

# Lecture 22:

## Ensembles: Bagging, Random Forests, & Intro to Boosting

Big Data and Machine Learning for Applied Economics  
Econ 4676

Ignacio Sarmiento-Barbieri

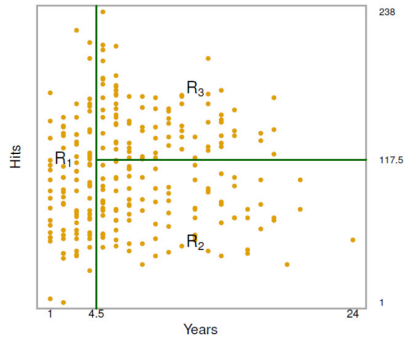
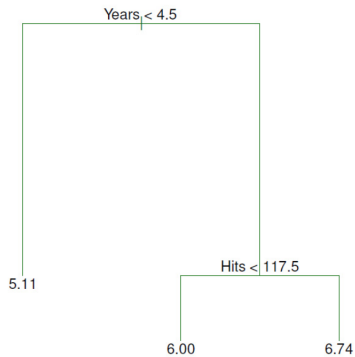
Universidad de los Andes

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# Agenda

- 1 Recap
- 2 Bagging and Random Forests
- 3 Comparisons: Lasso, CART, Random Forests
- 4 Boosting
  - AdaBoost
  - Causal Forests
- 5 Review & Next Steps
- 6 Further Readings

# CARTs



# CARTs

- Problem then boils down to searching the partition variable  $X_j$  and the partition point  $s$  such that

$$\min_{j,s} \left[ \min_{c_1} \sum_{x_i \in R_1(j,s)} (y - c_1)^2 + \min_{c_2} \sum_{x_i \in R_2(j,s)} (y - c_2)^2 \right] \quad (1)$$

- For each partition variable, and partition point, the internal minimization is the mean of each region

$$\hat{c}_m = \frac{1}{n_m} \sum (y_i | x_i \in R_m) \quad (2)$$

- Process is repeated inside each region.

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- Process is repeated inside each region.
- If the final tree has  $M$  regions then the prediction is

$$\hat{f}(x) = \sum_{m=1}^M \hat{c}_m I(x \in R_m) \quad (3)$$

- ▶ Cost complexity of tree  $T$

$$C_\alpha(T) = \sum_{m=1}^{[T]} n_m Q_m(T) + \alpha [T] \quad (4)$$

- ▶ where  $Q_m(T) = \frac{1}{n_m} \sum_{x_i \in R_m} (y_i - \hat{c}_m)^2$  for regression trees
- ▶  $Q_m(T)$  penalizes heterogeneity (impurity) within each region, and the second term the number of regions.
- ▶ Objective: for a given  $\alpha$ , find the optimal pruning that minimizes  $C_\alpha(T)$

# CARTs

- ▶ Smart way to represent nonlinearities. Most relevant variables on top.
- ▶ Very easy to communicate.
- ▶ Reproduces human decision-making process.
- ▶ Trees are intuitive and do OK, but
  - ▶ They are not very good at prediction
  - ▶ If the structure is linear, CART does not work well.
  - ▶ Not very robust

# Bagging

- ▶ We can improve performance a lot using either bootstrap aggregation (bagging), random forests, or boosting.
- ▶ Bagging & Random Forests:
  - ▶ Repeatedly draw bootstrap samples  $(X_i^b, Y_i^b)_{i=1}^N$  from the observed sample.



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  - ▶ Repeatedly draw bootstrap samples  $(X_i^b, Y_i^b)_{i=1}^N$  from the observed sample.
  - ▶ For each bootstrap sample, fit a regression tree  $\hat{f}^b(x)$
  - ▶ Average across bootstrap samples to get the predictor

$$\hat{f}_{bag} = \frac{1}{B} \sum_{b=1}^B \hat{f}^b(x) \quad (5)$$

- ▶ Basically we are smoothing predictions.
- ▶ Idea: the variance of the average is less than that of a single prediction.

