An introduction to `grf`

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This vignette introduces forest-based causal inference using the `grf` package for `R`. Generalized random forests (Athey, Tibshirani, and Wager, 2019) extend Breiman (2001)'s random forest to enable data-driven estimation of heterogeneity in causal parameters, such as conditional average treatment effects under unconfoundedness. We outline core functionality, including doubly robust estimation of average treatment effects, best linear projections, tools for assessing heterogeneity using TOC curves, evaluation of treatment rules via Qini curves, and methods for learning interpretable tree-based policies. The vignette contains two applied examples illustrating treatment effect heterogeneity in practice. For further details, see the package documentation and additional applied tutorials by Athey & Wager (2019), and Sverdrup, Petukhova, and Wager (2024).

Vignette sections:

1) *A grf guided tour*
2) *Example: Treatment heterogeneity and financial proficiency*
3) *Example: The effect of poverty on attention*

A grf guided tour

We describe how modern machine learning can flexibly control for confounding when estimating average treatment effects, how Breiman's (2001) random forest can be adapted to target treatment effect heterogeneity, and how this approach extends to a broad class of settings, including instrumental variables and right-censored outcomes. Further info is provided in Athey & Wager (2019), Cui et al. (2023), Dandl et al. (2024), Friedberg et al. (2020), and Sverdrup et al. (2024). For a more detailed theoretical treatment, Wager (2024) provides a textbook-style discussion.

Machine learning for causal inference

Machine learning methods are designed for prediction. In causal applications, however, our goal is often inference rather than prediction. Although statistical *estimation* and statistical *prediction* are distinct tasks (Efron, 2020), predictive tools can support valid inference when used appropriately. This section outlines how modern machine learning, combined with ideas from semi-parametric statistics, can provide flexible and model-agnostic tools for causal inference.

Suppose we observe outcomes Y_i, a binary treatment indicator $W_i \in 0, 1$, and covariates X_i. Our goal is to estimate the average treatment effect $\tau = \mathrm{E}[Y_i(1) - Y_i(0)]$. In an observational study, we might run a

regression

$$Y_i \sim \tau W_i + \beta X_i,$$

and interpret $\hat{\tau}$ as an estimate of the ATE. This interpretation relies on:

1) W being unconfounded given X (e.g., Rosenbaum and Rubin, 1983).
2) The confounders X having a linear effect on Y.
3) The treatment effect τ being constant.

Assumption 1) is an *identifying* assumption and something we choose to live with depending on domain knowledge. Assumptions 2) and 3) are *modeling* choices that we may wish to relax.

Relaxing assumption 2). Instead of assuming linearity, consider the partially linear model

$$Y_i = \tau W_i + f(X_i) + \varepsilon_i, \quad \mathrm{E}[\varepsilon_i \mid X_i, W_i] = 0,$$

where f is an unknown function. The challenge is that $f(X_i)$ is not specified.

Robinson (1988) shows that via the following intermediary objects, the *propensity score*

$$e(x) = \mathrm{E}[W_i \mid X_i = x],$$

and the *conditional mean*,

$$m(x) = \mathrm{E}[Y_i \mid X_i = x] = f(x) + \tau e(x),$$

we can rewrite the model in centered form:

$$Y_i - m(x) = \tau\Big(W_i - e(x)\Big) + \varepsilon_i.$$

This means that τ can be estimated by regressing the residualized outcome on the residualized treatment. Importantly, Robinson (1988) showed that this estimator is root-n consistent even if $m(x)$ and $e(x)$ are estimated more slowly. This robustness is often called orthogonality: small errors in *nuisance* estimates do not substantially bias the target parameter.

How do we flexibly estimate $m(x)$ and $e(x)$? These are standard regression problems: predict Y_i from X_i, and W_i from X_i. Modern machine learning methods, such as boosting or random forests, are well suited for this task (James et al., 2013). A wrinkle is that directly plugging machine learning estimates into this estimator can lead to bias due to regularization. A standard technique in semi-parametric estimation is sample splitting, where separate data subsets are used to train and evaluate these components. Cross-fitting, alternating which folds are used for training and estimation, is a common implementation (Chernozhukov et al., 2018; Zheng & van der Laan, 2011). In `grf`, *out-of-bag* predictions provide a forest-native leave-one-out form of cross-fitting.

