Big Data and Economics

Bootstrapping, Functions, and Parallel Processing

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Software requirements

R packages

It's important to note that "base" R already provides all of the tools to implement a regression discontinuity design, but there are some great other tools to use!

- New: boot, parallel, tictoc, pbapply, future, future.apply, furrr
- · Already used: tidyverse, fixest, haven, sandwich, broom, stargazer, lmtest, AER

A convenient way to install (if necessary) and load everything is by running the below code chunk.

```
## Load and install the packages that we'll be using today
if (!require("pacman")) install.packages("pacman")
pacman::p_load(boot,doParallel,fixest,sandwich,haven,tidyverse,broom,stargazer,lmtest,tictoc,parallel,p
## My preferred ggplot2 plotting theme (optional)
theme_set(theme_minimal())
```

Dataset

In this activity, we are going to work through bootstrapping to get standard errors for a regression model. We will use the ResumeNames dataset, which is in the AER package. The data come from the Resume Audit study by Bertrand and Mullainathan (2004). The data contain a row that corresponds to a ficitious resume that the authors sent to businesses in Boston and Chicago. The resumes were randomly assigned either a "white-sounding" name or a "black-sounding" name, among other characteristics. The data contain information on whether the business called the resume back for an interview. The outcome of interest way the average callback gap.

```
call=call=='yes') %>%
select(female, afam,call) %>%
as_tibble()

cleaned_resumes
```

```
## # A tibble: 4,870 x 3
##
     female afam call
      <lgl> <lgl> <lgl> <lgl>
##
   1 TRUE
           FALSE FALSE
##
   2 TRUE
           FALSE FALSE
##
  3 TRUE
           TRUE FALSE
##
  4 TRUE
           TRUE FALSE
## 5 TRUE
           FALSE FALSE
## 6 FALSE FALSE FALSE
## 7 TRUE
          FALSE FALSE
## 8 TRUE
            TRUE FALSE
## 9 TRUE
            TRUE FALSE
## 10 FALSE TRUE FALSE
## # i 4,860 more rows
```

Functions

Functions are a great way to make your code more readable and reusable. They are also a great way to make sure you don't make mistakes when you're running the same code over and over again.

Basic function to sample data

Let's start with a basic function. Here is a function that creates a random sample of the data with the same number of observations. This is the first step in bootstrapping.

```
set.seed(1000) # replication
random_sample <- function(data) {
    # Get the number of observations
    n <- nrow(data)
    # Sample the data
    data[sample(1:n, n, replace = TRUE), ]
}
random_sample(data=cleaned_resumes)</pre>
```

```
## # A tibble: 4,870 x 3
##
     female afam call
      <lgl> <lgl> <lgl> <lgl>
##
   1 TRUE
           FALSE FALSE
##
   2 FALSE TRUE FALSE
##
  3 TRUE
##
            TRUE FALSE
## 4 TRUE
            FALSE FALSE
## 5 FALSE TRUE FALSE
## 6 FALSE FALSE FALSE
## 7 TRUE
            TRUE FALSE
            TRUE FALSE
## 8 TRUE
## 9 TRUE
            FALSE FALSE
            TRUE FALSE
## 10 TRUE
## # i 4,860 more rows
```

Can you think of some way to improve on this sampling procedure? What is true in the experiment that we will lose through a random sample? Try to write up a better version that accounts for this.

Now test yourself by amending the function to run a regression of callback on ethnicity. Test it out!

```
set.seed(1000) # replication
random_sample_lm <- function(data) {
    # Get the number of observations
    n <- nrow(data)
    # Sample the data
    m1 <- lm(call ~ afam, data[sample(1:n, n, replace = TRUE), ])
    # Return the coefficients
    return(coef(m1)[2])
}
random_sample_lm(cleaned_resumes)</pre>
```

```
## afamTRUE
## -0.03381552
```

Iteration

Iteration is a great way to run the same code over and over again. It's especially useful for bootstrapping because bootstrapping is just running the same code over and over again with different data.

For loop

The most basic way to iterate is with a for loop. This is a great way to start because it's easy to understand and you can see what's going on. Let's write a for loop that iterates through the random_sample_lm() function 100 times. Make sure to create a list to store the results in.

```
results <- vector("list",100)
set.seed(1000) # replication
for (i in 1:100) {
   results[[i]] <- random_sample_lm(cleaned_resumes)
}</pre>
```

Apply

You can do the same thing using the *apply family. Here it is with lapply():

```
set.seed(1000) # replication
results <- lapply(1:100, function(x) random_sample_lm(cleaned_resumes))</pre>
```

Also, the pbapply package has a progress bar, which is nice for long simulations.

```
set.seed(1000) # replication
results <- pbapply::pblapply(1:100, function(x) random_sample_lm(cleaned_resumes))</pre>
```

Map

The map() function from the **purrr** package is a great way to iterate over a list. It's a bit more complicated than lapply(), but it's worth learning because it's very powerful. Here's an example with map_df().

```
set.seed(1000)
results <- map_df(1:100, function(x) random_sample_lm(cleaned_resumes))</pre>
```

Parallel Processing

Parallel processing is a pretty handy trick to speed up your code. It's especially useful for computationally intensive simulations that can be broken up into smaller chunks. As a warning, parallel programming can get quite complex, so we're sticking with the basics today. For more advanced coverage, check work by Grant McDermott.