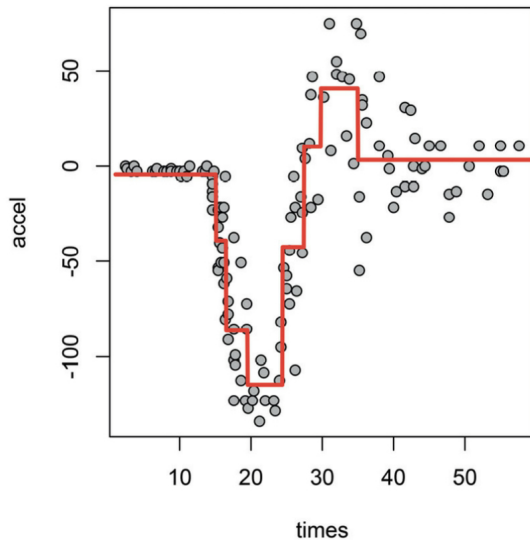
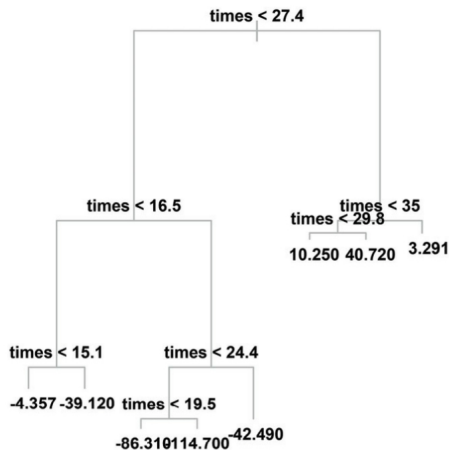
# Bagging



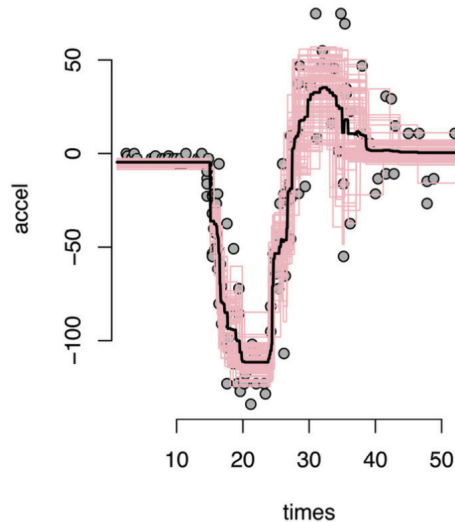
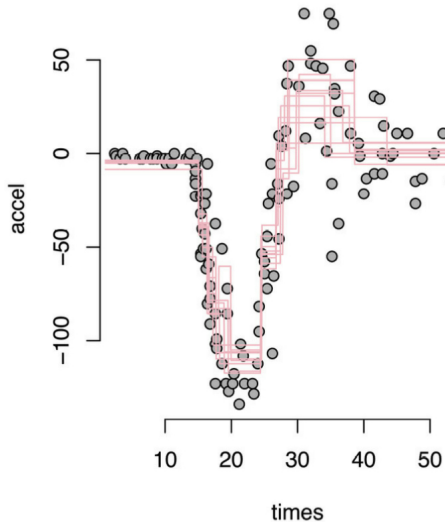
# Random Forests

- ▶ Problem with bagging: if there is a strong predictor, different trees are very similar to each other. High correlation. Is the variance really reduced?
- ▶ Forests: lower the correlation between the trees in the bootstrap.
- ▶ If there are  $p$  predictors, in each partition use only  $m < p$  predictors, chosen randomly
- ▶ Bagging is forest with  $m = p$  (use all predictors in each partition).
- ▶ Typically  $m = \sqrt{p}$

