

The background of the entire image is a dark blue field filled with a pattern of red dots. These dots are arranged in a way that they form a large, faint, stylized 'H' shape that encompasses the central text. The dots are of varying sizes and are more densely packed in certain areas, creating a textured, digital effect.

# HUST

**ĐẠI HỌC BÁCH KHOA HÀ NỘI**  
HANOI UNIVERSITY OF SCIENCE AND TECHNOLOGY

ONE LOVE. ONE FUTURE.



ĐẠI HỌC  
BÁCH KHOA HÀ NỘI  
HANOI UNIVERSITY  
OF SCIENCE AND TECHNOLOGY

# Machine Learning

IT3190E

Lecture: Linear Regression

ONE LOVE. ONE FUTURE.

# Contents

---

- Lecture 1: Introduction to Machine Learning
- **Lecture 2: Linear regression**
- Lecture 3+4: Clustering
- Lecture 5: Decision tree and Random forest
- Lecture 6: Neural networks
- Lecture 7: Support vector machines
- Lecture 8: Performance evaluation
- Lecture 9: Probabilistic models
- Lecture 10: Ensemble learning
- Lecture 11: Reinforcement learning
- Lecture 12: Regularization
- Lecture 13: Discussion on some advanced topics

# Regression

- There is an *unknown* function  $y^*$  that maps each  $\mathbf{x}$  to a number  $y^*(\mathbf{x})$ 
  - In practice, we can collect some pairs:  $(\mathbf{x}_i, y_i)$ , where  $y_i = y^*(\mathbf{x}_i)$ .
  - Each observation of  $\mathbf{x}$  is represented by a vector in an n-dimensional space, e.g.,  $\mathbf{x}_i = (x_{i1}, x_{i2}, \dots, x_{in})^T$ . Each dimension represents an *attribute (thuộc tính) or feature (đặc trưng) or variate*.
  - Bold characters denote vectors or matrices.
- **Regression problem:** learn a function  $y = f(\mathbf{x})$  from a given training set  $\mathbf{D} = \{(\mathbf{x}_1, y_1), (\mathbf{x}_2, y_2), \dots, (\mathbf{x}_M, y_M)\}$  such that  $y_i \cong f(\mathbf{x}_i)$  for every  $i$

# Linear regression (Hồi quy tuyến tính)

- **Linear model:** assume that  $y^*(\mathbf{x})$  can be well approximated by

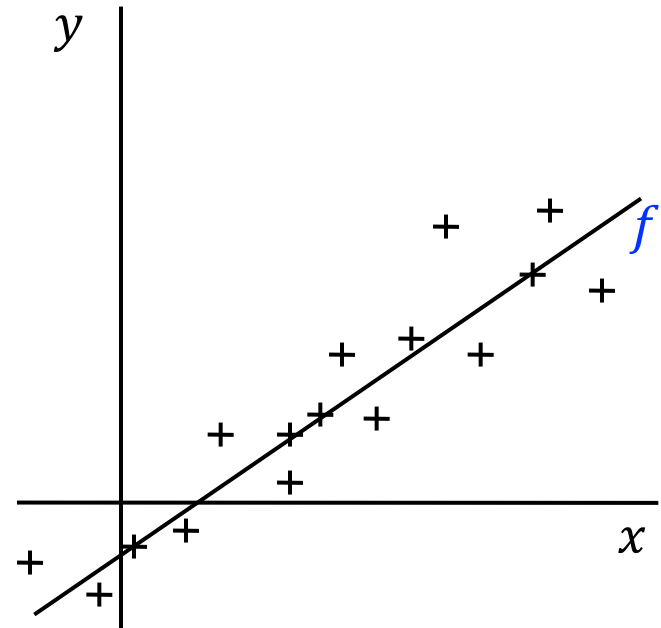
$$f(\mathbf{x}, \mathbf{w}) = w_0 + w_1x_1 + \dots + w_nx_n$$

- $w_0, w_1, \dots, w_n$  are the regression coefficients/weights.  
 $w_0$  sometimes is called “*bias*”.
  - In other words, we use a hyperplane to approximate the unknown function.
  - $f(\mathbf{x}, \mathbf{w})$  may not be linear in  $\mathbf{x}$ .
- **Note:** learning a linear model is equivalent to finding the weight vector  $\mathbf{w} = (w_0, w_1, \dots, w_n)^T$ .

