### 8. Regression discontinuity

LPO 8852: Regression II

Sean P. Corcoran

LPO 8852 (Corcoran)

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### RD - introduction

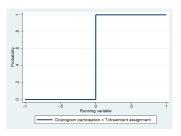
RD can be used when a **precise** rule based on a **continuous** characteristic determines treatment assignment. Examples:

- Test scores: can determine school admission, financial aid, summer school, remediation, graduation
- Income or poverty score: eligibility for income assistance or benefits, community eligibility for a means-tested anti-poverty program
- Date: age cutoff for retirement benefits, health insurance, school enrollment (KG or PK)
- Elections: fraction that voted for a particular candidate or ballot measure (e.g., school bond)

The continuous characteristic is typically called a **running variable**, **forcing variable**, or **index**.

#### RD - introduction

**Sharp RD**: treatment *assignment* goes from  $0 \rightarrow 1$  at a threshold c. Treatment *receipt* goes from 0% to 100% at c (full compliance).

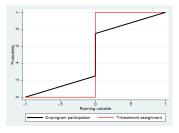


Re-center the running variable so that the threshold value is 0 (X - c).

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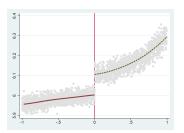
### RD - introduction

**Fuzzy RD**: treatment assignment goes from  $0 \rightarrow 1$  at a threshold c. Treatment receipt increases sharply at c but there is partial compliance.



#### RD - introduction

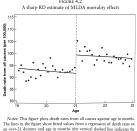
If there is a discrete change in treatment and program participation at c (and the program has a treatment effect) one would expect to see a discrete change in the mean outcome at c.



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#### RD - introduction

Under certain assumptions, this change can be interpreted as the (local) causal effect of the treatment. The challenge is estimating this change. There is often a relationship between the running variable and Y, even in the absence of treatment. We need to carefully estimate this relationship on either side of c, since this is what estimates the treatment effect.



### Potential outcomes and sharp RD

Each unit has potential outcomes  $Y_i(0)$  and  $Y_i(1)$ , which vary with the running variable  $X_i$ . We only observe  $Y_i(1)$  for those above c and  $Y_i(0)$  for those below c. The observed  $Y_i$ :

$$Y_i = (1 - D_i)Y_i(0) + D_iY_i(1) = \begin{cases} Y_i(0) & \text{if } X_i < c \\ Y_i(1) & \text{if } X_i \ge c \end{cases}$$

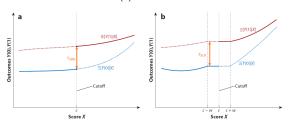
Here assume assignment to treatment and actual treatment are the same (i.e., sharp RD).

We never observe treated and untreated units with the same X—there is a complete lack of common support. RD estimation requires extrapolation!

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# Potential outcomes and sharp RD

There are two main frameworks for RD: the continuity framework (a) and local randomization framework (b).



The RD treatment effect at c is  $\tau_{SRD}$ . Figures: Cattaneo & Titiunik (2022)

# Potential outcomes and sharp RD

Under the **continuity assumption**, mean potential outcomes E[Y(0)|X] and E[Y(1)|X] are continuous near c. We need to model the relationship between Y and X below and above c to estimate  $\tau_{SRD}$ .

Under the **local randomization assumption**, treatment assignment—being above or below c—is random within a neighborhood  $(\pm w)$  of c. Mean potential outcomes do not vary within this neighborhood.

We will focus first and primarily on the continuity framework since it is more common. But there are cases where the randomization assumption arguably holds.

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### RD effects are local

Under certain assumptions our estimate of  $au_{SRD}$  can be considered causal, but at a single point in the distribution of X. It may have limited external validity. How representative au is depends on the context:

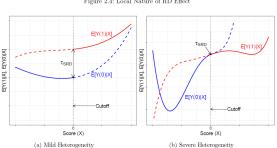


Figure 2.4: Local Nature of RD Effect

### RD plots

How should we model the relationship between Y and X?

- Linear function?
- Quadratic?
- Higher order polynomial?

A good place to start is a plot of the data, with X centered at c.

It can be difficult, however, to see systematic relationships in a scatterplot of the raw data. The user-written rdplot command can help (See also binscatter).