First, let's check how many cores you have on your machine. R runs things in parallel by splitting them up between cores!

```
# future::availableCores() ## Another option
detectCores()
```

```
## [1] 8
```

My computer has 8 cores, yours likely, but not necessarily, has fewer. That's okay! You can still split things into chunks and run them in parallel.

I'll show several different approaches.

future.apply

```
set.seed(123L)
tic()
future_lapply(1:1e3,
  function(i) random sample lm(cleaned resumes)) %>%
 bind rows()
## Warning: UNRELIABLE VALUE: One of the 'future.apply' iterations
## ('future lapply-1') unexpectedly generated random numbers without declaring so.
## There is a risk that those random numbers are not statistically sound and the
## overall results might be invalid. To fix this, specify 'future.seed=TRUE'. This
## ensures that proper, parallel-safe random numbers are produced via the
### L'Ecuyer-CMRG method. To disable this check, use 'future.seed = NULL', or set
## option 'future.rng.onMisuse' to "ignore".
## Warning: UNRELIABLE VALUE: One of the 'future.apply' iterations
## ('future_lapply-2') unexpectedly generated random numbers without declaring so.
## There is a risk that those random numbers are not statistically sound and the
## overall results might be invalid. To fix this, specify 'future.seed=TRUE'. This
## ensures that proper, parallel-safe random numbers are produced via the
## L'Ecuyer-CMRG method. To disable this check, use 'future.seed = NULL', or set
## option 'future.rng.onMisuse' to "ignore".
## Warning: UNRELIABLE VALUE: One of the 'future.apply' iterations
## ('future lapply-3') unexpectedly generated random numbers without declaring so.
## There is a risk that those random numbers are not statistically sound and the
## overall results might be invalid. To fix this, specify 'future.seed=TRUE'. This
## ensures that proper, parallel-safe random numbers are produced via the
### L'Ecuyer-CMRG method. To disable this check, use 'future.seed = NULL', or set
## option 'future.rng.onMisuse' to "ignore".
## Warning: UNRELIABLE VALUE: One of the 'future.apply' iterations
## ('future_lapply-4') unexpectedly generated random numbers without declaring so.
## There is a risk that those random numbers are not statistically sound and the
## overall results might be invalid. To fix this, specify 'future.seed=TRUE'. This
## ensures that proper, parallel-safe random numbers are produced via the
## L'Ecuyer-CMRG method. To disable this check, use 'future.seed = NULL', or set
## option 'future.rng.onMisuse' to "ignore".
## Warning: UNRELIABLE VALUE: One of the 'future.apply' iterations
```

```
## ('future_lapply-5') unexpectedly generated random numbers without declaring so.
## There is a risk that those random numbers are not statistically sound and the
## overall results might be invalid. To fix this, specify 'future.seed=TRUE'. This
## ensures that proper, parallel-safe random numbers are produced via the
## L'Ecuyer-CMRG method. To disable this check, use 'future.seed = NULL', or set
## option 'future.rng.onMisuse' to "ignore".
## Warning: UNRELIABLE VALUE: One of the 'future.apply' iterations
## ('future lapply-6') unexpectedly generated random numbers without declaring so.
## There is a risk that those random numbers are not statistically sound and the
## overall results might be invalid. To fix this, specify 'future.seed=TRUE'. This
## ensures that proper, parallel-safe random numbers are produced via the
## L'Ecuyer-CMRG method. To disable this check, use 'future.seed = NULL', or set
## option 'future.rng.onMisuse' to "ignore".
## Warning: UNRELIABLE VALUE: One of the 'future.apply' iterations
## ('future_lapply-7') unexpectedly generated random numbers without declaring so.
## There is a risk that those random numbers are not statistically sound and the
## overall results might be invalid. To fix this, specify 'future.seed=TRUE'. This
## ensures that proper, parallel-safe random numbers are produced via the
## L'Ecuyer-CMRG method. To disable this check, use 'future.seed = NULL', or set
## option 'future.rng.onMisuse' to "ignore".
## Warning: UNRELIABLE VALUE: One of the 'future.apply' iterations
## ('future lapply-8') unexpectedly generated random numbers without declaring so.
## There is a risk that those random numbers are not statistically sound and the
## overall results might be invalid. To fix this, specify 'future.seed=TRUE'. This
## ensures that proper, parallel-safe random numbers are produced via the
## L'Ecuyer-CMRG method. To disable this check, use 'future.seed = NULL', or set
## option 'future.rng.onMisuse' to "ignore".
## # A tibble: 1,000 x 1
##
      afamTRUE
##
         <dbl>
## 1 -0.0412
## 2 -0.0391
## 3 -0.0428
## 4 -0.0290
## 5 -0.0414
## 6 -0.0284
## 7 -0.0317
## 8 -0.0243
## 9 -0.0381
## 10 -0.0343
## # i 990 more rows
toc(log=TRUE)
## 4.12 sec elapsed
set.seed(123L)
tic()
sim_pblapply = pblapply(1:1e3, function(i) random_sample_lm(cleaned_resumes), c1 = parallel::detectCore
```

