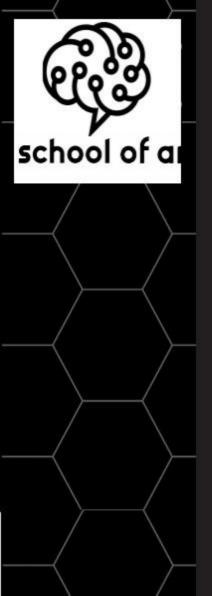


### **Ensemble Learning: About Ensemble Learning**

**AAA-Python Edition** 



### Plan

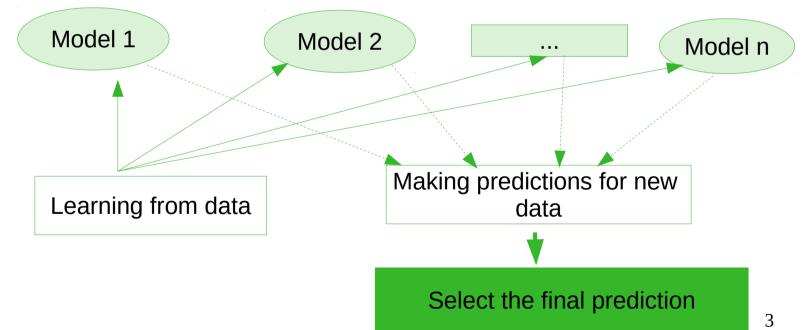
- 1- Ensemble Learning
- 2- Bagging & Pasting
- 3- Features sampling
- 4- Boosting : Adaptive Boosting
- 5- Boosting: Gradient boosting
- 6- Stacking or Blending



# Leal Ensembl

#### Concept

- In Ensemble Learning, we combine several models to build a better model. The algorithm used in Ensemble learning is called: an Ensemble Method.
- We can combine classifiers or regressors.
- The models can be all the same type, or different.





# -earn Ensembl

#### Voting

- Its about: how to select the final prediction.
- In Classification
  - Hard Voting: For each sample, a classifier will make a prediction: a class for that sample
    - Select the **most** predicted class by all the classifiers, for that sample .
  - Soft Voting: available when the classifiers can predict class probabilities.
    - Select the class with the **highest averaged** probability
- In Regression
  - The average of the predicted values.



# Leal Ensembl

#### **Ensemble Methods**

- The ensemble methods can vary by:
  - Varying or not the types of the models: use the same or different models.
  - Select the same sample only once or several times in the same model.
  - Whether or not the model use all the features or only a subset of features.
  - The models learn in parallel or sequentially.
  - The type of mechanism used to make a prediction.



# Learnin Ensemble

#### Example

```
1 from sklearn.tree import DecisionTreeClassifier
 2 from sklearn.ensemble import VotingClassifier
                                                            The will combine the models
 3 from sklearn.linear model import LogisticRegression
 4 from sklearn.svm import SVC
                                                         The 3 models used in
 6 logReg = LogisticRegression()
                                                         this ensemble method
  decTree = DecisionTreeClassifier()
 8 SVMClass = SVC()
 9 myEnsembleMethd = VotingClassifier(estimators=[('lr',logReg),('dt',decTree),('svc',SVMClass)]
                                    ,voting='hard')
  myEnsembleMethd.fit(x train, y train)
12 y pred= myEnsembleMethd.predict(x test.reshape(-1,4))
  myConfMat = confusion_matrix(y_test,y_pred)
14 print( myConfMat)
                                                            A string representing
Hard voting
                                                            the type of the model:
was used to
                                                            'dt' for : decision tree
determine
these
classes
```



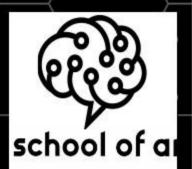
### Pastir 6 gin 9 -Ba

#### Definition

- Bagging (or Bootstrap aggregating) and pasting are both ensemble methods that combine same type of models. They both train the models on different random sub sets. All the models run in parallel. They use the voting mechanism for the final prediction
- Both Bagging and Basting apply random sampling:
  - The training set for each model is a subset of the original data randomly selected.
  - The same sample can be found in different models (different subsets).
- The difference is:

[By Amina Delali]

- Bagging: random sampling with replacement <==>
  - One sample can be found several times in the same model (same subset).
- Pasting: random sampling without replacement <==>
  - One sample can be found only once in the same model (same subset).