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# RD plots

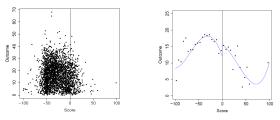


Figure 3.1: Scatter Plot—Meyersson Data 
Figure 3.2: RD Plot for Meyersson Data Using 40 Bins of Equal Length

Figures: Cattaneo et al. (2020) Foundations

### RD plots

#### A typical RD plot includes:

- Global polynomial fit (solid line) based on 4th or 5th order polynomial, fitted separately below and above c
- Local sample means (dots) based on binning X and plotting the mean for each bin at the midpoint of the bin

The plot can help you to see:

- Nonlinear relationship between Y and X
- Local variability around the fitted line
- Any evidence of a discontinuity at c
- Any other discontinuities away from c

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### rdplot

rdplot  $y \ x$  [, c(#) nbins(# #) p(#) binselect(binmethod) kernel(kernelfn) otheroptions]

- c(#) specifies the cutpoint c (default is 0)
- ullet nbins(# #) allows you to specify the number of bins on the L and R
- binselect(binmethod) specifies a procedure to determine the number of bins (if not set manually)
- p(#) is the order of the polynomial fit (default is 4)
- kernel(kernelfn) allows you to select a kernel: weights that depend on distance from c (default is uniform)
- Other options include confidence intervals, shading, covariates, etc.

### RD plots: deciding how bin width is determined

### Evenly-spaced bins (ES)

- All bins are equal width
- Number of observations per bin varies (thus, precision varies across bins)

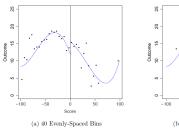
#### Quantile-spaced bins (QS)

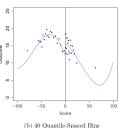
- All bins contain roughly the same number of observations
- Bin width varies
- Advantage: easier to see the density of the running variable

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# RD plots: deciding how bin width is determined

Figure 3.3: RD Plots-Meversson Data





Figures: Cattaneo et al. (2020) Foundations

### RD plots: choosing number of bins

#### Manual

- Can choose your own values nbins(# #)
- This is an ad hoc decision, however.

#### Integrated mean squared error (IMSE) method

- Chooses a number of bins that balances bias and variance when estimating local means
- A larger number of bins reduces bias, but there are fewer observations per bin, which leads to more sampling variance.
- · Results in means that roughly trace out the global polynomial fit
- Good for seeing general shape or other discontinuities away from c

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# RD plots: choosing number of bins

#### Mimicking variance (MV) method

- Chooses a number of bins so that local means have variability that approximates that of the data
- Usually leads to more bins than IMSE
- Provides a better picture of variability—it is less smoothed.

Note IMSE and MV methods will generally select a different number of bins on the L and R of c. If manually choosing the # of bins, you can also choose different values on the L and R.

Cattaneo et al. (2020) recommend starting with MV method, ideally comparing ES and QS for bin widths to show density of scores. Selecting IMSE method is preferable for global features of the regression function.

### rdplot bin options

Use these options with rdplot to select the number and type of bins:

nbins(# #) binselect(es): manual # of bins, equally spaced

nbins(# #) binselect(qs): manual # of bins, quantile spaced

binselect(es): IMSE method, equally spaced

binselect(qs): IMSE method, quantile spaced

binselect(esmv): MV method, equally spaced

binselect(qsmv): MV method, quantile spaced

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### Sharp RD estimation

The global polynomial fit in the RD plot can provide a good approximation overall, but a poor approximation at c, which is critical to RD estimation. Additionally, outlying observations far from c may be overly influential.

- When the full range of data are used to fit the relationship between Y and X, this is called a global or parametric approach.
- When a limited range of data around c are used, this is called a local, flexible, or non-parametric approach.

Current state of the art recommends the non-parametric approach: choosing a bandwidth around  $\boldsymbol{c}$  and fitting a local polynomial of low order.

