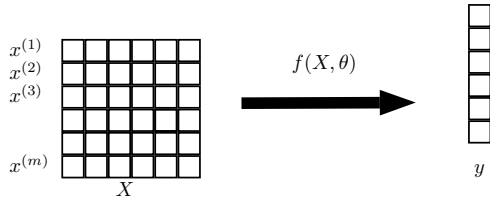
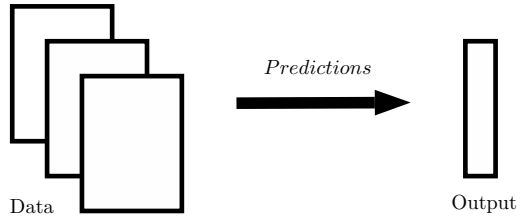


**POLIMI** GRADUATE  
SCHOOL OF **MANAGEMENT**

# MACHINE LEARNING IN PYTHON - LAB SESSION

Andrea Mor - [andrea.mor@polimi.it](mailto:andrea.mor@polimi.it) - <https://github.com/mrondr/BABD2023>

# SUPERVISED LEARNING



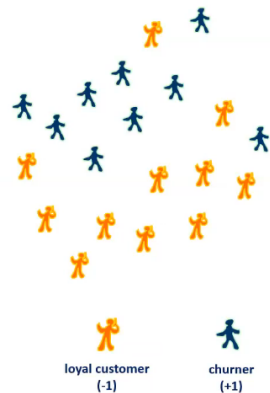
# THE PROBLEM: BANK TELEMARKETING<sup>1</sup>

Attribute		Type	Description/Values
Personal	age	num	Age of the potential client
	job	cat	admin., blue- collar, entrepreneur, housemaid, ... , unknown
	marital_status	cat	divorced, married, single, unknown
	education	cat	basic.4y, basic.6y, basic.9y, high.school, ... , unknown
Bank	default	cat	The client has credit in default: no,yes,unknown
	housing loan	cat	The client has a housing loan contract: no,yes,unknown
	loan	cat	The client has a personal loan: no,yes,unknown
Campain	contact	cat	Communication type: cellular,telephone
	month	cat	Last month contacted: jan, feb ,..., dec
	day_of_week	cat	Last contact day : mon, tue,..., fri
	duration	num	Last contact duration (in seconds)
	campaign	num	Number of contacts performed during this campaign
	pdays	num	Number of days that passed by after last contact
	previous poutcome	cat	Number of contacts performed before this campaign Outcome of the prev. marketing campaign: failure,nonexistent,success
Economical	emp.var.rate	num	Employment variation rate in the last quarter
	cons.price.idx	num	Consumer price index in the last month
	cons.conf.idx	num	Monthly consumer confidence index
	euribor3m	num	Dayly Euro Interbank Offered Rate
	nr.employed	num	Number of employed citizens in the last quarter (thousands)
<b>Target</b>	<b>success</b>	<b>target</b>	<b>O: no, 1: yes</b>

<sup>1</sup> A data-driven approach to predict the success of bank telemarketing. S. Moroa, P. Cortez, P. Rita. Decision Support Systems, 62:22-31, 2014.

# CLASSIFICATION PROBLEM

attributes						
customers	Area	Pothers	Pmob	...	NumSMS	Class
	2	0.14	0.59	...	18	1
	3	0.26	0.35	...	9	-1
	1	0.37	0.23	...	1	1
	⋮	⋮	⋮	⋮	⋮	⋮
	4	0.41	0.27	...	64	-1
	← past data →					
Area	Pothers	Pmob	...	NumSMS	Class	
1	0.27	0.67	...	36	?	
4	0.44	0.22	...	50	?	
4	0.31	0.47	...	14	?	
⋮	⋮	⋮	⋮	⋮	⋮	
2	0.31	0.14	...	49	?	
← future data →						

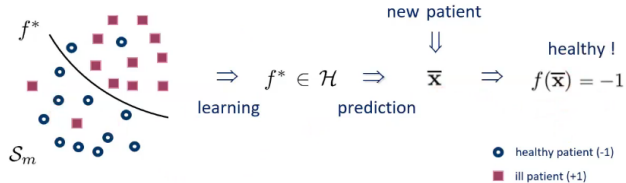


# CLASSIFICATION FORMULATION

$S_m = \{(\mathbf{x}_i, y_i), i \in \mathcal{M}\}$  : **training set, where**  $\mathbf{x}_i \in \mathbb{R}^n$  **and**  $y_i \in \mathcal{D}$

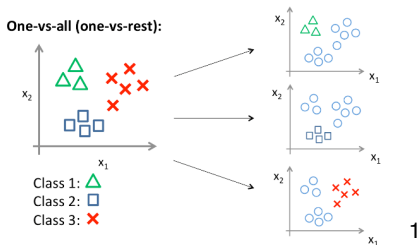
$\mathcal{H}$  **denotes a set of functions**  $f(\mathbf{x}) : \mathbb{R}^n \mapsto \mathcal{D}$

**Classification problem: define a hypotheses space**  $\mathcal{H}$  **and a function**  $f^* \in \mathcal{H}$  **which optimally describes the relationship between**  $\mathbf{x}_i$  **and**  $y_i$



# MULTI-CLASS CLASSIFICATION

1. **One-vs-Rest** We perform  $|H|$  different binary classifications: one for every class.



