## Lecture 11: Machine Learning

Generalized Random Forests

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# Machine Learning for Causal Treatment Effect Estimation

# Machine Learning for Causal Treatment

The goal of Random Forests is to estimate a conditional mean,

$$\mu(x_0) = E[Y \mid X = x_0]$$

But what if instead we wanted to estimate... **any** conditional function  $heta(x_0)$  ?

- Conditional linear regression coefficients
- Conditional causal treatment effects
- Conditional local average treatment effects using instruments
- ullet Quantiles of the conditional distribution of  $Y \mid X = x_0$
- Conditional class probabilities

Enter Athey, Tibshirani, and Wager (2019).

#### Generalized Random Forests

**Generalized Random Forests** by <u>Athey, Tibshirani, and Wager (2019)</u> extends the ensemble methods of Random Forests to essentially any conditional function  $\theta(x)$  that can be identified by local moment conditions.

They recast forests as an adaptive locally weighted estimator

- 1. Use the forest to calculate weighted set of neighbors for each observation (sound familiar?)
  - $\circ$  Employ problem-specific splitting rules to maximize the likelihood of capturing heterogeneity in heta(x)
- 2. Use those neighbors to estimate heta(x) as the solution to a local moment equation

# **GRF Setup**

#### **More formally:**

Suppose you have an IID sample with N observations.

You observe:

- $oldsymbol{\cdot}$   $O_i$  an observable quantity that encodes information relevant to estimating  $heta(\cdot)$ 
  - $\circ$  For nonparametric regression, this is just our our outcome  $O_i = \{Y_i\}$
  - $\circ$  For treatment effect estimation,  $O_i = \{Y_i, W_i\}$
  - Generally can contain more info
- ullet Auxiliary covariates Xi

#### **GRF**

Goal: Estimate solutions to local estimating equations of the form

$$E[\psi_{ heta(x),\ 
u(x)}(O_i\mid X_i=x)]=0$$

- heta(x) the parameter we care about (i.e. conditional mean, treatment effect, regression slope)
- ullet u(x) a (optional) nuisance parameter

--

ullet If conditional mean without noise, then  $\psi_{\mu(x)}(Y_i)=Y_i-\mu(x)$ 

#### **GRF Forest-Based Local Estimation**

One approach to estimating functions like  $\theta(x)$ :

• Find  $\hat{ heta}(x),\hat{
u}(x)$  that solve

$$\sum_{i=1}^n lpha_i \psi_{\hat{ heta}(x),\hat{
u}(x)}(O_i) = 0$$

- $oldsymbol{lpha}_i$ : weights measuring relevance of the  $i^{th}$  observation for fitting  $heta(\cdot)$  at x
  - Traditionally obtained using a kernel function (and sometimes a bandwidth)
  - Traditionally sensitive to curse of dimensionality

#### **GRF Forest-Based Local Estimation**

**GRF Approach:** use forests to derive  $\alpha_i$ .

- 1. Grow  $b=1,\ldots,B$  trees
- 2. Choose weights  $\alpha_i(x)$  to capture how frequently the  $i^{th}$  training example falls into the same leaf as x.

Weights within a given tree,  $lpha_{bi}(x)$ , are calculated as

$$lpha_{bi}(x) = rac{I[X_i \in L_b(x)]}{|L_b(x)|}.$$

ullet  $L_b(x)$  the set of observations falling within the same leaf as x

Overall weights for observation i are then the average across all trees:

$$lpha_i(x) = rac{1}{B} \sum_{b=1}^B lpha_{bi}(x)$$

# **GRF Weights**

For example, we can reframe random forests and our estimate of the conditional mean function as

$$\hat{\mu}(x) = \sum_{i=1}^n rac{1}{B} \sum_{b=1}^B Y_i rac{I[X_i \in L_b(x)]}{|L_b(x)|}$$

Or as the weighted sum of single tree predictions  $\hat{\mu}_b(\cdot)$ :

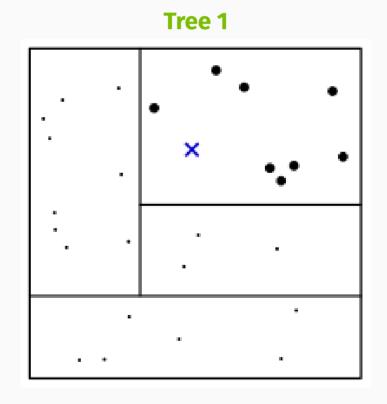
$$\hat{\mu}(x) = rac{1}{B} \sum_{b=1}^B \hat{\mu}_b(x)$$

Or more simply, as

$$\hat{\mu}(x) = \sum_{i=1}^n Y_i lpha_i(x)$$

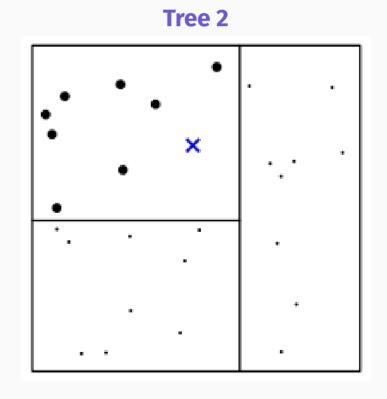
Let's see how these weights are developed with a simple example: a forest with three trees and two covariates.