### Sharp RD estimation

Figure 4.1: RD Estimation with local polynomial

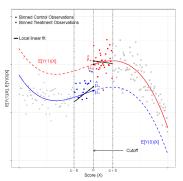


Figure: Cattaneo et al. (2020) Foundations

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# Sharp RD estimation: basic steps

Basic steps for sharp RD estimation under the continuity assumption:

- **1** Choose a polynomial order p and a kernel function K()
- Choose a bandwidth h
- Fit a weighted least squares regression of Y on  $(X_i c)$ ,  $(X_i c)^2$ , ...,  $(X_i c)^p$  up to the order p using K() function as weights. Do this for observations below and above the cutoff c. The difference in the two intercepts is  $\hat{\tau}_{SRD}$ . Note these can be done in the same regression (see next slides).

### Sharp RD estimation

Set aside the kernel and bandwidth for now and let  $\tilde{X}_i = X_i - c$ . The two regression models fit for observations below and above c are:

$$Y_{0i} = \alpha_0 + \beta_{01}\tilde{X}_i + \beta_{02}\tilde{X}_i^2 + \dots + \beta_{0p}\tilde{X}_i^p + u_i \text{ for } X_i < c$$

$$Y_{1i} = \alpha_1 + \beta_{11}\tilde{X}_i + \beta_{12}\tilde{X}_i^2 + \dots + \beta_{1p}\tilde{X}_i^p + u_i \text{ for } X_i \ge c$$

Notice the slope coefficients differ on either side of c. Let  $D_i = 1$  if  $X_i \ge c$ . Then we can pool the data and estimate one regression:

$$Y_{i} = \alpha_{0} + \beta_{01}\tilde{X}_{i} + \beta_{02}\tilde{X}_{i}^{2} + \dots + \beta_{0p}\tilde{X}_{i}^{p}$$
$$+ \rho D_{i} + \beta_{1}^{*} D_{i}\tilde{X}_{i} + \beta_{2}^{*} D_{i}\tilde{X}_{i}^{2} + \beta_{p}^{*} D_{i}\tilde{X}_{i}^{p} + u_{i}$$

where  $\rho=\alpha_1-\alpha_0$  (the difference in intercepts at c—this is our  $\hat{\tau}_{SRD}$ ) and  $\beta_j^*=(\beta_{1j}-\beta_{0j})$  (the difference in slope coefficients above and below c).

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## Sharp RD estimation

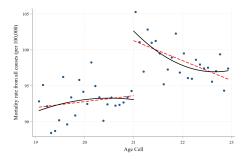
As an example, let p = 2 (quadratic):

$$Y_{i} = \alpha + \beta_{01}\tilde{X}_{i} + \beta_{02}\tilde{X}_{i}^{2} + \rho D_{i} + \beta_{1}^{*}D_{i}\tilde{X}_{i} + \beta_{2}^{*}D_{i}\tilde{X}_{i}^{2} + u_{i}$$

This is very easily estimated using OLS. See in-class example based on *Mastering Metrics* chapter 4 and Carpenter & Dobkin (2009): estimating the mortality effects of legal access to alcohol using the discontinuity in treatment at age 21. Note these data are already binned and limited to ages 19-23.

### In-class example: Carpenter & Dobkin 2009

Linear and quadratic fits with different slopes below and above c.



Note: uses agecell for x-axis instead of centered version.

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## Sharp RD estimation

You can improve your estimate of  $\hat{\tau}_{SRD}$  by choosing an appropriate bandwidth (h), polynomial order (p) and weighting function, or kernel K().

A **kernel function** assigns weights to the data based on their distance from the cutoff c. It may make sense to give greater weight to observations closer to c. Examples:

- uniform: equal weight
- triangular: weight declines symmetrically with distance from c (usually recommended)
- Epanechnikov: quadratic decline with distance from c

### Kernel functions

Figure 4.2: Different Kernel Weights for RD Estimation

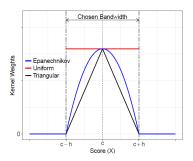


Figure: Cattaneo et al. (2020) Foundations

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## Polynomial order

Choice of polynomial order (p) is usually more consequential than K().

- For a given h, increasing p improves fit but also increases variability
  of the estimator
- Higher order polynomials tend to overfit, and have less reliable results near c (where it matters)

The current state of the art recommends a (local) **linear estimator**, but it is common to test sensitivity to higher orders (e.g., quadratic).

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## Choosing a bandwidth

The most consequential decision in RD is usually the bandwidth h.

