

IFT6390 Fondements de l'apprentissage machine

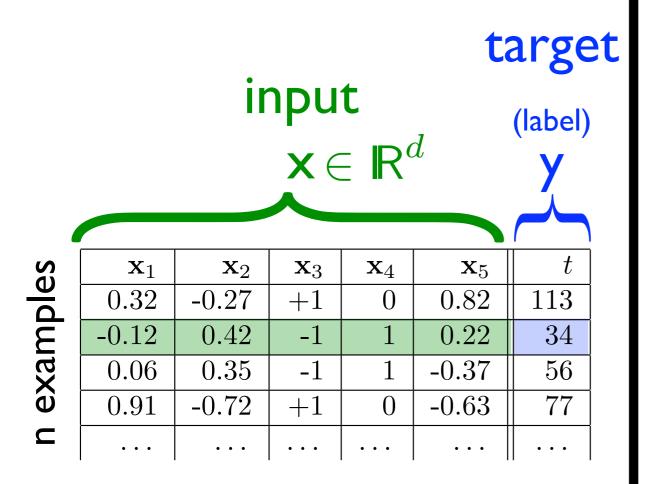
Linear regression and Regularized linear regression

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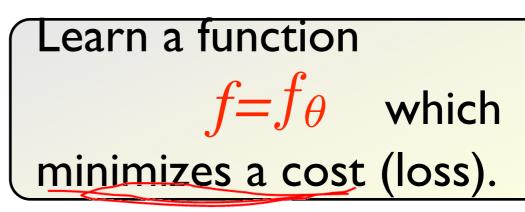
Supervised task

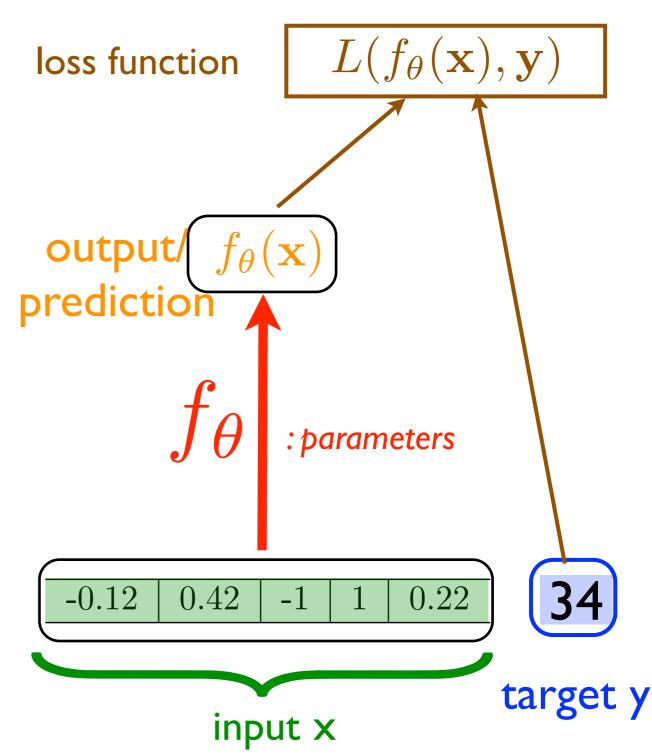
predict y from x



Training set Dn

$$D_n = \{(x^{(1)}, t^{(1)}), \dots, (x^{(n)}, t^{(n)})\}$$





Empirical risk minimization

We must specify:

- ullet A parametric form for our functions, $f_{ heta}$
- ullet A specific cost (loss) function $\ L(y,t)$

So we define the empirical risk as:

$$\hat{R}(f_{\theta}, D_n) = \sum_{i=1}^{n} L(f_{\theta}(\mathbf{x}^{(i)}), t^{(i)})$$

i.e. total loss on the training set

Learning amounts to finding the optimal values for the parameters:

$$\theta^* = \underset{\theta}{\operatorname{arg\,min}} \hat{R}(f_{\theta}, D_n)$$

It is the principle of empirical risk minimization.

