# Lecture 16: Big data and machine learning

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# From last time: fuzzy regression discontinuity

As usual, we'd like to run:

$$Y_i = \alpha + \tau D_i + \varepsilon_i$$

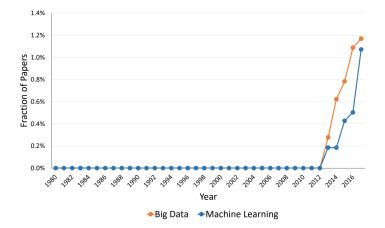
The regression discontinuity:

- Suppose  $D_i$  is determined by whether or not  $X_i$  lies above a cutoff, c
- Idea: Having X<sub>i</sub> just above or just below c is as good as random...
- ... And there is a discontinuous change in  $D_i$  as a result of crossing c
- $\rightarrow$  We can compare  $Y_i$  for units with  $X_i$  just above c to  $Y_i$  for units with  $X_i$  just below c
  - With incomplete changes in  $D_i$  from  $X_i < c$  to  $X_i \ge c$ :

$$D_i = \alpha + \gamma \mathbf{1}[X_i \ge c] + f(X_i) + \varepsilon_i \text{ for } c - h < X_i < c + h$$

$$Y_i = \alpha + \tau \hat{D}_i + f(X_i) + \varepsilon_i$$
 for  $c - h < X_i < c + h$ 

# And now for something completely different...



A computer scientist might say:

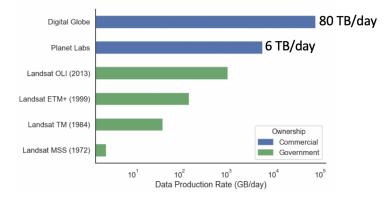
• "Too big to fit on your computer"

An economist might say:

• "I dunno, 15 GB?"

 $\rightarrow$  All of this is a bit fuzzy

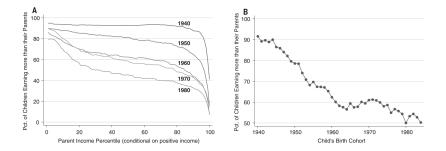
# Big data are getting bigger every day



The era of Big Data isn't just good because of small standard errors:

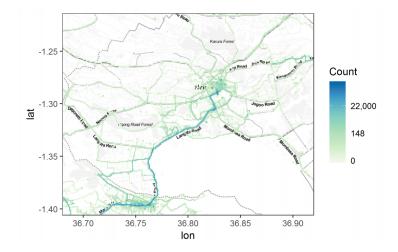
- New data collection methods present opportunities
- We can study previously unanswerable questions
- Sometimes this requires getting a bit creative

### Thinking outside of the box to get new data



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# Thinking outside of the box to get new data



(Especially) with Big Data, we have to be careful:

- Just because a dataset is large doesn't mean it's unbiased
- Large data can also have errors
- And come with additional concerns too (privacy, etc)

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- $\rightarrow\,$  It's important to understand what we're using
- $\rightarrow$  (The following slides owe credit to Tamma Carleton)

We typically interact with three types of data:

- **1** Raw, out of the source
- Processed "in house"
- 8 Processed "out of the house"
- $\rightarrow$  All of these data can be used as Y, D, or X (or even Z)
- $\rightarrow\,$  Each has its own pros and cons

Major pros of raw data:

- We know what we're dealing with
- We get a fighting chance to understand measurement error and bias

Major cons of raw data:

- Raw data are often not exactly what we want
- We have to be careful when we use them as a proxy

# Raw data



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#### Program Evaluation

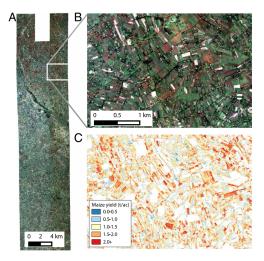
Major pros of home-grown data:

- We know what we're dealing with
- We get a fighting chance to understand measurement error and bias

Major cons of home-grown data:

