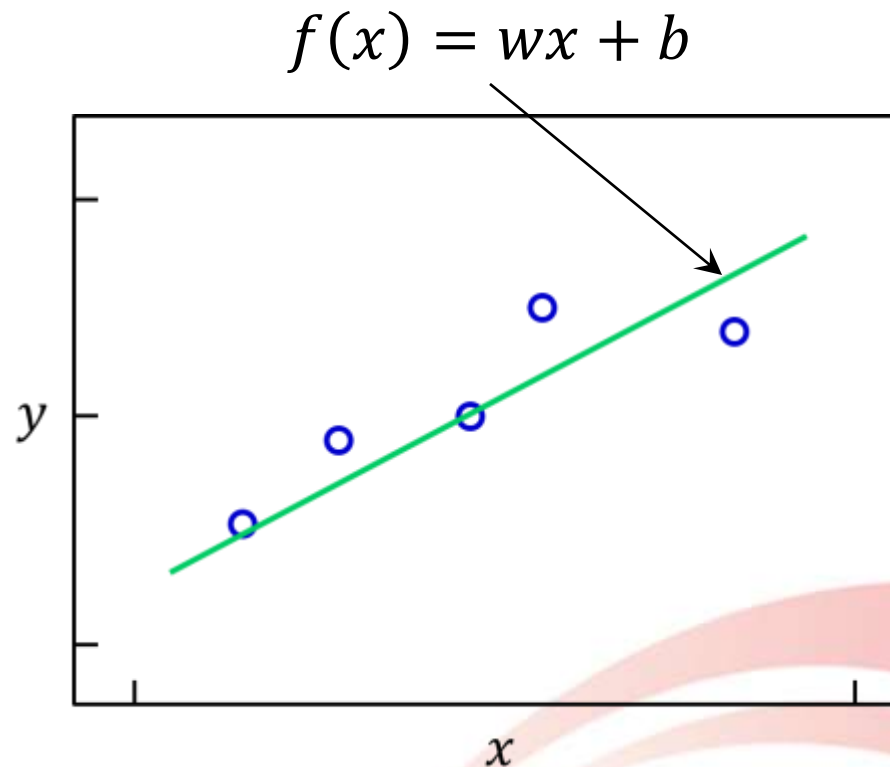
- For example,

$$h(z) = \begin{cases} +1 & \text{if } z \geq 0 \\ -1 & \text{if } z < 0 \end{cases}$$

**Next Lecture**

# Linear Regression: One-Dimension

- Each instance is represented by only one input feature
- To learn a linear function  $f(x)$  in terms of  $w$  and  $b$  (both are scalars) from  $\{x_i, y_i\}, i = 1, \dots, N$



# 1D Linear Regression (cont.)

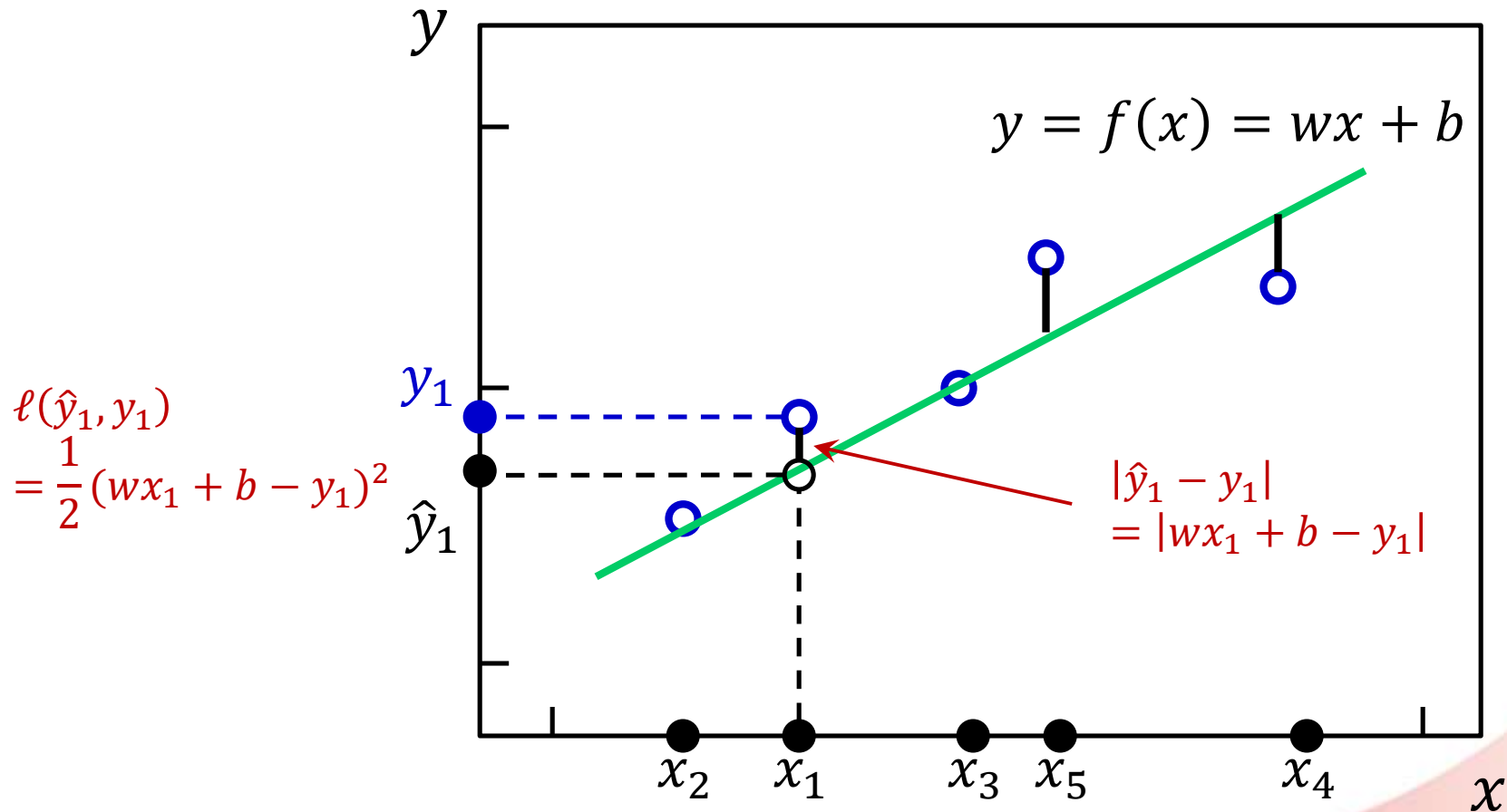
$$[\hat{w}, \hat{b}] = \arg \min_{[w, b]} \sum_{i=1}^N \ell(\hat{y}_i, y_i)$$

