Data Science for Causal Inference

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2024-07-15

Table of contents I

Introductions

Data Science in Causal Inference

Heterogeneous Treatment Effects

Variable Selection



About Me

- Associate Prof of Government (American University)
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- Senior Social Scientist (The Lab @ DC)
- ➤ Fellow in Methodology (US Office of Evaluation Sciences: "OES")

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- Research agenda: political methodology, causal inference, experimental design, experiments in public policy

Name?

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- ▶ Olympic sport you look forward to?

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- ▶ Data Science in Causal Inference
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- Sensitivity

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- Data Science in Causal Inference
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- Sensitivity
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- ▶ Data Science in Causal Inference
 - Models
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 - Staggered adoption

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 - Calloway-Sant'Anna approach

Data Science in Causal Inference

The "potential outcomes" framework: $% \left(1\right) =\left(1\right) \left(1\right) \left($

		Would Enroll if	Would Enroll if	
Citizen	Canvass?	Canvass?	No Canvass?	Enroll
1	Yes	Yes		Yes
2	Yes			Yes
3	No			No
4	No			No

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2	Yes	Yes	(No)	Yes
3	No	(Yes)	No	No
4	No	(No)	No	No

The "potential outcomes" framework, more abstractly:

					True τ
Unit i	Treatment T	Y(1)	Y(0)	$Y^{ m obs}$	Y(1) - Y(0)
1	1	10		10	
2	1	20		20	
3	0		15	15	
4	0		5	5	

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3	0	(40)	15	15	25
4	0	(20)	5	5	15

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				$\widehat{ATE} = \hat{\bar{\tau}} =$	15 - 10 = 5

The "potential outcomes" framework, notation:

- \triangleright Units indexed by i
- Treatment T_i or D_i or Z_i
- \triangleright Outcome if treated $Y_i(1)$
- \triangleright Outcome if control $Y_i(0)$
- ightharpoonup True treatment effect $\tau_i = Y_i(1) Y_i(0)$
- True average treatment effect
 - $\bar{\tau} = \frac{1}{n} \sum_{i=1}^{n} (Y_i(1) Y_i(0))$
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$$\bar{\tau} = \frac{1}{n} \sum_{i=1}^{n} (Y_i(1) - Y_i(0))$$

▶ Pre-treatment covariates X

(and we'll draw some DAG's, too)

Data Science Approaches

Three tasks of data science:

Description

Three tasks of data science:

- Description
- ▶ Prediction

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Models/algorithms central to all three.

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Models/algorithms central to all three.

Hernán, Hsu, and Healy (2019)

Description

▶ Identifying patterns, etc.

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- ► E.g., clustering to discover groups

Prediction

► Components

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 - ► Inputs/outputs (predictors/outcomes, features/responses, ...)

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 - Metric for evaluating mapping
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- ▶ E.g., regression, random forests, neural networks, ...

Causal Inference

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 - (the more knowledge, the better!)
 - (alternative: solve fundamental problem of causal inference!)
- ► E.g., experiments, observational causal designs, ...

Causal Inference with Machine Learning

Causal Inference with Machine Learning



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I finally found it in real life: the consultant who runs OLS in Excel and calls it machine learning

9:17 AM · Jan 31, 2019 · Twitter for iPhone

54 Retweets	7 Quote Tweets	511 Likes		
\Diamond	↑	\bigcirc	riangle	

Causal Inference with Machine Learning



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(OK, not "machine learning", perhaps, but models at least ...)