- *x* the point of interest
  - lines our splits
  - darker dots the points
     within the same leaf



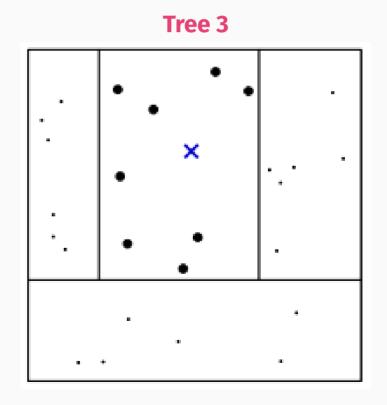
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Let's see how these weights are developed with a simple example: a forest with three trees and two covariates.

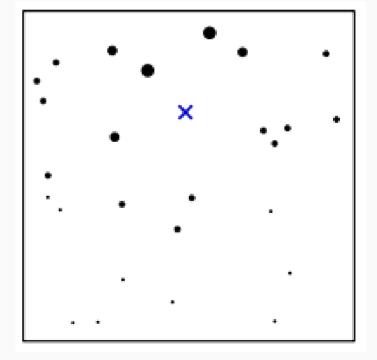
- *x* the point of interest
  - lines our splits
  - darker dots the points
     within the same leaf



Let's see how these weights are developed with a simple example: a forest with three trees and two covariates.

Then we average across all the individual tree-based weights to obtain the overall  $\alpha_i(x)$ .

#### Forest Weights for x



## Honesty

One additional benefit of using GRF: growing **honest** forests

**Honesty** aims to reduce bias in tree predictions by using **distinct subsamples** for 1. Building the tree, and 2. Making predictions

#### **Classic Random Forest**

 Draws a single subsample, use that subsample for both choosing a tree's splits and making predictions in that tree

#### **Honest Forests**

- Draw a training sample
- Use part of that sample to choose the tree's splits
- Use the rest of the training sample to make predictions
- Prune away any remaining empty leaves

#### **GRF Functions**

Conveniently for us, the **grf** package contains tons of features and extensive documentation to estimate a wide range of forests, including...

#### **Causal Treatment Effects**

Approach	Forest Function
Causal Treatment Effects	<pre>causal_forest()</pre>
Causal Treatment Effects with right-censored outcomes	<pre>causal_survival_forest()</pre>
Multi-arm/multi-outcome causal treatment effects	<pre>multi_arm_causal_forest()</pre>

## **GRF Functions**

Conveniently for us, the **grf** package contains tons of features and extensive documentation to estimate a wide range of forests, including...

#### **Moments of Conditional Distributions**

Approach	Forest Function
Conditional Mean	<pre>regression_forest()</pre>
Multi-outcome conditional means	<pre>multi_regression_forest()</pre>
Right-censored Survival	<pre>survival_forest()</pre>
Conditional Quantiles	<pre>quantile_forest()</pre>
Conditional Class Probabilities	<pre>probability_forest()</pre>

#### **GRF Functions**

Conveniently for us, the **grf** package contains tons of features and **extensive documentation** to estimate a wide range of forests, including...

#### **Regression Outcomes**

Approach	Forest Function
Conditional Linear Model	<pre>lm_forest()</pre>
Local Average Treatment Effects (IV)	<pre>instrumental_forest()</pre>

#### Causal Forests

When  $W_i \in \{0,1\}$  and **unconfoundedness holds**, we can estimate heterogeneous causal treatment effects with **causal forests**.