Drop the regularization  
term for simplicity at first

where  $\hat{y}_i = wx_i + b$

- The loss function  $\ell(\hat{y}_i, y_i)$  is to measure the difference between  $\hat{y}_i$  and  $y_i$ 
  - For regression, the magnitude of the difference, i.e.,  $|\hat{y}_i - y_i|$
- To make the resultant optimization problem easier to solve
  - We expect the loss function has some good properties, e.g., differentiable everywhere
  - The square of magnitude,  $|\hat{y}_i - y_i|^2 = (\hat{y}_i - y_i)^2$  or  $\frac{1}{2}(\hat{y}_i - y_i)^2$

# Regression Loss Function



$$\frac{1}{2} \sum_{i=1}^5 \ell(\hat{y}_i, y_i) = \frac{1}{2} \sum_{i=1}^5 (wx_i + b - y_i)^2$$

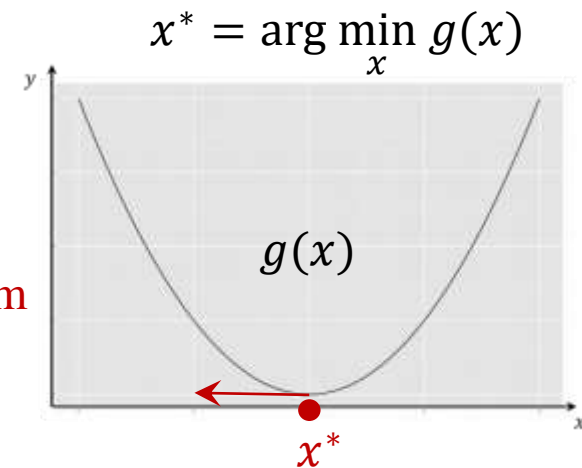
# Optimization

- Learn  $w$  and  $b$  by minimizing the square loss

$$[\hat{w}, \hat{b}] = \arg \min_{[w, b]} \frac{1}{2} \sum_{i=1}^N (wx_i + b - y_i)^2$$

- The objective of the optimization problem
- The objective is convex

- Unconstrained optimization problem
- Set the derivatives of the objective w.r.t.  $w$  and  $b$  to zero, respectively
- $\hat{w}$  and  $\hat{b}$  can be obtained by solving the equations



# Closed-form Solution

$$\frac{\partial \left( \frac{1}{2} \sum_{i=1}^N (wx_i + b - y_i)^2 \right)}{\partial w} = 0$$

$$\frac{\partial \left( \frac{1}{2} \sum_{i=1}^N (wx_i + b - y_i)^2 \right)}{\partial b} = 0$$

$$\sum_{i=1}^N (wx_i + b - y_i)x_i = 0$$
$$\sum_{i=1}^N (wx_i + b - y_i) = 0$$

Chain rule of calculus

$$y = g(x)$$

$$z = f(y) = f(g(x))$$

$$\frac{\partial z}{\partial x} = \frac{\partial z}{\partial y} \frac{\partial y}{\partial x}$$

$$z_i = wx_i + b - y_i$$

$$\begin{aligned} \frac{\partial z_i^2}{\partial w} &= \frac{\partial z_i^2}{\partial z_i} \frac{\partial z_i}{\partial w} \\ &= 2z_i \frac{\partial (wx_i + b - y_i)}{\partial w} \\ &= 2z_i x_i \end{aligned}$$

# Closed-form Solution (cont.)

$$\sum_{i=1}^N (wx_i + b - y_i)x_i = 0$$

$$\sum_{i=1}^N (wx_i + b - y_i) = 0$$

$$\sum_{i=1}^N (wx_i + b - y_i)x_i = 0$$

$$b = \frac{1}{N} \sum_{i=1}^N (y_i - wx_i)$$

$$\sum_{i=1}^N (wx_i - y_i) + Nb = 0$$



# Closed-form Solution (cont.)

$$\sum_{i=1}^N (wx_i + b - y_i)x_i = 0$$

$$b = \frac{1}{N} \sum_{i=1}^N (y_i - wx_i)$$

$$\sum_{i=1}^N (wx_i + b - y_i)x_i = 0$$

$$b = \frac{1}{N} \sum_{i=1}^N y_i - w \frac{1}{N} \sum_{i=1}^N x_i = \bar{y} - w\bar{x}$$

The diagram illustrates the derivation of the closed-form solution for the bias term  $b$ . A large blue curved arrow on the left indicates the flow from the first equation to the third. Red boxes and arrows highlight the substitution of sample means into the equation for  $b$ . The first equation is  $\sum_{i=1}^N (wx_i + b - y_i)x_i = 0$ . The second equation is  $b = \frac{1}{N} \sum_{i=1}^N (y_i - wx_i)$ . The third equation is  $b = \frac{1}{N} \sum_{i=1}^N y_i - w \frac{1}{N} \sum_{i=1}^N x_i = \bar{y} - w\bar{x}$ . Red boxes are drawn around the terms  $\frac{1}{N} \sum_{i=1}^N y_i$ ,  $\frac{1}{N} \sum_{i=1}^N x_i$ ,  $\bar{y}$ , and  $\bar{x}$ . Red arrows show the substitution of  $\bar{y}$  for  $\frac{1}{N} \sum_{i=1}^N y_i$  and  $\bar{x}$  for  $\frac{1}{N} \sum_{i=1}^N x_i$ .

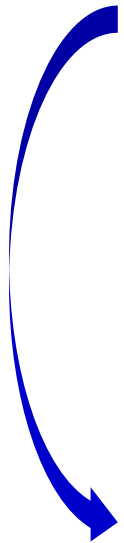
# Closed-form Solution (cont.)

$$\sum_{i=1}^N (wx_i + b - y_i)x_i = 0$$

$$b = \bar{y} - w\bar{x}$$

$$\sum_{i=1}^N (wx_i + \bar{y} - w\bar{x} - y_i)x_i = 0$$

$$b = \bar{y} - w\bar{x}$$



# Closed-form Solution (cont.)

$$\sum_{i=1}^N (wx_i + \bar{y} - w\bar{x} - y_i)x_i = 0$$

$$b = \bar{y} - w\bar{x}$$

$$w \sum_{i=1}^N (x_i - \bar{x})x_i = \sum_{i=1}^N (y_i - \bar{y})x_i$$

$$b = \bar{y} - w\bar{x}$$

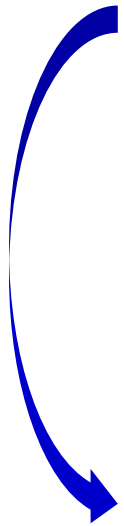
# Closed-form Solution (cont.)