Loaded two datasets:

str(df1)

str(df2)

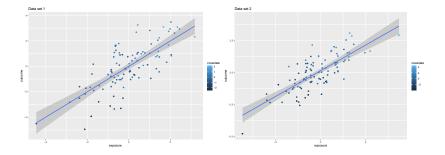
```
tibble [100 x 3] (S3: tbl_df/tbl/data.frame)
$ covariate: num [1:100] -0.622 1.137 -0.238 1.529 -0.154
$ exposure : num [1:100] 0.0332 0.3627 0.2422 1.4633 0.779
$ outcome : num [1:100] -0.429 2.675 -0.647 2.238 1.044
```

```
tibble [100 x 3] (S3: tbl_df/tbl/data.frame)

$ exposure : num [1:100] 0.4862 0.0653 -1.4021 -0.546 -0.4

$ outcome : num [1:100] 1.706 0.669 -1.597 -1.733 0.617
```

\$ covariate: num [1:100] 2.24 0.924 -0.999 -2.343 0.207 .



Model each

```
lm_df1 <- lm(outcome ~ exposure, data = df1)
lm_df2 <- lm(outcome ~ exposure, data = df2)</pre>
```

```
# A tibble: 4 x 4
data term estimate std.error
<chr> <chr> <chr> <chr> 0.00671 0.120
df1 (Intercept) -0.00671 0.120
df1 exposure 0.996 0.0927
df2 (Intercept) 0.133 0.0890
df2 exposure 1.00 0.0841
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▶ Both cases: effect of exposure ≈ 1 .

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```

- ▶ Both cases: effect of exposure ≈ 1 .
- ▶ Is this good?
- ▶ What if we adjust for covariate?

```
lm_df1_adj <- lm(outcome ~ exposure + covariate, data = df:
lm_df2_adj <- lm(outcome ~ exposure + covariate, data = df:</pre>
```

▶ Both cases: effect of exposure ≈ 0.5 .

```
lm_df1_adj <- lm(outcome ~ exposure + covariate, data = df:
lm_df2_adj <- lm(outcome ~ exposure + covariate, data = df:</pre>
```

```
# A tibble: 4 x 4
data term estimate std.error
<chr> <chr> <chr> <chr> dbl> cdbl>
1 df1 exposure 0.501 0.108
2 df1 covariate 0.970 0.147
3 df2 exposure 0.554 0.0990
4 df2 covariate 0.385 0.0598
```

- ▶ Both cases: effect of exposure ≈ 0.5 .
- ▶ Is this good?

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- ▶ Both cases: effect of exposure ≈ 0.5 .
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- Which is correct? $\beta = 1$? $\beta = 0.5$?

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```

- ▶ Both cases: effect of exposure ≈ 0.5 .
- ▶ Is this good?
- Which is correct? $\beta = 1$? $\beta = 0.5$?
- ► Should we adjust for covariate?

There is nothing in the data that tells us.

There is nothing in the data that tells us. ©

There is nothing in the data that tells us. \odot Here are the true structures:





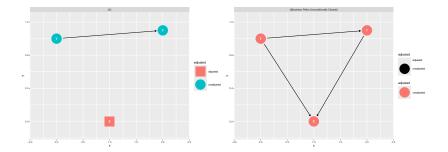
When know structures, adjustment sets for unbiasedness differ:

- ▶ df1: confounding \Rightarrow adjust for X
- ▶ df2: collider \Rightarrow do not adjust for X