We decide based on a majority vote.

2. **One-vs-One** We perform  $|H|(|H| - 1)/2$  binary classifications: one for every pair of classes. We decide based on a majority vote.

# CLASSIFICATION MODELS

- ▶ Heuristics Methods
  - Nearest Neighbours
  - Classification Trees
- ▶ Probabilistic Methods
  - Bayesian Methods
- ▶ Regression Methods
  - Logistic regression
- ▶ Separation Methods
  - Support vector machine
  - Perceptron
  - Neural Networks



# EVALUATION DIMENSIONS

- ▶ Prediction accuracy
- ▶ Speed
- ▶ Robustness
- ▶ Scalability
- ▶ Interpretability
- ▶ Rules effectiveness

# QUALITY MEASURES - CONFUSION MATRIX

		Prediction outcome	
		0	1
Actual value	0	True Negative	False Positive
	1	False Negative	True Positive

► **Accuracy** =  $\frac{TP+TN}{TP+FP+TN+FN}$

**When:** balanced problems, equal costs of FN and FP.

**When NOT:** imbalanced problems.

► **Recall** (True Positive rate) =  $\frac{TP}{FN+TP}$

*“proportion of true positives among actual positive”*

**When:** you want to be sure to catch all the 1s, at the cost of having false alerts.

► **Precision** (Positive Predictive value) =  $\frac{TP}{TP+FP}$

*“proportion of true positives among positive predictions”*

**When:** you want to be sure that all that you predict as 1, are indeed 1, e.g., false alerts are costly.

► Precision and Recall are typically evaluated together.

# QUALITY MEASURES - CONFUSION MATRIX

		Prediction outcome	
		0	1
Actual value	0	True Negative	False Positive
	1	False Negative	True Positive

- ▶ **Geom. mean** =  $\sqrt{\text{Precision} \times \text{Recall}}$
- ▶ **F-score** =  $\frac{(1+\beta^2)}{\beta^2} \frac{1}{\frac{1}{\text{Precision}} + \frac{1}{\text{Recall}}} = (1 + \beta^2) \frac{\text{Precision} \cdot \text{Recall}}{(\beta^2 \cdot \text{Precision}) \cdot \text{Recall}}$
- ▶ **F1**: harmonic mean of Prec. and Recall  
$$= \frac{2TP}{2TP + FP + FN}$$

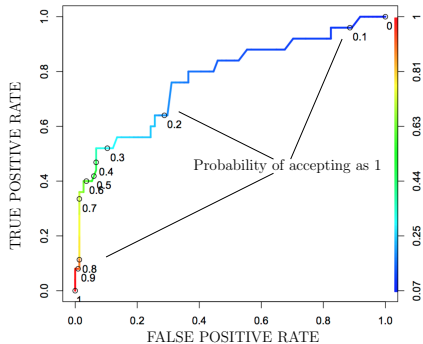
Frequently the to-go metric.
- ▶ **F2**: 2x emphasis on Recall  
**When:** False Negatives are more costly.

# QUALITY MEASURES - CONFUSION MATRIX (AUXILLIARY)

		Prediction outcome	
		0	1
Actual value	0	True Negative	False Positive
	1	False Negative	True Positive

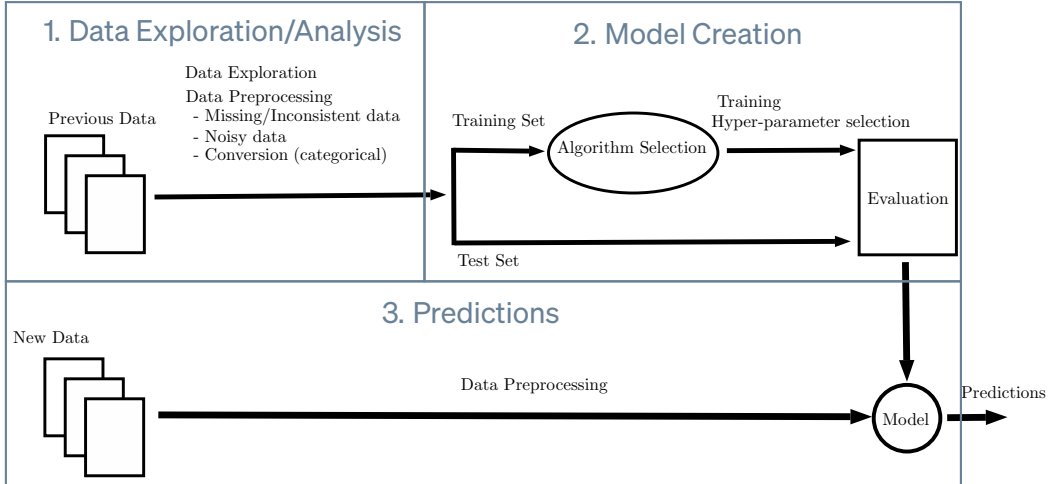
- ▶ **False Positive rate** =  $\frac{FP}{FP+TN}$   
“proportion of false positives among actual negatives”  
**When:** cost of dealing with positive prediction is high (e.g., false alarms).
- ▶ **True Negative rate (Specificity)** =  $\frac{TN}{TN+FP}$   
**When:** e.g., you want to be sure when you say something is safe for health.