# Linear regression: **example**

- What is the best function?

x	y
0.13	-0.91
1.02	-0.17
3.17	1.61
-2.76	-3.31
1.44	0.18
5.28	3.36
-1.74	-2.46
7.93	5.56
...	...



# Prediction

- For each observation  $\mathbf{x} = (x_1, x_2, \dots, x_n)^T$ 
  - The *true output*:  $y^*(\mathbf{x})$  (but unknown for future data)
  - *Prediction* by our linear model:

$$y_x = w_0 + w_1x_1 + \dots + w_nx_n$$

- We often expect  $y_x \cong y^*(\mathbf{x})$ .
- Prediction for a future observation  $\mathbf{z} = (z_1, z_2, \dots, z_n)^T$ 
  - Use the learned function to make prediction

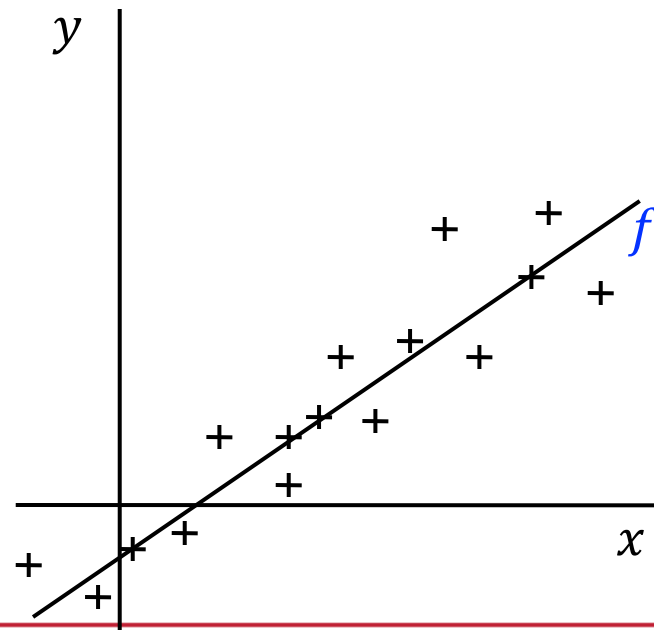
$$f(\mathbf{z}, \mathbf{w}) = w_0 + w_1z_1 + \dots + w_nz_n$$

# Learning a regression function

- **Learning goal:** *learn a function  $f^*$  such that its prediction in the future is the best.*
  - Its generalization is the best.
- **Difficulty:** infinite number of functions

$$H = \{ f(x, \mathbf{w}): \mathbf{w} = (w_0, w_1, \dots, w_n) \in \mathbb{R}^{n+1} \}$$

- How can we learn?
- Is function  $f$  better than  $g$ ?
- Use a measure
  - *Loss function* is often used to guide learning.





# Loss function

- The *error/loss* of the prediction for an example  $\mathbf{x} = (x_1, x_2, \dots, x_n)^T$ :

$$r(f, \mathbf{x}) = [y^*(\mathbf{x}) - f(\mathbf{x}, \mathbf{w})]^2$$

Cost, risk

- The *expected loss* (*risk*) of  $f$  over the whole space:

$$E = E_x[r(f, \mathbf{x})] = E_x[y^*(\mathbf{x}) - f(\mathbf{x}, \mathbf{w})]^2$$

( $E_x$  is the expectation over  $\mathbf{x}$ )

- About the loss/cost:  $r(f, \mathbf{x})$ 
  - Square loss is used above. Other loss functions can be used, e.g.
    - *Absolute loss*:  $|y^*(\mathbf{x}) - f(\mathbf{x}, \mathbf{w})|$
    - *Hinge loss*:  $\max\{0, 1 - y^*(\mathbf{x}) f(\mathbf{x}, \mathbf{w})\}$
    - ...

- The goal of learning is to find  $f^*$  that minimizes the expected loss:

$$f^* = \arg \min_{f \in H} E_x[r(f, x)]$$

- For linear model:  $H$  is the space of functions of linear form.
- But we cannot work directly with this problem during the learning phase.  
(Why?)