```
g_conf <- dagitty("dag{ x -> y ; x <- c -> y }")
g_coll <- dagitty("dag{ x -> y ; x -> c <- y }")
adjustmentSets(g_conf, "x", "y")
{ c }
adjustmentSets(g_coll, "x", "y")</pre>
```

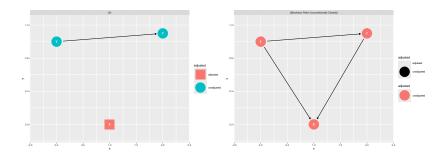
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(Data from D'Agostino McGowan (2023))

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- ► Importance of identifying "pre-treatment covariates", "proper covariates"; doing "design before analysis"
- ► Importance of experiments: strong knowledge about (part of) causal structure
- ➤ Causal inference is critical to scientific questions, and separate from prediction
- ➤ Though, methods from prediction can aid causal inference
- (A perspective on "causal euphimisms": Hernán (2018))

Approaches of Prediction and Causal Inference

Two Cultures, (Breiman 2001)

- ▶ Data Models: our "social science modeling"
- ▶ Algorithmic Models: our "data science algorithms"

Methods for Prediction and Causal Inference

- ► Cross-validation
- ▶ Regression/Decision trees
- ▶ Random forests

James et al. (2021)

Cross-validation

k-fold cross-validation

- \triangleright Randomly partition data into k groups
- \blacktriangleright Apply method to k-1 groups
- ▶ Use result to predict for left-out group
- ► Calculate $MSE_i = \frac{1}{n} \sum_{i=1}^{n} (y_i \hat{y}_i)^2$
- \triangleright Calculate test error as average of the k MSE's:

$$CV_{(k)} = \frac{1}{k} \sum_{i=1}^{k} MSE_i$$

 \blacktriangleright Select model that minimises $CV_{(k)}$

```
library(tidyverse)
## Make data
mk_data \leftarrow function(n = 90, n_folds = 10){
  df <- tibble(
    x1 = rnorm(n),
    x2 = rnorm(n),
    x3 = rnorm(n).
    y = 0.1 * x1 + 0.2 * x2 + 0.5 * x3 + rnorm(n),
    cv_fold = sample(rep(1:n_folds, (n / n_folds)))
df <- mk data()</pre>
```

head(df)

```
# A tibble: 6 x 5
    x1    x2    x3    y cv_fold
    <dbl>    <db
```

head(df)

```
# A tibble: 6 x 5
    x1    x2    x3    y cv_fold
    <dbl>    <dbl>    <dbl>    <dbl>    <dbl>    <dbl>    <int>
    1    1.81    0.827    -0.0417    -0.841    2
    2    0.336    -1.42    0.314    0.197    7
    3    0.372    0.602    -0.601    0.424    1
    4    0.648    1.27    -0.122    1.22    2
    5    -1.62    -0.670    1.25    0.158    4
    6    0.841    -0.318    -0.104    0.561    6
```

table(df\$cv_fold)

```
1 2 3 4 5 6 7 8 9 10
9 9 9 9 9 9 9 9 9 9
```

```
cv lm <- function(data, fmla){</pre>
 n folds <- max(data$cv fold)</pre>
  store_mses <- vector("numeric", length = n_folds)</pre>
  for(idx in 1:n folds){
    df_train <- data |> filter(cv_fold != idx)
    df_test <- data |> filter(cv_fold == idx)
    lm_out <- lm(fmla, data = df train)</pre>
    predictions <- predict(lm_out, newdata = df_test)</pre>
    store mses[idx] <- mean((df test$y - predictions)^2)}
  test_error_cv_k <- mean(store_mses)</pre>
  return(test error cv k)
```