# QUALITY MEASURES - ROC CURVE & AUC

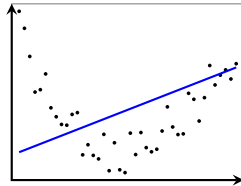
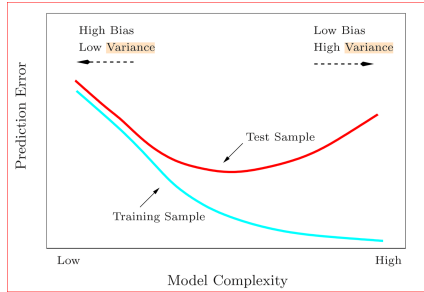


- ▶ If we accepting even with small probability then  $TPR = FPR = 1$
- ▶ If we accepting just with high probability then  $TPR = FPR = 0$
- ▶ The perfect classifier is the the point (0, 1)
- ▶  $AUC \in [0.5, 1]$  area under the curve is a quality measure of our algorithm.

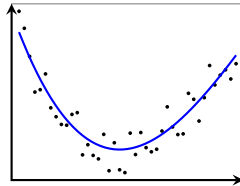
# WORKFLOW



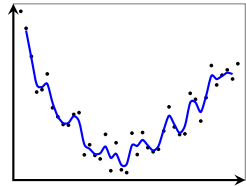
# UNDER/OVER-FITTING



Underfitting

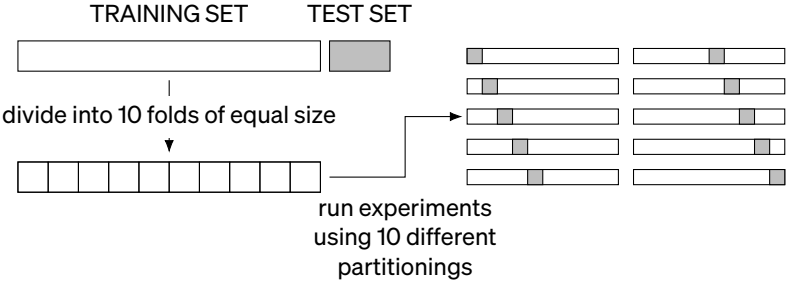


Balance



Overfitting

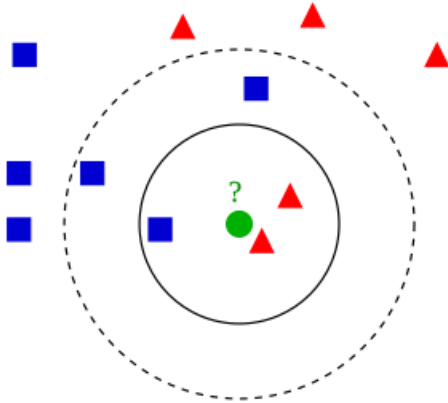
# CROSS VALIDATION





# THE ALGORITHMS

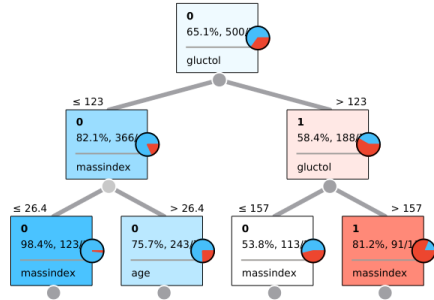
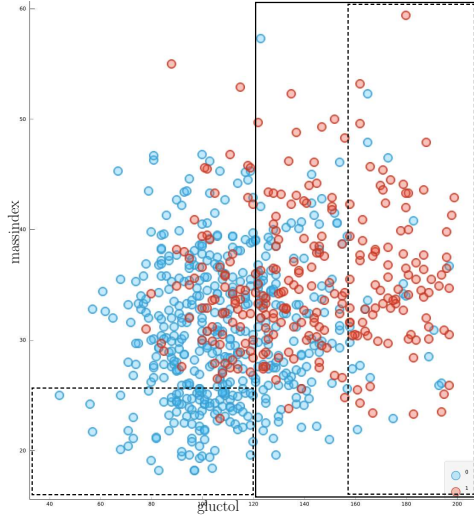
# K-NEAREST NEIGHBORS



