## Ch 8.2.1, 8.2.2: Bagging and Random Forests

Lecture 24 - CMSE 381

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Mon. Nov 14, 2022

#### Announcements

#### Last time:

• 8.1 Decision Trees

#### This lecture:

- 8.2.1 Bagging
- 8.2.2 Random forest.

#### **Announcements:**

• Some rearranging:

			carranging.		
20	F	Nov 4	Polynomial & Step Functions.	7.1,7.2	
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	W	Dec 7	Review		
	F	Dec 9	Midterm #3	non-interr	neat sheet and a let-connected culator

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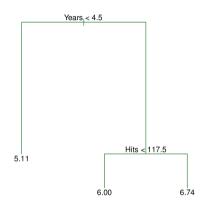
## Section 1

Last time

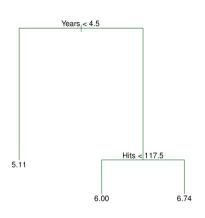
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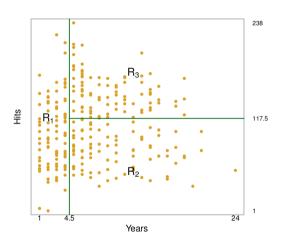
## First decision tree example

	Hits	Years	LogSalary
1	81	14	6.163315
2	130	3	6.173786
3	141	11	6.214608
4	87	2	4.516339
5	169	11	6.620073
317	127	5	6.551080
318	136	12	6.774224
319	126	6	5.953243
320	144	8	6.866933
321	170	11	6.907755



# Viewing Regions Defined by Tree



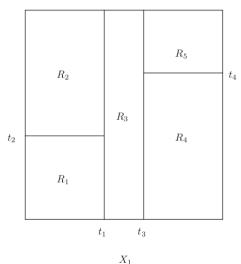


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# How do we actually get the tree? Two steps

- We divide the predictor space that is, the set of possible values for  $X_1, X_2, \cdots, X_p$  — into J distinct and non-overlapping regions,  $R_1, R_2, \cdots, R_L$
- For every observation that falls into the region  $R_i$ , we make the same prediction = the mean of the response values for the training observations in  $R_i$ .



# Recursive binary splitting

#### Goal:

Find boxes  $R_1, \dots, R_J$  that minimize

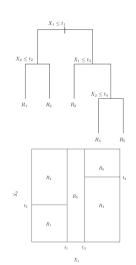
$$\sum_{j=1}^J \sum_{i \in R_j} (y_i - \hat{y}_{R_j})^2$$

 $\hat{y}_{R_j}$  = mean response for training observations in *j*th box

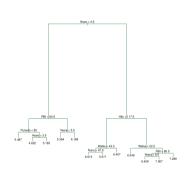
Pick s so that splitting into  $\{X \mid X_j < s\}$  and  $\{X \mid X_j \geq s\}$  results in largest possible reduction in RSS:

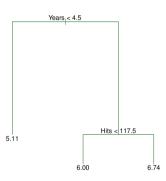
$$R_1(j,s) = \{X \mid X_j < s\}$$
  
 $R_2(j,s) = \{X \mid X_j \ge s\}$ 

$$\sum_{i|x_i \in R_1(j,s)} (y_i - \hat{y}_{R_1})^2 + \sum_{i|x_i \in R_2(j,s)} (y_i - \hat{y}_{R_2})^2$$



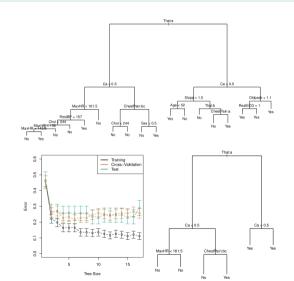
# Pruning





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#### Classification version

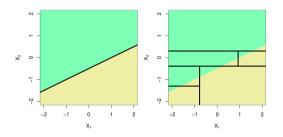


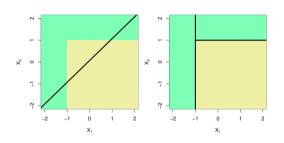
#### **Evaluating the splits:**

- $\hat{p}_{mk}$  = proportion of training observations in  $R_m$  from the kth class
- Error:  $E = 1 \max_k(\hat{p}_{mk})$
- Gini index:

$$G = \sum_{k=1}^K \hat{
ho}_{mk} (1-\hat{
ho}_{mk})$$

## Linear models vs trees





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# $\mathsf{Pros}/\mathsf{Cons}$

Pros: Cons:

## Section 2

8.2.1 Bagging

## Recall: The bootstrap

## Want to do (but can't):

Build separate models from independent training sets, and average resulting predictions:

- $\hat{f}^1(x), \dots, \hat{f}^B(x)$  for B separate training sets
- Return the average

$$\hat{f}_{avg}(x) = \frac{1}{B} \sum_{b=1}^{B} \hat{f}^b(x)$$

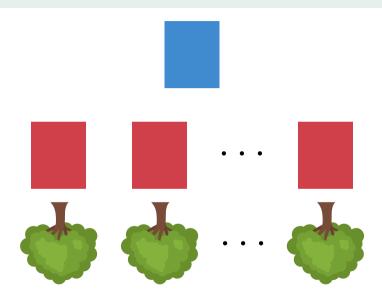
#### **Boostrap modification:**

- Work with fixed data set
- Take B samples from this data set (with replacement)
- Train method on *b*th sample to get  $\hat{f}^{*b}(x)$
- Return average of predictions (regression)

$$\hat{f}_{bag}(x) = \frac{1}{B} \sum_{b=1}^{B} \hat{f}^{*b}(x)$$

or majority vote (classification)

## Tree version



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# Prediction on new data point



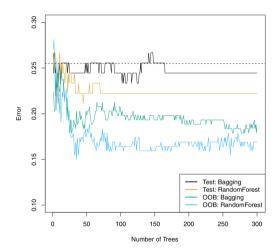
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## Bagging vs Bootstrap

**Bootstrap** 

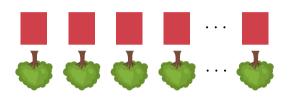
Bagging

# Example: Heart classification data



## Out of Bag Error Estimation

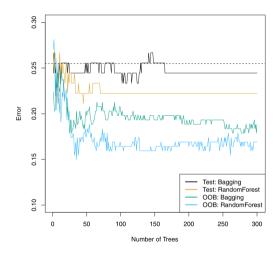
- On average, bootstrap sample uses about 2/3 of the data
- Remaining observations not used are called out-of-bag (OOB) observations
- For each observation, run through all the trees where it wasn't used for building
- Return the average (or majority vote) of those as test prediction



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# Error using OOB



# Bagging code example

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## Section 3

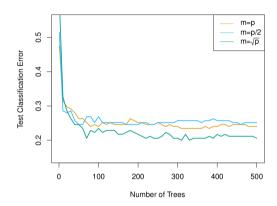
## Random Forests

#### The idea

- Goal is to decorrelate the bagged trees:
  - If there is a strong predictor, the first split of most trees will be the same
  - Most or all trees will be highly correlated
  - Averaging highly correlated quantities doesn't decrease variance as much as uncorrelated

- The random forrest fix:
  - Each time a split is considered, only use a random subset of m the predictors
  - Fresh sample taken every time
  - ▶ Typically  $m \approx \sqrt{p}$
  - ▶ On average, (p m)/p of splits won't consider strong predictor
  - ightharpoonup m = p gives back bagging

## Example on gene expression



Coding example for random forests

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## TL:DR

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