```
cv_lm(data = df, fmla = y \sim x1 + x2)
```

[1] 1.323677

[1] 1.097002

```
cv_lm(data = df, fmla = y ~ x1 + x2)
[1] 1.323677

df <- mk_data()
cv_lm(df, y ~ x1 + x2)</pre>
```

[1] 0.9505413

```
cv lm(data = df, fmla = y \sim x1 + x2)
[1] 1.323677
df <- mk data()</pre>
cv lm(df, y \sim x1 + x2)
[1] 1.097002
df <- mk data()</pre>
cv lm(df, y \sim x1 + x2 + x3)
```

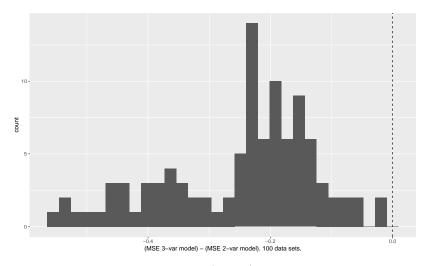


Figure 1: MSE always less (better) for 3-variable model.

- Partition predictor space into regions R_1, R_2, \dots, R_J .
- ▶ If unit falls in region R_j , use average outcome in R_j as predicted value: \hat{y}_{R_j}
- \blacktriangleright (For decision on discrete outcome, count votes in R_j)
- Goal: minimise residual sum of squares (RSS), just like LS regression:

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

How to define regions R_j ?

How to define regions R_i ?

- ➤ Top-down, greedy recursive binary split
- At each step, find predictor and cut-point that minimise

$$\sum_{i:x \in R_1(j,s)} \left(y_i - \hat{y}_{R_1(j,s)}\right)^2 + \sum_{i:x \in R_2(j,s)} \left(y_i - \hat{y}_{R_2(j,s)}\right)^2$$

▶ Overfitting is a potential problem

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- "Pruning"

Pruning

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Pruning

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Sum squared pred. error (plus penalty that grows with tree size) across units in region, then regions.

But, how to choose α ? (Use cross-validation.)

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- 4. Using that α , select best subtree from Step 2

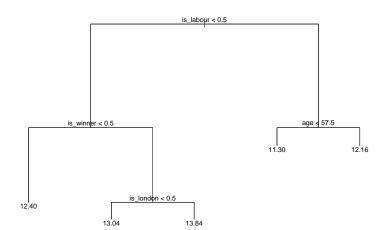
Example: Regression Tree library(qss) library(rsample) library(tree) data("MPs") mps <- MPs |> mutate(age = yod - yob, is labour = if else(party == "labour" is_london = if_else(region == "Greater is_winner = if_else(margin > 0, 1, 0)) select(ln.net, age, is_labour, is_london, is_winner) |> na.omit()

mp split <- initial split(mps, prop = 0.7)</pre>

mp_train <- training(mp_split)
mp test <- testing(mp split)</pre>

set.seed(765076184)

```
plot(tree_mp)
text(tree_mp)
```



Would pruning help?

```
cv_mps <- cv.tree(tree_mp, K = 10)
plot(cv_mps$size, cv_mps$dev, type = "b")</pre>
```

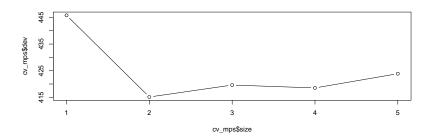


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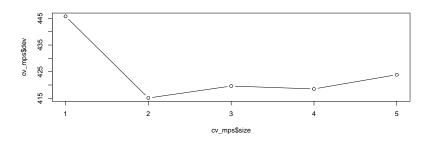


Figure 3: Subtree size 2 minimises SSR

```
prune_mps <- prune.tree(tree_mp, best = 2)

plot(prune_mps)
text(prune_mps)</pre>
```



Figure 4: The pruned tree

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► MSE for pruned: 1.922

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(Pretty good for 1 split!?)

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Bagging: bootstrap aggregation

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- Linear regression: lower variance)

Random forests: decorrelated, bagged trees

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- ightharpoonup (Often choose $m \approx \sqrt{p}$)
- So, different splits consider different predictors
- So, trees will look very different to each other

```
library(randomForest)
# Full bag:
bag mps <- randomForest(ln.net ~ ., data = mp train,</pre>
                          ntree = 500, mtry = 4,
                          importance = TRUE)
# Decorrelate:
rf mps <- randomForest(ln.net ~ ., data = mp train,</pre>
                          ntree = 500, mtry = 2,
                         importance = TRUE)
```

Predict:

```
preds_bag <- predict(bag_mps, newdata = mp_test)
preds_rf <- predict(rf_mps, newdata = mp_test)</pre>
```

- MSE for RF: 1.995
- ▶ MSE for full bag: 2.536

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Treatment + ϵ

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```
lm_out <- lm(ln.net ~ is_winner, data = mps)
lm_out</pre>
```

```
Call:
lm(formula = ln.net ~ is_winner, data = mps)
Coefficients:
(Intercept) is_winner
    12.2464    0.5176
```