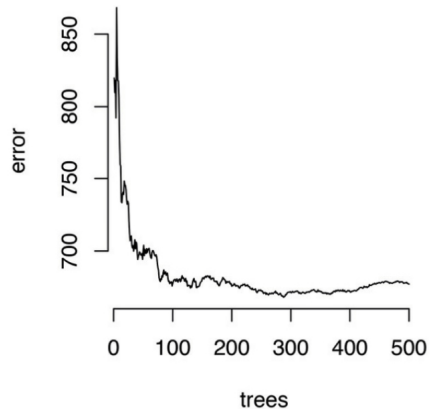
# Random Forests



# Random Forests



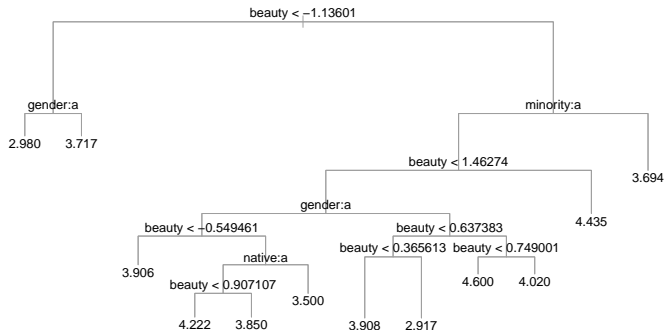
# Random Forests



# Random Forests and Trees

## Trees

```
pstcut <- prune.tree(pstree, best=12)  
plot(pstcut, col=8)  
text(pstcut)
```





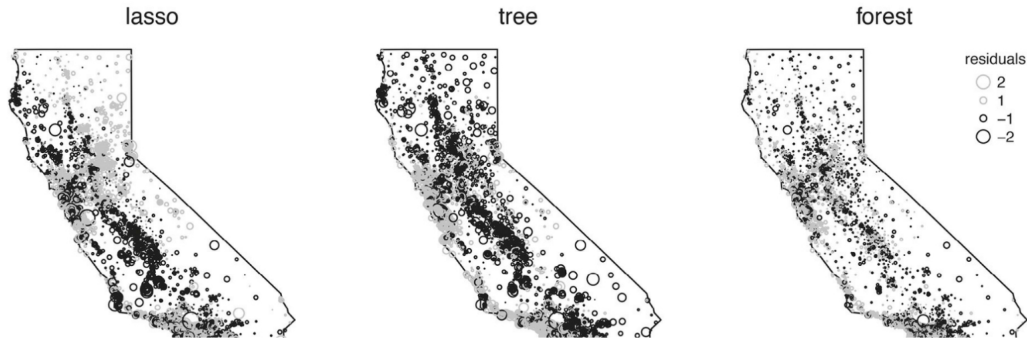
# Random Forests and Trees

```
require("ranger")

rf_tree <- ranger(eval ~beauty + gender + minority + native +
  tenure + division, data=tr,
  write.forest=TRUE,num.tree=200,min.node.size=25,
  importance="impurity")
sort(rf_tree$variable.importance, decreasing = TRUE)
```

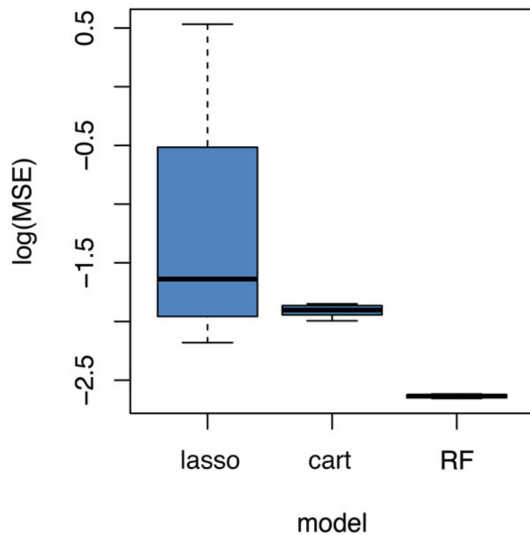
```
##      beauty  minority    gender    native    tenure  division
## 22.881176   3.089366   2.608295   2.104095   2.062075   1.627261
```

# In sample residuals



```
# Model Matrix for Lasso  
XXca <- model.matrix(logMedVal~.*longitude*latitude,  
  data=data.frame(scale(CAhousing)))[, -1]
```