Non-constant treatment effects. To relax Assumption 3), suppose treatment effects vary with covariates:

$$Y_i = \tau(X_i) W_i + f(X_i) + \varepsilon_i, \ \mathrm{E}[\varepsilon_i \mid X_i, W_i] = 0,$$

where $\tau(x)$ is the conditional average treatment effect $\mathrm{E}[Y_i(1) - Y_i(0) \mid X_i = x]$. If we had access to a neighborhood $\mathcal{N}(x)$ where τ were approximately constant, we could proceed exactly as before, by performing a residual-on-residual regression on the samples belonging to $\mathcal{N}(x)$, i.e.:

$$\tau(x) := \mathrm{lm}\left(Y_i - \hat{m}^{(-i)}(X_i) \sim W_i - \hat{e}^{(-i)}(X_i), \mathrm{weights} = 1\{X_i \in \mathcal{N}(x)\}\right),$$

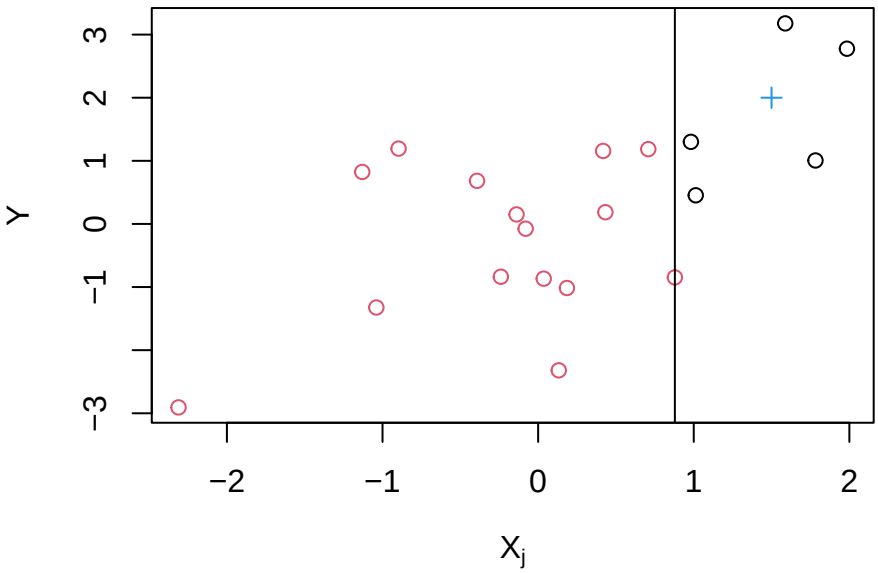
where $^{(-i)}$ denotes out-of-bag estimates. Conceptually, causal forest estimates $\tau(x)$ by running a locally weighted residual-on-residual regression using samples predicted to have similar treatment effects.

The key question is how to construct these weights. The next section reviews Breiman's random forest and explains how it serves as an adaptive neighborhood finder.

Random forest as an adaptive neighborhood finder

Breiman's random forest estimates the conditional mean $\mu(x) = \mathrm{E}[Y_i \mid X_i = x]$ in two phases:

1) **Building phase**: Grow B trees by greedily choosing covariate splits that maximize the squared difference in subgroup means $n_L n_R (\bar{y}_L - \bar{y}_R)^2$.
2) **Prediction phase**: For a target sample x, average the outcomes Y_i in the same terminal leaf $L_b(x)$ across trees: $\hat{\mu}(x) = \frac{1}{B} \sum_{b=1}^{B} \sum_{i=1}^{n} Y_i \frac{1\{Xi \in L_b(x)\}}{|L_b(x)|}$.



For a single tree and split, Figure 1 shows the chosen split point: the vertical line that maximizes the difference in leaf means. A target point x falls into one leaf, and prediction averages outcomes in that leaf. Repeating over trees and averaging yields the final estimate.

Rewriting the prediction as

$$\begin{aligned}\hat{\mu}(x) &= \sum_{i=1}^{n} \frac{1}{B} \sum_{b=1}^{B} Y_i \frac{1\{X_i \in L_b(x)\}}{|L_b(x)|} \\ &= \sum_{i=1}^{n} Y_i \alpha_i(x),\end{aligned}$$

shows that a random forest prediction is a weighted average of outcomes. The weights $\alpha_i(x)$ reflect how often training sample i shares a leaf with x. They can be interpreted as adaptive similarity weights. This weighting view of random forests appears frequently in statistical applications, see for example Zeileis, Hothorn & Hornik (2003).