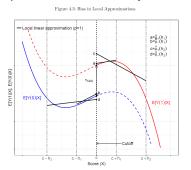


Figure: Cattaneo et al. (2020) Foundations

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## Choosing a bandwidth

A smaller h reduces the risk of misspecification ("smoothing bias") but will tend to increase the estimator's variability due to fewer observations. There is a **bias-variance** tradeoff in the choice of h.

You can choose your own bandwidth, but the decision is  $ad\ hoc$  and could lead to specification searching. Need a data-driven way to select h.

The most popular (and recommended) optimal bandwidth is **minimum MSE** (e.g., Imbens & Kalyanaraman, 2012). This minimizes the MSE of a local polynomial RD estimator, given a choice of p and K(). See Cattaneo et al. (2020) for a technical discussion.

### Sharp RD estimation using rdrobust

You could choose an h and K() and implement these using OLS (as in the in-class example above). There are several issues with this, however:

- It will involve multiple steps, and you would also want to use an optimal bandwidth selector first to choose h for you.
- The standard errors will be invalid. The MSE method allows for some bias in the choice of optimal bandwidth. OLS standard errors do not account for the effects of this approximation.

The user-written command rdrobust is very flexible and can handle all of these things and more.

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#### rdrobust

rdrobust y x [, c(#) p(#) h(# #) bwselect(bwmethod)
kernel(kernelfn) covs(covars) otheroptions]

- c(#) specifies the cutpoint c (default is 0)
- bwselect(bwmethod) specifies a procedure to determine the bandwidth (if not set manually)
- ullet p(#) is the order of the polynomial fit (default is 1)
- kernel(kernelfn) (default is triangular)
- covs(covars) allows you to include covariates in your model
- Other options

### Sharp RD estimation using rdrobust

See in the in-class example using data from Carpenter & Dobkin (2009), using rdrobust instead of regress.

Example using fixed bandwidth of  $\pm 2,$  uniform kernel, and either linear or quadratic fit:

```
rdrobust all age, c(0) h(2) kernel(uniform) p(1) rdrobust all age, c(0) h(2) kernel(uniform) p(2)
```

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# Optimal bandwidth selection using rdbwselect

The related command rdbwselect will perform optimal bandwidth selection methods. (Note this can also be called by rdrobust in one step).

#### Two common options:

- mserd: MSE method, same bandwidth on both sides of c
- msetwo: MSE method, different bandwidths on each side

There are other options, see Cattaneo et al. (2020) for details.

#### rdbwselect

rdbwselect y x [, c(#) p(#) bwselect(bumethod)
kernel(kernelfn) covs(covars) otheroptions]

- c(#) specifies the cutpoint c (default is 0)
- bwselect(bumethod) specifies a procedure to determine the bandwidth (default is mserd)
- p(#) is the order of the polynomial fit (default is 1)
- kernel(kernelfn) (default is triangular)
- covs(covars) allows you to include covariates in your model
- Other options

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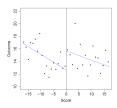
### Optimal bandwidth selection

You can pass through the optimal bandwidth (uses rdbwselect) to the rdrobust and rdplot commands.

In rdrobust: rdrobust y x, bwselect(mserd)

See the in-class example code for this and the RD plot code.

Figure 4.4: Local Polynomial RD Effect Illustrated with rdplot—Meyersson Data



#### Standard errors and confidence intervals

As noted earlier, the usual OLS standard errors are invalid. See Cattaneo et al. (2020) for an extended discussion of inference in sharp RD. Their recommendation is to use the **robust bias-corrected** approach to inference (i.e., confidence intervals). Can add all option to see inference using different methods.

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# Including covariates

You can include covariates in your model to improve precision, <u>but</u> if the continuity assumption holds, one would not expect covariates to be systematically different below/above c. Obtaining materially different results with covariates is a red flag.

In rdrobust: rdrobust y x, covs(covars)

Note the optimal bandwidth will likely change with the new model specification (w/covariates).

Another way to use covariates: splitting the sample by subgroups. This is commonly done, and requires no modification of the steps above (just subset on the group when using rdrobust).

### RD assumptions

To assess the validity of the sharp RD, consider the assumptions:

- Treatment assignment (and receipt) occurs at a known threshold c.
- The relationship between potential outcomes Y(1), Y(0) and X is continuous in the neighborhood of c. There is no reason to expect a sharp break in Y in the absence of treatment.