$$w \sum_{i=1}^N (x_i - \bar{x})x_i = \sum_{i=1}^N (y_i - \bar{y})x_i$$

$$b = \bar{y} - w\bar{x}$$

$$w = \frac{\sum_{i=1}^N (y_i - \bar{y})x_i}{\sum_{i=1}^N (x_i - \bar{x})x_i}$$

$$b = \bar{y} - w\bar{x}$$



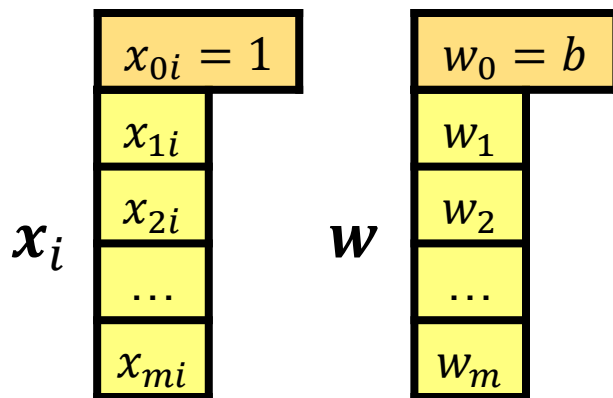
# Multi-Dimension Case

- Each instance has  $m$  dimensions, a linear function  $f(\mathbf{x})$  is defined as

$$f(\mathbf{x}) = \mathbf{w} \cdot \mathbf{x} + b$$

- By defining  $w_0 = b$ , and  $x_0 = 1$ ,  $f(\mathbf{x})$  can be rewritten as

$$f(\mathbf{x}) = \mathbf{w} \cdot \mathbf{x}$$



Both  $\mathbf{w}$  and  $\mathbf{x}$  have  
 $m + 1$  dimensions

$$= \sum_{k=0}^m x_{ki} w_k$$

$$= \sum_{k=1}^m x_{ki} w_k + x_{0i} w_0$$

# Optimization

- Learn  $\mathbf{w}$  by minimizing the total square loss

$$\hat{\mathbf{w}} = \arg \min_{\mathbf{w}} \frac{1}{2} \sum_{i=1}^N (\mathbf{w} \cdot \mathbf{x}_i - y_i)^2$$

- A closed-form solution can be obtained by setting the derivative of the objective w.r.t.  $\mathbf{w}$  to zero, and solving the resultant equations

$$\frac{\partial \left( \frac{1}{2} \sum_{i=1}^N (\mathbf{w} \cdot \mathbf{x}_i - y_i)^2 \right)}{\partial \mathbf{w}} = \mathbf{0}$$

# Brief Linear Algebra Review

- Linear algebra plays a crucial role in deriving solutions for various machine learning methods
- You are highly recommended to refer to Part I of the Deep Learning book at <https://www.deeplearningbook.org/>
- Transpose of a vector or matrix

$$\mathbf{x} (m \times 1)$$

$x_1$
$x_2$
$\dots$
$x_m$

$$\mathbf{x}^T (1 \times m)$$

$x_1$	$x_2$	$\dots$	$x_m$
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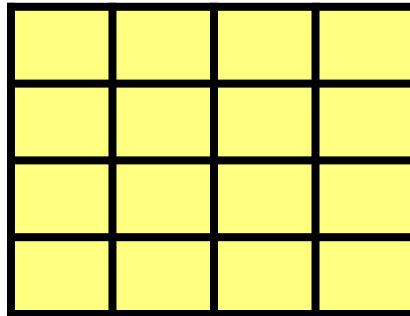
$$\mathbf{A} (m \times N)$$


$$\mathbf{A}^T (N \times m)$$


# Matrix/Vector Concepts

- Square matrix
  - If a matrix  $\mathbf{A}$  has the same number of rows and columns, then it is said to be square matrix

$\mathbf{A} (m \times m)$



- Symmetric matrix
  - If a **square** matrix  $\mathbf{A}$  satisfies  $\mathbf{A} = \mathbf{A}^T$



# Matrix Multiplication

- Matrix multiplication is associative
  - $(\mathbf{AB})\mathbf{C} = \mathbf{A}(\mathbf{BC})$
- Matrix multiplication is distributive
  - $\mathbf{A}(\mathbf{B} + \mathbf{C}) = \mathbf{AB} + \mathbf{AC}$
- Matrix multiplication is NOT commutative in general
  - $\mathbf{AB} \neq \mathbf{BA}$
- The identity matrix,  $\mathbf{I} (m \times m)$ , is a symmetric matrix with ones on the diagonal and zeros everywhere else
  - If  $\mathbf{A}$  is a square matrix  $(m \times m)$ :  $\mathbf{AI} = \mathbf{IA} = \mathbf{A}$
  - If  $\mathbf{A}$  is  $(N \times m)$ :  $\mathbf{AI} = \mathbf{A}$
  - If  $\mathbf{A}$  is  $(m \times N)$ :  $\mathbf{IA} = \mathbf{A}$

$\mathbf{I} (m \times m)$

1	0	0	0
0	1	0	0
0	0	1	0
0	0	0	1

# Matrix Operations

- The transpose of  $\mathbf{A}^T$ :

$$(\mathbf{A}^T)^T = \mathbf{A}$$

- The transpose of  $\mathbf{AB}$ :

$$(\mathbf{AB})^T = \mathbf{B}^T \mathbf{A}^T$$

- The transpose of  $\mathbf{Ax}$

$$(\mathbf{Ax})^T = \mathbf{x}^T \mathbf{A}^T$$

- The transpose of  $\mathbf{x}^T \mathbf{y}$

$$(\mathbf{x}^T \mathbf{y})^T = \mathbf{y}^T \mathbf{x}$$


- The transpose of a scalar is the scalar itself

$$a^T = a$$


# Matrix Operations (cont.)

- For a square matrix  $\mathbf{A}$  ( $m \times m$ ), if it is invertible, then there exists a unique matrix, denoted by  $\mathbf{A}^{-1}$ , such that

$$\mathbf{A}\mathbf{A}^{-1} = \mathbf{A}^{-1}\mathbf{A} = \mathbf{I}$$

- If it is not invertible, then such a matrix  $\mathbf{A}^{-1}$  does not exist
  - Non-square matrices do not have inverses by definition
  - Properties of the inverse ( $\mathbf{A}$  and  $\mathbf{B}$  are invertible)
    - $(\mathbf{A}^{-1})^{-1} = \mathbf{A}$
    - $(\mathbf{AB})^{-1} = \mathbf{B}^{-1}\mathbf{A}^{-1}$
    - $(\mathbf{A}^{-1})^T = (\mathbf{A}^T)^{-1}$
- 

# Linear System

- Given the following system of linear equations

$$\begin{aligned}2x_1 + x_3 &= 5 \\3x_1 - 4x_2 + 2x_3 &= 4 \\2x_2 - 3x_3 &= -3 \\-x_1 + 2x_2 - 5x_3 &= 1\end{aligned}$$