Eg: Linear regression

A very simple learning algorithm

We select

A linear (affine) form for the function:

$$f_{\theta}(\mathbf{x}) = \langle \mathbf{w}, \mathbf{x} \rangle + b$$

scalar product (inner product)

Cost: quadratic error:

$$L(y,t) = (y-t)^2$$

t---> v^

We look for the parameters that minimize the empirical risk

Principle of empirical risk minimization (ERM)

$$\theta^* = \underset{\theta}{\operatorname{arg\,min}} \, \hat{R}(f_{\theta}, D_n)$$

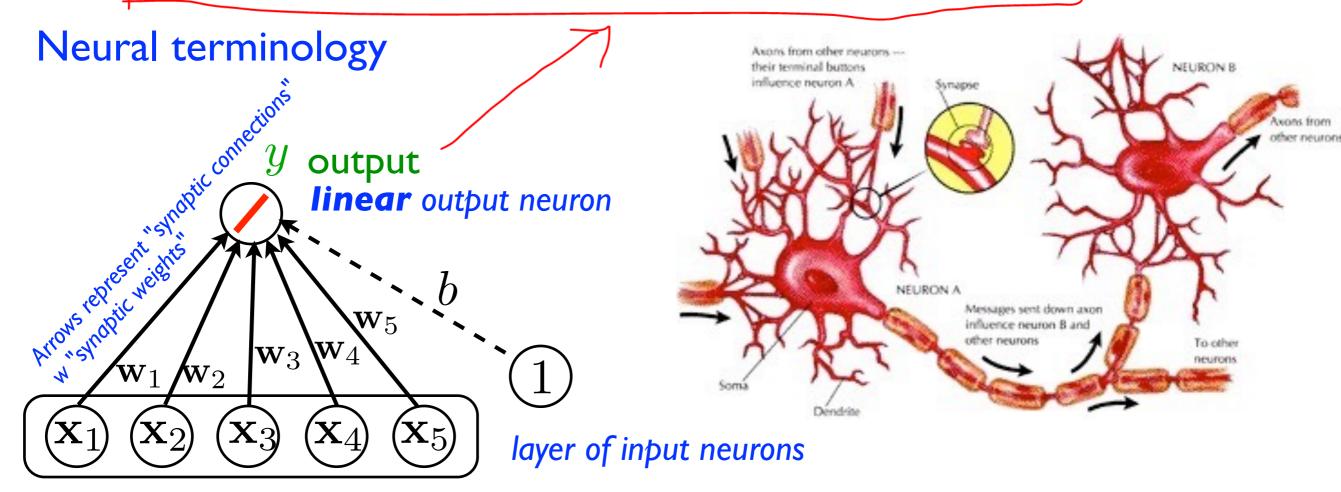
$$\theta = \{\mathbf{w}, b\}, \mathbf{w} \in \mathbb{R}^d, b \in \mathbb{R}$$
 "weight vector" bias

Linear regression

Neural inspiration

Intuitive understanding of the scalar product each component of x has a weighted influence on the output y

$$y = f_{\theta}(\mathbf{x}) = \mathbf{w}_1 \mathbf{x}_1 + \mathbf{w}_2 \mathbf{x}_2 + \ldots + \mathbf{w}_d \mathbf{x}_d + b$$



input (observation) X

Regularized empirical risk

It is often necessary to induce a "preference" for some parameter values rather than others to avoid overfitting

We define regularized empirical risk as follows:

$$\hat{R}_{\lambda}(f_{\theta}, D_{n}) = \underbrace{\left(\sum_{i=1}^{n} L(f_{\theta}(\mathbf{x}^{(i)}), t^{(i)})\right)}_{\text{empirical risk}} + \underbrace{\lambda\Omega(\theta)}_{\text{regularization term (penalty)}}$$

If we go toward, the complex model more, then the penalize alswo will be more

When we dont have additional data, we can use reg. to avoid overfitting

 Ω penalizes more or less the different parameter values. $\lambda \ge 0$ the importance of this regularization term

(in relation to the empirical risk)

Eg: Ridge Regression

= linear regression + quadratic (L2) regularization

We penalize the large weights

$$\Omega(\theta) = \Omega(\mathbf{w}, b) = \|\mathbf{w}\|^2 = \sum_{j=1}^{a} \mathbf{w}_j^2$$

Neural terminology:

"weight decay" penalty

$$\hat{R}_{\lambda}(f_{\theta}, D_{n}) = \underbrace{\left(\sum_{i=1}^{n} L(f_{\theta}(\mathbf{x}^{(i)}), t^{(i)})\right)}_{\text{regularization term (penalty)}} + \underbrace{\lambda\Omega(\theta)}_{\text{regularization term (penalty)}}$$

Eg: Ridge regression

= linear regression + quadratic (L2) regularization

L2 also known as weight decay(in neural network)