#### These imply:

- X has **not been manipulated** to affect who receives treatment.
- There are no other programs or services with the same eligibility rule (to avoid confounding with some other treatment).

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# Some common RD validity tests

The following are commonly performed as validity tests with RD:

- Test for effects at c on pre-treatment covariates or placebo outcomes. Would expect a null effect.
- Test for continuity in the density of the running variable around c ("manipulation test").
- Test for discontinuities elsewhere in the distribution of X (i.e., artificial cutoffs). A "smoothness" test.
- Exclusion of observations near c ("donut hole" approach).
- Sensitivity tests for bandwidth choice.

### Test for effects on pre-treatment covariates

Use the same rdrobust code but swapping in pre-treatment covariates (or placebo outcomes) for y. Note the optimal bandwidth will change (this is OK). Can also show RD plots:

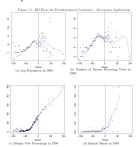


Figure: Cattaneo et al. (2020) Foundations

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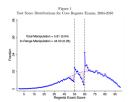
## Manipulation test

**Manipulation** occurs whenever units have their value of X altered in order to affect their treatment status. For example, a teacher might adjust a test score in order to help a student pass or become eligible for a program.

- This may be visible in a histogram, or not if the manipulation goes both ways.
- If manipulation is <u>random</u> or <u>uninformed</u>, such that potential outcomes in the absence of treatment are no different on average for those whose X has been manipulated, then manipulation will not pose a problem. Manipulation is not usually random, however.
- Manipulation may lead one to find an "effect" where there is none.
- Note manipulation is <u>not</u> the same thing as non-compliance (the "fuzzy RD" case).

# Manipulation test

#### Sometimes manipulation is clear from inspecting densities or histograms:



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Figure: Dee et al. (2011)

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# Manipulation test

Figure 5.3: Histogram of Score

Cutoff

Score (X)

(a) No Sorting

(b) Sorting

Figure: Cattaneo et al. (2020) Foundations

## Manipulation test using rddensity

You should present both a figure and a formal statistical test. The user-written rddensity command can help with this.

#### rddensity x, plot otheroptions

The idea of the statistical test is to fit a local polynomial to the density of X on the L and R of c and testing for a discontinuity at c. This general procedure is commonly referred to as a McCrary test (McCrary, 2008) although details of the test vary.

 $H_0$  is "no manipulation," or no discontinuity at c.

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# Manipulation test using rddensity

Figure 5.4: Histogram and Estimated Density of the Score

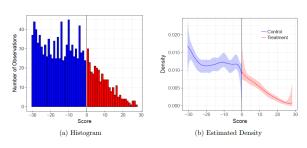
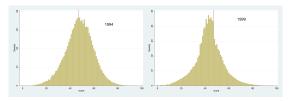


Figure: Cattaneo et al. (2020) Foundations

### Example: Camacho and Conover (2011)

In 1998, Colombia set eligibility threshold for social welfare benefits at a poverty index of 47.



There is evident manipulation in 1999 but the manipulation test also fails in 1994. This test can over-reject with a **discrete** running variable (here, an integer poverty index), even more so when *N* is large.

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#### Tests at alternative cutoffs

Can use rdrobust to test for discontinuities at other points. To avoid "contamination" from real cutpoints, include *only* the treated  $(X \geq c)$  or untreated (X < c) observations, depending on where alternative cutoff is located. Example:

rdrobust y x if  $x \ge 0$ , c(1)

Cattaneo et al. (2020): "evidence of continuity away from the cutoff is neither necessary nor sufficient for continuity at the cutoff, but the presence of discontinuities away from the cutoff can be interpreted as potentially casting doubt on the RD design, at the very least in cases where such discontinuities can not be explained by substantive knowledge of the specific application."