## Main Parameters

- ▶  $k$  : number of neighbors
- ▶ neighbor weights
- ▶ distances

# CLASSIFICATION TREE

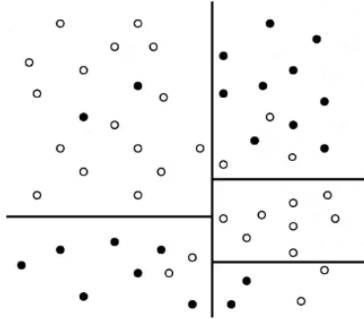


Tree

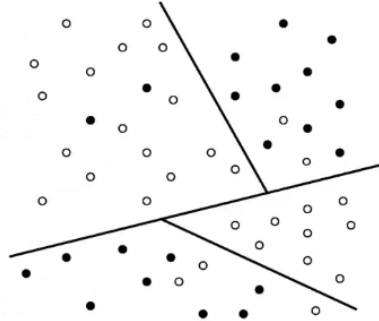
## types

- ▶ Binary tree (zero/two descendants)
- ▶ General trees
- ▶ Uni-variate tree ( $X_j < b$ )
- ▶ Multi-variate tree ( $\sum_{j=1}^n w_j x_j < b$ )

# CLASSIFICATION TREE

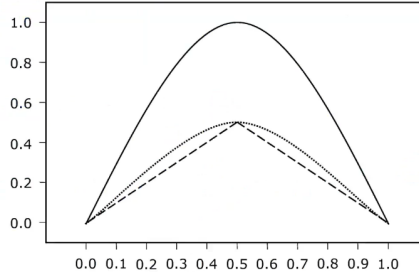
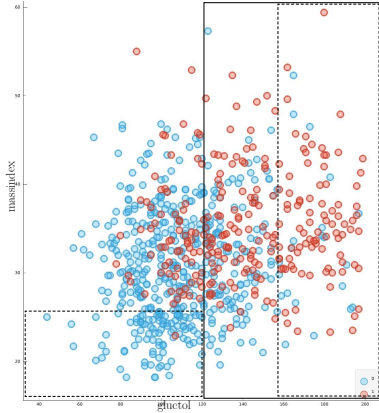


classification by an  
axis parallel tree



classification by an  
oblique tree

# CLASSIFICATION TREE



- ▶ **Entropy index:**
- ▶ **Gini index:**
- ▶ **Miss-classification index:**

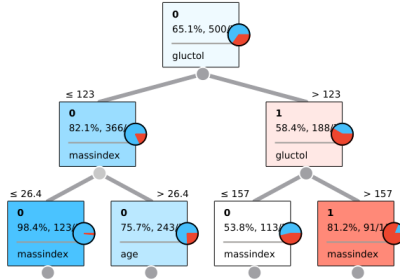
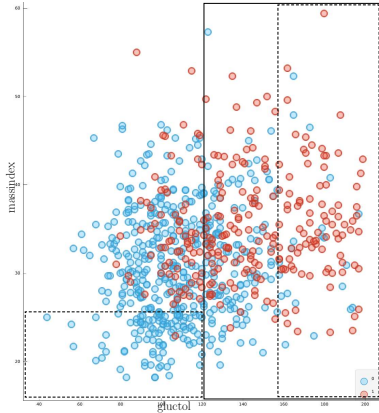
## Split criteria

$$-\sum_{h=1}^H f_h \log_2 f_h$$

$$1 - \sum_{h=1}^H f_h^2$$

$$1 - \max_h f_h$$

# CLASSIFICATION TREE



## Main Parameters

- ▶ impurity measure: “gini”, “entropy”
- ▶ max\_depth
- ▶ min\_samples\_split: minimum number of samples to split an internal node
- ▶ min\_sample\_leaf: minimum number of samples required to be at a leaf node

# NAIVE BAYESIAN CLASSIFIER

- Bayes Theorem

$$P(y|\mathbf{x}) = \frac{P(\mathbf{x}|y)P(y)}{P(\mathbf{x})} = \frac{P(\mathbf{x}|y) P(y)}{\sum_{l=1}^H P(\mathbf{x}|y) P(y)}$$

- Maximum a posteriori hypothesis

$$y_{MAP} = \arg \max_{y \in \mathcal{H}} P(y|\mathbf{x}) = \arg \max \frac{P(\mathbf{x}|y)P(y)}{P(\mathbf{x})}$$

- Independence (Naive)

$$P(\mathbf{x}|y) = P(x_1|y) \cdot P(x_2|y) \cdot \dots \cdot P(x_n|y) = \prod_{j=1}^n P(x_j|y)$$