- This takes a lot of time and effort
- And we don't always have the right toolkit

# Processed in house



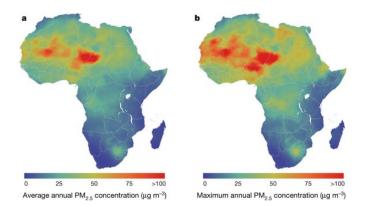
Major pros of outsourced data:

- We leverage external expertise
- We potentially have less measurement error than the in-house version
- This is a lot less work than the

Major cons of outsourced data:

- We don't know exactly what we're measuring
- We can't look "under the hood" to uncover bias

# Processed out of the house



Just like with small data, you need to know what you've got:

- Big Data allow for new possibilities
- But require additional processing tools and time
- A careful combination of in-house and out-of-house work can yield benefits

# Estimation vs. prediction

This class has been about asking:

- $\rightarrow$  What is the causal effect of D on Y?
- $\rightarrow$  Aka, in

$$Y_i = \alpha + \tau D_i + \varepsilon_i$$

what is  $\hat{\tau}$ ?

• Focus is on **unbiasedness** 

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what is  $\hat{\tau}$ ?

• Focus is on **unbiasedness** 

Machine learning instead asks:

- $\rightarrow$  What is the best guess of some outcome?
- $\rightarrow$  What is  $\hat{Y}$ ?
  - Want to consider a bias-variance tradeoff

An estimation problem:

"What is the causal effect of my rain dance on rainfall today?"

 $\rightarrow$  Estimation problem: rain dances (maybe) affect rainfall

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→ Estimation problem: rain dances (maybe) affect rainfall A prediction problem:

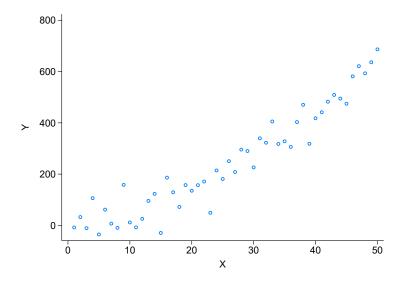
"Do I need an umbrella today?"

 $\rightarrow$  Prediction problem: rainfall doesn't depend on umbrellas

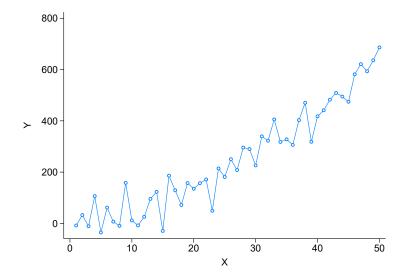
Machine learning is just methods trying to generate predicions:

- Given a dataset with outcome Y and covariates X, what function f(X) best predicts Y?
- Note the difference between this and causal inference!









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### In-sample prediction:

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#### Oross-validation:

- Instead of using the whole sample for step (1)...
- Split the sample into pieces...
- Do step (1) on one part, and predict  $\hat{Y}$  on the other part
- Record how well the model fits (eg  $Y \hat{Y}$ )

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8 Repeat:

- Do this several times over different sample splits
- Pick the final model that does best

#### Program Evaluation

It's worth unpacking this a bit further:

- **Goal:** Produce the best guess at  $\hat{f}(\mathbf{X})$
- This typically involves being very flexible: interactions between Xs
- We know we want to avoid over-fitting
- Just running a ton of OLS regressions is a slow way to do this