**1. Build Phase:** greedily choose splits to maximize the **squared difference** in subgroup treatment effects

$$(n_L n_R (\hat{ au}_L - \hat{ au})_R)^2$$

•  $\hat{ au}$ 's chosen through a centered residual-on-residual regression (Robinson 1988)

### Causal Forests

**2. Estimate**  $\tau(x)$ , where

$$au(x) := lm \ \left( Y_i - \hat{m}^{-1}(X_i) \sim W_i - \hat{e}^{-1}(X_i), ext{ weights} = lpha_i(x) 
ight)$$

- $ullet \left(Y_i \hat{m}^{-1}(X_i), \; W_i \hat{e}^{-1}(X_i)
  ight)$ : orthogonalized
  - outcome/treatment
    - $\circ$  Residual after "regressing out" the main effect of  $X_i$  on  $Y_i$  (W\_i)
    - $\hat{m}^{-1}(X_i)$ ,  $\hat{e}^{-1}(X_i)$  obtained from separate regression forests

Let's load in some data to see how we can run causal forests, using a sample from the **National Study of Learning Mindsets**.



```
nslm ← read.csv("data/NSLM.csv")
```

Taking a look at the data, we have the following variables:

- schoolid the ID of randomly selected US public high schools
- Y: continuous measure of achievement
- z: treatment status
- s3: students' self-reported expectations for future success
- c1: student race (categorical)
- c2 : student gender (categorical)
- c3: first-generation status (categorical)
- xc: school urbanicity (categorical)

- X1: school-level mean of students' fixed mindsets
- x2: school achievement level (test scores + college prep)
- x3: school racial minority composition (% black, latino, native american)
- X4: school poverty (% in familes below FPL)
- x5: school size

We want to answer the following research questions:

- 1. Was a nudge-like mindset intervention designed to instill a "growth mindset" in students effective at improving achievement?
- 2. Do schools' prior achievement levels  $(x_2)$  and pre-existing mindset norms  $(x_1)$  effect the magnitude of this effect?

Let's find out!

We could just run an (interacted) OLS regression...

```
 reg1 \leftarrow feols(Y \sim Z + S3 + C1 + C2 + C3 + XC + X1 + X2 + X3 + X4 + X5, \ data = nsl \\ reg2 \leftarrow feols(Y \sim Z + Z:X2 + Z:X3 + S3 + C1 + C2 + C3 + XC + X1 + X2 + X3 + X4 + X \\ etable(reg1, reg2, keep = c("Z", "X2", "X1"))
```

	reg1	reg2
Dependent Var.:	Υ	Υ
Z	0.2549*** (0.0115)	0.2539*** (0.0115)
X1	-0.0954*** (0.0073)	-0.0954*** (0.0073)
X2	-0.0163. (0.0087)	-0.0263** (0.0098)
Z x X2		0.0313* (0.0141)
Z x X3		0.0152 (0.0134)

...which gives us an estimate of the mean treatment effect and mean mediating effect of  $x_1/x_2$ .

Instead, let's use causal\_forest() to obtain heterogeneous treatment effects for each school.

Let's first build some data objects we'll use as arguments.

- 1. Convert c1 and xc to dummies:
  - C1 has 15 levels, so rather than code each one by hand let's use model.matrix
    - Argument: formula

```
# expand C1 categorical variable into dummies (with intercept)
C1_exp ← model.matrix(~ factor(nslm$C1) + 0)
```

Let's first build some data objects we'll use as arguments.

Repeat for xc:

```
XC_exp ← model.matrix(~ factor(nslm$XC) + 0)
```

Next, build the predictor matrix

```
# first select all covariates except for XC and C1
X ← select(nslm, -schoolid:-Y, -XC, - C1)
# Add in the dummies for XC and C1
Xmat ← cbind(X, C1_exp, XC_exp)
```

And get the outcome and treatment vectors along with the school clusters:

```
Y ← nslm$Y # our outcome
Z ← nslm$Z # our treatment indicator
schoolid ← nslm$schoolid
```

To perform the residual-on-residual regression step, our forest wants predictions for  $\hat{Y}$  and  $\hat{W}$ . We can either

- 1. Omit them from the forest setup
  - The forest will estimate them for us in a separate regression forest
- 2. Run the initial regression forests manually and supply the predictions

Let's do the latter, but first an important consideration...

As with random forests, by default GRF methods assume **independent observations**.

Here, we know that treatment assignment occurred at the **school level**, so students' treatment is dependent on which school that they're at.

The good news is, we can account for this **clustering** within our forest to draw our random units at the school level.