Homogeneous effects:

```
t.test(ln.net ~ is_winner, data = mps)
```

Welch Two Sample t-test

```
data: ln.net by is_winner

t = -3.9552, df = 287.65, p-value = 9.636e-05

alternative hypothesis: true difference in means between the second second
```

Homogeneous and Heterogeneous Effects: Estimation Homogeneous effects:

$$\text{Outcome} = \beta_0 + \beta_1 \text{Treatment} + \sum \beta_j X_j + \epsilon$$

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Outcome =
$$\beta_0 + \beta_1$$
Treatment + $\sum \beta_j X_j + \epsilon$

lm(formula = ln.net ~ is_winner + is_labour + is_london + a
 data = mps)

Coefficients:

Homogeneous effects:

lm lin(ln.net ~ is winner, covariates = ~ is labour + is lo

	Estimate	Std. Error	t valı
(Intercept)	1.226687e+01	0.078894901	155.4836617
is_winner	3.459885e-01	0.131207672	2.6369536
is_labour_c	-1.613663e-01	0.152608515	-1.0573873
is_london_c	2.427360e-01	0.250214401	0.9701118
age_c	4.740367e-03	0.007031323	0.6741786
<pre>is_winner:is_labour_c</pre>	-9.104022e-01	0.264395760	-3.4433313
<pre>is_winner:is_london_c</pre>	-8.847770e-02	0.426241818	-0.2075763
is_winner:age_c	-4.778657e-05	0.012753800	-0.0037468
	CI Lower	CI Upper	DF
(Intercept)	12.111785723	12.42195044	416

0.088075873 0.60390123 416

-0.461346226 0.13861367 416

-0.249106208 0.73457813 416

is_winner

is_labour_c

is london c



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- Inference: not "evidence against TE = 0?", but "evidence against $CATE_1 = CATE_2$?"

Heterogeneous effects:

 $\label{eq:outcome} \text{Outcome} = \beta_0 + \beta_1 \text{Treatment} + \beta_2 \text{Group} + \beta_3 \text{Treatment} \cdot \text{Group} + \epsilon$

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Heterogeneous effects:

 $Outcome = \beta_0 + \beta_1 Treatment + \beta_2 Group + \beta_3 Treatment \cdot Group + \epsilon$

- \triangleright β_1 gives TE for Group == 0
- \triangleright $\beta_1 + \beta_3$ gives TE for Group == 1

Heterogeneous effects:

```
(Intercept) is_winner is_labour
11.959 0.780 -0.165
age is_winner:is_labour
0.005 -0.914
```

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$$Y(0), Y(1) \perp \!\!\!\perp T \mid \mathbf{X}$$

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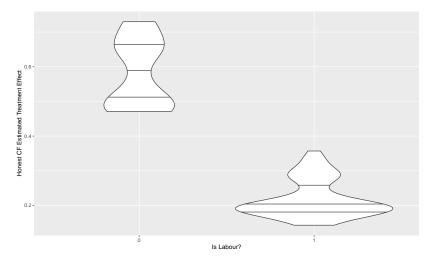
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- \blacktriangleright Build a random forest (decorrelated deep trees picking from m predictors) of causal trees

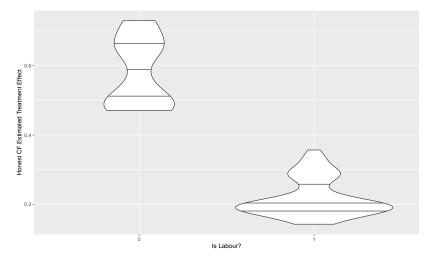
Example: Causal Forests

```
library(grf)
X <- mp_train |> select(age, is_labour, is_london)
W <- mp_train |> select(is_winner) |>
  unlist() |> as.numeric()
Y <- mp_train |> select(ln.net) |> unlist()
cf out <- causal forest(X, Y, W)
```