- They can be written in a more compact form as

$$\mathbf{Ax} = \mathbf{b}$$

$\mathbf{A} (4 \times 3)$

2	0	1
3	-4	2
0	2	-3
-1	2	-5

$\mathbf{x} (3 \times 1)$

$x_1$
$x_2$
$x_3$

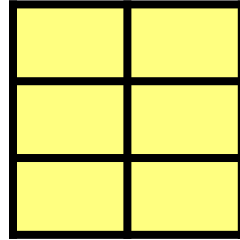
$\mathbf{b} (4 \times 1)$

5
4
-3
1

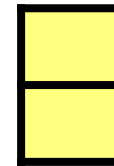
# Linear System (cont.)

$$\mathbf{A}x = b$$

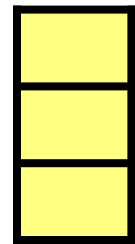
$$\mathbf{A} \ (k \times d)$$



$$x \ (d \times 1)$$



$$b \ (k \times 1)$$



- If  $\mathbf{A}$  is square, and invertible, then we multiply both sides of the equation by  $\mathbf{A}^{-1}$  to obtain a **unique** solution

$$\mathbf{A}^{-1}\mathbf{A}x = \mathbf{A}^{-1}b \implies x = \mathbf{A}^{-1}b$$

# Linear System (cont.)

$$\mathbf{A}\mathbf{x} = \mathbf{b}$$

$\mathbf{A} (k \times d)$


$\mathbf{x} (d \times 1)$


$\mathbf{b} (k \times 1)$


- If  $\mathbf{A}$  is not invertible, solutions are **not** unique, we can find a solution by using the pseudo inverse (also known as generalized inverse) of  $\mathbf{A}$  instead, denoted by  $\mathbf{A}^\dagger$

$$\mathbf{x} = \mathbf{A}^\dagger \mathbf{b}$$

- Special case: when  $\mathbf{A}$  is square ( $k \times k$ ) but not invertible

$\mathbf{A}^\dagger \mathbf{A} = \mathbf{A} \mathbf{A}^\dagger = \text{diag}(\lambda_1, \dots, \lambda_k)$ , where  $\lambda_i \in \{0, 1\}$ , at least one  $\lambda_i = 0$

$$\mathbf{A} \mathbf{A}^\dagger = \mathbf{A}^\dagger \mathbf{A}$$

1	0	0	0
0	1	0	0
0	0	0	0
0	0	0	0

# Closed-form Solution for Linear Regression

$$\frac{\partial \left( \frac{1}{2} \sum_{i=1}^N (\mathbf{w} \cdot \mathbf{x}_i - y_i)^2 \right)}{\partial \mathbf{w}} = \mathbf{0}$$

$$\frac{1}{2} \sum_{i=1}^N \frac{\partial (\mathbf{w} \cdot \mathbf{x}_i - y_i)^2}{\partial \mathbf{w}} = \mathbf{0}$$

$$\sum_{i=1}^N (\mathbf{w} \cdot \mathbf{x}_i - y_i) \mathbf{x}_i = \mathbf{0}$$

$$\sum_{i=1}^N (\mathbf{w} \cdot \mathbf{x}_i) \mathbf{x}_i - \sum_{i=1}^N y_i \mathbf{x}_i = \mathbf{0}$$

$$z_i = \mathbf{w} \cdot \mathbf{x}_i - y_i$$

$$\frac{\partial z_i^2}{\partial \mathbf{w}} = \frac{\partial z_i^2}{\partial z_i} \frac{\partial z_i}{\partial \mathbf{w}} = 2z_i \frac{\partial (\mathbf{w} \cdot \mathbf{x}_i - y_i)}{\partial \mathbf{w}}$$

$$\frac{\partial (\mathbf{w} \cdot \mathbf{x}_i - y_i)}{\partial \mathbf{w}} = \frac{\partial (\mathbf{w} \cdot \mathbf{x}_i)}{\partial \mathbf{w}} - 0 = \mathbf{x}_i$$

The Matrix Cookbook

<https://www.math.uwaterloo.ca/~hwolkowi/matrixcookbook.pdf>

# Closed-form Solution (cont.)

$$a\mathbf{x} = \mathbf{x}a$$

$$\sum_{i=1}^N \boxed{\text{scalar } (\mathbf{w} \cdot \mathbf{x}_i)} \mathbf{x}_i - \sum_{i=1}^N y_i \mathbf{x}_i = \mathbf{0}$$

$$\sum_{i=1}^N \boxed{\mathbf{x}_i (\mathbf{w} \cdot \mathbf{x}_i)} - \sum_{i=1}^N y_i \mathbf{x}_i = \mathbf{0}$$

$$\mathbf{x}_i (\mathbf{w} \cdot \mathbf{x}_i) = \mathbf{x}_i (\mathbf{w}^T \mathbf{x}_i) = \mathbf{x}_i (\mathbf{x}_i^T \mathbf{w}) = (\mathbf{x}_i \mathbf{x}_i^T) \mathbf{w}$$

$$\left( \sum_{i=1}^N \boxed{\mathbf{x}_i \mathbf{x}_i^T} \right) \mathbf{w} - \sum_{i=1}^N y_i \mathbf{x}_i = \mathbf{0}$$

$\mathbf{x}_i \in \mathbb{R}^{(m+1) \times 1}$  and  $\mathbf{x}_i^T \in \mathbb{R}^{1 \times (m+1)}$ , thus  $\mathbf{x}_i \mathbf{x}_i^T$  is a  $(m+1)$  by  $(m+1)$  matrix



# Closed-form Solution (cont.)

$$\left( \sum_{i=1}^N \boxed{x_i x_i^T} \right) \mathbf{w} - \sum_{i=1}^N y_i \mathbf{x}_i = \mathbf{0}$$

$\mathbf{x}_i \in \mathbb{R}^{(m+1) \times 1}$  and  $\mathbf{x}_i^T \in \mathbb{R}^{1 \times (m+1)}$ , thus  $\mathbf{x}_i \mathbf{x}_i^T$  is a  $(m+1)$  by  $(m+1)$  matrix

$(m+1) \times 1$

 $\mathbf{x}_i$ 

$x_{0i}$
$x_{1i}$
...
$x_{mi}$

$\mathbf{x}_i^T$

$x_{0i}$	$x_{1i}$	...	$x_{mi}$
----------	----------	-----	----------

$1 \times (m+1)$

$\mathbf{x}_i^T \mathbf{x}_i$  Inner product, scalar

$\mathbf{x}_i \mathbf{x}_i^T$   $(m+1)$  by  $(m+1)$  matrix

$x_{0i}x_{0i}$	$x_{0i}x_{1i}$	...	$x_{0i}x_{mi}$
$x_{1i}x_{0i}$	$x_{1i}x_{1i}$	...	$x_{1i}x_{mi}$
...	...	...	...
$x_{mi}x_{0i}$	$x_{mi}x_{1i}$	...	$x_{mi}x_{mi}$