- clusters the clustering variable (here our schoolid factor)
- equalize.cluster.weights = TRUE ensures we draw the same fraction
   frmo each cluster

Running the initial regression forests for our  $\hat{Y}$  and  $\hat{W}$  predictions:

Setting up the causal forest:

• x: the covariate matrix

```
# Define the random forest
cf \leftarrow causal\_forest(X = Xmat,
                     Y = Y
                     Y.hat = Y_hat,
                     W = Z,
                     W.hat = W_hat,
                     clusters = schoolid,
                     equalize.cluster.weights =
                     num.trees = 200,
                     honesty = TRUE,
                     honesty.fraction = 0.5,
                     tune.parameters = "all"
```

- x: the covariate matrix
- Y: the outcome vector (our measure of student achievement)
- w: the treatment vector

```
# Define the random forest
cf \leftarrow causal\_forest(X = Xmat,
                     Y = Y.
                     W = Z.
                     Y.hat = Y_hat,
                     W.hat = W_hat,
                     clusters = nslm$schoolid,
                     equalize.cluster.weights =
                     num.trees = 2000.
                     honesty = TRUE,
                     honesty.fraction = 0.5,
                     tune.parameters = "all"
```

- x: the covariate matrix
- Y: the outcome vector (our measure of student achievement)
- w: the treatment vector
- Y.hat: our prediction of Y as a function of X
- W.hat: our prediction of W as a function of X

```
# Define the random forest
cf \leftarrow causal\_forest(X = Xmat,
                     Y = Y.
                     W = Z,
                     Y.hat = Y hat,
                     W.hat = W hat,
                     clusters = nslm$schoolid,
                     equalize.cluster.weights =
                     num.trees = 200.
                     honesty = TRUE,
                     honesty.fraction = 0.5,
                     tune.parameters = "all"
```

- clusters: the cluster identifier vector
- equalize.cluster.weights: whether to draw equal sample sizes from each cluster

```
# Define the random forest
cf \leftarrow causal\_forest(X = Xmat,
                     Y = Y.
                     W = Z.
                     Y.hat = Y hat,
                     W.hat = W hat,
                     clusters = nslm$schoolid,
                     equalize.cluster.weights =
                     num.trees = 200,
                     honesty = TRUE,
                     honesty.fraction = 0.5,
                     tune.parameters = "all"
```

- clusters: the cluster identifier vector
- equalize.cluster.weights: whether to draw
   equal sample sizes
   from each cluster
- num.trees: size of forest to grow (default is 2000)

```
# Define the random forest
cf \leftarrow causal\_forest(X = Xmat,
                     Y = Y.
                     W = Z.
                     Y.hat = Y hat,
                     W.hat = W hat,
                     clusters = nslm$schoolid,
                     equalize.cluster.weights =
                     num.trees = 200,
                     honesty = TRUE,
                     honesty.fraction = 0.5,
                     tune.parameters = "all"
```

- honesty: whether or not to grow honestly (default to TRUE)
- honesty.fraction:
   the subsample
   portion used for
   growing trees
   (defaults to 50%)

```
# Define the random forest
cf \leftarrow causal\_forest(X = Xmat,
                     Y = Y.
                     W = Z.
                     Y.hat = Y hat,
                     W.hat = W hat,
                     clusters = nslm$schoolid,
                     equalize.cluster.weights =
                     num.trees = 200.
                     honesty = TRUE,
                     honesty.fraction = 0.5,
                     tune.parameters = "all"
```

## Causal Forests Example

#### Setting up the causal forest:

tune.parameters: which/whether to tune parameters

```
o none, all,
sample.fraction,
mtry,
min.node.size,
honesty.fraction,
honesty.prune.leaves,
alpha,
imbalance.penalty
```

```
# Define the random forest
cf \leftarrow causal\_forest(X = Xmat,
                     Y = Y.
                     W = Z.
                     Y.hat = Y hat,
                     W.hat = W hat,
                     clusters = nslm$schoolid,
                     equalize.cluster.weights =
                     num.trees = 200.
                     honesty = TRUE,
                     honesty.fraction = 0.5,
                     tune.parameters = "all"
```

## Variable Importance

Looking at variable importance with variable\_importance(forest):

• Scaled count of how frequently variables were chosen to split on (i.e. that they maximized treatment effect heterogeneity)

covar	imp
X1	0.214
X2	0.149
Х3	0.137
X4	0.131
X5	0.0915
S3	0.0656

## Split Frequencies

Alternatively, we can obtain the raw split frequencies with...

```
split_frequencies(forest, max.depth)
```

```
# getting split frequencies 4 branches deep
splitfreq ← split_frequencies(cf, max.depth = 4)
colnames(splitfreq) ← colnames(Xmat)
rownames(splitfreq) ← paste("Depth", 1:4)
splitfreq[,1:8]
```

```
## Depth 1 9 9 5 50 33 28 29 15 ## Depth 2 38 23 19 57 45 59 38 59 ## Depth 4 112 74 65 70 64 63 70 80
```

#### Variable Selection

Our variable importance and split frequencies reveals useful information about the covariates: some **really don't matter** 

Since we're only considering a random subset of covariates at each potential split, by including covariates without explanatory power for treatment effect heterogeneity, we're likely **adding noise** and **diluting** the performance of our forest.