# Out of sample MSE



# Boosting

- ▶ Problem with CART: high variance. Instability
- ▶ Weak classifier: marginally better classifier than flipping a coin (error rate slightly better than .5)
- ▶ E.g.: CART with few branches ('stump', two branches)
- ▶ Boosting: weighted average of a succession of weak classifiers.
- ▶ Vocab
  - ▶  $y \in -1, 1$  (for simplicity),  $X$  vector of predictors.
  - ▶  $y = G(X)$  (classifier)
  - ▶  $err = \frac{1}{N} \sum_i^N I(y_i \neq G(x_i))$

# AdaBoost

- 1 Start with weights  $w_i = 1/N$
- 2 For  $m = 1$  through  $M$ :
  - 1 Adjust  $G_m(x)$  using weights  $w_i$ .
  - 2 Compute prediction

$$err_m = \frac{\sum_{i=1}^N I(y_i \neq G_m(x_i))}{\sum_{i=1}^N w_i} \quad (6)$$

- 3 Compute  $\alpha_m = \ln \left[ \frac{(1-err_m)}{err_m} \right]$
- 4 Update weights:  $w_i \leftarrow w_i c_i$

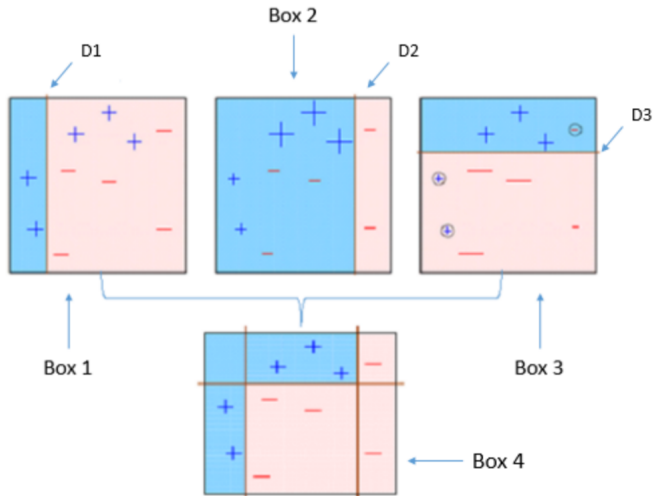
$$c_i = \exp [\alpha_m I(y_i \neq G_m(x_i))] \quad (7)$$

- 3 Output:  $G(x) = \text{sgn}[\sum_{m=1}^M \alpha_m G_m(x)]$

# AdaBoost

- ▶  $c_i = \exp[\alpha_m I(y_i \neq G_m(x_i))]$
- ▶ If it was correctly predicted,  $c_i = 1$ . No issue.
- ▶ Otherwise,  $c_i = \exp(\alpha_m) = \frac{(1 - \text{err}_m)}{\text{err}_m} > 1$
- ▶ At each step the method gives more relative importance to the predictions that were wrong.
- ▶ Final step: weighted average of predictions at each step.

# AdaBoost



Source: <https://www.analyticsvidhya.com/blog/2015/11/quick-introduction-boosting-algorithms-machine-learning/>

*American Economic Review: Papers & Proceedings* 2017, 107(5): 546–550  
<https://doi.org/10.1257/aer.p20171000>

## *LABOR MARKETS AND CRIME<sup>‡</sup>*

### Using Causal Forests to Predict Treatment Heterogeneity: An Application to Summer Jobs<sup>†</sup>

By JONATHAN M.V. DAVIS AND SARA B. HELLER\*



# Idle hands are the devil's workshop

- ▶ The application uses two large scale RCTs of Chicago's One Summer Plus (OSP) program conducted in 2012 and 2013. OSP provides disadvantaged youth ages 14 to 22 with 25 hours a week of employment, an adult mentor, and some other programming.
- ▶ Participants are paid Chicago's minimum wage (\$8.25 at the time).
- ▶ Find a 43 percent reduction in violent crime arrests in the 16 months after random assignment.