# Closed-form Solution (cont.)

$$\left( \sum_{i=1}^N (x_i x_i^T) \right) w - \sum_{i=1}^N y_i x_i = 0$$

$$x_i x_i^T$$

$(m + 1)$  by  $(m + 1)$  matrix

$x_{0i}x_{0i}$	$x_{0i}x_{1i}$	...	$x_{0i}x_{mi}$
$x_{1i}x_{0i}$	$x_{1i}x_{1i}$	...	$x_{1i}x_{mi}$
...	...	...	...
$x_{mi}x_{0i}$	$x_{mi}x_{1i}$	...	$x_{mi}x_{mi}$

$$\sum_{i=1}^N (x_i x_i^T)$$

$\sum_{i=1}^N x_{0i}x_{0i}$	$\sum_{i=1}^N x_{0i}x_{1i}$	...	$\sum_{i=1}^N x_{0i}x_{mi}$
$\sum_{i=1}^N x_{1i}x_{0i}$	$\sum_{i=1}^N x_{1i}x_{1i}$	...	$\sum_{i=1}^N x_{1i}x_{mi}$
...	...	...	...
$\sum_{i=1}^N x_{mi}x_{0i}$	$\sum_{i=1}^N x_{mi}x_{1i}$	...	$\sum_{i=1}^N x_{mi}x_{mi}$

# Closed-form Solution (cont.)

$$\sum_{i=1}^N (\mathbf{x}_i \mathbf{x}_i^T)$$

$\sum_{i=1}^N x_{0i} x_{0i}$	$\sum_{i=1}^N x_{0i} x_{1i}$	...	$\sum_{i=1}^N x_{0i} x_{mi}$
$\sum_{i=1}^N x_{1i} x_{0i}$	$\sum_{i=1}^N x_{1i} x_{1i}$	...	$\sum_{i=1}^N x_{1i} x_{mi}$
...	...	...	...
$\sum_{i=1}^N x_{mi} x_{0i}$	$\sum_{i=1}^N x_{mi} x_{1i}$	...	$\sum_{i=1}^N x_{mi} x_{mi}$

$$\sum_{i=1}^N (\mathbf{x}_i \mathbf{x}_i^T) = \mathbf{X} \mathbf{X}^T$$

$(m + 1)$  by  $N$

$N$  by  $(m + 1)$

$\mathbf{X}$

$\mathbf{X}^T$

$x_{01}$	$x_{02}$	...	$x_{0N}$
$x_{11}$	$x_{12}$	...	$x_{1N}$
...	...	...	...
$x_{m1}$	$x_{m2}$	...	$x_{mN}$

$x_{01}$	$x_{11}$	...	$x_{m1}$
$x_{02}$	$x_{12}$	...	$x_{m2}$
...	...	...	...
$x_{0N}$	$x_{1N}$	...	$x_{mN}$

# Closed-form Solution (cont.)

$$\left( \sum_{i=1}^N (\mathbf{x}_i \mathbf{x}_i^T) \right) \mathbf{w} - \sum_{i=1}^N y_i \mathbf{x}_i = \mathbf{0}$$

$$\mathbf{X} \mathbf{X}^T \mathbf{w} - \sum_{i=1}^N y_i \mathbf{x}_i = \mathbf{0}$$

$$\mathbf{X} \mathbf{X}^T \mathbf{w} - \mathbf{X} \mathbf{y} = \mathbf{0}$$

$(m+1) \text{ by } N$

$\mathbf{X}$

$x_{01}$	$x_{02}$	$\dots$	$x_{0N}$
$x_{11}$	$x_{12}$	$\dots$	$x_{1N}$
$\dots$	$\dots$	$\dots$	$\dots$
$x_{m1}$	$x_{m2}$	$\dots$	$x_{mN}$


$N \text{ by } 1$

$\mathbf{y}$

$y_1$
$y_2$
$\dots$
$y_N$

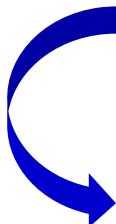
$$\sum_{i=1}^N y_i \mathbf{x}_i = \mathbf{X} \mathbf{y}$$

# Closed-form Solution (cont.)


$$\mathbf{X}\mathbf{X}^T \mathbf{w} - \mathbf{X}\mathbf{y} = \mathbf{0}$$

$$\mathbf{X}\mathbf{X}^T \mathbf{w} = \mathbf{X}\mathbf{y}$$

When  $\mathbf{X}\mathbf{X}^T$  is invertible


$$(\mathbf{X}\mathbf{X}^T)^{-1} \mathbf{X}\mathbf{X}^T \mathbf{w} = (\mathbf{X}\mathbf{X}^T)^{-1} \mathbf{X}\mathbf{y}$$


$$\cancel{\mathbf{I}} \mathbf{w} = (\mathbf{X}\mathbf{X}^T)^{-1} \mathbf{X}\mathbf{y}$$

What if  $\mathbf{X}\mathbf{X}^T$  is NOT invertible?



# When $\mathbf{XX}^T$ is NOT Invertible

- We can use the pseudo inverse (also known as generalized inverse) of  $\mathbf{XX}^T$  instead, i.e.,  $(\mathbf{XX}^T)^\dagger$
- In this case

$$\mathbf{w} = (\mathbf{XX}^T)^\dagger \mathbf{X}\mathbf{y}$$


# Regularized Linear Regression

- Recall structural risk minimization

$$\hat{f} = \arg \min_f \sum_{i=1}^N \ell(f(\mathbf{x}_i), y_i) + \lambda \Omega(f)$$

$\lambda > 0$  tradeoff hyper-parameter

$$\hat{\mathbf{w}} = \arg \min_{\mathbf{w}} \frac{1}{2} \sum_{i=1}^N (\mathbf{w} \cdot \mathbf{x}_i - y_i)^2 + \frac{\lambda}{2} \|\mathbf{w}\|_2^2$$

$\|\mathbf{w}\|_2^2 = \mathbf{w} \cdot \mathbf{w}$

A regularization term to control the complexity of the model

Also known as Ridge regression

# Optimization

$$\hat{\mathbf{w}} = \arg \min_{\mathbf{w}} \frac{1}{2} \sum_{i=1}^N (\mathbf{w} \cdot \mathbf{x}_i - y_i)^2 + \frac{\lambda}{2} \|\mathbf{w}\|_2^2$$

- Still an unconstrained optimization problem, and the objective is still convex
- A closed-form solution can be obtained by setting the derivative of the objective w.r.t.  $\mathbf{w}$ , and solving the resultant equation

$$\frac{\partial \left( \frac{1}{2} \sum_{i=1}^N (\mathbf{w} \cdot \mathbf{x}_i - y_i)^2 + \frac{\lambda}{2} \|\mathbf{w}\|_2^2 \right)}{\partial \mathbf{w}} = \mathbf{0}$$