Solution: re-run the tree, keeping only the most important variables

• Athey and Wager (2019): keep only variables greater than mean importance

## Variable Selection

#### Selecting the "important" covariates:

```
varimp ← variable_importance(cf)
X_sel ← Xmat[,which(varimp> mean(varimp))]
colnames(X_sel)

## [1] "S3" "C2" "X1" "X2"

## [5] "X3" "X4" "X5" "factor(nslm$C1
```

# Causal Forests Example

And rerunning the forest with just these variables:

Use the average\_treatment\_effect() function to get one of several average treatment effects:

- target.sample = "all": The ATE
  - $\circ E[Y(1)-Y(0)]$
- target.sample = "treated": The ATT

$$\circ \ E[Y(1) - Y(0) \mid W_i = 1]$$

Use the average\_treatment\_effect() function to get one of several average treatment effects:

target.sample = "control": The average treatment effect on the
 control

$$\circ \ E[Y(1) - Y(0) \mid W_i = 0]$$

- target.sample = "overlap": The overlap-weighted average treatment effect on the control
  - $\circ E[e(X)(1-e(X))(Y(1)-Y(0))]/E[e(X)(1-e(X))]$
  - $\circ$  Where  $e(X) = P[W_i = 1 \mid X_i = x]$

Use the average\_treatment\_effect() function to get one of several average treatment effects:

```
ate ← average_treatment_effect(cf_sel, target.sample = "all")
ate

## estimate std.err
## 0.25446427 0.01082919
```

Retrieving the out-of-bag predictions for au(X) with predict():

```
tauhat_oob ← predict(cf_sel, estimate.variance = TRUE)
head(tauhat_oob)
```

predictions	variance.estimates	debiased.error	excess.error
0.238	0.00167	0.0151	1.13e-05
0.196	0.00194	0.00903	1.52e-05
0.241	0.00155	0.0344	1.12e-05
0.245	0.00196	0.161	1.16e-05
0.226	0.00213	0.355	1.15e-05
0.242	0.00249	0.0934	1.07e-05

#### Causal Forest Error

In addition to our predicted treatment effects and variance estimate, we get **two types of error**:

#### **Debiased Error**

- Estimate of the "R-loss"Criterion
  - i.e. the **predictive fit** of our forest

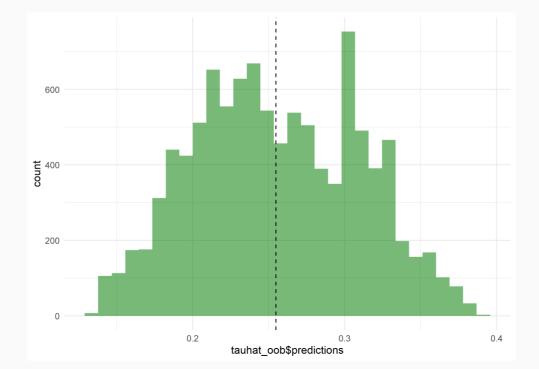
#### **Excess Error**

- Error resulting from random nature of forests
  - i.e. how unstable our estimates would be if we re-grew forests with the same data/setup
  - We want this to be very small relative to estimate's variance

### **Treatment Effects**

Plotting the distribution of individual treatment effects (and relative to the OLS mean estimate:)

```
ggplot() +
  geom_histogram(aes(tauhat_oob$predictions), fill = "forestgreen", alpha = 0.6) +
  geom_vline(aes(xintercept = reg1$coefficients["Z"]), linetype = "dashed") +
  theme_minimal()
```



## **ATE Subset Comparison**

We can also use subset to see differences in top/bottom treatment effects

```
# Getting boolean vectors of top/bottom quartile treatment effects high_effect \leftarrow tauhat_oob$predictions > quantile(tauhat_oob$predictions, 0 low_effect \leftarrow tauhat_oob$predictions < quantile(tauhat_oob$predictions, 0
```

## **ATE Subset Comparison**

We can also use subset to see differences in top/bottom treatment effects

```
# getting ATE for each group
ate_high = average_treatment_effec
ate_high

# getting ATE for each group
ate_low = average_treatment_effect
ate_low

## estimate std.err
## 0.35006924 0.05311834

## [1] "95% CI for difference in ATE: 0.117 +/- 0.124"
```

## Causal Forests Example

Re-estimating the forest **without clusters** shows the importance of accounting for the cluster assignment:

#### Test Calibration

Looking at the **Test Calibration** results for each forest shows the importance of clustering.

test\_calibration() Regresses the forest error for prediction of Y on **two terms:** 