Regularized empirical risk

$$\hat{R}_{\lambda}(f_{\theta}, D_{n}) = \underbrace{\left(\sum_{i=1}^{n} L(f_{\theta}(\mathbf{x}^{(i)}), t^{(i)})\right)}_{\text{in Empirical risk}} + \underbrace{\lambda\Omega(\theta)}_{\text{regularization term (penalty)}}$$

We are looking for the parameter values that minimize this objective

$$\{\mathbf{w}^*, b^*\} = \boldsymbol{\theta}^* = \underset{\boldsymbol{\theta}}{\operatorname{argmin}} \hat{R}_{\lambda}(f_{\boldsymbol{\theta}}, D_n)$$

Eg: Ridge Regression

- = linear regression + quadratic (L2) regularization
 - For linear regression or ridge regression a little linear algebra gives us an analytical solution

we solve for $\theta = \{b, \mathbf{w}\}$:

$$\frac{\partial \hat{R}_{\lambda}(f_{\theta}, D_n)}{\partial \theta} = 0$$

we obtain:
$$\begin{pmatrix} b^* \\ \mathbf{w}^* \end{pmatrix} = (\check{X}^T\check{X} + \lambda\check{I})^{-1}\check{X}^T\mathbf{t}$$

This is a closed form solution

$$\text{où } \check{X} = \begin{pmatrix}
 1 & \mathbf{x}_{1}^{(1)} & \dots & \mathbf{x}_{d}^{(1)} \\
 \vdots & \vdots & \ddots & \vdots \\
 1 & \mathbf{x}_{1}^{(n)} & \dots & \mathbf{x}_{d}^{(n)}
 \end{pmatrix}, \mathbf{t} = \begin{pmatrix}
 t^{(1)} \\
 \vdots \\
 t^{(n)}
 \end{pmatrix}
 \quad
 \check{I} = \begin{pmatrix}
 0 & 0 & & & \\
 0 & 1 & 0 & & \\
 0 & 1 & 0 & & \\
 0 & \ddots & 0 & \\
 0 & 1 & 0 & \\
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- Most of the time (other choices of f and L) we do not have an analytical solution.
- More generally, we can use a gradient descent method.

$$f(x) = x^{T}\Theta \qquad \theta = w \quad , \text{ no blas}$$

$$D_{n} = \left\{ (x_{1}, x_{1}), \dots (x_{n}, v_{n}) \right\} \quad x_{1} \in \mathbb{R}^{d}$$

$$X = \left[(x_{1}, x_{2}, \dots x_{n}) \right]^{T} \in \mathbb{R}^{d}$$

$$Y = \left[(v_{1}, v_{2}, \dots v_{n}) \right]^{T} \in \mathbb{R}^{d}$$

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 $X^TX\Theta^* = X^Ty$ $\Theta^* = (X^TX)^{-1}X^Ty$

Other possibility:

optimization by gradient descent

 \hat{R}_{λ}

$$\hat{R}_{\lambda}(f_{ heta}, D_n) = \underbrace{\left(\sum_{i=1}^n L(f_{ heta}(\mathbf{x}^{(i)}), t^{(i)})\right)}_{\text{regularization term}} + \lambda \Omega(\theta)$$

- we initialize the parameters randomly
- we update them iteratively following the gradient

Either batch gradient descent (whole dataset):

Loop:
$$\theta \leftarrow \theta - \eta \frac{\partial \hat{R}_{\lambda}}{\partial \theta}$$

$$= \left(\sum_{i=1}^{n} \frac{\partial}{\partial \theta} L(f_{\theta}(\mathbf{x}^{(i)}), t^{(i)}) \right) + \lambda \frac{\partial}{\partial \theta} \Omega(\theta)$$

Or stochastic gradient descent:

Loop:

For i in 1...n
$$\theta \leftarrow \theta - \eta \frac{\partial}{\partial \theta} \left(L(f_{\theta}(\mathbf{x}^{(i)}), t^{(i)}) + \frac{\lambda}{n} \Omega(\theta) \right)$$

Or other variants of the gradient descent idea (conjugate gradient, Newton's method, natural gradient, ...)