Causal forests combines this forest weighting with Robinson's approach:

1) **Building phase**: Choose splits that maximize the squared difference in estimated subgroup treatment effects, $n_L n_R (\hat{\tau}_L - \hat{\tau}_R)^2$, where $\hat{\tau}$ is computed via residual-on-residual regression.
2) **Prediction phase**: Use the resulting forest weights $\alpha(x)$ to estimate

$$\tau(x) := \mathrm{lm}\left(Y_i - \hat{m}^{(-i)}(X_i) \sim W_i - \hat{e}^{(-i)}(X_i),\ \text{weights} = \alpha_i(x) \right).$$

Thus, causal forest performs a forest-localized version of Robinson's regression. In `causal_forest`, the arguments `Y.hat` and `W.hat` provide estimates of $m(x)$ and $e(x)$, by default obtained from separate regression

forests. If we are in a randomized setting where we know the propensity score, we can simply supply these via the `W.hat` argument. Robinson's transformation also motivates a general loss function for heterogeneous treatment effects (Nie & Wager, 2021), which `causal_forest` uses to guide the selection of tuning parameters.

The same idea extends beyond treatment effects and residual-on-residual regressions. In the generalized random forest framework, the forest is tailored to target heterogeneity in any parameter defined by an estimating equation $\psi_\theta(\cdot)$. Given an estimating equation that identifies a parameter θ without covariates, `grf` then constructs weights $\alpha(x)$ that yield covariate-specific estimates $\theta(x)$, using Breiman's random forest as a blueprint. For causal forests, this estimating equation corresponds to Robinson's regression.

Orthogonal estimating equations offers practical advantages by separating the statistical tasks. Other forests, such as `instrumental_forest` and `causal_survival_forest`, use similar constructions. For example, `causal_survival_forest` fits two separate survival forests to construct right-censoring adjustments. These are then used to construct an orthogonal estimating equation, and `grf` trains a final random forest targeting heterogeneity in the treatment effect parameter.

For computational and implementation reasons, `grf` does not re-estimate the target parameter at every candidate split in the *building phase*. Instead, it uses an influence-function-based approximation computed once in the parent node. `grf` then reuses the standard random forest machinery (Breiman-style tree growth and regression-based splitting) but replaces the raw outcomes with pseudo-outcomes derived from the orthogonal score. From an implementation perspective, this means that new forest estimators can be built by specifying the appropriate score and pseudo-outcome, while leaving the underlying random forest algorithm unchanged.

> **Note**: The target parameter θ may be vector-valued. In that case, splits are based on the squared ℓ_2 distance $n_L n_R \|\hat{\theta}_L - \hat{\theta}_R\|^2$, as in `grf`'s `multi_arm_causal_forest` and `lm_forest`.

Efficiently estimating summaries of the CATEs

Robinson's residual-on-residual approach is orthogonal, which makes it well suited for estimates that converge slowly. But what if we want summaries of $\tau(x)$, such as the ATE or a best linear projection (BLP), that have guaranteed root-n rate of convergence along with exact confidence intervals?

Simply averaging the estimates $\hat{\tau}(X_i)$ is not optimal: prediction errors are too large and do not cancel out. Robins, Rotnitzky & Zhao (1994) proposed the Augmented Inverse Probability Weighted (AIPW) estimator, which combines propensity scores and conditional mean estimates in a way that is robust to estimation errors in either component. It can be expressed as a one-step debiasing correction, taking the following form

$$\hat{\Gamma}_i = \hat{\tau}^{(-i)}(X_i) + \frac{W_i - \hat{e}^{(-i)}(X_i)}{\hat{e}^{(-i)}(X_i)[1 - \hat{e}^{(-i)}(X_i)]} \Big(Y_i - \hat{\mu}^{(-i)}(X_i, W_i)\Big),$$

where $\mu(x, w) = E[Y_i \mid X_i = x, W_i = w]$ (this can be backed out from a causal forest fit). The forest-based ATE is then given by

$$\hat{\tau}_{\text{AIPW}} = \frac{1}{n} \sum_{i=1}^{n} \hat{\Gamma}_i.$$

This construction yields **doubly robust** estimating equations and forms the basis of summary methods in `grf`, including `average_treatment_effect` and `best_linear_projection`. The associated doubly robust scores $\hat{\Gamma}_i$ are accessible through the function `get_scores`. As a concrete example, the ATE is an average of $\hat{\Gamma}_i$'s, while the best linear projection is a regression of the $\hat{\Gamma}_i$'s on a set of covariates X_i.

Assessing heterogeneity with TOC curves

Combined with techniques such as honesty and subsampling (Wager & Athey, 2018), the generalized random forest algorithm supports pointwise asymptotic guarantees. Results like these are reassuring, as they characterize the large-sample behavior of the forest. For practical purposes, however, there are more direct and transparent ways to assess whether meaningful heterogeneity is present in a given dataset.