#### Tests at alternative cutoffs

A graphical summary of tests at alternative cutoffs, including the real cutoff at X = c as a reference:

Figure 5.5: RD Estimation for True and Placebo Cutoffs

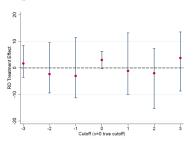


Figure: Cattaneo et al. (2020) Foundations

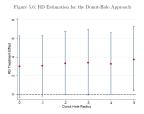
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#### Donut hole robustness test

Units closest to c get the most weight and are also those most susceptible to manipulation (in some applications). Can exclude observations within a certain radius of c to see if/how results change. Example:

rdrobust  $y x \text{ if } abs(x) \ge 0.3$ 



### Sensitivity to bandwidth

It is common to see estimates using different bandwidth choices as a robustness check. Could use multiples of the original bandwidth (E.g.,  $\times 0.5$ ,  $\times 1.5$ ,  $\times 2$ )

Results should be interpreted with caution: wider bandwidths can introduce bias, while narrower bandwidths introduce more sampling variability.

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# Sharp RD under local randomization

Thus far we have relied on the **continuity** framework, where mean potential outcomes are continuous near c. We modeled the relationship between Y and X below and above c in order to estimate  $\tau_{SRD}$ .

Ideally we would not have to rely on functional form assumptions at all. Under **local randomization**, treatment assignment—being above or below c—is <u>random</u> within a neighborhood ( $\pm w$ ) of c. Mean potential outcomes do not vary within this neighborhood.

RD is often described as "as good as random assignment" around c. But this is not strictly true if we are fitting a polynomial to estimate  $\tau_{SRD}$ . If local randomization held, we could just compare means on either side.

### Sharp RD under local randomization

With local randomization, potential outcomes are not related to the running variable x in a window  $(\pm w)$  around c:

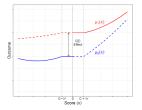


Figure 2.2: Local Randomization RD

Figure: Cattaneo et al. (2023) Extensions

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# Sharp RD under local randomization

The assumption of local randomization is stronger than the assumption of continuity. So when might you want to use this approach?

- If you have a large sample and a continuous running variable where you believe random factors play an important role in whether a unit is just above or below c (e.g., a test, voter turnout).
- If you have a small sample near c, making it difficult to reliably approximate the function on either side of c.
- If you have a discrete running variable, such that the continuity assumption is not conceptually appropriate.

The local randomization approach is often used as a robustness check to a continuity-based analysis when the running variable is continuous.

#### Local randomization: estimation and inference

Cattaneo et al. (2023) provide a full description of sharp RD estimation and inference under the local randomization assumption. See their paper for more details. It boils down to comparing means on either side of c.

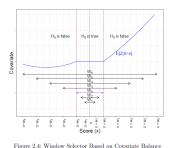
The most consequential decision: within what window  $(\pm w)$  can we assume units are "as good as randomly assigned?"

They recommend testing for balance in pre-determined covariates, starting with the smallest possible window around c and getting progressively wider until the null of no difference is rejected. The chosen window is the largest one such that  $H_0$  fails to be rejected inside that window (and all windows within it).

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#### Window selection under local randomization

 $W_1$  is the first window tested, then  $W_2$ , etc.  $W_0$  is the widest window where  $H_0$  is not rejected. Can use rdwinselect for this.



#### rdwinselect

rdwinselect runvar covs [, c(#) wobs(#) wstep(#) seed(#)
otheroptions]

- c(#) specifies the cutpoint c (default is 0)
- wobs (#) tells Stata to increase window width by a fixed number of observations each time.
- wstep(#) instead tells Stata to increase window width by a fixed amount each time
- seed(#) sets a random seed, for small-sample randomization inference.
- Other options

By default the smallest window is the one with  $\geq$  10 obs on each side.

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#### Local randomization: estimation and inference

Once the window is selected, can use rdrandinf to estimate the difference in means above and below c, construct confidence intervals, etc.

Inference is based on either small sample (exact) randomization inference, or large sample approximations. See Cattaneo et al. (2023) for details.

Validity tests are important here as well, and are carried out with some modification. See Cattaneo et al. (2023) for details.

- Test for effects on pre-treatment covariates and placebo outcomes
- Manipulation test
- Test for effects at artificial cutoffs
- Sensitivity to window width

#### rdrandinf

rdrandinf y x [, c(#) wl(#) wr(#) seed(#) otheroptions]

- c(#) specifies the cutpoint c (default is 0)
- wl(#) and wr(#) are the lower and upper bounds of the window.
   You can choose these yourself or use rdwinselect.
- seed(#) sets a random seed, for small-sample randomization inference.
- Other options

You can select the window and perform estimation all in one step, but it is advisable to make decisions about the window before seeing the results.