- 1. Mean Prediction: indication of how well the forest captures the mean ATE
  - $\circ$  Coefficient indistinguishable from 1  $\rightarrow$  accurate mean prediction
- 2. Differential Prediction: how well the forest captures heterogeneity in the ATF
  - $\circ$  Coefficient distinguishable from 0  $\to$  confirms existence of heterogeneous treatment effects
  - Coefficient indistinguishable from 1 → accurately capture heterogeneity

### Test Calibration

First for the non-clustered forest:

```
##
## Best linear fit using forest predictions (on held-out data)
## as well as the mean forest prediction as regressors, along
## with one-sided heteroskedasticity-robust (HC3) SEs:
##
## Estimate Std. Error t value Pr(>t)
## mean.forest.prediction 0.997121 0.042448 23.4903 < 2.2e-16 ***
## differential.forest.prediction 1.148856 0.202310 5.6787 6.969e-09 ***
## ---
## Signif. codes: 0 '***' 0.001 '**' 0.01 '*' 0.05 '.' 0.1 ' ' 1</pre>
```

#### Test Calibration

Looking at the **Test Calibration** results for each forest shows the importance of clustering.

Next for the unclustered forest:

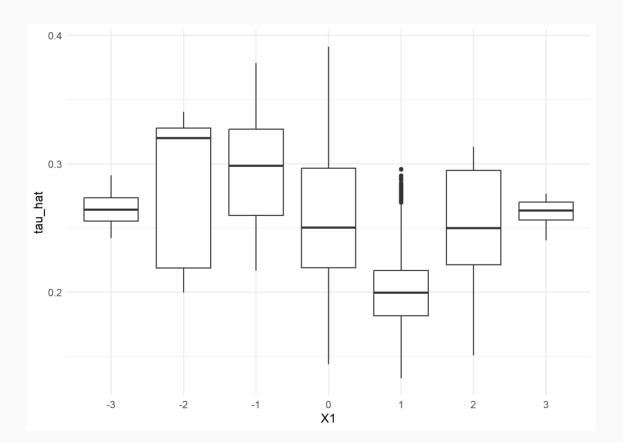
```
test_calibration(cf_noclust)
```

```
##
## Best linear fit using forest predictions (on held-out data)
## as well as the mean forest prediction as regressors, along
## with one-sided heteroskedasticity-robust (HC3) SEs:
##
## Estimate Std. Error t value Pr(>t)
## mean.forest.prediction 1.015750 0.045235 22.4549 < 2.2e-16 ***
## differential.forest.prediction 0.531436 0.122606 4.3345 7.374e-06 ***
## ---
## Signif. codes: 0 '***' 0.001 '**' 0.05 '.' 0.1 ' ' 1</pre>
```

# Treatment Effect Heterogeneity

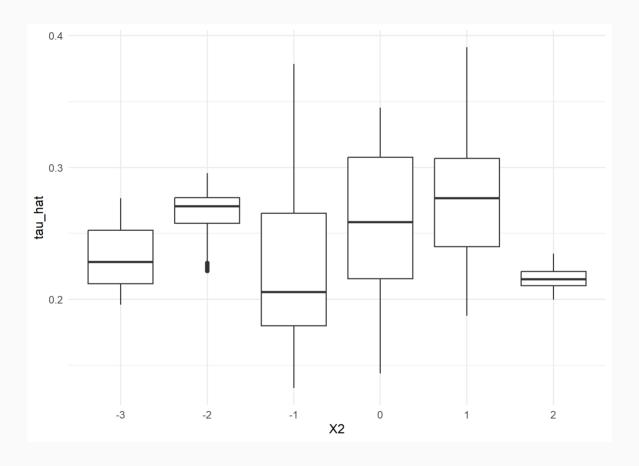
It looks like there's a clear effect of the intervention nudge on achievement, but what about the mediating effects of x1 and x2?

First, let's look at the variation in treatment effects by the values of x1:



## Treatment Effect Heterogeneity

Repeating for pre-existing achievement levels (x2):



It appears that the nudge had the least impact in schools with the highest preexisting achievement level ( $x_2$ )

#### **Covariate Correlation**

One potential limitation: there are a lot of factors at the school-level associated with pre-existing achievement levels.

```
x2\_reg \leftarrow feols(X2 \sim X1 + X3 + X4 + X5 + C2 + C3 \mid XC + C1, data = nslm) etable(x2_reg)
```

	x2_reg
Dependent Var.:	X2
X1	-0.2858 (0.1592)
Х3	-0.2635 (0.1368)
X4	-0.0902 (0.0859)
X5	0.2635. (0.1117)
C2	0.0023 (0.0155)

#### Simulations

**Solution:** use a new "simulated" covariate matrix to predict treatment effects if schools had different achievement levels, holding other characteristics fixed.