# Empirical loss

- We can observe a data set  $\mathbf{D} = \{(\mathbf{x}_1, y_1), (\mathbf{x}_2, y_2), \dots, (\mathbf{x}_M, y_M)\}$ , and have to learn  $f$  from  $\mathbf{D}$ .
- Residual sum of squares:

$$RSS(f) = \sum_{i=1}^M (y_i - f(\mathbf{x}_i, \mathbf{w}))^2 = \sum_{i=1}^M (y_i - w_0 - w_1 x_{i1} - \dots - w_n x_{in})^2$$

- **Empirical loss** (lỗi thực nghiệm):  $L(f, \mathbf{D}) = \frac{1}{M} RSS(f)$ 
  - $L(f, \mathbf{D})$  is an approximation of  $\mathbf{E}_x[r(\mathbf{x})]$ .
- $|L(f, \mathbf{D}) - \mathbf{E}_x[r(\mathbf{x})]|$  is often known as **generalization error** (lỗi tổng quát hoá) of  $f$ .
- Many learning algorithms base on this RSS or its variants.

# Methods: **ordinary least squares (OLS)**

- Given  $\mathbf{D}$ , we find  $f^*$  that minimizes RSS:

$$f^* = \arg \min_{f \in H} \text{RSS}(f) \quad (1)$$
$$\Leftrightarrow \mathbf{w}^* = \arg \min_{\mathbf{w}} \sum_{i=1}^M (y_i - w_0 - w_1 x_{i1} - \cdots - w_n x_{in})^2$$

- This method is often known as *ordinary least squares (OLS, bình phương tối thiểu)*.
- Find  $\mathbf{w}^*$  by taking the gradient of RSS and solving the equation  $\text{RSS}'=0$ . We have:

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

- Where  $\mathbf{A}$  is the data matrix of size  $M \times (n + 1)$ , where the  $i^{\text{th}}$  row is  $\mathbf{A}_i = (1, x_{i1}, x_{i2}, \dots, x_{in})$ ;  $\mathbf{B}^{-1}$  is the inversion of matrix  $\mathbf{B}$ ;  $\mathbf{y} = (y_1, y_2, \dots, y_M)^T$ .
- Note: we assume that  $\mathbf{A}^T \mathbf{A}$  is invertible (ma trận  $\mathbf{A}^T \mathbf{A}$  khả nghịch).**

# Methods: OLS

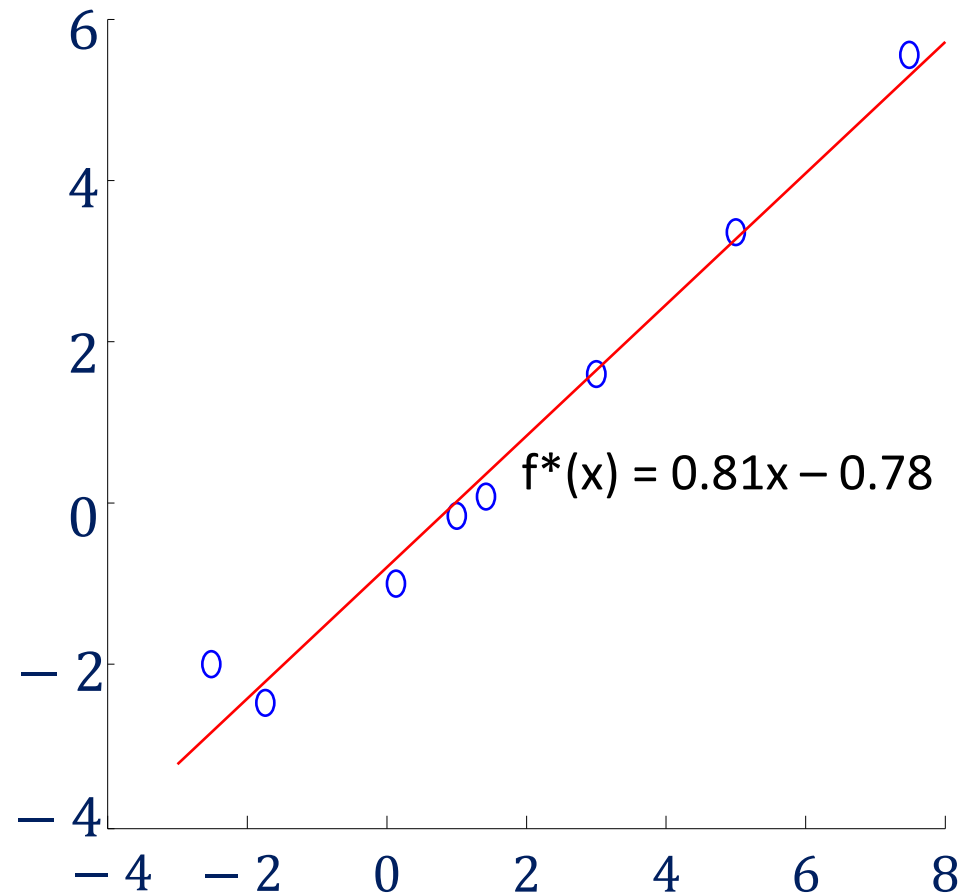
- Input:  $\mathbf{D} = \{(\mathbf{x}_1, y_1), (\mathbf{x}_2, y_2), \dots, (\mathbf{x}_M, y_M)\}$
- Output:  $\mathbf{w}^*$
- Learning: compute

