# MLRF Lecture 06 J. Chazalon, LRDE/EPITA, 2021

# Some classifiers – part 2

Lecture 06 part 02

#### How to build non-linear classifiers?

2 solutions:

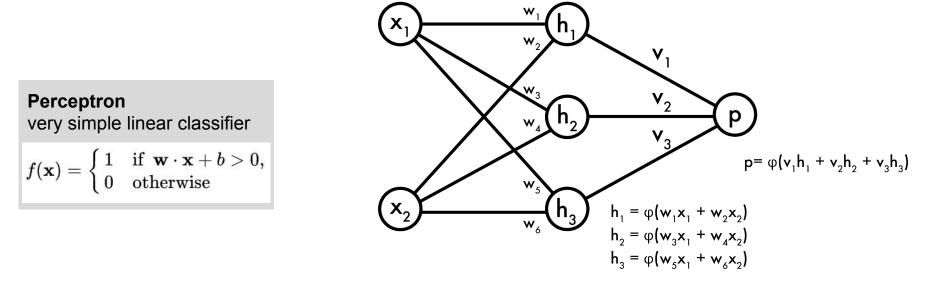
- Preprocess the data seen last time Ex.: explicit embedding, kernel trick...
  ⇒ change the input to make it linearly separable
- 2. **Combine** multiple **linear classifiers** into nonlinear classifier *current topic Ex.: boosting, neural networks...*

 $\Rightarrow$  split the input space into linear subspaces

# Non-linear classification using combinations of linear classifiers

#### **Multi-layer Perceptron**

Combine features linearly, apply a linear activation function  $\varphi$ , repeat.



#### Universal approximation theorem

What if φ not linear?

Universal approximation theorem (Cybenko 89, Hornik 91)

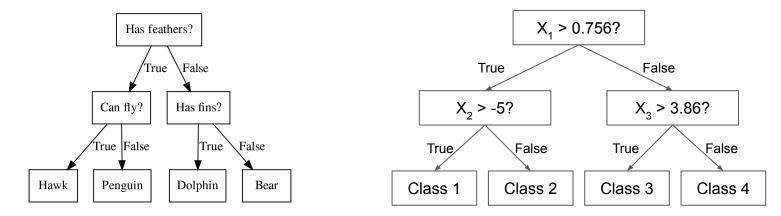
φ: any nonconstant, bounded, monotonically increasing function  $I_m$ : m-dimensional unit hypercube (interval [0-1] in m dimensions) Then 1-layer neural network with φ as activation can model any continuous function f:  $I_m → R$ (no bound on size of hidden layer)

By extension, works on f: bounded  $\mathsf{R}^m \to \mathsf{R}$ 

What can we learn? What can't we? UAT just says it's possible to model, not how.

#### **Decision trees**

Works on categorical (like "red", "black") and numerical (both discrete and continuous) random variables.



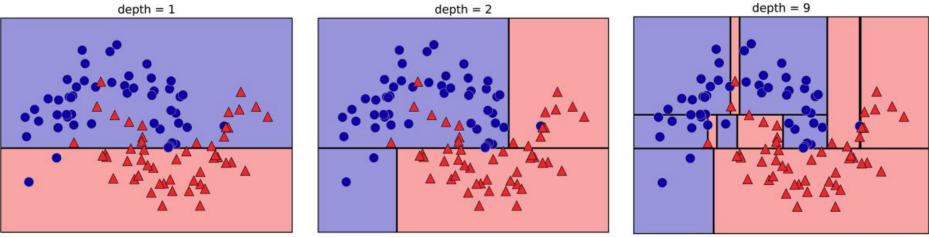
Train by optimizing classification "purity" at each decision (threshold on a particular dimension in numerical case).

#### **Decision trees**

Very fast training and testing. Non parametric.

No need to preprocess the features.

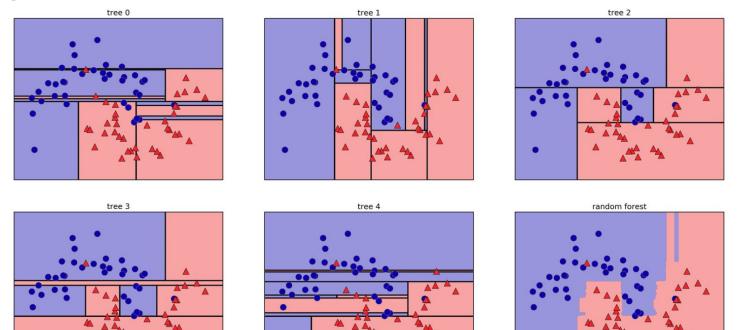
BUT: Very prone to **overfitting** without strong limits on depth.



