### Outline



#### 2 Survey

- for
- apply
- lapply
- lapply: Extended Example #1
- lapply: Extended Example #2
- mapply: a secret weapon

### 3 Bootstrapping

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# Iteration: for, appply, etc Efficiency and Clarity

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2013

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# R Frame of Mind

- Iteration is commonly needed
  - repeat the same thing over and over with new samples
  - process several subgroups of data (compare cities)
  - apply various functions to one data set
- Some idioms make code faster.
- Some idioms make code more understandable.

#### Fit These Notes Into Context

- Use of iterators requires the ability to write small functions.
- If you have never written a small function for R, please review the lecture functions-1 before tackling this material.
- This lecture was once part of functions-1. In fact, it was the major motivation for functions-1, because I had to teach people how to write functions before using R apply statements.

# Clarity and Understandability

- Especially in the early years of R, people who used for loops were ridiculed and urged to use apply() instead.
- Some ridicule was justified because code based on for() often makes heavy use of '[' to access data, and that is a very slow operator.
- I have examples of silly/slow code using for()
- However, if you have only a few situations to loop through, there is not usually a substantial speedup by recoding from for() to apply() (see Chambers, *Software for Data Analysis*)
- On the other hand, for() loops, especially nested loops, are prone to user-error and miscalculations, and they will be more difficult to read.

# Bootstrapping is at the End

- Difficult to be sure bootstrapping should be included in this lecture
- It is included here because people who are frustrated with R's apply concepts are also usually frustrated with bootstrapping in R.
- Why this makes a difference: Efficiency! People who do bootstrapping in the literal, obvious way, are generally wasting memory and time.

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R has lots of ways to do things over and over

- for loop: process by "i" or by "element"
- apply: process rows and/or columns in a matrix
- lapply: process each element in a list
- sapply: attempts to simplify output from lapply
- replicate: shorthand for sapply for simple simulations
- mapply: for functions that need several arguments, separately drawn from separate vectors or lists

# for looping

- First, I initialize x1, then
- loop over elements to set their values

```
doubleMe <- function(input = 0){
    newval <- 2 * input
}
x1 <- vector(mode = "numeric", length = 57)
for(i in 1:57) {x1[i] <- doubleMe(i)}</pre>
```

integers i from 1 to 57 are sent to double me, results collect

 Note, it is not necessary to actually do this for loop in R, because R is vectorized.

```
x2 <- doubleMe(1:57)
all.equal(x1,x2)
```

[1] TRUE

• Using vectorized code is much faster.

# "apply()"

- useRs are urged to avoid "for loops" when possible
- Why? Accessing particular values with "[" (vector or matrix indexes) is SLOW. Better to exploit R's "vectorization"
- apply() is one of a family of functions that can replace a for loop.
- apply() takes a matrix, and does "the same FUN" to all of its rows or columns (or both).
- Definition: MARGIN=1 means "work row by row", MARGIN=2 means "column by column"

# Example of "apply()" With a Built-In FUN

- Given a matrix xyz with columns "x", "y", and "z"
- On the columns, MARGIN=2, apply the R "mean" function.

```
xyz <- matrix( rnorm(9), ncol=3)
xyz</pre>
```

× y z 0.39523051 -0.55518856 0.02325157

• If there is no "built in" function that does what you want, then you have to write your own.

# Write your own Function for apply

- Suppose you want the second-highest score from each column.
- Write a little function called "second()"

```
second <- function(acol = NULL){
    sort(acol)[2]
}
print(xyz)</pre>
```

	х	У	Z
[1,]	0.5855288	-0.4534972	0.6300986
[2,]	0.7094660	0.6058875	-0.2761841
[3,]	-0.1093033	-1.8179560	-0.2841597

apply(xyz, MARGIN = 2, FUN = second)

x	у	Z
0.5855288	-0.4534972	-0.2761841

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# Apply the normedEntropy function to rows

• The normedEntropy() function is presented in the lecture functions-1. I reproduce it for completeness here

```
divr <- function(p = 0){
    ifelse ( p > 0 & p < 1, -p * log2(p), 0)
}
entropy <- function(p){
    sum( divr(p) )
}
maximumEntropy <- function(N) - log2(1 / N)
normedEntropy <- function(x) entropy(x) /
    maximumEntropy(length(x))</pre>
```

• First, create a matrix in which the sum of each row is 1.0

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 Bootstrapping

### Apply the normedEntropy function to rows ...

```
xmat <- matrix(rmultinom(6, size = 20, prob = c
(1,2,3,4,5)), byrow = T, ncol = 5)
xmat <- prop.table(xmat, 1)
print(round(xmat,3))
```

	[,1]	[,2]	[,3]	[,4]	[,5]
[1,]	0.00	0.30	0.15	0.20	0.35
[2,]	0.20	0.15	0.20	0.20	0.25
[3,]	0.10	0.15	0.10	0.30	0.35
[4,]	0.10	0.00	0.15	0.40	0.35
[5,]	0.05	0.10	0.30	0.35	0.20
[6,]	0.10	0.05	0.30	0.25	0.30

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### Entropy for each row!