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

- Where  $\mathbf{A}$  is the data matrix of size  $M \times (n + 1)$ , where the  $i^{\text{th}}$  row is  $\mathbf{A}_i = (1, x_{i1}, x_{i2}, \dots, x_{in})$ ;  $\mathbf{B}^{-1}$  is the inversion of matrix  $\mathbf{B}$ ;  
 $\mathbf{y} = (y_1, y_2, \dots, y_M)^T$ .
- Note: we assume that  $\mathbf{A}^T \mathbf{A}$  is invertible.
- Prediction for a new  $\mathbf{x}$ :  $y_x = w_0^* + w_1^* x_1 + \dots + w_n^* x_n$

# Methods: OLS example

x	y
0.13	-1
1.02	-0.17
3	1.61
-2.5	-2
1.44	0.1
5	3.36
-1.74	-2.46
7.5	5.56



# Methods: **limitations of OLS**

- OLS cannot work if  $\mathbf{A}^T\mathbf{A}$  is not invertible
  - If some columns (attributes/features) of  $\mathbf{A}$  are dependent, then  $\mathbf{A}$  will be singular and therefore  $\mathbf{A}^T\mathbf{A}$  is not invertible.  
(Nếu một vài cột của  $\mathbf{A}$  phụ thuộc tuyến tính thì  $\mathbf{A}$  sẽ không khả nghịch)
- OLS requires considerable computation due to the need of computing a matrix inversion.
  - Intractable for the very high dimensional problems.
- OLS likely tends to overfitting, because the learning phase just focuses on minimizing the error of the training data.

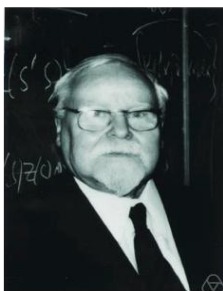
# Methods: Ridge

- Given  $\mathbf{D} = \{(\mathbf{x}_1, y_1), (\mathbf{x}_2, y_2), \dots, (\mathbf{x}_M, y_M)\}$ , we solve for:

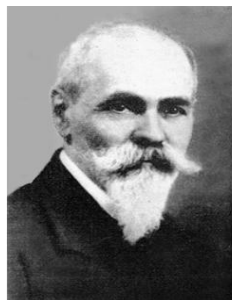
$$f^* = \arg \min_{f \in H} RSS(f) + \lambda \|\mathbf{w}\|_2^2$$

$$\Leftrightarrow \mathbf{w}^* = \arg \min_{\mathbf{w}} \sum_{i=1}^M (y_i - \mathbf{A}_i \mathbf{w})^2 + \lambda \sum_{j=0}^n w_j^2 \quad (2)$$

- Where  $\lambda$  is a regularization constant ( $\lambda > 0$ ),  $\|\mathbf{w}\|_2$  is the  $L^2$  norm.