Various regularizers

«Ridge»: regularization, L_2

In Bayesian terms: corresponds to a Gaussian prior on the weights

$$\Omega(\theta) = \Omega(\mathbf{w}, b) = \|\mathbf{w}\|_2^2 = \sum_{j=1}^{\infty} \mathbf{w}_j^2$$

«Lasso»: regularization, L_1

In Bayesian terms: corresponds to a Laplacian prior on the weights

$$\Omega(\theta) = \Omega(\mathbf{w}, b) = ||\mathbf{w}||_1 = \sum_{j=1}^{n} |\mathbf{w}_j|$$

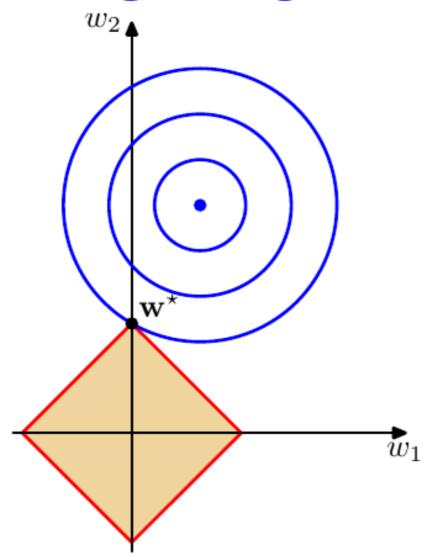
=> automatic selection of components (a number of weights will be zero)

«Elastic net»: combination of the two

$$\Omega(\theta) = \Omega(\mathbf{w}, b) = \lambda_1 \|\mathbf{w}\|_1 + \lambda_2 \|\mathbf{w}\|_2^2$$

Etc...

Visualizing L_1 regularization

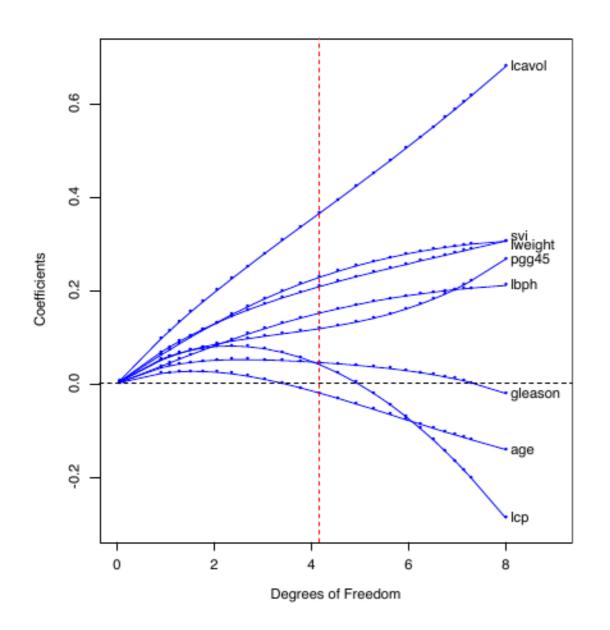


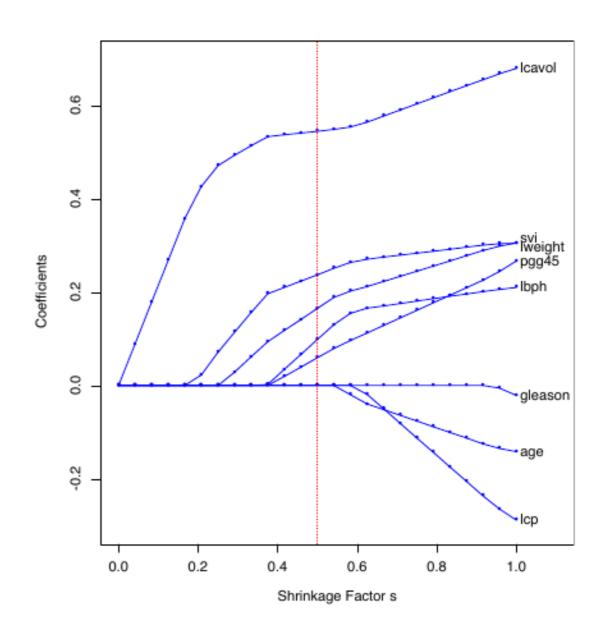
- ullet If λ is big enough, the circle is very likely to intersect the diamond at one of the corners
- This makes L_1 regularization much more likely to make some weights exactly 0

Pros and cons of L_1 regularization

- If there are irrelevant input features, Lasso is likely to make their weights 0, while L_2 is likely to just make all weights small
- Lasso is biased towards providing sparse solutions in general
- ullet Lasso optimization is computationally more expensive than L_2
- More efficient solution methods have to be used for large numbers of inputs (e.g. least-angle regression, 2003).
- \bullet L_1 methods of various types are very popular
- ullet One can combine L_1 and L_2 regularization (elastic-net)

Example of L1 vs L2 effect





- ullet Note the sparsity in the coefficients induces by L_1
- ullet Lasso is an efficient way of performing the L_1 optimization