# NAIVE BAYESIAN CLASSIFIER

- Categorical/discrete attributes

$$P(x_j|y) = P(x_j = r_{jk}|y = v_h)$$

→ empirical frequency of the observed value on the class  $v_h$

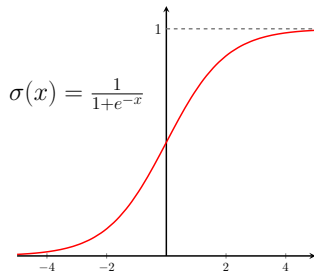
- Numerical attribute

$$P(x_j|y) \sim N(\mu_{jh}, \sigma_{jh})$$

→ assuming Gaussian density with empirical parameters



# LOGISTIC REGRESSION



The model postulates that

$$\log \left[ \frac{P(y = 1|x)}{P(y = 0|x)} \right] = w^T x$$

then if  $p = P(y = 1|x)$ :

$$\frac{p}{1-p} = e^{w^T x}$$

$$\Rightarrow p = P(y = 1|x) = \frac{e^{w^T x}}{1 + e^{w^T x}} = \frac{1}{1 + e^{-w^T x}},$$

$$P(y = 0|x) = 1 - p = \frac{1}{1 + e^{w^T x}} = \frac{e^{-w^T x}}{1 + e^{-w^T x}}$$

# LOGISTIC REGRESSION

Therefore if we maximize the likelihood

$$L(w) := P(Y_1 = y_1, \dots, Y_m = y_m | w, x_1, \dots, x_m) = \prod_{i=1}^n P(Y_i = y_i | w, x_1, \dots, x_m).$$

Assuming independence:

$$L(w) := \prod_{i|y_i=1} p(x_i) \cdot \prod_{i|y_i=0} (1 - p(x_i))$$

$$L(w) := \prod_{i=1}^n p(x_i)^{y_i} (1 - p(x_i))^{1-y_i}$$

is equivalent to maximize the log likelihood

$$\begin{aligned} l(w) &= \sum_{i=1}^n (y_i \log(p(x_i)) + (1 - y_i) \log(1 - p(x_i))) \\ &= \sum_{i=1}^n (y_i \log\left(\frac{p(x_i)}{1 - p(x_i)}\right) + \log(1 - p(x_i))) \end{aligned}$$

Given the quantities defined above

$$l(w) = \sum_{i=1}^n (y_i w^T x_i - \log(1 + e^{-w^T x_i}))$$

Therefore we aim to resolve the following optimization problem

$$\max_w l(w)$$

# LOGISTIC REGRESSION

Regulation term(s) can be considered (recall that  $\max l(w)$  is equivalent to  $\min -l(w)$ ):

$$\min_w \frac{1}{2} ||w||^2 - C \cdot l(w)$$

## Main Parameters

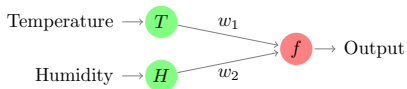
- ▶  $C$ : Inverse of regularization strength
- ▶ Resolution algorithm parameters:
  - solver: lbfgs, newton-cg, liblinear, sag, saga.
  - tol: Tolerance for stopping criteria.
  - max\_iter: max. number of iterations
  - n\_jobs: Number of CPU cores

# MULTI-LAYER PERCEPTRON - SMALL EXAMPLE

<b>Temp. [C]</b>	20	31	15	18	21
<b>Humidity [%]</b>	40	36	23	45	30
<b>Prob. Rain</b>	0.70	0.52	0.55	0.73	0.60

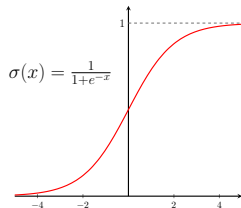
# MULTI-LAYER PERCEPTRON - SMALL EXAMPLE

Temp. [C]	20	31	15	18	21
Humidity [%]	40	36	23	45	30
Prob. Rain	0.70	0.52	0.55	0.73	0.60



$$f(T, H, w_1, w_2) = \underbrace{\sigma}_{\text{activation function}}(w_1 \cdot T + w_2 \cdot H)$$

$$\min_{w_1, w_2} \sum_{i=1}^5 [P_i(\text{rain}) - f(T_i, H_i, w_1, w_2)]^2$$



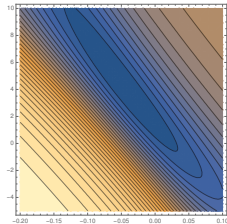
For a classification problem we can use the Likelihood as cost function.

# MULTI-LAYER PERCEPTRON - SMALL EXAMPLE

$$\begin{aligned} & \min_{w_1, w_2} \sum_{i=1}^5 [P_i(\text{rain}) - f(T_i, H_i, w_1, w_2)]^2 \\ &= \min \left[ 0.7 - 1/(1 + e^{-(w_1 \cdot 20 + w_2 \cdot 0.4)}) \right]^2 + \left[ 0.52 - 1/(1 + e^{-(31 \cdot w_1 + w_2 \cdot 0.36)}) \right]^2 + \dots \end{aligned}$$