Tikhonov,  
smoothing an ill-  
posed problem



Zaremba, model  
complexity  
minimization



Bayes: priors  
over parameters



Andrew Ng: need no  
maths, but it prevents  
overfitting!



- Problem (2) is equivalent to the following:

$$\mathbf{w}^* = \arg \min_{\mathbf{w}} \sum_{i=1}^M (y_i - \mathbf{A}_i \mathbf{w})^2 \quad \text{Subject to } \sum_{j=0}^n w_j^2 \leq t \quad (3)$$

- for some constant  $t$ .

- The **regularization/penalty** term:  $\lambda \|\mathbf{w}\|_2^2$

- Limits the magnitude/size of  $\mathbf{w}^*$  (i.e., reduces the search space for  $\mathbf{f}^*$ ).
- Helps us to trade off between *the fitting of  $f$  on  $\mathbf{D}$*  and *its generalization* on future observations.

- We solve for  $\mathbf{w}^*$  by taking the gradient of the objective function in (2), and then zeroing it. Therefore we obtain:

$$\mathbf{w}^* = (\mathbf{A}^T \mathbf{A} + \lambda \mathbf{I}_{n+1})^{-1} \mathbf{A}^T \mathbf{y}$$

- Where  $\mathbf{A}$  is the data matrix of size  $M \times (n + 1)$ , where the  $i^{\text{th}}$  row is  $\mathbf{A}_i = (1, x_{i1}, x_{i2}, \dots, x_{in})$ ;  $\mathbf{y} = (y_1, y_2, \dots, y_M)^T$ ;  $\mathbf{I}_{n+1}$  is the identity matrix of size  $n + 1$ .
- Compared with OLS, Ridge can
  - Avoid the cases of singularity, unlike OLS. Hence Ridge always works.
  - Reduce overfitting.
  - Increase the error for the training set.
- **Note:** *the predictiveness of Ridge depends heavily on the choice of  $\lambda$ .*

# Methods: Ridge

- Input:  $\mathbf{D} = \{(\mathbf{x}_1, y_1), (\mathbf{x}_2, y_2), \dots, (\mathbf{x}_M, y_M)\}$  and  $\lambda > 0$
- Output:  $\mathbf{w}^*$
- Learning: compute

$$\mathbf{w}^* = (\mathbf{A}^T \mathbf{A} + \lambda \mathbf{I}_{n+1})^{-1} \mathbf{A}^T \mathbf{y}$$

- Prediction for a new  $\mathbf{x}$ :

$$y_x = w_0^* + w_1^* x_1 + \dots + w_n^* x_n$$

- **Note:** to avoid some negative effects of the magnitude of  $y$  on covariates  $\mathbf{x}$ , one should remove  $w_0$  from the penalty term in (2). In this case, the solution of  $\mathbf{w}^*$  should be modified slightly.