### Pastin Ŋ 9 6 -Ba 7

#### Bagging example using scikit-learn

```
# load the data
myIris = load_iris()
X = myIris.data
y = myIris.target
x_train,x_test,y_train,y_test= train_test_split(X, y, test_size=0.25)

from sklearn.ensemble import BaggingClassifier
from sklearn.svm import SVC
theModel2 = SVC()
myBagClass2 = BaggingClassifier(
theModel2, n_estimators=300,max_samples=90, myBagClass2.fit(x_train, y_train)
y_pred = myBagClass2.predict(x_test)
```

Bootstrap =True ==> Ensemble method : bagging
TheModel2 = SVC == > all the models are support vector machine classifiers
max\_samples=90 ==> the size of the subsets (bags) == 90 sample
n\_jobs=-1 ==> use all the available cores (to compute in parallel)
n\_estimators= 300 ==> use 300 SVC

```
myConfMat = confusion_matrix(y_test,y_pred)
print( myConfMat)
myBagClass.score(x_test,y_test)

[[11 0 0]
[0 8 3]
[0 0 16]]
```

The data



#### stin Pa Ŋ 6 ļ 6 6 -Ba 7

#### Pasting example using sckit-learn

```
from sklearn.svm import SVC
myBagClass2P = BaggingClassifier(
theModel2, n estimators=300, max samples=90, bootstrap=False, n jobs=-1)
myBagClass2P.fit(x_train, y_train)
v pred = myBagClass2P.predict(x test)
Bootstrap =False ==> Ensemble method : pasting
The remaining parameter are the same as the
previous ones
                                myConfMat2P = confusion matrix(y test,y pred)
                                print( myConfMat2P)
from sklearn.linear model import LogisticRegression
theModel = LogisticRegression()
 The model used is a LogisticRegression classifier
                                                                            1 12]]
```



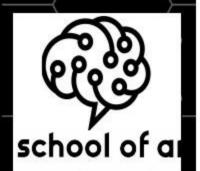
# Features sampling

M

#### Definition

- All the following methods use features sampling: each model will be trained in a random subset of features.
- Sampling features can be with or without replacement.
- Random patches method
  - Sampling both "training instances" and "features"

- Random subspaces method
  - Keeping all "training" instances but "sampling" features



# 3- Features sampling

[By Amina Delali]

#### Random Patches method

```
# import the tool for generating the data
from sklearn.datasets import make regression
# generate a random regression problem
xr, yr= make regression( n features = 100)
xr train,xr test,yr train,yr test= train test split(xr, yr, test size=0.25)
print(xr train.shape)
from sklearn.ensemble import BaggingRegressor
from sklearn.linear model import LinearRegression
theModelR = LinearRegression()
myBagReg = BaggingRegressor(
               theModelR, n estimators=1100, max samples=50, max features=80,
              bootstrap=True,
                              bootstrap features= True, n jobs=-1)
myBagReg.fit(xr train, yr train)
yr_pred = myBagReg.predict(xr test)
myBagReg.score(xr test,yr test)
                               Bagging
     0.5057404688880153
                        Number of selected samples = 50 <
                        75 ==> instances sampling
```

The data: 75 samples (instance), with 100 features

(75, 100)

Number of selected features = 80 < 100 features = > features sampling

Sampling features with replacement

1



samb

Features

3

#### Random Subspaces method

Pasting

**Features** 

sampling

replacement

without

Sampling features :

0.6483102156919015

80 < 100

Since all samples are used without replacement ==> the instances are **not sampled** ==> all the training data is used (75 samples)

Since max\_samples = 1.0 (a float value)==> max\_samples

training data (100% \*75 = 75)

=100% of the

[By Amina Delali]



#### Definition

- Boosting: Ensemble method, that combines several weak learners into a stronger learner.
- This is done by training the models sequentially ==> each model correct (boost) its predecessor.
- Uses the same models on the same data each time.
- The most known boosting methods are: Adaptive Boosting and Gradient Boosting.
  - Adaptive Boosting: each new predictor focus on the training samples that its predecessor underfitted (for example: misclassified in a classification problem) by modifying the instances weight.
  - Gradient Boosting: the new predictor tries to fit to the residual errors made by the previous predictor.