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### Fuzzy RD

**Fuzzy RD**: treatment assignment goes from  $0 \rightarrow 1$  at c. Treatment receipt (e.g., program participation) increases sharply at c but there is some non-compliance.

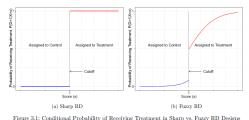


Figure 3.1: Conditional Probability of Receiving Treatment in Snarp vs. Fuzzy RD Designs

Figure: Cattaneo et al. (2023) Extensions

### Fuzzy RD: intent-to-treat effects

Treatment assignment changes sharply at c. If the continuity (and/or local randomization) assumption holds, we still can estimate the causal effect of treatment assignment at c using the sharp RD methods described thus far.

This is <u>not</u> the treatment effect at c, but rather the **intent-to-treat**,  $\tau_{ITT}$ .

In a fuzzy RD context, one should also estimate the effect of treatment assignment on the probability of receiving the treatment,  $\tau_D$ , using sharp RD methods. This is the **first stage**, for reasons that will become clear.

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## Fuzzy RD: nonrandom take-up

The challenge for estimating the treatment effect at c is that units may strategically decide to take up (or not) the treatment based on their expected gains from it. There is no reason to believe potential outcomes are unrelated to the decision to take up treatment.

### Potential treatment take-up

Potential treatment take-up, analogous to potential outcomes:

 $D_i(0)$  is i's treatment take-up/receipt when  $T_i = 0$ 

 $D_i(1)$  is i's treatment take-up/receipt when  $T_i = 1$ 

$D_i(T_i)$		Types:
$T_i = 0$	$T_i = 1$	
0	1	Compliers
1	1	Always-takers
0	0	Never-takers
1	0	Defiers

There are now <u>four</u> potential outcomes that depend on both treatment assignment and take-up:  $Y_i(T_i, D_i)$ .

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## Fuzzy RD as IV

Treatment assignment  $T_i$  at c is potentially a good instrumental variable for treatment receipt  $D_i$ . If so, we know from IV that the (local) average treatment effect of  $D_i$  on the outcome Y would be:

$$\tau_{\mathit{FRD}} = \frac{\tau_{\mathit{ITT}}}{\tau_{\mathit{D}}}$$

The numerator and denominator are bout sharp RD parameters that we know how to estimate.

Call  $\tau_{FRD}$  the fuzzy RD parameter. Under what conditions does it provide the local average treatment effect (LATE) at c?

### Fuzzy RD as IV

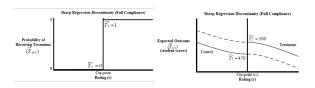
#### Conditions under which $\tau_{FRD}$ provides the LATE:

- Exclusion restriction: treatment assignment (being above/below c)
  cannot affect potential outcomes (or potential treatments) except
  through treatment received.
- Monotonicity: we assume there are no defiers at or near c (or inside the window W if using local randomization).

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# Example from Bloom (2009, 2012)

#### Sharp RD:



$$\tau_{RD} = 500 - 470 = 30$$

## Example from Bloom (2009, 2012)

#### Fuzzy RD (one-sided non-compliance):



$$\tau_{ITT} = 495 - 470 = 25$$
 $\tau_{FS} = 0.80 - 0 = 0.80$ 
 $\tau_{FRD} = \tau_{ITT}/\tau_{FS} = 25/0.80 = 31.25$ 

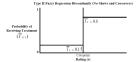
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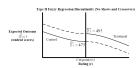
# Example from Bloom (2009, 2012)

Above: there are no "always-takers" (else we would see them below c); there are only compliers and never-takers. Above c, it is clear that 80% are compliers and 20% are never-takers. The ITT is a weighted average of  $\tau_{FRD}$  (for compliers) and 0 (for never-takers). The fuzzy RD parameter "scales up" the ITT to reflect the effect on the compliers.