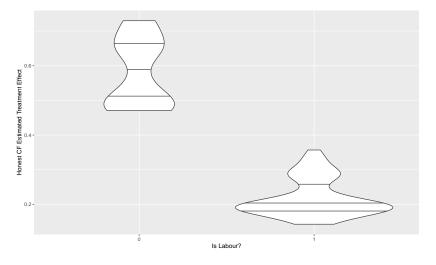
Example: Causal Forests

```
X test <- mp test |> select(age, is labour, is london)
cf pred est var <- predict(cf out, X test,
                           estimate.variance = TRUE)
cf preds <- cf pred est var$predictions
df cf <- tibble(X test,
                cf te = cf preds,
                cf se = sqrt(cf pred est var$variance.
                te 1se lower = cf te - cf se,
                te 1se upper = cf te + cf se)
```

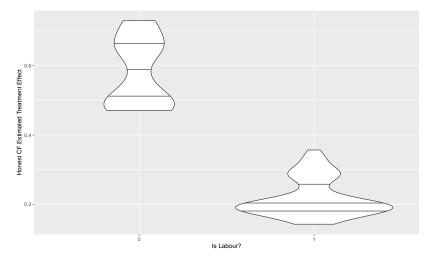




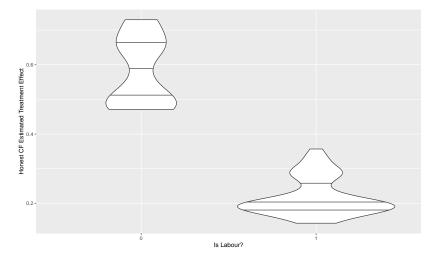
▶ Mean CF TE, Tory: 0.58



▶ Mean CF TE, Tory: $0.58 \rightsquigarrow £192,000$

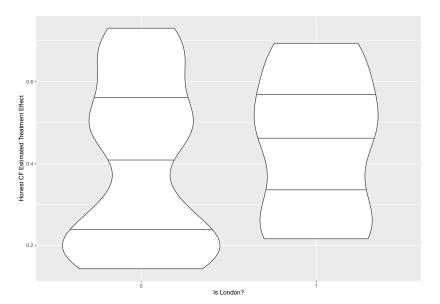


- ► Mean CF TE, Tory: 0.58 → £192,000
- ▶ Mean CF TE, Labour: 0.219

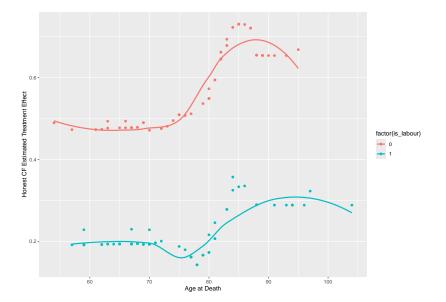


- Mean CF TE, Tory: $0.58 \rightsquigarrow £192,000$
- Mean CF TE, Labour: $0.219 \rightsquigarrow £60,000$

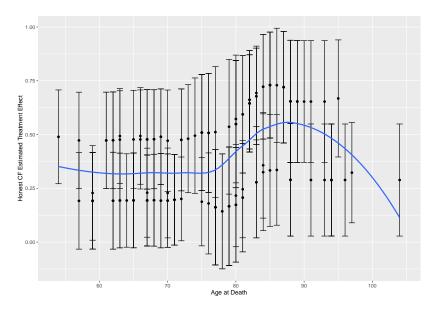
Example: Causal Forests Results, London



Example: Causal Forests Results, Age



Example: Causal Forests Results, Age





Feature Selection

▶ Wrappers: pick subset of covars, train on data (estimate model), test on hold-out, score predictions. Keep best-scoring subset.

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- ▶ Wrappers: pick subset of covars, train on data (estimate model), test on hold-out, score predictions. Keep best-scoring subset.
- Filters: correlate covars with outcome. Keep strongest.
- ▶ Embeds: select features and estimate model at same time. Penalize using more predictors.