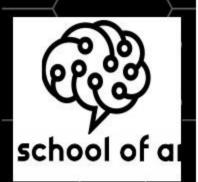
#### AdaBoost: training

- Weighting samples ==> each sample value will be multiplied by its weight.
- AdaBoost is Applicable in binary classification.
- The steps of the algorithm are as follow:
  - initialize the samples weight **w**<sup>i</sup> (for the first predictor ) by 1/m. m is the number of the training samples.
  - for each predictor j compute:
    - $\rightarrow$  the weighted error rate :  $\mathbf{r_i} = \sum \mathbf{W^i}$  (whre the prediction is wrong) /  $\sum \mathbf{W^i}$
    - $\rightarrow$  compute the j predictor's weight:  $\alpha_j = \eta \log(1 r_{j}) / r_{j}$ .  $\eta$  is the learning rate parameter.
    - Compute the new weights (to be used by the following new predictor j+1): wi=wi,if,y<sub>true</sub>(i)=y<sub>pred</sub>(i)

$$w^i = w^i exp(\alpha_i), if, y_{true}(i) \neq y_{pred}(i)$$

- → Normalize the new weights w' by: w'=1/∑w'
- → The process is repeated until the perfect predictor is found, or the maximum number of predictors is reached.

[By Amina Delali]



#### AdaBoost: Predicting

- To make a prediction:
  - make a prediction with each predictor j from the resulting N predictors.
  - attribute a weight to each prediction by the predictor's  $\boldsymbol{j}$  weight  $\boldsymbol{\alpha}_i$
  - $\rightarrow$  for each sample  $\mathbf{x}$  select the class  $\mathbf{k}$  that receives the majority of weighted votes: for each predicted class  $\mathbf{k}$  sum up the corresponding  $\mathbf{\alpha}_{\mathbf{j}}$  weights, then select the class  $\mathbf{k}$  with the biggest sum.

#### AdaBoost: SAMME

- SAMME: Stagewise Additive Modeling using a Multi-class Exponential loss function
- Enhanced version of AdaBoost, applicable in multiclass classification.
- Same steps as AdaBoost, just the  $\alpha$  weight is computed differently:  $\alpha_i = \eta * (log[(1-rj)/rj] + log(K-1))$ . K is the number of classes.



#### Example

```
from sklearn.ensemble import AdaBoostClassifier

# we will use our a decision tree classifier

from sklearn.tree import DecisionTreeClassifier

theModel4 = DecisionTreeClassifier(max_depth=1)

mySAMME = AdaBoostClassifier(theModel4, n_estimators=200,

algorithm="SAMME", learning_rate=1

}
```

```
mySAMME.fit(x_train, y_train)
y_pred = mySAMME.predict(x_test)

myConfMat3 = confusion_matrix(y_test,y_pred)
print(myConfMat3)
mySAMME.score(x_test,y_test)

[[12 0 0]
        [ 0 11 2]
        [ 0 1 12]]

0.9210526315789473
```

A weak classifier: a decision tree with 1 level: the 2 leafs, the split of the root node. This Tree is called a: Decision Stump

In sckit-learn to apply the adaptive boosting to regression, the weights are adjusted according to the error of the predictions



5- Boosting: Gradient boosting

#### The concept

- 1. The predictor will be first trained on a set of data: x,y
- 2. The residual errors are computed from its prediction:  $r = y y_{pred}$
- 3.A new predictor will be trained with the new set of data: x,r
- 4. The residual errors are computed again as follow:  $r^2 = r r_{pred}$
- 5. The steps 3 and 4 are repeated until: you predict using all the predefined number of predictors, or you determine the optimal consecutive predictors (the least generated error) and you select those predictors as your final model. Or, you continue adding predictors until the errors will not diminish
- 6. The final prediction will be the sum of all the predictions.
- Scikit learn implements gradient tree boosting: the models used are decision trees.