# Closed-form Solution

$$\frac{\partial \left( \frac{1}{2} \sum_{i=1}^N (\mathbf{w} \cdot \mathbf{x}_i - y_i)^2 + \frac{\lambda}{2} \|\mathbf{w}\|_2^2 \right)}{\partial \mathbf{w}} = \mathbf{0}$$

$$\frac{\partial \frac{1}{2} \sum_{i=1}^N (\mathbf{w} \cdot \mathbf{x}_i - y_i)^2}{\partial \mathbf{w}} + \frac{\partial \frac{\lambda}{2} \|\mathbf{w}\|_2^2}{\partial \mathbf{w}} = \mathbf{0}$$

$$\left( \sum_{i=1}^N (\mathbf{x}_i \mathbf{x}_i^T) \right) \mathbf{w} - \sum_{i=1}^N y_i \mathbf{x}_i + \lambda \mathbf{w} = \mathbf{0}$$

$$\frac{\partial \|\mathbf{w}\|_2^2}{\partial \mathbf{w}} = \frac{\partial (\mathbf{w} \cdot \mathbf{w})}{\partial \mathbf{w}} = 2\mathbf{w}$$

The matrix cookbook

# Closed-form Solution (cont.)

$$\left( \sum_{i=1}^N (\mathbf{x}_i \mathbf{x}_i^T) \right) \mathbf{w} - \sum_{i=1}^N y_i \mathbf{x}_i + \lambda \mathbf{w} = \mathbf{0}$$

$$\mathbf{X}\mathbf{X}^T \mathbf{w} - \mathbf{X}\mathbf{y} + \lambda \mathbf{I} \mathbf{w} = \mathbf{0}$$

$$(\mathbf{X}\mathbf{X}^T + \lambda \mathbf{I}) \mathbf{w} - \mathbf{X}\mathbf{y} = \mathbf{0}$$

Always invertible as  
long as  $\lambda$  is positive

$$(\mathbf{X}\mathbf{X}^T + \lambda \mathbf{I}) \mathbf{w} = \mathbf{X}\mathbf{y} \quad \longrightarrow \quad \mathbf{w} = (\mathbf{X}\mathbf{X}^T + \lambda \mathbf{I})^{-1} \mathbf{X}\mathbf{y}$$

# Why $\mathbf{XX}^T + \lambda \mathbf{I}$ Invertible?

- A square matrix is invertible if and only if it does not have a zero eigenvalue
- If a symmetric matrix  $\mathbf{A}$  is positive semidefinite, then all of its eigenvalues are non-negative ( $\geq 0$ )
  - When a symmetric matrix  $\mathbf{A}$  ( $d \times d$ ) is said to be positive semidefinite iif for any non-zero column vector  $\mathbf{x}$  ( $d \times 1$ ),  $\mathbf{x}^T \mathbf{A} \mathbf{x} \geq 0$
- If a symmetric matrix  $\mathbf{A}$  is positive definite, then all of its eigenvalues are positive ( $> 0$ )
  - When a symmetric matrix  $\mathbf{A}$  ( $d \times d$ ) is said to be positive definite iif for any non-zero column vector  $\mathbf{x}$  ( $d \times 1$ ),  $\mathbf{x}^T \mathbf{A} \mathbf{x} > 0$
- A positive definite matrix is invertible (all eigenvalues  $> 0$ , and of course non-zero)

# Why $\mathbf{XX}^T + \lambda \mathbf{I}$ Invertible? (cont.)

- $\mathbf{XX}^T$  is a positive semidefinite
  - Need to prove for any non-zero  $(m + 1)$ - dimensional column vector  $\mathbf{z}$ ,  $\mathbf{z}^T \mathbf{XX}^T \mathbf{z} \geq 0$
  - Proof: Denote  $\mathbf{y} = \mathbf{X}^T \mathbf{z}$ . Then  $\mathbf{z}^T \mathbf{XX}^T \mathbf{z} = \boxed{\mathbf{y}^T \mathbf{y}} \quad \|\mathbf{y}\|_2^2 \geq 0$
- $\mathbf{XX}^T + \lambda \mathbf{I}$  is positive definite if  $\lambda > 0$   $\|\mathbf{y}\|_2^2 = 0$  iff  $\mathbf{y} = \mathbf{X}^T \mathbf{z} = \mathbf{0}$ 
  - Need to prove for any non-zero  $(m + 1)$ - dimensional column vector  $\mathbf{z}$ ,  $\mathbf{z}^T (\mathbf{XX}^T + \lambda \mathbf{I}) \mathbf{z} = \mathbf{z}^T \mathbf{XX}^T \mathbf{z} + \lambda \mathbf{z}^T \mathbf{z} > 0$
  - Proof:  $\mathbf{z}^T (\mathbf{XX}^T + \lambda \mathbf{I}) \mathbf{z} = \boxed{\mathbf{z}^T \mathbf{XX}^T \mathbf{z}} + \boxed{\lambda \mathbf{z}^T \mathbf{z}} > 0$ 

$\geq 0$

$> 0$  since  $\lambda > 0$   
and  $\mathbf{z}$  is non-zero

# Large-scale Issue

$(m + 1) \times (m + 1)$

$$\mathbf{w} = (\mathbf{X}\mathbf{X}^T + \lambda\mathbf{I})^{-1}\mathbf{X}\mathbf{y}$$

- The computation complexity of computing an inverse of a  $(m + 1) \times (m + 1)$  matrix is  $O((m + 1)^3)$
- When  $m$  is large, it is time consuming
- Rather than computing the inverse to obtain a closed-form solution, consider the following linear system

$$(\mathbf{X}\mathbf{X}^T + \lambda\mathbf{I})\mathbf{w} = \mathbf{X}\mathbf{y} \quad \text{Linear system: } \mathbf{Ax} = \mathbf{b}$$

- We can solve it by using various numerical methods, e.g., Gaussian elimination, etc.

# Large-scale Issue (cont.)

$$\min_{\mathbf{w}} \frac{1}{2} \sum_{i=1}^N (\mathbf{w} \cdot \mathbf{x}_i - y_i)^2 + \frac{\lambda}{2} \|\mathbf{w}\|_2^2$$

- Alternatively, rather than trying to derive an analytical solution, we can apply numerical methods to iteratively minimize the objective, e.g., gradient descent
- Denote the objective by

$$E(\mathbf{w}) = \frac{1}{2} \sum_{i=1}^N (\mathbf{w} \cdot \mathbf{x}_i - y_i)^2 + \frac{\lambda}{2} \|\mathbf{w}\|_2^2$$

The diagram illustrates the gradient descent update rule. The equation  $\mathbf{w}_{t+1} = \mathbf{w}_t - \rho \frac{\partial E(\mathbf{w})}{\partial \mathbf{w}}$  is shown. A red arrow points from the text "Objective to be minimized" to the term  $E(\mathbf{w})$  in the numerator of the gradient. Another red arrow points from the text "Learning rate  $\rho \in (0,1]$ " to the variable  $\rho$  in the denominator. The entire equation is set against a background of red wavy lines.