# MULTI-LAYER PERCEPTRON - SMALL EXAMPLE

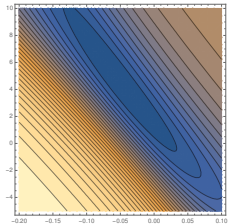
$$\begin{aligned} \min_{w_1, w_2} \sum_{i=1}^5 [P_i(\text{rain}) - f(T_i, H_i, w_1, w_2)]^2 \\ = \min \left[ 0.7 - 1/(1 + e^{-(w_1 \cdot 20 + w_2 \cdot 0.4)}) \right]^2 + \left[ 0.52 - 1/(1 + e^{-(31 \cdot w_1 + w_2 \cdot 0.36)}) \right]^2 + \dots \end{aligned}$$



$$(w_1^*, w_2^*) = (-0.044, 4.147)$$

# MULTI-LAYER PERCEPTRON - SMALL EXAMPLE

$$\begin{aligned} \min_{w_1, w_2} \sum_{i=1}^5 [P_i(\text{rain}) - f(T_i, H_i, w_1, w_2)]^2 \\ = \min \left[ 0.7 - 1/(1 + e^{-(w_1 \cdot 20 + w_2 \cdot 0.4)}) \right]^2 + \left[ 0.52 - 1/(1 + e^{-(31 \cdot w_1 + w_2 \cdot 0.36)}) \right]^2 + \dots \end{aligned}$$

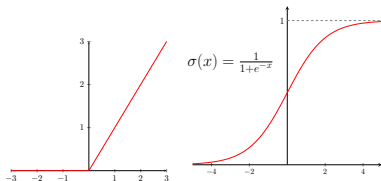
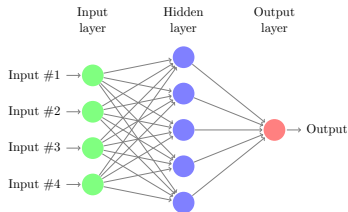


$$(w_1^*, w_2^*) = (-0.044, 4.147)$$

<b>Temp. [C]</b>	20	31	15	18	21
<b>Humidity [%]</b>	40	36	23	45	30
<b>Prob. Rain</b>	0.70	0.52	0.55	0.73	0.60
<b>Predicted</b>	0.70	0.56	0.58	0.75	0.60
<b>Error</b>	<b>0.0</b>	<b>-0.04</b>	<b>-0.03</b>	<b>-0.02</b>	<b>0.0</b>



# MULTI-LAYER PERCEPTRON



ReLU

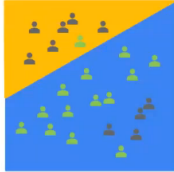
Sigmoid

## Main Parameters

- ▶ hidden\_layer\_sizes:  $(n_1, n_2, \dots, n_L)$
- ▶ activation: identity, logistic, tanh, relu
- ▶ alpha regularization term parameter
- ▶ Resolution algorithm parameters: solver, tol, batch\_size, learning\_rate, max\_iter.

# NEURAL NETWORK

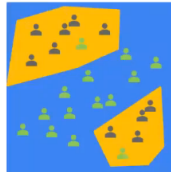
1 LAYER



2 LAYER



3+ LAYER



1 LAYER, 3 NODES

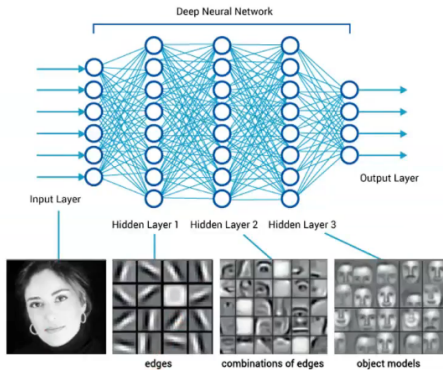
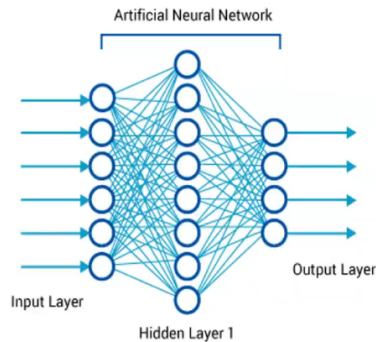


1 LAYER, 6 NODES



1 LAYER, 20 NODES

# DEEP LEARNING



# CLASSIFICATION AND STATISTICAL LEARNING THEORY

Given an hypothesis space  $\mathcal{F}$ , a predictive function  $f \in \mathcal{F}$ , and a loss function  $V(y, f(\mathbf{x}))$ , we define

- ▶ the **empirical risk** on a train set  $\mathcal{T}$  as

$$R_{emp}(f) = \frac{1}{m} \sum_{i=1}^m V(y_i, f(\mathbf{x}_i))$$

- ▶ and the **expected risk** as

$$R(f) = \int V(y, f(\mathbf{x})) dP(\mathbf{x}, y)$$

Overfitting arises when the difference between these errors is large.