A simple and pragmatic approach is **train/test evaluation**. We fit a CATE model on a training set, then on a held-out test set, we form subgroups based on predicted treatment effects and examine whether average treatment effects differ systematically across these groups. This approach is quite general: we may evaluate any machine-learned targeting function, including alternative effect measures, such as predicted risk or other quantities that may be aligned with treatment benefit in a given application. As long as higher values correspond to higher expected treatment gains, the same evaluation logic applies.

The `rank_average_treatment_effect` function provides an automated implementation of this exercise by stratifying the test sample according to quantiles of the estimated targeting function. The *Targeting Operator Characteristic* (TOC) curve (Yadlowsky et al., 2024) compares the average treatment effect among units with predicted CATEs in the top q-th quantile to the overall ATE.

$$\text{TOC}(q) = \text{E}[Y_i(1) - Y_i(0) \mid \text{Estimated CATE}(X_i) \geq (1-q)\text{-th quantile}] - \text{ATE}.$$

Given a trained evaluation forest, `grf` provides doubly robust estimates of the TOC and the area under the TOC, along with standard errors that enable a simple and valid t-test for the presence of heterogeneity. Under the null of constant treatment effects, the area under the TOC curve is zero in expectation.

Policy evaluation with Qini curves

We are often interested in data-driven rules to guide treatment allocation decisions, also referred to as *policies*. Under a binary treatment, a policy is a rule $\pi(X_i) \in \{0, 1\}$ that maps covariates to a treatment decision. The value of a policy is the expected gain from following it,

$$G(\pi) = \text{E}[\pi(X_i)\{Y_i(1) - Y_i(0)\}].$$

Given doubly robust scores $\hat{\Gamma}_i$, we can estimate the value of any fixed policy by averaging $\hat{\Gamma}_i$ over the sample for which $\pi(X_i) = 1$. An estimated CATE function $\hat{\tau}(\cdot)$ naturally induces targeting rules. A simple example is the threshold policy

$$\pi_0(X_i) = \begin{cases} 1, & \hat{\tau}(X_i) \geq 0 \\ 0, & \text{otherwise.} \end{cases}$$

If the true CATE were known, the optimal policy would treat units with $\tau(X_i) \geq 0$. Since we only have access to estimates $\hat{\tau}(\cdot)$, we instead evaluate how well such data-driven policies perform by estimating their value $G(\pi_0)$.

Often, treatment is costly or limited, and we may only treat a fraction of the population. For example, we might target the top 20% of units ranked by predicted effect by thresholding at the 80-th percentile. Varying this threshold produces a family of policies indexed by the fraction treated budget $B \in (0, 1]$. The *Qini curve* summarizes their performance by plotting

$$G(\pi_B) = \text{E}[\pi_B(X_i)\{Y_i(1) - Y_i(0)\}],$$

where

$$\pi_B(X_i) = \begin{cases} 1, & \hat{\tau}(X_i) > 0 \text{ and } \hat{\tau}(X_i) \geq 1 - B\text{-th quantile of positive predicted CATEs} \\ 0, & \text{otherwise.} \end{cases}$$

Using doubly robust scores and appropriate resampling methods, we can obtain confidence intervals for $G(\pi_B)$ across the curve. This approach is implemented in the sister package `maq`, which also supports multiple treatment arms with variable costs by solving an intermediate linear program that maps predicted effects and costs into budget-constrained policies.

Learning interpretable policies

In some applications, directly assigning treatment based on complex, black-box CATE predictions may be undesirable. Decision makers often require transparent and interpretable decision rules that assign treatment

based on a small number of observable characteristics. These policies are easy to communicate and implement (e.g., via a few conditional statements in a database query), while still leveraging heterogeneity in treatment effects.

Given experimental or observational data, we can learn such interpretable policies by maximizing the policy value

$$V(\pi) = \mathrm{E}[Y_i(\pi(X_i))],$$

where we restrict the policy π to belong to a class of pre-specified rules, for example shallow decision trees. Because we never observe both potential outcomes, we need suitable reward proxies. It turns out (Athey & Wager, 2021) that we can solve the empirical counterpart of this problem efficiently by using doubly robust scores for $\mathrm{E}[Y_i(1)]$ and $\mathrm{E}[Y_i(0)]$ as reward proxies. The `grf` package constructs these scores automatically, and the sister package `policytree` solves the optimization problem for shallow depth-k decision trees. This yields simple treatment rules that can be evaluated on held-out data and compared to natural benchmarks such as treating everyone or no one.