# The in-sample prediction step

Function class $\mathcal F$ (and its parametrization)	Regularizer $R(f)$
Global/parametric predictors	
Linear $\beta' x$ (and generalizations)	Subset selection $  \beta  _0 = \sum_{j=1}^{k} 1_{\beta_j \neq 0}$
	LASSO $  \beta  _1 = \sum_{j=1}^k  \beta_j $
	Ridge $  \beta  _2^2 = \sum_{i=1}^k \beta_i^2$
	Elastic net $\alpha   \beta  _1 + (1 - \alpha)   \beta  _2^2$
Local/nonparametric predictors	
Decision/regression trees	Depth, number of nodes/leaves, minimal lead size, information gain at splits
Random forest (linear combination of trees)	Number of trees, number of variables used in each tree, size of bootstrap sample, complexity of trees (see above)
Nearest neighbors	Number of neighbors
Kernel regression	Kernel bandwidth
Mixed predictors	
Deep learning, neural nets, convolutional neural networks	Number of levels, number of neurons per level, connectivity between neurons
Splines	Number of knots, order
Combined predictors	
Bagging: unweighted average of predictors from bootstrap draws	Number of draws, size of bootstrap samples (and individual regularization parameters)
Boosting: linear combination of predictions of residual	Learning rate, number of iterations (and individual regularization parameters)
Ensemble: weighted combination of different predictors	Ensemble weights (and individual regularization parameters)

ML is a great tool, but we have to be careful!

- Do NOT try to interpret the function! 222
- The ML model gives you  $\hat{Y}$ ...but not  $\hat{\tau}$ !
- We're not recovering causal effects
- And we don't get standard errors
- And the models are typically unstable

Machine learning is not designed for  $\hat{\tau}$ :

- We can't directly use ML for what we want to estimate
- But does this mean ML is useless for us?

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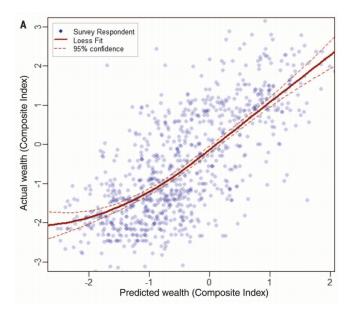
 $\rightarrow$  No.

 $\rightarrow\,$  We just need to be a little bit creative!

There are three main ways to use ML for causal inference:

- **1** Data generation
- **2** Heterogeneity analysis
- **8** Estimating  $\hat{\tau}$

#### ML for data generation



We often want to find heterogeneous effects:

- The FPCI gets in our way for this too
- We need to compare treated vs. untreated units...
- ... conditional on  $X_i = x$
- But there might be many X<sub>i</sub>
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- We need to compare treated vs. untreated units...
- ... conditional on  $X_i = x$
- But there might be many X<sub>i</sub>
- And they might even be continuous
- $\rightarrow$  Which  $X_i$  have interesting heterogeneity in  $\tau_i(X_i)$ ?

Reframe this into a prediction question:

- What is predicted  $\hat{Y}_i(X_i)$ ?
- That is, which  $X_i$ s give you different  $\hat{Y}_i$ ?
- For this to be the same as heterogeneity in  $\tau_i(X_i)...$
- ... we need random assignment to treatment
- Under random assignment, there is a 1:1 mapping between  $\hat{Y}_i$  and  $\hat{\tau}_i$

The most common approach is the **causal tree**:

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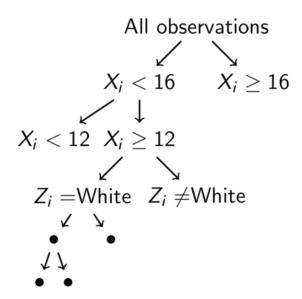
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- $\rightarrow$  Repeat with different training samples to construct a causal forest

#### Causal trees



Now we're really pushing the frontier:

- 1 ML with selection on observables
- **2** ML with selection on unobservables
- $\rightarrow\,$  We need to reframe our questions as prediction problems

Consider an underlying model:

$$Y_i = \alpha + \tau D_i + f(\mathbf{X}_i) + \varepsilon_i$$

where  $E[\varepsilon|D, \mathbf{X}] = 0$ : Conditional on **X**, *D* is as good as random

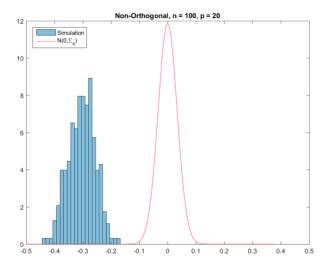
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- But which **X**s matter?
- And what is the right f(X)?
- We can use ML to help us figure this out
- → Simple guess: simply predict  $\hat{Y}$  based on D and X; interpret coefficient on D as  $\tau$



The simple guess doesn't work!