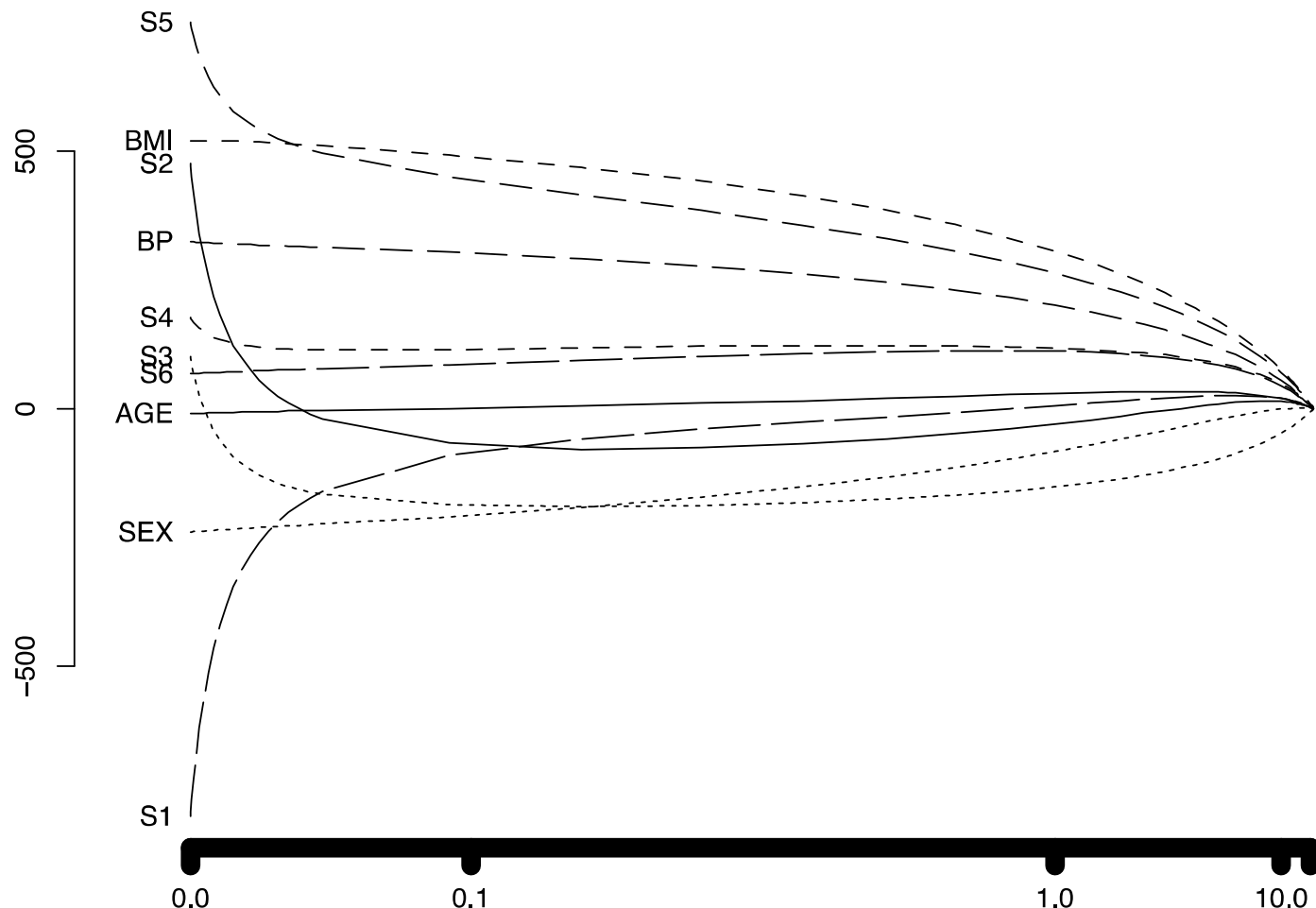
# An example of using Ridge and OLS

- The training set **D** contains 67 observations on prostate cancer, each was represented with 8 attributes. Ridge and OLS were learned from **D**, and then predicted 30 new observations.

w	Ordinary Least Squares	Ridge
0	2.465	2.452
lcavol	0.680	0.420
lweight	0.263	0.238
age	-0.141	-0.046
lbph	0.210	0.162
svi	0.305	0.227
lcp	-0.288	0.000
gleason	-0.021	0.040
pgg45	0.267	0.133
<b>Test RSS</b>	<b>0.521</b>	<b>0.492</b>

# Effects of $\lambda$ in Ridge regression

- $\mathbf{W}^* = (w_0, S1, S2, S3, S4, S5, S6, \text{AGE}, \text{SEX}, \text{BMI}, \text{BP})$  changes as the regularization constant  $\lambda$  changes.



- Ridge regression use  $L^2$  norm for regularization:

$$\mathbf{w}^* = \arg \min_{\mathbf{w}} \sum_{i=1}^M (y_i - \mathbf{A}_i \mathbf{w})^2, \text{ subject to } \sum_{j=0}^n w_j^2 \leq t \quad (3)$$

- Replacing  $L^2$  by  $L^1$  norm will result in LASSO:

$$\mathbf{w}^* = \arg \min_{\mathbf{w}} \sum_{i=1}^M (y_i - \mathbf{A}_i \mathbf{w})^2$$

Subject to  $\sum_{j=0}^n |w_j| \leq t$

- Equivalently:

$$\mathbf{w}^* = \arg \min_{\mathbf{w}} \sum_{i=1}^M (y_i - \mathbf{A}_i \mathbf{w})^2 + \lambda \|\mathbf{w}\|_1 \quad (4)$$

- This problem is non-differentiable  $\rightarrow$  the training algorithm should be more complex than Ridge.