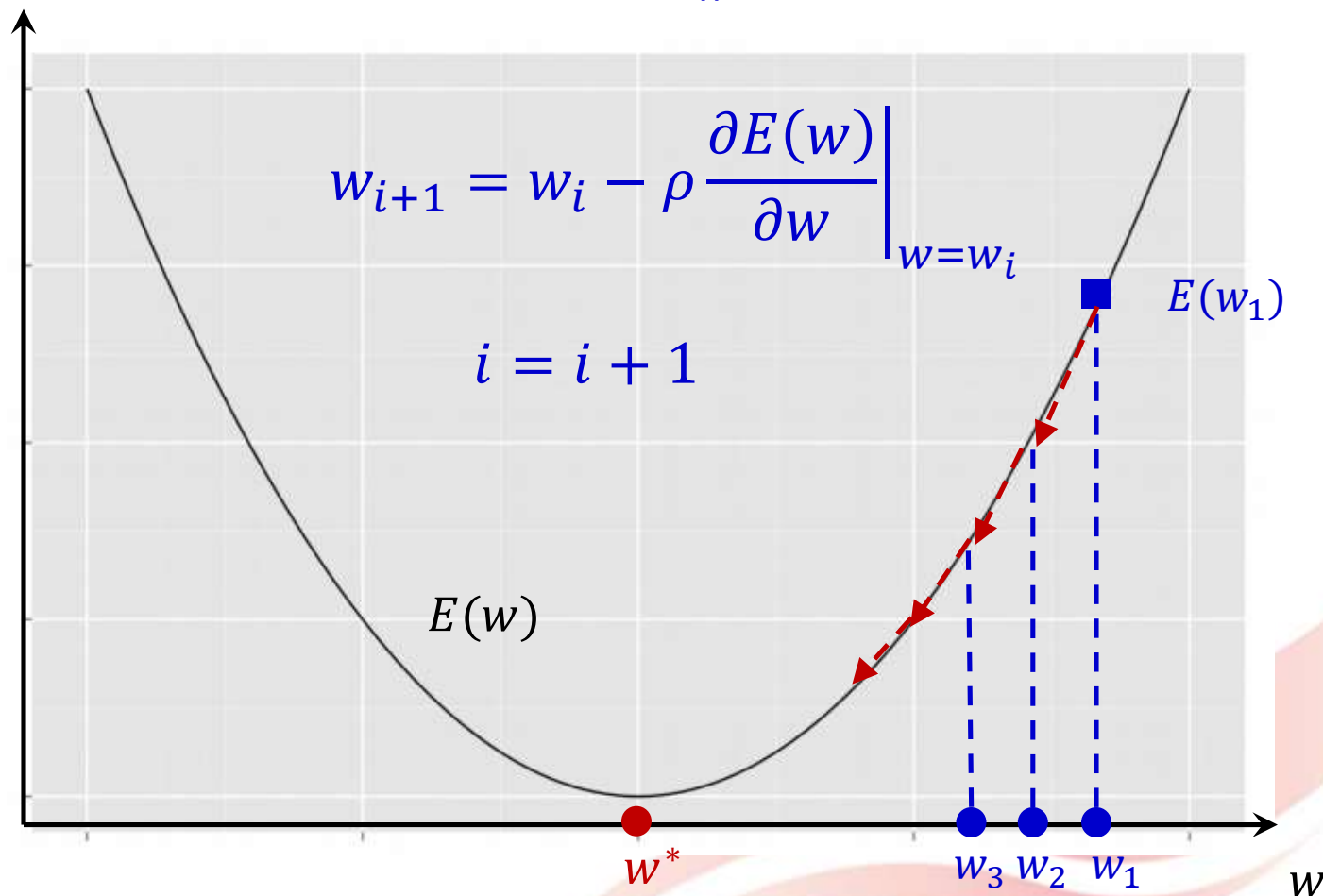
$$\mathbf{w}_{t+1} = \mathbf{w}_t - \rho \frac{\partial E(\mathbf{w})}{\partial \mathbf{w}}$$

Objective to be minimized

Learning rate  $\rho \in (0,1]$

# Gradient Descent

$$w^* = \arg \min_w E(w)$$



# Implementation using scikit-learn

- API: `sklearn.linear_model`: Linear Models

[https://scikit-learn.org/stable/modules/classes.html#module-sklearn.linear\\_model](https://scikit-learn.org/stable/modules/classes.html#module-sklearn.linear_model)

- Classical linear regressors
- `linear_model.LinearRegression` → linear regression without regularization
- `linear_model.Ridge` → regularized linear regression

## Classical linear regressors

<code>linear_model.LinearRegression(*[, ...])</code>	Ordinary least squares Linear Regression.
<code>linear_model.Ridge([alpha, fit_intercept, ...])</code>	Linear least squares with l2 regularization.
<code>linear_model.RidgeCV([alphas, ...])</code>	Ridge regression with built-in cross-validation.
<code>linear_model.SGDRegressor([loss, penalty, ...])</code>	Linear model fitted by minimizing a regularized empirical loss with SGD



# Example

```
>>> from sklearn.linear_model import LinearRegression
```

```
>>> from sklearn.linear_model import Ridge
```

```
>>> import numpy as np
```

```
>>> n_samples, n_features = 10, 5
```

```
>>> rng = np.random.RandomState(0)
```

```
>>> y = rng.randn(n_samples)
```

```
>>> X = rng.randn(n_samples, n_features)
```

```
>>> rr = Ridge(alpha=0.1)
```

```
>>> rr.fit(X, y)
```

```
>>> pred_train_rr= rr.predict(X)
```

```
>>> lr = LinearRegression()
```