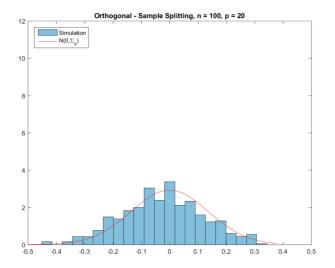
- The overall best fit ignores the SOO assumption
- Some X<sub>i</sub> that are important for D<sub>i</sub> may be left out
- $\rightarrow$  ML will choose  $X_i$  that are important for  $Y_i$
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We can do better!

- **1** Predict  $\hat{Y}$  as a function of X
- **2** Also predict  $\hat{D}$  as a function of X
- **8** Estimate treatment effects using both sets of covariates
- $\rightarrow$  This only works when you do steps (1) and (2) with LASSO

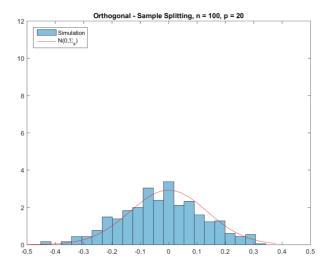


The most up-to-date approach is:

- **1** Predict  $\hat{Y}$  as a function of X
- **2** Predict  $\hat{D}$  as a function of X
- **3** Compute residuals:  $Y^R = Y \hat{Y}$  and  $D^R = D \hat{D}$
- **4** Recover  $\hat{\tau}$  by regressing:

$$Y_i^R = \alpha + \tau D^R + \varepsilon_i$$

- $\rightarrow$  You can do steps (1) and (2) with any ML method
- $\rightarrow\,$  Note that this approach still needs the SOO assumptions



Just like with SOO, ML can be useful to pick covariates:

- ML helps pick  $X_i$  when we don't need them for identification
- This works for RCTs, DD, IV, etc with exogenous covariates only

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- ML helps pick  $X_i$  when we don't need them for identification
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- We can use ML when we want a better fit

Great application: first stage of IV:

- Conditional on (a) good instrument(s), we just want a good fit
- Nothing wrong with using ML to improve the first stage...
- As long as you're only using exogenous covariates

We need to reframe as a prediction exercise:

- What is a(n estimated) counterfactual?
- Just a guess at what would've happened without treatment
- This is a simple prediction exercise
- We can potentially use ML to help us generate this counterfactual
- (Subject to all of the standard selection / identification issues)

# An example: Energy efficiency in California schools

Policy issue:

- Lots of money is being spent on EE upgrades
- But are they effective?

Approach:

- Look at hourly data from 2,000 public K-12 California schools
- Some schools decided to implement EE upgrades
- This was not randomized, so we use an FE approach
- $\rightarrow\,$  Leverage high-resolution data for an ML-augmented FE method

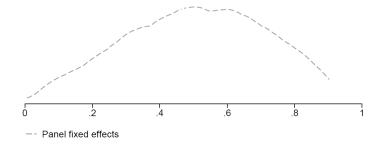
# Estimating the effects of EE upgrades: version A

- We compare:
  - Consumption at schools that retrofitted to those that didn't
  - Consumption before and after retrofits
- We progressively add a series of control variables (school, hour and month-of-sample fixed effects, plus interactions):

$$Y_{ith} = \tau D_{it} + \alpha_i + \kappa_h + \gamma_t + \varepsilon_{ith}$$

Interpretation of  $\tau$ : Average reduction in KWh at treated schools.

#### Panel FE results are unstable



- Panel FE models aren't properly specified.
- Schools are very heterogeneous (e.g., climate, size, school calendar).
  - Ideally, introduce school-specific coefficients and trends in a very flexible manner.
- We easily came up with ~6,000,000 candidate control variables by making them school-hour specific!
- No clear ex ante optimal choice.