#### **Random Forests**

[Breiman 2001]

**Average** the decision of multiple decision trees.



from A. Müller

#### **Random Forests**

Randomize in two ways:

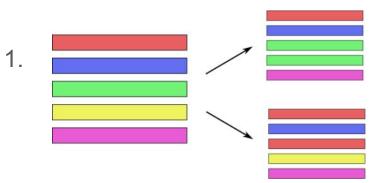
1. For each tree,

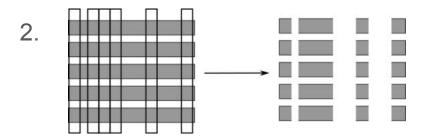
pick a **bootstrap** sample of data

2. For each split,

pick random sample of features

More trees are always better.





## **Ensemble methods**

### "Bagging" or "bootstrap aggregating"

[Breiman 96]

Underlying idea: part of the variance is due to the specific choice of the training data set.

- 1. Let us create many similar training data sets using **bootstrap**.
- 2. For each of them, train a new classifier.
- 3. The final function will be the **<u>average</u>** of each function outputs.

 $\Rightarrow$  If generalization error is decomposed into bias and variance terms then bagging reduces variance. (average of large number of random errors  $\approx$  0)

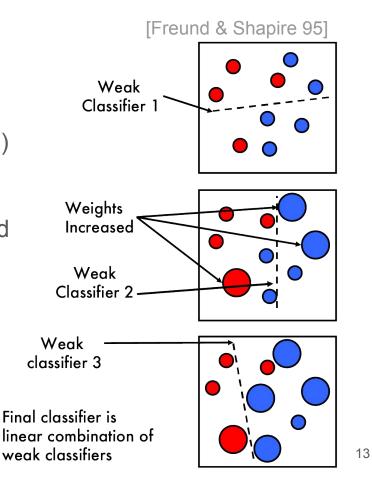
Random forest = a way of bagging trees.

#### "Boosting", AdaBoost variant

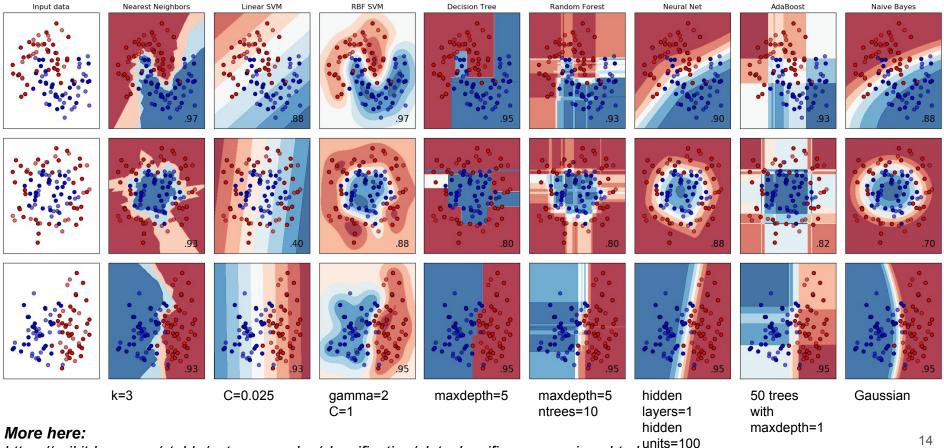
Combinaison of weak classifiers  $\sum_{m} \alpha_{m} G_{m}(x)$ 

 $\alpha_m$  increases with precision (less errors, bigger  $\alpha_m$ )

The classifier  $G_m$  is trained with **increased error cost** for the observations which were misclassified by  $G_{m-1}$ .



#### A quick comparison



https://scikit-learn.org/stable/auto\_examples/classification/plot\_classifier\_comparison.html "

# More tricks

#### Data augmentation

Add realistic deformations to your input in order to improve domain coverage.

For **image data**, depending on what is possible in production: rotations, horizontal & vertical flips, scaling, translation, illumination change, warping, noise, etc.



For **vector data**: interesting problem. Possible approach: train/fit PCA then add random noise in low-energy features.

#### Reject

Several options:

- Improve the model of class boundary In 1-vs-all training, add noise to the "others" samples.
- 2. Adjust the decision function depending on your application Look at the prediction probability of your classifier, and threshold it as per your need using a ROC curve.
- 3. Model the noise

Add a "none" class to your classifier, with samples for real life cases of negative samples.