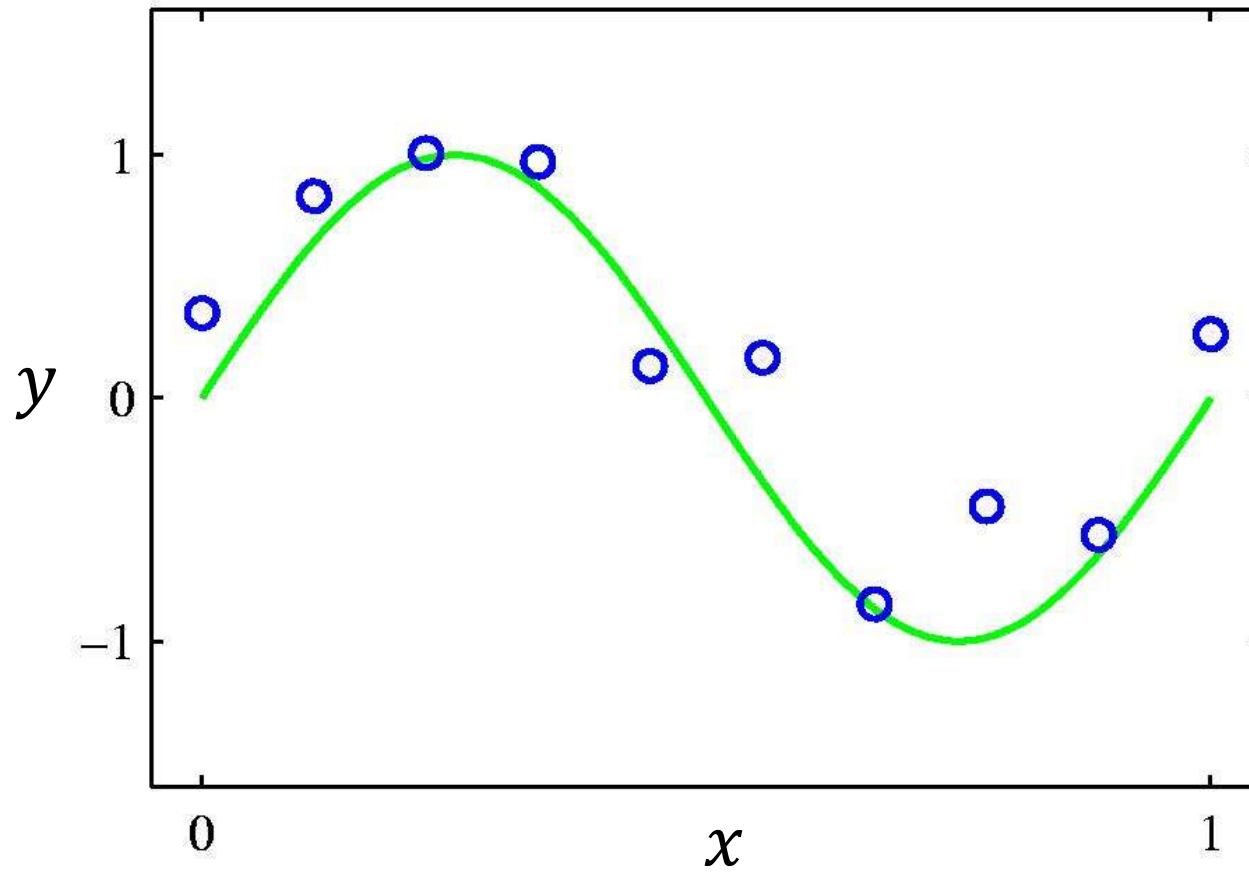
```
>>> lr.fit(X, y)
```

```
>>> pred_train_lr= lr.predict(X)
```

Model training and testing

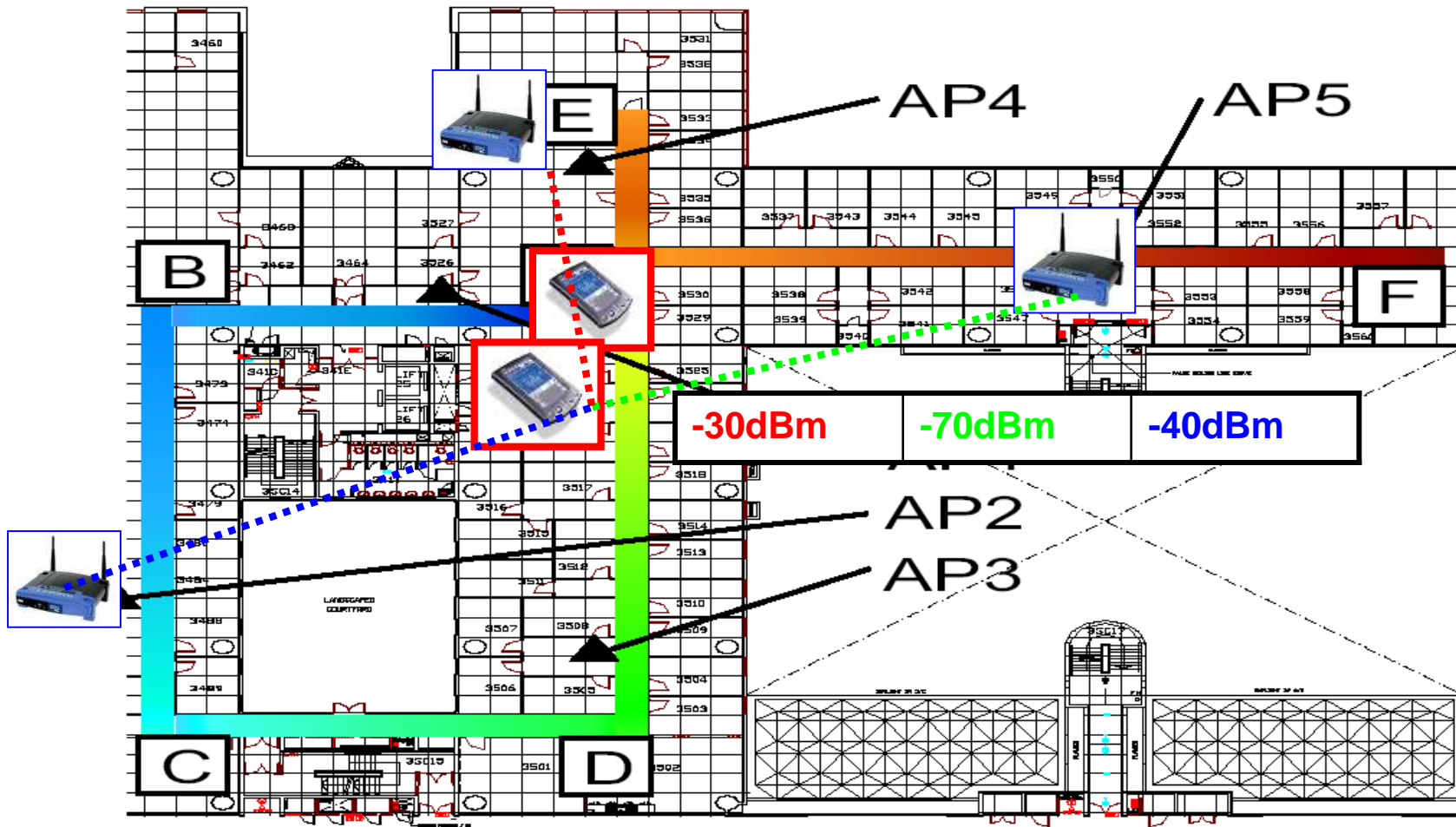
# Nonlinear Regression

Kernel methods (L5)



# Regression: Real-world Example

## Indoor WiFi localization



**Thank you!**