Let's write a function that will let us obtain treatment effect predictions if we swap the value of x2 with a particular value in all rows:

### Simulations

Next, build a vector of X2 values to predict for:

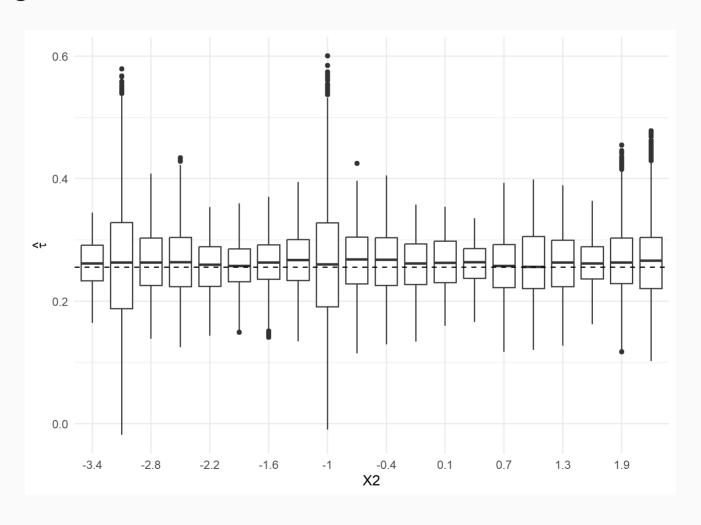
```
# get range of X2 to predict over
min_x2 ← min(Xmat$X2)%>% round(2)
max_x2 ← max(Xmat$X2) %>% round(2)
x2_seq ← seq(min_x2, max_x2, length.out = 20)
```

#### And running:

```
# Vectorizing and saving out
sim_x2 ← map_dfr(x2_seq, sim_x2)
saveRDS(sim_x2, "output/x2_sim_df.rds")
```

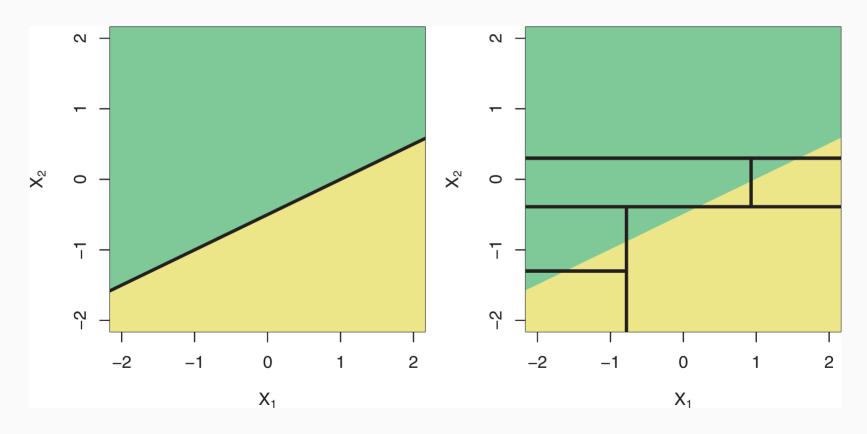
# Simulations

#### Plotting the simulation results:



## **Linear Boundaries**

What if I want to use these methods to estimate heterogeneous treatment effects but I'm concerned about replicating **linear boundaries**?



A: local linear forests

<u>Developed in Friedberg et al. (2021)</u>, local linear forests combine the local smoothness benefits of local linear regression with the high-dimensional flexibility of random forests.

#### **Random Forests**

- Highly data-adaptive
- Struggle to capture smoothness in conditional mean function

#### **Local Linear Regression**

- Great at capturing locally smooth signals
- Quickly deteriorates due to curse of dimensionality
  - Relies on Euclidean distance, which loses locality even in 4-5 dimensions

Local linear forests function as a Two-stage adaptive kernel method:

- 1. Grow a random forest with splits chosen to minimize the sum of squared errors between the resulting two branches
  - Uses a standard splitting procedure but splitting on the residuals
     from a ridge regression
  - Partitions the covariate space based on local effects, captures existing linear signals
- 2. Use the forest to obtain weights  $lpha_i(x_0)$
- 3. Rather than use the weights to fit the local average at  $x_0$  as in random forests, Use the  $lpha_i(x_0)$  weights to fit a **local linear regression** 
  - Includes a ridge penalty for regularization

#### **More formally:**

Consider two parameters:

- 1.  $\mu(x_0)$ : the local average at  $x_0$
- 2.  $heta(x_0)$ : the slope of the local line
  - $\circ$  Correction for local trend in  $X_i-x$