Example: Treatment heterogeneity and financial proficiency

As a simple empirical illustration, we use the `causal_forest` function to analyze a randomized controlled trial measuring the effect of financial education. Bruhn et al. (2016) randomly assigned Brazilian high schools to a financial education program and measured outcomes such as financial proficiency. Randomization at the school level avoids interference between students. We use a processed copy of the data included in `grf`, containing student-level information for around 17,000 students at 850 high schools. The outcome is post-treatment financial literacy, and there are 13 pre-treatment covariates, including survey responses and two indices measuring students' ability to save and their financial autonomy.

Between 4% and 10% of the covariates contain missing values (`NA`). We assume missingness is random and retain them, as `grf` handles missing data via "missingness incorporated in attributes" (MIA) splitting (Mayer et al., 2020; Twala et al., 2008).

```
data("schoolrct")
Y <- schoolrct$outcome
W <- schoolrct$treatment
school <- schoolrct$school
X <- schoolrct[-(1:3)]
```

Here, we thought it would be interesting to define the CATE in terms of the following pre-treatment covariates

```
colnames(X)
#>  [1] "is.female"
#>  [2] "mother.attended.secondary.school"
#>  [3] "father.attened.secondary.school"
#>  [4] "failed.at.least.one.school.year"
#>  [5] "family.receives.cash.transfer"
#>  [6] "has.computer.with.internet.at.home"
#>  [7] "is.unemployed"
#>  [8] "has.some.form.of.income"
#>  [9] "saves.money.for.future.purchases"
#> [10] "intention.to.save.index"
#> [11] "makes.list.of.expenses.every.month"
#> [12] "negotiates.prices.or.payment.methods"
#> [13] "financial.autonomy.index"
```

Next, we fit a causal forest using default settings. Since this is an RCT, we pass the trial randomization probabilities via `W.hat = 0.5`. We also pass `clusters = school` to account for clustering at the school level in variance estimates.

Remark: On large datasets, setting `options(grf.verbose = TRUE)` can display progress during training and prediction.

Remark: We use `num.threads = 2` here to limit CPU usage. The default (`NULL`) uses all available cores.

```
c.forest <- causal_forest(X, Y, W, W.hat = 0.5, clusters = school, num.threads = 2)
```

Computing the ATE:

```
ate <- average_treatment_effect(c.forest)
ate
#> estimate  std.err
#>     4.35     0.52
ate[1] / sd(Y) # effect in SD units
#> estimate
#>      0.3
```

The effect of the program is positive and significant, increasing proficiency by roughly a quarter of a standard deviation.

Next, we summarize $\tau(X_i)$ by estimating a best linear projection on selected covariates:

$$\tau(X_i) \sim a + \beta_1 \text{is.female}_i + \beta_2 \text{family.receives.cash.transfer}_i + \beta_3 \text{is.unemployed}_i + \beta_4 \text{financial.autonomy.index}_i.$$

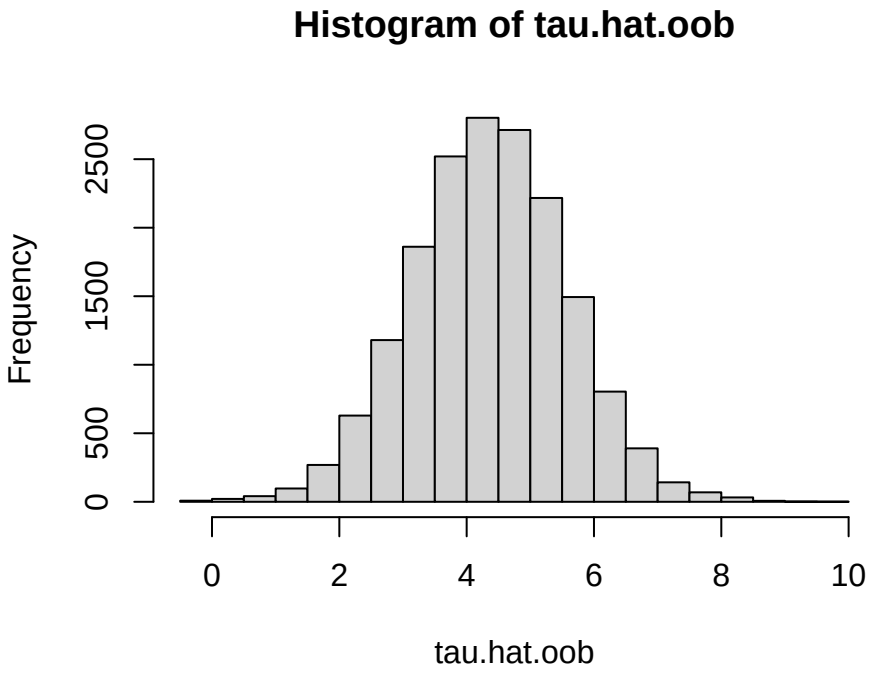
We get a doubly robust estimate of the coefficients in the model via

```
best_linear_projection(c.forest, X[c("is.female",
                                     "family.receives.cash.transfer",
                                     "is.unemployed",
                                     "financial.autonomy.index")])
#>
#> Best linear projection of the conditional average treatment effect.
#> Confidence intervals are cluster- and heteroskedasticity-robust (HC3):
#>
#>                               Estimate Std. Error t value Pr(>|t|)
#> (Intercept)                     6.7392     0.9777    6.89  5.7e-12 ***
#> is.female                      -0.7268     0.5307   -1.37   0.1709
#> family.receives.cash.transfer  -1.1992     0.6526   -1.84   0.0661 .
#> is.unemployed                   0.3206     0.6277    0.51   0.6095
#> financial.autonomy.index       -0.0329     0.0128   -2.58   0.0099 **
#> ---
#> Signif. codes:  0 '***' 0.001 '**' 0.01 '*' 0.05 '.' 0.1 ' ' 1
```