pbapply

toc(log=TRUE)

```
## 4.22 sec elapsed
```

furrr

```
tic()
future_map_dfr(1:1e3,
 function(i) random sample lm(cleaned resumes),
  .options = furrr_options(seed=123L)) %>%
 bind rows()
## # A tibble: 1,000 x 1
     afamTRUE
##
##
        <dbl>
  1 -0.0266
##
## 2 -0.0198
## 3 -0.0420
## 4 -0.0461
## 5 -0.0310
##
  6 -0.0372
## 7 -0.0226
## 8 -0.0328
## 9 -0.0177
## 10 -0.0241
## # i 990 more rows
toc(log=TRUE)
```

5.28 sec elapsed

This is the best way to make sure you understand the bootstrapping process.

Bootstrapping out of the box

sandwich package

The sandwich package is great for estimating all sorts of funny standard errors in R. It is often combined with the coeftest function from the **Imtest** package to substitute new standard errors into the model after estimation and specify the number of replications.

```
#library(sandwich); library(lmtest) #already loaded
# Run a regression assuming normal (iid) errors
m <- lm(call ~ afam, data = cleaned_resumes)</pre>
# Obtain the boostrapped SEs
coeftest(m, vcov = vcovBS(m,
 R=1000 # Number of replications
 )
)
##
## t test of coefficients:
##
##
               Estimate Std. Error t value Pr(>|t|)
## (Intercept) 0.0965092 0.0060210 16.0287 < 2.2e-16 ***
## afamTRUE
             ## ---
```

```
## Signif. codes: 0 '*** 0.001 '** 0.01 '* 0.05 '.' 0.1 ' ' 1
```

sandwich can be run in parallel by specifying the "cores" to use. This is useful for computationally intensive simulations.

```
coeftest(m, vcov = vcovBS(m,
  R=1000, # Number of replications
  cores=2 # Number of cores to use
  )
)
```

```
##
## t test of coefficients:
##
## Estimate Std. Error t value Pr(>|t|)
## (Intercept) 0.0965092 0.0059297 16.2755 < 2.2e-16 ***
## afamTRUE -0.0320329 0.0079433 -4.0327 5.599e-05 ***
## ---
## Signif. codes: 0 '***' 0.001 '**' 0.05 '.' 0.1 ' ' 1</pre>
```

One challenge is **sandwich** only works for standard error calculations and does not give you as much information on the bootstrap needed to use it other ways!

boot package

The boot package is one of the most flexible packages for bootstrapping in R. It allows you to bootstrap any function that you can write in R. It also allows you to parallelize your bootstrap simulations, which can be very useful for computationally intensive simulations.

Here's that again, but now in parallel:

There are a number of handy visualization tools and table shortcuts to use with the **boot** package too:

```
# See the coefficient estimates and their bootstrap standard errors plot(results)
```

```
-0.045 -0.040 -0.035 -0.030 -0.025 -0.020 -0.015
```

```
# Optional: print regression table with the bootstrap SEs
# This uses stargazer, but the method is similar
# with other table-making packages,
# see /Presentation/export_a_formatted_regression_table.html
# library(broom) # already loaded
tidy_results <- tidy(boot_results)

#library(stargazer) # already loaded
m1 <- lm(call~afam, data = cleaned_resumes)
stargazer(m1, se = list(tidy_results$std.error), type = 'text')</pre>
```

##
% Error: Argument 'se' must be NULL (default), or a list of numeric vectors.

That said, it can be a bit cumbersome to use. For example, if you want to bootstrap a function that takes more than two arguments, you need to use the . . . argument and define extraPar and numCenter. This can be a bit confusing and easy to screw up.

The code below does not work. Can you figure out why?

```
statFun <- function(funData, indices, addPars, centerMean)
{
    # Check to see if extra parameters
    if(addPars)
    {
        result <- mean(funData[indices] - centerMean)
    }else
    {
        result <- mean(funData[indices])
    }

# Return the value
    return(result)</pre>
```

```
boot(testData, statFun, R = 100, extraPar = TRUE, numCenter = mean(testData))
```

Other resources

- For more info on functions, check out Grant McDermott and Ed Rubin's introductory and advanced material.
- For more info on paralell programming, check out Grant McDermott and Ed Rubin's chapter on all the nitty gritty details.
- For more explanations on bootstrapping in sections, see 13.3.2 and 15.5 of *The Effect* by Nick Huntington-Klein or The Library of Statistical Techniques also provides a concise summary. A much longer explanation is available in Data Science for R.