**Goal:** solve for estimates  $\hat{\mu}(x_0)$ ,  $\hat{ heta}(x_0)$  that minimize

$$\sum_{i=1}^n \underbrace{lpha_i(x_0)}_{ ext{Weights}} (Y_i - \mu(x_0) - \underbrace{(X_i - x_0) heta(x_0)}_{ ext{Local Trend Correction}})^2 + \underbrace{\lambda|| heta(x_0)||_2^2}_{RidgePenalty}$$

Local linear forests also extend to causal forests:

1. Use local linear forests to obtain the estimates for

$$\left(\hat{m}^{-1}(X_i),\; \hat{e}^{-1}(X_i))
ight)$$

ullet i.e. the leave-one-out initial predictions for Y and W obtained from auxiliary forests

Local linear forests also extend to causal forests:

2. Estimate conditional average treatment effect  $\hat{ au}$  as the arg min of

$$egin{aligned} \sum_{i=1}^n lpha_i(x_0) \Big( Y_i - \hat{m}^{-1}(x_i) - a - (X_i - x_0) heta_a \ &- ( au + heta_t(X_i - x_0)) (W_i - \hat{e}^{-1}(X_i)) \Big)^2 \ &+ \lambda_ au || heta_t||_2^2 + \lambda_a || heta_a||_2^2 \end{aligned}$$

- a intercept (0 if  $\left(\hat{m}^{-1}(X_i),\;\hat{e}^{-1}(X_i)
  ight)$  estimates are accurate)
- ullet  $\lambda_{ au}$  and  $\lambda_a$ : ridge penalties for local trends in treatment and outcome
  - As these get large, we recover a standard causal forest

To estimate a local linear causal forest:

1. Use <code>ll\_regression\_forest()</code> instead of <code>regression\_forest()</code> to obtain the initial estimates of  $\left(\hat{m}_{ll}^{-1}(X_i),\;\hat{e}_{ll}^{-1}(X_i)\right)$ 

To estimate a local linear causal forest:

2. Run Causal forest using the local linear forest estimates from 1.

To estimate a local linear causal forest:

3. Use local linear arguments of predict() when generating predictions for au

Comparing overall fit (Root MSE) between the two:

```
 rmse\_cf \leftarrow sqrt(sum((cf\_sel\$Y.orig - cf\_sel\$Y.hat)^2)/length(cf\_sel\$Y.ori\xi rmse\_ll \leftarrow sqrt(sum((cf\_sel\_ll\$Y.orig - cf\_sel\_ll\$Y.hat)^2)/length(cf\_sel\_data.frame(Forest = c("Causal Forest", "Local Linear Causal Forest"), \\ RMSE = c(rmse\_cf, rmse\_ll) %>% round(4))
```

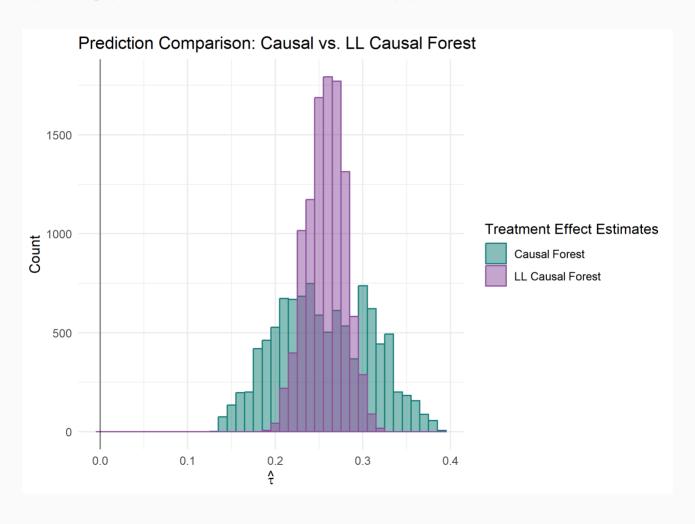
Forest	RMSE
Causal Forest	0.533
Local Linear Causal Forest	0.551

Comparing test calibration results:

```
##
## Best linear fit using forest predictions (on held-out data)
## as well as the mean forest prediction as regressors, along
## with one-sided heteroskedasticity-robust (HC3) SEs:
##
## Estimate Std. Error t value Pr(>t)
## mean.forest.prediction 0.997121 0.042448 23.4903 < 2.2e-16 ***
## differential.forest.prediction 1.148856 0.202310 5.6787 6.969e-09 ***
## ---
## Signif. codes: 0 '***' 0.001 '**' 0.01 '*' 0.05 '.' 0.1 ' ' 1</pre>
```

Comparing test calibration results:

And comparing predictions from the two approaches:



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