We see that students with higher financial autonomy, on average, benefit less from the training program.

Next, we estimate out-of-bag CATE predictions along all 13 covariates:

```
tau.hat.oob <- predict(c.forest)$predictions
hist(tau.hat.oob)
```



The histogram shows most students have effects near the overall ATE, with some variation. To assess whether these capture meaningful heterogeneity, we estimate the area under the TOC:

```
rate.oob <- rank_average_treatment_effect(c.forest, tau.hat.oob)
t.stat.oob <- rate.oob$estimate / rate.oob$std.err
pnorm(t.stat.oob, lower.tail = FALSE)  # one-sided p-value Pr(>t)
#> [1] 0.65
```

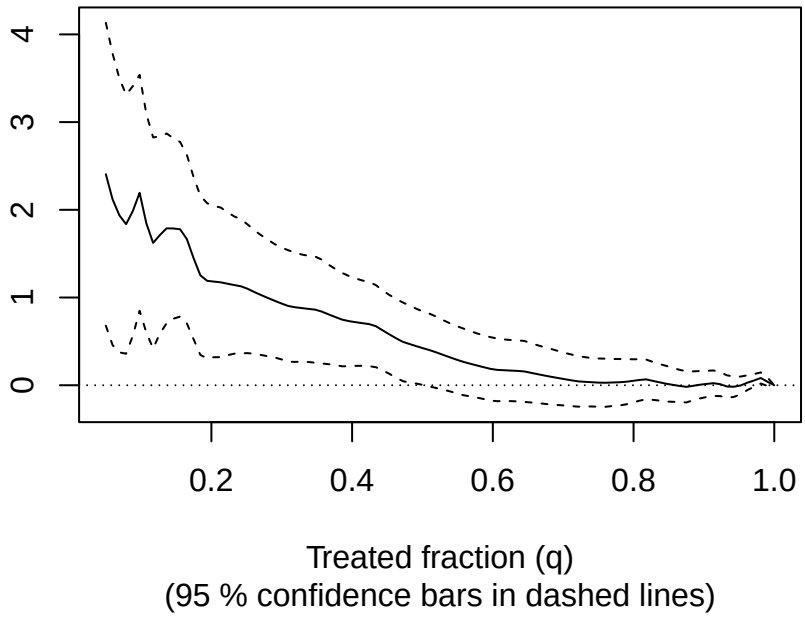
Here we use out-of-bag predictions both to construct $\hat{\tau}(\cdot)$ and to evaluate the TOC. Since the same data are reused, exact finite-sample guarantees for the resulting t-statistic are not available. Instead of performing an explicit train/test split, we adopt a pragmatic heuristic and report a one-sided test in the direction of interest (AUTOC > 0). In this example, the one-sided test rejects the null of no heterogeneity. This should be interpreted as suggestive evidence rather than a formally validated inference procedure based on sample splitting.

However, the TOC remains useful for evaluating alternative prioritization strategies. For example, students with low financial autonomy may benefit more. The causal forest variable importance confirms `financial.autonomy.index` is a key predictor:

```
var.imp <- variable_importance(c.forest)
head(colnames(X)[order(var.imp, decreasing = TRUE)])
#> [1] "financial.autonomy.index"          "intention.to.save.index"
#> [3] "family.receives.cash.transfer"     "father.attened.secondary.school"
#> [5] "is.female"                         "has.computer.with.internet.at.home"
```

We can construct a TOC prioritizing students with lower financial autonomy:

```
rate.autonomy <- rank_average_treatment_effect(
  c.forest,
  -1 * X$financial.autonomy.index, # Negate priorities to rank by lowest first.
  q = seq(0.05, 1, length.out = 100),
  subset = !is.na(X$financial.autonomy.index))
plot(rate.autonomy, main = "TOC: By increasing financial autonomy")
```



This reveals that students in the lower 20% of financial autonomy have an ATE roughly 1.1 points above the overall average, with a confidence interval excluding zero.