# Causal Tree: Theory Details

- ▶ Work well in RCTs
- ▶ Issue: we do not observe the ground truth
- ▶ Honest estimation (Innovation):
  - ▶ One sample to choose partition
  - ▶ One sample to estimate leaf effects
- ▶ Why is the split critical?
- ▶ Fitting both on the training sample risks overfitting: Estimating many “heterogeneous effects” that are really just noise idiosyncratic to the sample.
- ▶ We want to search for true heterogeneity, not noise

# Heterogeneous Treatment Effects Assumptions

- ▶ There are a couple of assumptions that are key
- ▶ Assumption 1: Unconfoundedness

$$Y_i(1), Y_i(0) \perp W_i \mid X_i \quad (8)$$

- ▶ The *unconfoundedness* assumption states that, once we condition on observable characteristics, the treatment assignment is independent to how each person would respond to the treatment.
- ▶ i.e., the rule that determines whether or not a person is treated is determined completely by their observable characteristics.
- ▶ This allows, for example, for experiments where people from different genders get treated with different probabilities,
- ▶ **rules out** experiments where people self-select into treatment due to some characteristic that is not observed in our data.

# Heterogeneous Treatment Effects

## ► Assumption 2: Overlap

$$\forall x \in \text{supp}(X), \quad 0 < P(W = 1 \mid X = x) < 1 \quad (9)$$

- The *overlap* assumption states that at every point of the covariate space we can always find treated and control individuals.
- i.e., in order to estimate the treatment effect for a person with particular characteristics  $X_i = x$ , we need to ensure that we are able to observe treated and untreated people with those same characteristics so that we can compare their outcomes.

# The Honest Target: Athey and Imbens Innovation

- ▶ The ultimate goal is to construct and assess an algorithm  $\pi(\cdot)$  that maximizes the honest criterion

$$\max Q^H(\pi) = -E_{S^{te}, S^{est}, S^{tr}} [MSE_{\mu}(S^{te}, S^{est}, S^{tr}, \pi(S^{tr}))] \quad (10)$$

- ▶ In CART the target is different (adaptive target)

$$\max Q^C(\pi) = -E_{S^{te}, S^{tr}} [MSE_{\mu}(S^{te}, S^{tr}, \pi(S^{tr}))] \quad (11)$$

# From Causal Trees to Causal Forest

The implementation steps are as follows in Davis and Heller (2017):

- ▶ (1) Draw a subsample  $b$  without replacement containing  $n_b = 0.2N$  observations from the  $N$  observations in the dataset
- ▶ (2) Randomly split the  $n_b$  observations in half to form a training sample ( $tr$ ) and an estimation sample ( $e$ ) such that  $n_{tr} = n_e = \frac{n_b}{2}$ . Using just the training sample, start with a single leaf containing all  $n_{tr}$  observations.

# From Causal Trees to Causal Forest

The implementation steps are as follows in Davis and Heller (2017):

- ▶ (3) For each value of each covariate,  $X_j = x$ , form candidate splits of the observations into two groups based on whether  $X_j \leq x$ . Consider only splits where there are at least ten treatment and ten control observations in both new leaves.
- ▶ Choose the single split that maximizes an objective function  $O$  capturing how much the treatment effect estimates vary across the two resulting subgroups, with a penalty for within leaf variance. If this split increases  $O$  relative to no split, implement it and repeat this step in both new leaves. If no split increases  $O$ , this is a terminal leaf.

$$O = (n_T + n_C) \hat{\tau}_l^2 - 2 \left( \frac{\hat{Var}(Y_{Tl})}{n_T} + \frac{\hat{Var}(Y_{Cl})}{n_C} \right) \quad (12)$$

# From Causal Trees to Causal Forest

- ▶ (4) Once no more splits can be made in step 3, the tree is defined for subsample  $b$ . Move to the estimation sample, and group the new observations into the same tree based on their  $X$ s.
- ▶ (5) Using just the estimation sample, calculate  $\hat{\tau} = \bar{y}_{Tl} - \bar{y}_{Tc}$  within each terminal leaf. This step makes the tree honest, since treatment effect estimates are made using different observations than the ones that determined the splits.
- ▶ (6) Return to the full sample of  $N$  observations. Assign  $\hat{\tau}_{l,b} = \hat{\tau}_l$  to each observation whose  $X$ s would place it in leaf  $l$ , and save this prediction.
- ▶ (7) Repeat steps (i) to (vi)  $B = 25,000$  times