Regularization Methods

OLS reminder

Minimize SSR:

$$\begin{aligned} & \operatorname{argmin}_{\beta} \sum_{i=1}^{n} \left(y_{i} - \hat{y}_{i} \right)^{2} \\ & \operatorname{argmin}_{\beta} \sum_{i=1}^{n} \left(\mathbf{y} - \mathbf{X} \hat{\beta} \right)^{2} \end{aligned}$$

L1 regularization: the LASSO (Least Absolute Shrinkage and Selection Operator)

$$\operatorname{argmin}_{\beta} \left[\sum_{i=1}^{n} \left(y_i - \mathbf{X} \hat{\beta} \right)^2 + \lambda \sum_{j=1}^{k} |\hat{\beta}_j| \right]$$

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L2 regularization: Ridge regression

$$\operatorname{argmin}_{\beta} \left[\sum_{i=1}^{n} \left(y_i - \mathbf{X} \hat{\beta} \right)^2 + \lambda \sum_{j=1}^{k} \hat{\beta}_j^2 \right]$$

Mix L1 and L2: Elastic net

$$\operatorname{argmin}_{\beta} \left(\frac{\sum\limits_{i=1}^{n} \left(y_i - \mathbf{X} \hat{\beta} \right)^2}{2n} + \lambda \left[\alpha \sum\limits_{j=1}^{k} |\hat{\beta}_j| + \frac{1-\alpha}{2} \sum\limits_{j=1}^{k} \hat{\beta}_j^2 \right] \right)$$

Mix L1 and L2: Elastic net

$$\operatorname{argmin}_{\beta} \left(\frac{\sum\limits_{i=1}^{n} \left(y_i - \mathbf{X} \hat{\beta} \right)^2}{2n} + \lambda \left[\alpha \sum\limits_{j=1}^{k} |\hat{\beta}_j| + \frac{1-\alpha}{2} \sum\limits_{j=1}^{k} \hat{\beta}_j^2 \right] \right)$$

Regularized trees, ...

How to choose λ , α ?

How to choose λ , α ?

Cross-validation for λ :

```
df_lasso <- read_csv("~/Desktop/lasso.csv")

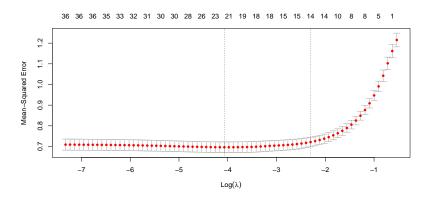
X <- as.matrix(df_lasso[, 2:ncol(df_lasso)])

Y <- as.matrix(df_lasso[, "y"])

library(glmnet)

cv_lasso <- cv.glmnet(X, Y, alpha = 1)</pre>
```

plot(cv_lasso)



cv_lasso\$lambda.min

[1] 0.0170891

Implement:

```
Call: glmnet(x = X, y = Y, alpha = 1, lambda = cv_lasso$16
```

Df %Dev Lambda 1 21 45.32 0.01709

Coefficients:

```
coef_lasso <- coef(lasso_out)</pre>
round(coef_lasso, 3)
37 x 1 sparse Matrix of class "dgCMatrix"
                 s0
(Intercept)
             0.000
x1
              0.112
              0.095
x2
xЗ
              0.086
x4
              0.147
x5
              0.002
              0.063
x6
x7
              0.051
8x
              0.074
x9
              0.042
x10
```

Coefficients:

(Intercept)

round(coef_lasso[,], 3)

x18

x24

x30

x36 0.048

0.010

0.000

0.000

0.147	0.086	0.095	0.112	0.000
x10	x9	x8	x7	x6
0.000	0.042	0.074	0.051	0.063
x16	x15	x14	x13	x12
0.000	0.000	0.026	0.000	0.039

x2

x20

x26

x32

-0.015

0.000

0.032

xЗ

x21

x27

x33

0.030

0.000

0.000

x2:

x28

x34

0.000

-0.010

-0.04

x1

x19

x25

x31

0.127

0.000

0.028

x1

x2

x3

x4

x5 x6

x7

8x

Implement, alternative λ :