Training programs like these typically require resources to implement and scale to many students. The analysis here suggests that targeting students with low financial autonomy may be a desirable strategy. If we estimate the area under this TOC we get a t-value above 1.96. We would reject a null at conventional levels, leading us to conclude this prioritization strategy is effective.

```
rate.autonomy
#>  estimate std.err             target
#>      0.69    0.22 priorities | AUTOC
rate.autonomy$estimate / rate.autonomy$std.err
#> [1] 3.2
```

Next, we consider learning a simple policy that maps covariates `X` to a treatment decision using the `policytree` package. We perform this task on the observations without missing covariate values. First, we retrieve doubly robust scores for the quantities $\mathrm{E}[Y_i(0)]$ and $\mathrm{E}[Y_i(1)]$ using the function `double_robust_scores` from the `policytree` package.

```
library(policytree)

Gamma.hat <- double_robust_scores(c.forest)
no.missing <- complete.cases(X)
X.nm <- X[no.missing, ]
Gamma.nm <- Gamma.hat[no.missing, ]
```

A doubly robust estimate of the average financial proficiency score in the two treatment states is

```
colMeans(Gamma.nm)
#> control treated
#>      58      62
```

In this application, it is hard to imagine that some students do *not* benefit from the treatment, and the *treat everyone* policy likely increases the overall financial proficiency of the population. Treating everyone is costly;

hence, it is natural to ask if there is a policy that delivers an overall population benefit on a cost-adjusted basis. Here, we imagine we are paying the ATE when we assign treatment and offset this in our objective:

```
Gamma.nm[, "treated"] <- Gamma.nm[, "treated"] - ate["estimate"]
```

Then, we let `policy_tree` perform the maximization

$$\underset{\pi \in \text{depth-1 trees}}{\text{argmax}} \frac{1}{n} \sum_{i=1}^{n} \hat{\Gamma}[\pi(X_i)],$$

where $\pi(X_i) \in \{1, 2\}$ corresponds to the control and treated column of the `Gamma` matrix.

```
ptree <- policy_tree(X.nm, Gamma.nm, depth = 1)
print(ptree)
#> policy_tree object
#> Tree depth:  1
#> Actions:  1: control 2: treated
#> Variable splits:
#> (1) split_variable: financial.autonomy.index  split_value: 41
#>   (2) * action: 2
#>   (3) * action: 1
```

Here, we arrive at the simple decision rule "assign financial education training program if financial autonomy score is ≤ 41", reaffirming our previous analysis that indicated the pre-treatment variable `financial.autonomy.index` could serve as a good candidate for treatment decisions (*Note*: in practice, we would have set aside a held-out test set to estimate the value of this policy).

Example: The effect of poverty on attention

In this section, we analyze experimental data originally studied by Carvalho et al. (2016). The authors conducted a randomized experiment in which low-income individuals completed a cognitive test either before payday or after payday. The outcome is the number of correct answers on a test designed to measure cognitive ability. The original study did not find a statistically significant average effect. However, subsequent analyses suggest that poverty may impair cognitive function for certain subgroups (Farbmacher et al., 2021). A processed version of the data is available in `grf`. It contains around 2,500 individuals and 24 pre-treatment characteristics (age, income, education, etc.). Education is coded ordinally (4: college graduate; 3: some college; 2: high school graduate; 1: less than high school), and may therefore be used directly in the random forest splitting.

We code treatment status so that $W_i = 1$ if the individual was observed after payday. With this coding, positive treatment effects correspond to improved cognitive performance (i.e., reduced impairment). This choice simplifies interpretation: for example, the TOC curve is then positive when prioritization successfully targets individuals who benefit most. Using potential outcomes notation, let $Y_i(1)$ denote the number of correct answers under "no poverty" (after payday) and $Y_i(0)$ the number under financial distress (before payday). The average treatment is

$$\mathrm{E}[Y_i(1) - Y_i(0)],$$

which measures the average improvement in cognitive performance when individuals are not financially constrained. The CATE then captures heterogeneity in this improvement.

```
data("attentionrct")
Y <- attentionrct$outcome.correct.ans.per.second
W <- 1 - attentionrct$treatment
X <- attentionrct[-(1:4)]
```