# From Causal Trees to Causal Forest

- Define observation  $i$ 's predicted CATE as  $\hat{\tau}_i^{CF}(x) = \frac{1}{B} \sum \hat{\tau}_{l,b}$

# From Causal Trees to Causal Forest

- ▶ Define observation  $i$ 's predicted CATE as  $\hat{\tau}_i^{CF}(x) = \frac{1}{B} \sum \hat{\tau}_{l,b}$
- ▶ The procedure requires the researcher to select three parameters: the number of trees, the minimum number of treatment and control observations in each leaf, and the subsample size.
- ▶ In the absence of formal criteria to guide our choices, we used a large number of trees (more trees reduce the Monte Carlo error introduced by subsampling; we found moving from 10,000 to 25,000 improved the stability of estimates across samples).
  - ▶ Increasing the minimum number of observations in each leaf trades off bias and variance; bigger leaves make results more consistent across different samples but predict less heterogeneity.
  - ▶ Smaller subsamples reduce dependence across trees but increase the variance of each estimate (larger subsamples made little difference in our application).

## From Causal Trees to Causal Forest

- ▶ We run the entire CF procedure using only  $S_{in}$ , then use the trees grown in  $S_{in}$  to generate predictions for all observations in  $S_{in}$  and  $S_{out}$ .
- ▶ This allows to assess the performance of the predictions in a hold-out sample (albeit with reduced statistical power) and to check whether heterogeneity is more distinct in  $S_{in}$  than  $S_{out}$ , which could be a sign of overfitting.
- ▶ Within each sample, we group youth by whether they are predicted to have a positive or negative treatment effect ( $\hat{\tau}_i^{CF} > 0$  is desirable for employment and adverse for arrests).
- ▶ We estimate separate treatment effects for these two subgroups by regressing each outcome on the indicator:

$$y_{ib} = \beta_1 I[\hat{\tau}_i^{CF} > 0] + \beta_2 T_i I[\hat{\tau}_i^{CF} > 0] + \beta_3 T_i (1 - I[\hat{\tau}_i^{CF} > 0]) + X\theta + \alpha_b + u_{ib} \quad (13)$$

# Causal Forests

TABLE 1—TREATMENT EFFECTS BY PREDICTED RESPONSE

Subgroup	No. of violent crime arrests	Any formal employment
<i>Panel A. In sample</i>		
$\hat{\tau}_i^{CF}(x) > 0$	0.22 (0.05)	0.19 (0.03)
$\hat{\tau}_i^{CF}(x) < 0$	-0.05 (0.02)	-0.14 (0.03)
$H_0$ : subgroups equal, $p =$	0.00	0.00
<i>Panel B. Out of sample</i>		
$\hat{\tau}_i^{CF}(x) > 0$	-0.01 (0.05)	0.08 (0.03)
$\hat{\tau}_i^{CF}(x) < 0$	-0.02 (0.02)	-0.01 (0.03)
$H_0$ : subgroups equal, $p =$	0.77	0.02
<i>Panel C. Adjusted in sample</i>		
$\hat{\tau}_i^{CF}(x) > 0$	-0.06 (0.04)	0.05 (0.03)
$\hat{\tau}_i^{CF}(x) < 0$	-0.02 (0.02)	-0.04 (0.03)
$H_0$ : subgroups equal, $p =$	0.41	0.02

# Review & Next Steps

- ▶ Bagging and Random Forests
- ▶ Comparisons: Lasso, CART, Random Forests
- ▶ AdaBoost
- ▶ Causal Forests
- ▶ Next class: More on boosting
- ▶ Questions? Questions about software?

## Further Readings

- ▶ Athey, S., & Imbens, G. (2016). Recursive partitioning for heterogeneous causal effects. *Proceedings of the National Academy of Sciences*, 113(27), 7353-7360.
- ▶ Davis, Jonathan M.V., and Sara B. Heller. 2017. "Using Causal Forests to Predict Treatment Heterogeneity: An Application to Summer Jobs." *American Economic Review*, 107 (5): 546-50.
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- ▶ Taddy, M. (2019). *Business data science: Combining machine learning and economics to optimize, automate, and accelerate business decisions*. McGraw Hill Professional.