# LASSO: regularization role

- The regularization types lead to different domains for  $\mathbf{w}$ .
- LASSO often produces **sparse** solutions, i.e., many components of  $\mathbf{w}$  are zero.
  - Shrinkage and selection at the same time

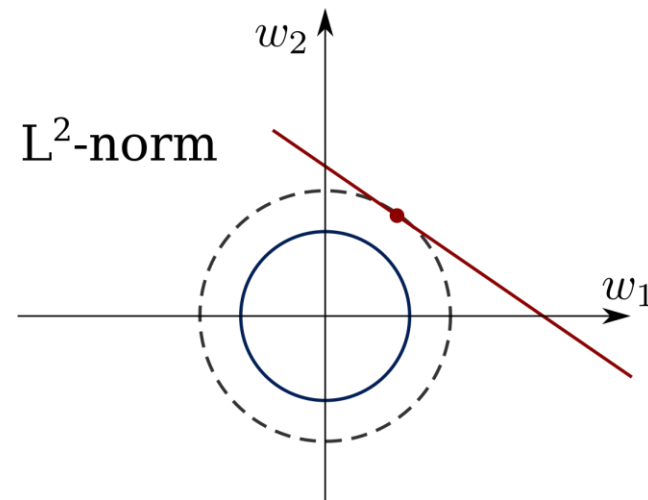
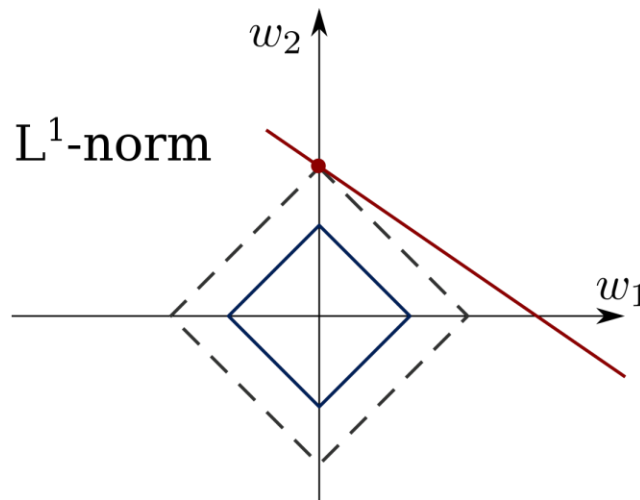


Figure by Nicoguaro - Own work, CC BY 4.0,  
<https://commons.wikimedia.org/w/index.php?curid=58258966>

# OLS, Ridge, and LASSO

- The training set **D** contains 67 observations on prostate cancer, each was represented with 8 attributes. OLS, Ridge, and LASSO were trained from **D**, and then predicted 30 new observations.

w	Ordinary Least Squares	Ridge	LASSO
0	2.465	2.452	2.468
lcavol	0.680	0.420	0.533
lweight	0.263	0.238	0.169
age	-0.141	-0.046	
lbph	0.210	0.162	0.002
svi	0.305	0.227	0.094
lcp	-0.288	0.000	
gleason	-0.021	0.040	
pgg45	0.267	0.133	
<b>Test RSS</b>	<b>0.521</b>	<b>0.492</b>	<b>0.479</b>

Some weights  
are 0  
→ some  
attributes  
may not be  
important



# References

- Hesterberg, T., Choi, N. H., Meier, L., & Fraley, C. (2008). Least angle and L1 penalized regression: A review. *Statistics Surveys*.
- Trevor Hastie, Robert Tibshirani, Jerome Friedman. *The Elements of Statistical Learning*. Springer, 2009.
- Tibshirani, Robert (1996). Regression Shrinkage and Selection via the lasso. *Journal of the Royal Statistical Society. Series B (methodological)*. Wiley. 58 (1): 267–88.

A decorative graphic on the left side of the slide. It features a dark blue background with a large, stylized circular pattern composed of many small red dots. The dots are arranged in a way that creates a sense of depth and movement, resembling a spiral or a series of concentric circles that are slightly offset from each other.

**HUST**

**THANK YOU !**