## Machine Learning: Advantages in this application

- Exogenous weather variation and predictable weekly and seasonal patterns drive variation in electricity consumption.
- Schools are relatively stable consumption units:
  - as opposed to single households that move around, unobservably buy a new appliance, expand family size, etc.
  - as opposed to businesses and manufacturing plants, exposed to macroeconomic shocks.

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Prediction can do well!

#### Step 1

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  - Use LASSO method (penalized regression).
    - Minimizing the sum of the squared errors plus  $\lambda \cdot \sum_{j=1}^{p} |\beta_j|$ .
    - Larger "tuning parameters" lead to fewer coefficients.
    - Use bootstrapped cross-validation with training and holdout samples within pre-treatment.

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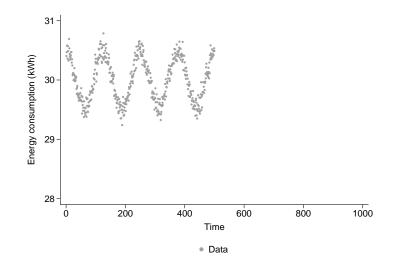
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    - Larger "tuning parameters" lead to fewer coefficients.
    - Use bootstrapped cross-validation with training and holdout samples within pre-treatment.
  - Include a wide range of school-specific variables, and also consumption at control schools (a la synthetic control).
  - Also consider other alternatives (random forests).

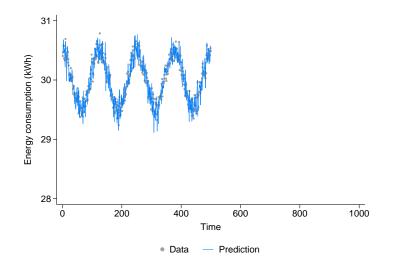
#### Step 2

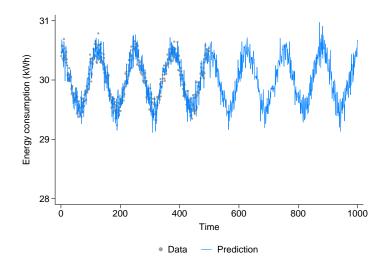
• Regress *prediction errors* on treatment and controls.

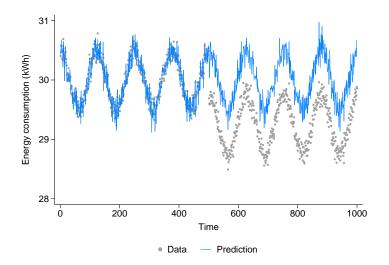
$$Y_{ith} = \tau D_{it} + \alpha_i + \kappa_h + \gamma_t + \varepsilon_{ith}$$

- Data pooled across schools
- Replicates diff-in-diff approach, but Y variable is now the prediction error

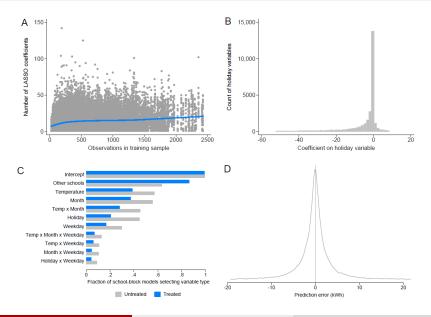






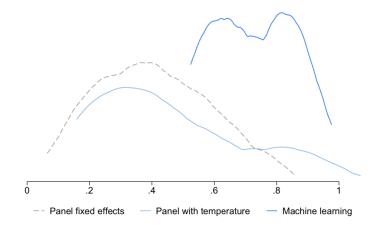


## ML diagnostics



Program Evaluation

#### ML results are stable across estimators



#### TL;DR:

- 1 New datasets open new questions
- 2 Machine learning offers opportunity
- **8** Both require some careful consideration or tweaks to be useful for us