0.1051188782

0.0898842045

0.0742522801

0.1513883536

0.0603811184

0.0389489143 0.0575738993

Coefficients:

0.029

0.004

0.000

0.000

x18

x24

x30

x36 0.030

round(coef(lasso	_1se)[,],	3)

	XS	XZ	XI	(Incercebe)
0.	0.074	0.090	0.105	0.000
	x9	x8	x7	х6
0.0	0.037	0.058	0.039	0.060
:	x15	x14	x13	x12

0.008

0.000

0.000

0.013

x20

x26

x32

0.000

0.000

0.000

0.000

x21

x27

x33

0.000

0.039

0.000

0.000

x19

x25

x31

15: x10 000 x10

0.000

0.000

0.000

0.000

x2:

x28

x34

The idea:

 \triangleright covariates may $\rightsquigarrow Y$ or $\rightsquigarrow T$

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- \triangleright covariates may $\rightsquigarrow Y$ or $\rightsquigarrow T$
- $\triangleright \approx$ "double robust", "AIPW" estimators
- (different to just "doing LASSO twice" for regularization + shrinkage)

1. Model Y = f(X) using LASSO

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rlasso out <- rlassoATE(primary2006 ~ age + is male + primary2006 ~ age + is male + primary2006

```
summary(rlasso_out)
```

```
Estimation and significance testing of the treatment effective. Type: ATE

Bootstrap: not applicable
    coeff. se. t-value p-value

TE 0.080091 0.002625 30.51 <2e-16 ***
---

Signif. codes: 0 '***' 0.001 '**' 0.05 '.' 0.1 '
```

```
X <- as.matrix(df_social[, c("age", "is_male", "primary2004
Y <- as.matrix(df_social[, "primary2006"])
D <- as.matrix(df_social[, "is_neighbors"])</pre>
summary(rlassoEffects(X, Y, method = "double selection"))
[1] "Estimates and significance testing of the effect of ta
            Estimate. Std. Error t value Pr(>|t|)
     0.0038449 0.0000681 56.456 < 2e-16 ***
age
is male 0.0086763 0.0018889 4.593 4.36e-06 ***
primary2004 0.1474364 0.0019924 74.000 < 2e-16 ***
hhsize 0.0004260 0.0012618 0.338 0.736
is_neighbors 0.0802361 0.0026278 30.534 < 2e-16 ***
Signif. codes: 0 '***' 0.001 '**' 0.01 '*' 0.05 '.' 0.1 '
```

R packages for Regularization, etc.

- ▶ glmnet
- caret

See also tidymodels, parsnip, \dots

${\bf Embedded} \ {\bf Regularization} \ {\bf Methods}$



Thanks!

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References I

- Breiman, Leo. 2001. "Statistical Modeling: The Two Cultures." *Statistical Science* 16 (3): 199–215. http://www.jstor.org/stable/2676681.
- D'Agostino McGowan, Lucy. 2023. quartets: Datasets to Help Teach Statistics. https://r-causal.github.io/quartets/.
- Hernán, Miguel A. 2018. "The c-Word: Scientific Euphemisms Do Not Improve Causal Inference from Observational Data." *American Journal of Public Health* 108 (5): 616–19.
- Hernán, Miguel A., John Hsu, and Brian Healy. 2019. "A Second Chance to Get Causal Inference Right: A Classification of Data Science Tasks." CHANCE 32 (1): 42–49. https://doi.org/10.1080/09332480.2019.1579578.
- James, Gareth, Daniela Witten, Trevor Hastie, and Robert Tibshirani. 2021. An Introduction to Statistical Learning with Applications in R. 2nd ed. New York, NY: Springer.