

# h2oEnsemble: Scalable Ensemble Learning in R

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July 2, 2015

# Introduction



# Overview

- Ensemble Learning
- Model Stacking (aka. Super Learning)
- H2O Machine Learning via h2o R package
- h2oEnsemble R package

# Ensemble Learning



In statistics and machine learning, **ensemble methods** use multiple models to obtain better predictive performance than could be obtained from any of the constituent models.

– *Wikipedia, 2015*

- Ensemble of weak learners (e.g. Random Forest)
- Generalized Model Stacking (combine the predictions from multiple models)

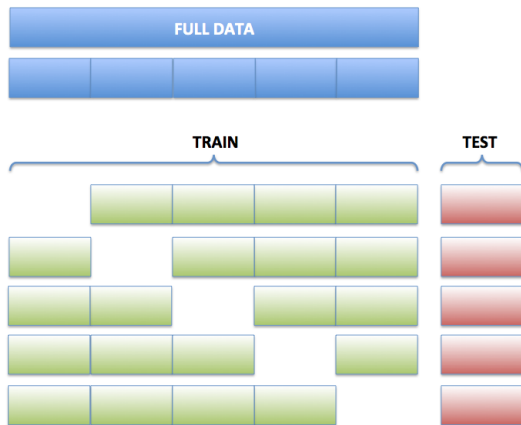
# Super Learner algorithm

The **Super Learner algorithm** is a loss-based supervised learning method that finds the optimal combination of a collection of prediction algorithms.



Super Learner performs asymptotically as well as best possible weighted combination of the base learners.

# K-fold Cross-validation



**Example:** 5-fold cross validation

## Super Learner: The setup

- 1 Define a base learner library of  $L$  learners,  $\Psi^1, \dots, \Psi^L$ .
- 2 Specify a metalearning method,  $\Phi$ .
- 3 Partition the training observations into  $V$  folds.

# Super Learner algorithm

## Super Learner: The algorithm

- 1 Generate a matrix  $Z$ , of dimension  $n \times L$ , of cross-validated predictions as follows: During cross-validation, we obtain fits,  $\hat{\Psi}'_{-v}$ , defined as fitting  $\Psi'$  on the observations that are not in fold  $v$ . Predictions are then generated for the observations in the  $v^{\text{th}}$  fold.
- 2 Find the optimal combination of subset-specific fits according to a user-specified metalearner algorithm,  $\hat{\Phi}$ , with a new design matrix,  $Z$ .
- 3 Fit  $L$  models (one for each base learner) on the original training set,  $X$ , and save the  $L$  individual model fit objects along with  $\hat{\Phi}$ . This ensemble model can be used to generate predictions on new data.



# Super Learning for Big Data

Practical solutions to this problem:

- 1 Develop alternative formulations of Super Learner that learn on subsets of data to overcome memory limitations.
- 2 Use candidate learners that can learn iteratively and thus do not require loading the entire training set into memory at once. (i.e., online learning)
- 3 Make use of distributed algorithms.
- 4 Rather than native R or Python, use a more “scalable” language (C++, Java, Scala, Fortran, Julia).

# H2O Machine Learning platform

H2O is an open source, distributed, Java machine learning library.



APIs available in:  
R, Python, Java, Scala and REST/JSON

# H2O Machine Learning platform

Distributed Supervised ML Algorithms available in H2O

- Generalized Linear Model with Elastic Net regularization
- Gradient Boosting Machines (w/ trees)
- Random Forest
- Deep Learning: Multi-Layer Feed-Forward Neural Networks



## h2o: How to start H2O & load data

### Example

```
library(h2o) # First install from CRAN
localH2O <- h2o.init() # Initialize the H2O cluster

# Data directly into H2O cluster (avoids R)
train <- h2o.importFile(path = "train.csv")

# Data into H2O from R data.frame
train <- as.h2o(my_df)
```

## h2o: How to train & test

### Example

```
y <- "Class"  
x <- setdiff(names(train), y)  
  
fit <- h2o.gbm(x = x, y = y, training_frame = train)  
pred <- h2o.predict(fit = fit, validation_frame = test)
```

# h2oEnsemble R package

h2oEnsemble: Set up the ensemble

## Example

```
learner <- c("h2o.randomForest.1",  
            "h2o.deeplearning.1",  
            "h2o.deeplearning.2")
```

```
metalearner <- "h2o.glm.wrapper"
```

```
family <- "binomial"
```

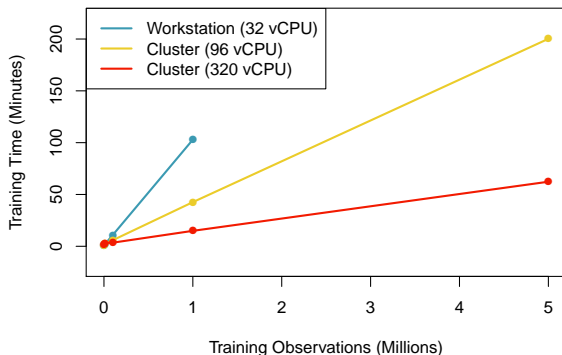
## h2oEnsemble: How to train & test

### Example

```
fit <- h2o.ensemble(x = x, y = y, training_frame = train,  
                  family = family,  
                  learner = learner,  
                  metalearner = metalearner)  
  
pred <- h2o.predict(fit = fit, validation_frame = test)
```

# H2O Ensemble: Performance

**Runtime Performance of H2O Ensemble**



R color palette: <https://github.com/karthik/wesanderson>



Thank you!

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