Given the moderate sample size relative to the number of predictors, fitting a CATE estimator on all covariates with train/test evaluation may be underpowered. To reduce dimensionality, we first estimate the conditional

mean $\mathrm{E}[Y_i \mid X_i]$ and use forest variable importance to select a subset of covariates for the CATE. In the example below, keeping the top 10 variables by importance is a simple heuristic.

```
train <- sample(nrow(X), nrow(X) * 0.6)
test <- -train

Y.forest <- regression_forest(X[train, ], Y[train], num.trees = 500, num.threads = 2)
varimp.Y <- variable_importance(Y.forest)
keep <- colnames(X)[order(varimp.Y, decreasing = TRUE)[1:10]]
X.cate <- X[, keep]
print(keep)
#>  [1] "age"                                "working"
#>  [3] "education"                          "current.income"
#>  [5] "pay.day.amount.fraction.of.income" "black"
#>  [7] "household.income"                   "white"
#>  [9] "disabled"                           "retired"
```

Next, we fit a CATE forest on the training sample and evaluate heterogeneity on the held-out test sample. Because the data come from a randomized experiment, we supply the known randomization probability through the `W.hat` argument.

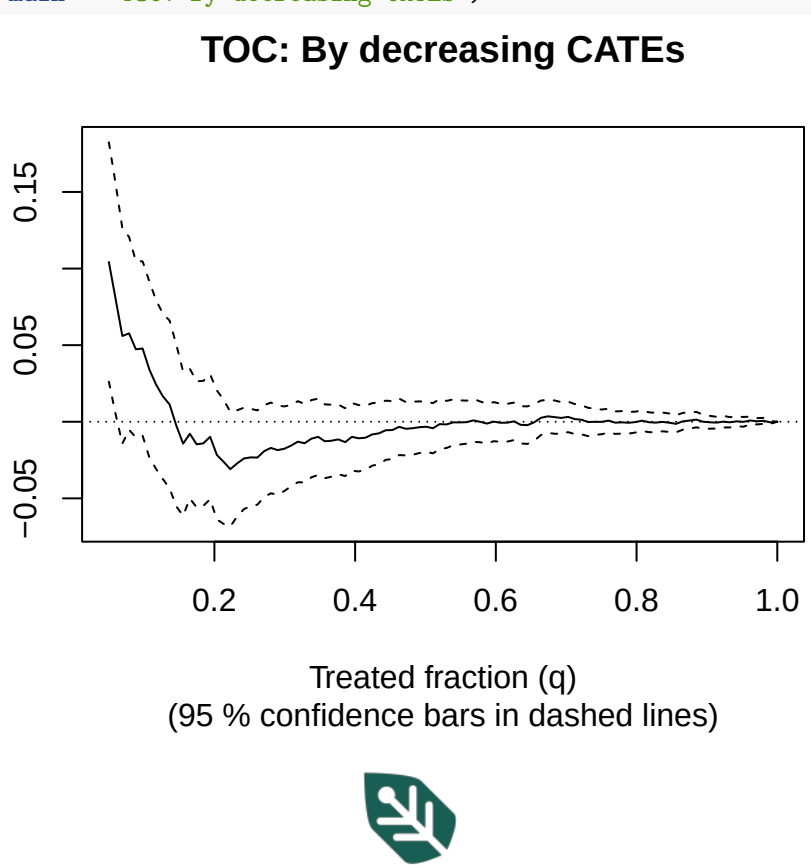
```
# Estimate CATE function.
Y.hat.train <- predict(Y.forest)$predictions
cate.forest <- causal_forest(
  X.cate[train, ],
  Y[train],
  W[train],
  Y.hat = Y.hat.train,
  W.hat = 0.5,
  num.threads = 2
)
tau.hat.eval <- predict(cate.forest, X.cate[test, ])$predictions

# Fit evaluation forest.
eval.forest <- causal_forest(
  X[test, ],
  Y[test],
  W[test],
  W.hat = 0.5,
  num.threads = 2
)
```

Next, we assess whether the estimated CATEs captures heterogeneity in cognitive impairment by estimating the TOC and area under the TOC.

```
rate.cate <- rank_average_treatment_effect(
  eval.forest,
  tau.hat.eval,
  q = seq(0.05, 1, length.out = 100)
)
print(rate.cate)
#>  estimate std.err             target
#>    0.0061  0.0092 priorities | AUTOC
rate.cate$estimate / rate.cate$std.err # t-statistic
#> [1] 0.67
```

```
plot(rate.cate, main = "TOC: By decreasing CATEs")
```



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