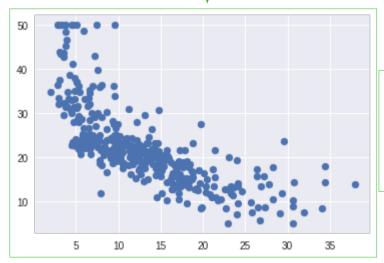
### 5- Boosting: Gradient boosting

#### Example: the data

```
# we will use load_boston utilities to load our data
# we will considere only the feature with the indices 12 (the last one): 'LSTAT'

from sklearn.datasets import load_boston
mydata = load_boston()
xr = mydata.data[:,12]
yr = mydata.target
xr_train,xr_test,yr_train,yr_test= train_test_split(xr, yr, test_size=0.25)
for x in [xr_train, xr_test, yr_train, yr_test]:
    x=x.reshape(-1,1)

plt.scatter(xr_train,yr_train)
```



- Boston House Prices dataset.
- The chosen feature represents the: % lower status of the population

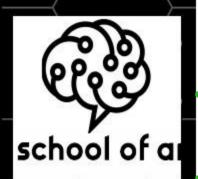


### 5- Boosting: Gradient boosting

#### Example: using a GBRT

GBRT for Gradient Boosted Regression Trees

```
1 from sklearn.ensemble import GradientBoostingRegressor
2 from sklearn.model_selection import train_test_split
          myGBRT = GradientBoostingRegressor(max depth=2, n estimators=45)
          myGBRT.fit(xr2 train, yr2 train)
        7 # xnew to draw the model
        8 xnew= np.linspace(4,40,2000).reshape(-1,1)
        9 ypred= myGBRT.predict(xnew)
        | | plt.plot(xnew, ypred,color="red")
          plt.scatter(xr2 train,yr2 train,marker=".")
30
                                    27.701804841167945
20
                  # testing the model
                  ypred= myGBRT.predict(xr2 test)
                  #plot the true values: in green and the predicted values: in black
                  plt.scatter(xr2_test,yr2_test,color="green",marker=".")
                   plt.scatter(xr2 test, vpred, color="black", marker="+")
                                                                                                            19
                  from sklearn.metrics import mean squared error as MSE
[By Amina Delali] | print (MSE(yr2 test,ypred))
```



### blending or 6 -Stackin

#### Concept

The idea here, is to **train** a model to learn how to **aggregate** the ensemble models **predictions**.

The method is composed of:

- Learner models : that will fit to the data, and make the predictions.
- Blender: the final model or meta learner, that will make the final prediction.

There are different methods to train the blender:

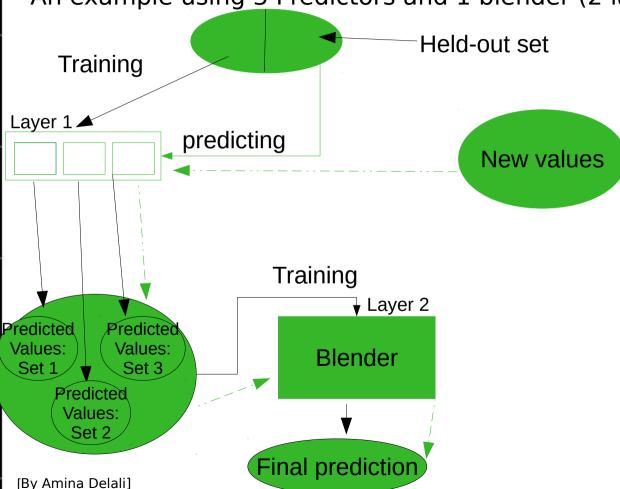
- Hold-out set: Blending
- Out-of-fold: Stacking



blending or 9 **6-Stackin** 

#### Hold-out set: principle

An example using 3 Predictors and 1 blender (2 layers)



- To make a prediction, the new instance will go through the first layer.
- The resulting predictions will serve as input for the second layer.
- The prediction made by this later one is the final result.