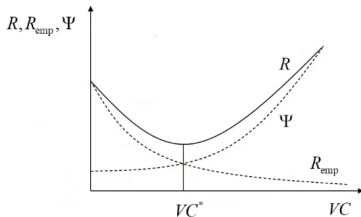
# VC-DIMENSION

The Vapnik–Chervonenkis (VC) dimension is the maximum number of points that can be correctly classified by classifiers in  $\mathcal{F}$ . We can prove that

$$R(f) - R_{emp}(f) \leq \sqrt{\frac{1}{t}(\gamma \log(2t/\gamma) - \log(\eta/4) + 1)} = \Psi(t, \gamma, \eta)$$

with probability  $1 - \eta$ , where

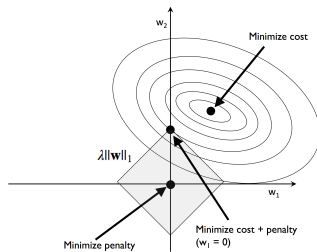
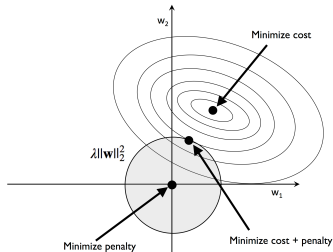
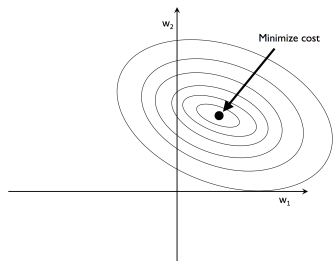
- ▶  $t$  is the number of training points
- ▶  $\gamma$  is the VC-dimension



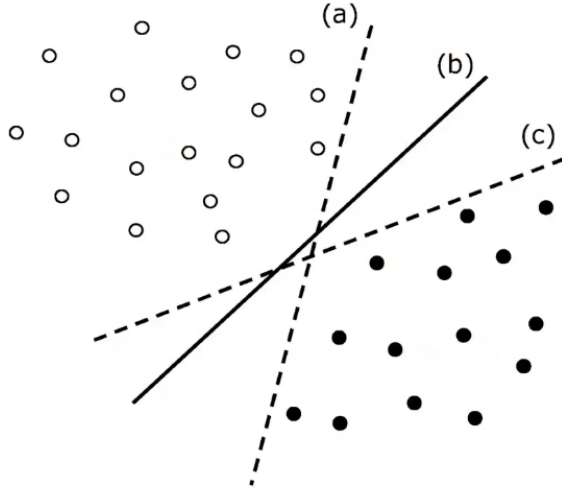
# REGULARIZATION

## Structural risk minimization (SRM)

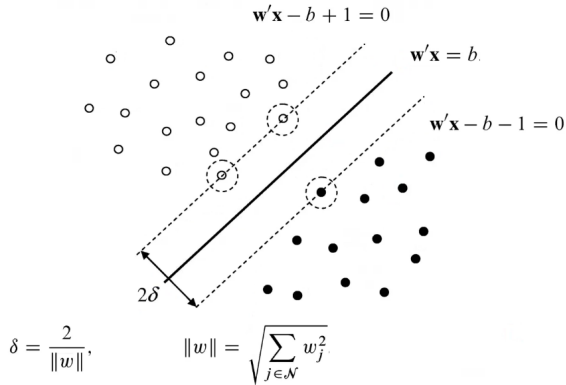
$$\hat{R}(f) = \frac{1}{t} \sum_{i=1}^t V(y_i, f(\mathbf{x}_i)) + \lambda \|f\|_K^2$$



# SUPPORT VECTOR MACHINE - LINEARLY SEPARABLE



# SVM - LINEARLY SEPARABLE

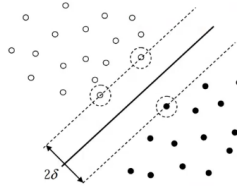


$$\begin{aligned} \min_{\mathbf{w}, b} \quad & \frac{1}{2} \|\mathbf{w}\|_2^2 \\ \text{s. t.} \quad & y_i(\mathbf{w}'\mathbf{x}_i - b) \geq 1 \quad i \in \mathcal{M} \end{aligned}$$

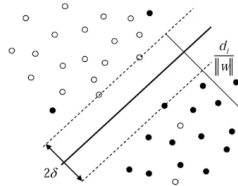


# SVM - GENERAL CASE

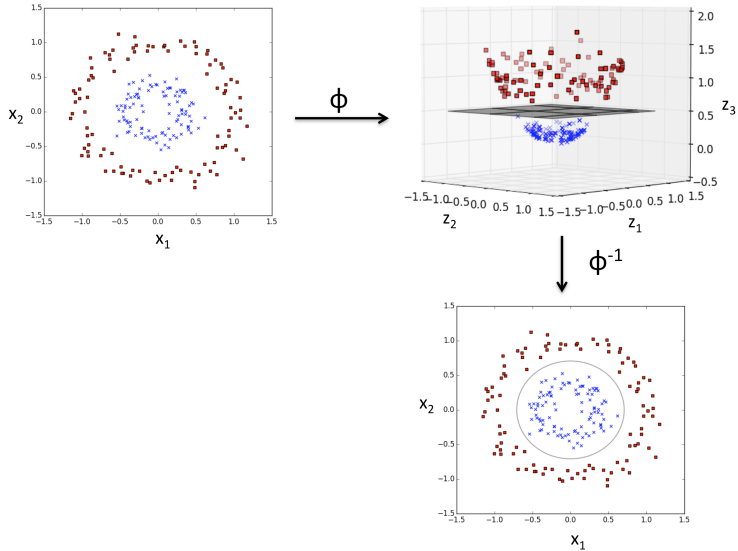
$$\begin{aligned} \min_{\mathbf{w}, b} \quad & \frac{1}{2} \|\mathbf{w}\|_2^2 \\ \text{s. t.} \quad & y_i(\mathbf{w}'\mathbf{x}_i - b) \geq 1 \quad i \in \mathcal{M} \end{aligned}$$