## Example from Bloom (2009, 2012)

#### Fuzzy RD (two-sided non-compliance):





$$\tau_{ITT} = 495 - 475 = 20$$
 $\tau_{FS} = 0.80 - 0.15 = 0.65$ 
 $\tau_{FRD} = \tau_{ITT}/\tau_{FS} = 20/0.65 = 38.46$ 

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# Example from Bloom (2009, 2012)

Above: there are both "always-takers" and "never-takers" (in addition to compliers). Assuming the proportion of always-takers and never-takers is constant in the neighborhood of c, we can estimate the never-taker rate to be 20% and always-taker rate to be 15%. The rest (65%) are compliers.

Again, the ITT is a weighted average of  $\tau_{FRD}$  (for compliers) and 0 (for never-takers). The fuzzy RD parameter "scales up" the ITT to reflect the effect on the compliers.

Note: in practice the proportion of non-compliant cases can vary with the running variable x, but it must be continuous through c.

### Fuzzy RD estimation

Since  $\tau_{FRD}$  is constructed as the ratio of two sharp RD parameters, one could just use rdrobust or rdrandinf to estimate those parameters. However:

- In rdrobust, the two bandwidths may (and probably will) differ. (If
  interested in the ITTs themselves, it's fine to use differing
  bandwidths). In rdrandinf, this is not an issue since the window is
  based on pre-determined covariates.
- You will lack a standard error, confidence intervals, p-value for  $\tau_{FRD}$

These two commands include a fuzzy option that (in the case of rdrobust) will choose <u>one</u> optimal bandwidth and report standard errors, etc., for the ratio.

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## Fuzzy RD: example

Access to financial aid program based on poverty index (SISBEN) in Colombia. First stage:

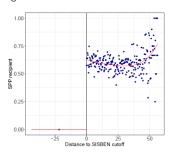


Figure 3.2: RD Plot: First Stage—SPP data

### Fuzzy RD: example

Reduced form (intent-to-treat) effect on enrollment in a high-quality college:

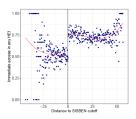


Figure 3.3: RD Plot: Intention-to-treat (p = 3)—SPP data

### Figure: Cattaneo et al. (2023) Extensions

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# Fuzzy RD: example

rdrobust with fuzzy option:



### Fuzzy RD validity tests

Validity tests should be carried out in the same manner as sharp RD, following under the continuity or local randomization approach. Focus on assignment to treatment, not actual treatment (which is nonrandom).

The fuzzy RD estimator is an IV estimator so a **weak instrument** is problematic. If the first stage—how assignment to treatment is related to actual treatment—is small, this can lead to unreliable estimates. Cattaneo et al. (2023) recommend an F-statistic of 20 or higher.

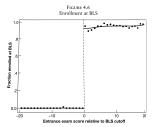
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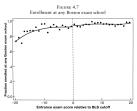
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# Abdulkadiroğlu, Pathak, & Roth (2014)



Notes: This figure plots enrollment rates at Boston Latin School (BLS) conditional on admissions test scores, for BLS applicants scoring near the BLS admissions cutoff. Solid lines show fitted values from a local linear regression estimated separately on either side of the cutoff (indicated by the vertical dashed line).

# Abdulkadiroğlu, Pathak, & Roth (2014)



Note: This figure plots enrollment rates at any Boston exam school, conditional on admissions test scores, for Boston Latin School (BLS) applicants scoring near the BLS admissions cutoff. Solid lines show fitted values from a local linear regression, estimated separately on either side of the cutoff (indicated by the vertical dashed line).

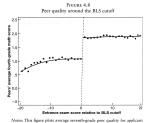
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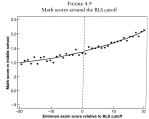
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# Abdulkadiroğlu, Pathak, & Roth (2014)



Notes: Ins ingute poist average seventir-grate peer quarry for application to Boston Litali School (BLS), conditional on admissions test scores, for BLS applicants scoring near the admissions cutoff. Peer quality is measured seventh-grade exchoolimates' fourth-grade math socres. Solid lines show fixed values from a local linear regression, estimated separately on either side of the cutoff (indicated by the vertical dashed line).

# Abdulkadiroğlu, Pathak, & Roth (2014)



Notes: This figure plots seventh- and eighth-grade math scores for applicants to the Boston Latin School (BLS), conditional on admissions test scores, for BLS applicants scoring near the admissions cutoff. Solid lines show fitted values from a local linear regression, estimated separately on either side of the cutoff (indicated by the verifical dashed line).

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