# -Stacking or blending

Ö

#### Hold-out set: Training

```
from sklearn.svm import SVR
from sklearn.tree import DecisionTreeRegressor

# predictors composing the first layer
myModel1 = SVR(kernel='poly')
myModel2 = LinearRegression()
myModel3 = DecisionTreeRegressor()

# the model choosen for the blender
myBlender = SVR(kernel='linear')

# subdividing the= training data into 2 subsets
size= int(xr2_train.shape[0]/2)

subXR1=xr2_train[:size]
subYR1= yr2_train[size:]

subYR1= yr2_train[size:]
subYR2= yr2 train[size:]
```

print(subXR1.shape,subXR2.shape,subYR1.shape,subYR2.shape)

Subset xr1 for training Subset xr2 for predicting

```
22
23 # training the first layer
24 myModell.fit(subXR1,subYR1)
```

myModel2.fit(subXR1,subYR1) myModel3.fit(subXR1,subYR1) # predicting with the second subset
pred1 = myModel1.predict(subXR2)

pred1 = myModel1.predict(subXR2)
pred2 = myModel2.predict(subXR2)
pred3 = myModel3.predict(subXR2)

Predicting with the first level

#### Training the first level

### Generate the new features

[By Amina Delali]

# constructing the new x values
xnew = np.append(pred1.reshape(-1,1), pred2.reshape(-1,1), axis= 1)
xnew= np.append(xnew, pred3.reshape(-1,1),axis = 1)
print(xnew.shape)

# train on the previous predicted values

myBlender.fit(xnew,subYR2)

Training the blender



### blending Or 5 -Stackin Ö

#### Hold-out set: Testing

```
The test data will be used
# predicting on test data
                                                   by the first level to predict
# predicting with the first layer
                                                  the values == features used
pred1 = myModel1.predict(xr2 test)
pred2 = myModel2.predict(xr2 test)
                                                      by later by the blender
pred3 = myModel3.predict(xr2 test)
# create the x vlaues
# constructing the new x values
xnew_test = np.append(pred1.reshape(-1,1), pred2.reshape(-1,1), axis= 1)
xnew test= np.append(xnew test, pred3.reshape(-1,1),axis = 1)
print(xnew test.shape)
# predicting with the second layer
myFP= myBlender.predict(xnew test)
                                                The final prediciton, will be
print(myBlender.score(xnew test,yr2 test))
                                                done by the blender.
# printing the mean square error
print (MSE(yr2 test,myFP))
                                                 0.6411442145726178
```

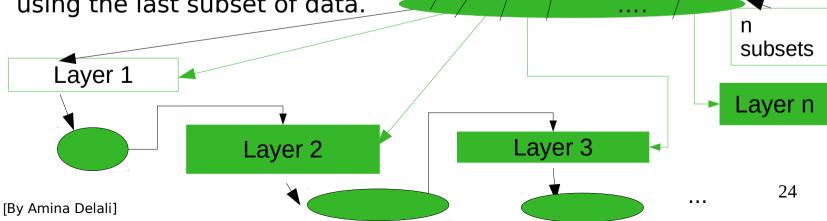
27.90487595475314

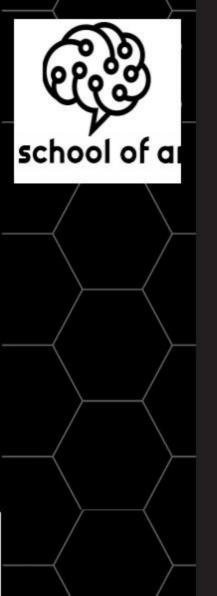


# blending -Stackin

#### Hold-out set: generalization

- Its possible to train several type of blenders. And , each one can be a set of models.
- The idea is to divide the original training set into several subsets: n subsets ==> n layers (n-1 blending phase)
- The first set of predictors will train from the first subset, and make prediction with the second one.
- The second set of predictors will train from the previous predictions.
   And then make new predictions using the third subset.
- The process is repeated until the last subset of predictors: it will train from the last predictions made by the previous predictors using the last subset of data.





### References

- Aurélien Géron. Hands-on machine learning with Scikit-Learn and Tensor-Flow: concepts, tools, and techniques to build intelligent systems. O'Reilly Media, Inc, 2017.
- Scikit-learn.org. scikit-learn, machine learning in python. On-line at https://scikit-learn.org/stable/. Accessed on 03-11-2018.



# Thank you!

FOR ALL YOUR TIME