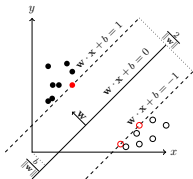
$$\begin{aligned} \min_{\mathbf{w}, b, d} \quad & \frac{1}{2} \|\mathbf{w}\|_2^2 + \lambda \sum_{i=1}^m d_i \\ \text{s. t.} \quad & y_i(\mathbf{w}'\mathbf{x}_i - b) \geq 1 - d_i \quad i \in \mathcal{M} \\ & d_i \geq 0 \quad i \in \mathcal{M} \end{aligned}$$



# SVM - KERNELS



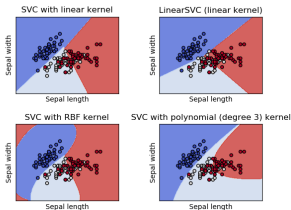
# SVM - GENERAL CASE



$$\min_{w,b,d} \quad \frac{1}{2} ||w||^2 + C \sum_{i=1}^m d_i$$

$$\text{subject to } y_i (w^T \underbrace{\phi(x_i)}_{\text{kernel}} - b) \geq 1 - d_i,$$

$$d_i \geq 0$$

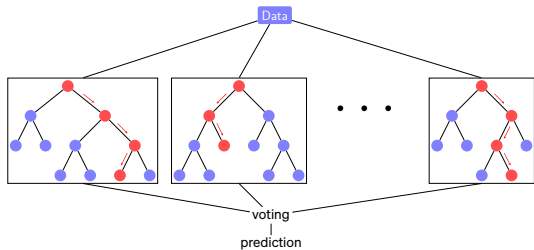


## Main Parameters

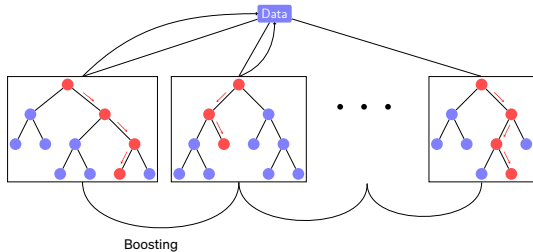
- ▶  $C$ : Inverse of regularization strength
- ▶ kernel:
  - linear:  $x'x$
  - poly:  $(\gamma x'x + r)^d$
  - rbf:  $\exp(-\gamma ||x - x'||^2)$
  - sigmoid:  $\tanh(\gamma x'x + r)$
- ▶  $\text{degree}(d)$ ,  $\text{gamma}(\gamma)$ ,  $\text{coef0}(r)$
- ▶ Resolution algorithm parameters

# ENSEMBLE METHODS

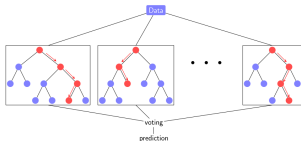
## Bagging



## Boosting



# RANDOM FOREST

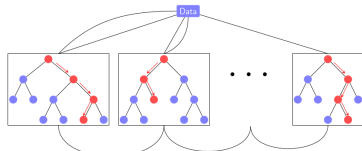


1. Create different (simple) tree models (stumps)
2. Each model is created with a subset of observation/features ( $\sim 2m/3$ )
3. We combine the prediction of all trees

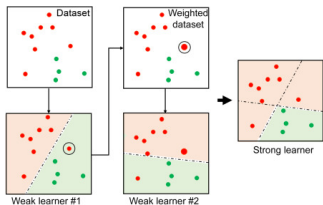
## Main Parameters

- ▶ `n_estimators`: Number of trees
- ▶ `max_features`: Number of features selected for the split
- ▶ `bootstrap=False`: Use all samples
- ▶ Tree parameters

# ADABOOST



Boosting



1. Assign equal weights to observations  $w_i^{(0)} = 1/m$
2. For  $k = 1, \dots, K$ 
  - Select a sample of observations based on the weights.
  - Create the  $k$ -th weak learner and compute predictions  $x^{(k)}$
  - Compute the model **weighted** error and assign its coefficient:

$$\alpha^{(k)} = \lambda \times \log((1 - \text{error})/\text{error})$$

- Update sample weights:

$$w_i^{(k+1)} \propto w_i^{(k)} \times \exp(-\alpha^{(k)} y_i \hat{x}_i^{(k)})$$

3. Final weighted prediction

## Main Parameters

- n\_estimators: Number of estimators ( $K$ )
- base\_estimator: Weak estimator type
- learning\_rate: weights of estimator in final decision ( $\lambda$ )

# THANK YOU