• apply normed Entropy to each Row with apply

apply(xmat, MARGIN = 1, FUN = normedEntropy)

[1] 0.8295351 0.9921503 0.9156704 0.7759110 0.8888583 0 .9003158

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"lapply()": Do same thing to all Elements of a List

 lapply() will take a list of things and apply a given function to each item, returning a new list. Generally,
 aNewlist <- lapply( some list ELIN - someEunction )</li>

aNewList <- lapply( someList, FUN = someFunction )

- someFunction MUST accept the elements from someList as the first argument
- Additional arguments to someFunction are allowed

# Example Use of lapply

• Create a list with 5 sets of random uniform normal variables

```
sampleList <- lapply(rep(1000,5), rnorm) sampleList [[1]][888]
```

[1] -0.3101479

#### • Same as

```
sampleList <- list() ## or <- vector(``list'', 5)
sampleList[[1]] <- rnorm(1000)
sampleList[[2]] <- rnorm(1000)
sampleList[[3]] <- rnorm(1000)
sampleList[[4]] <- rnorm(1000)
sampleList[[5]] <- rnorm(1000)</pre>
```

# Example Use of lapply

• Get the mean of sets 1 and 2 individually

mean(sampleList[[1]])

[1] 0.04081866

mean(sampleList[[2]])

[1] -0.02739241

• Grab means of all sets with lapply

```
(aNewList <- lapply(sampleList, mean))
```

```
[[1]]
[1] 0.04081866
[[2]]
[1] -0.02739241
[[3]]
[1] -0.0255273
```

# Why lapply, Not apply?

 Sometimes our "data" is not an even set of columns that fits in a data.frame or matrix

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### Why lapply, not apply?

```
for(i in 1:length(xlist)){
  cat("Given List")
  print(xlist[[i]])
  cat("Normed Entropy")
  print(round(elist[[i]],3))
  cat("\n")
 }
```

```
Given List [1] 1 1 1 2 3 3
Normed Entropy [1] 0.921
Given List [1] 3 2 5 2 5 2 1 6 2 4
Normed Entropy [1] 0.898
Given List [1] 101 101 100 101 100 99 101 100 100 102 100
102 100 99 100 101 100 101 100
Normed Entropy [1] 0.843
```

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# Example with additional arguments

• One NA wrecks mean (by default)

```
sampleList <- lapply( rep(1000,5), rnorm)
sampleList[[1]][77] <- NA
(aNewList <- lapply(sampleList, mean))</pre>
```

```
[[1]]
[1] NA
[[2]]
[1] -0.008354005
[[3]]
[1] -0.003276648
[[4]]
[1] -0.003438522
[[5]]
[1] 0.05110267
```

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# Example (cont.): Fix that Missing Value Problem

```
(aNewList <- lapply(sampleList, mean, na.rm = TRUE))
```

```
[[1]]
[1] -0.03336209
[[2]]
[1] -0.008354005
[[3]]
[1] -0.003276648
[[4]]
[1] -0.003438522
[[5]]
[1] 0.05110267
```

= nac

# Example: lapply to Simulate Regressions.

- The question:
  - Create 100 regression models from 100 data sets
  - Study the sampling distribution of the  $R^2$  statistic from those regressions.

# Step 1.

• The following generates 100 data frames in a list "mydatasets".

```
exs <- 10
exq <- 0.345
exstde <- 20
createOneDF <- function(run, s = NA, q = NA, stde = NA
    ){
    x <- 18 + 43*runif(1000)
    y <- s + q * x + rnorm(1000, mean = 0, sd = stde)
    mydf <- data.frame(run,x,y)
}
mydatasets <- lapply(1:100, createOneDF, exs, exq,
    exstde)
```

- Here the "list" is just a sequence 1,2,3,...
- lapply automatically gives each list element to function as first argument. (In this case, "run" number).

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# Step 2.

 Now apply a function to each data frame, make list "myregressions"

```
myregressions <- lapply(mydatasets, FUN = function(mydf) lm(y \sim x, data = mydf))
```

- Note: small functions can be written "inline"
- Could as well have written

```
calcReg <- function(adf = NULL){
  mod <- lm(y ~ x, data = adf)
  }
myregressions <- lapply(mydatasets, FUN = calcReg)</pre>
```

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# Take Stock of What We Have

• Each element in the list "mydatasets" really is a data frame:

head ( mydatasets [[33]])

	run	Х	у
1	33	41.47315	30.817774
2	33	48.78788	48.229489
3	33	31.71107	45.515414
4	33	50.28991	-22.129543
5	33	60.13310	33.632953
6	33	35.67771	9.532895

 Each element in "myregressions" really is a regression result object

```
myregressions [[33]]
```

Take Stock of What We Have ...

```
Call:

Im(formula = y \sim x, data = mydf)

Coefficients:

(Intercept) x

10.5261 0.3371
```

• Which can be summarized thus:

summary ( myregressions[[33]] )

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### Take Stock of What We Have ...

```
Call
lm(formula = y \sim x, data = mydf)
Residuals :
   Min 1Q Median
                            3Q
                                  Max
-56.643 - 11.595 0.873 12.462 57.854
Coefficients.
           Estimate Std. Error t value Pr(>|t|)
(Intercept) 10.52613 1.94869 5.402 8.26e-08 ***
            0.33713 0.04737 7.117 2.10e-12 ***
x
Signif. codes: 0 '***' 0.001 '**' 0.01 '*' 0.05 '.' 0.1 ' '
    1
Residual standard error: 18.79 on 998 degrees of freedom
Multiple R^2: 0.0483, Adjusted R^2: 0.04735
F-statistic: 50.66 on 1 and 998 DF, p-value: 2.101e-12
```

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Take Stock of What We Have ...

Note, the  $R^2$  value that we need is sitting there, in the middle of the summary output. We'll need that.

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# Step 3.

- Grab the  $R^2$  from each regression in the list.
- The estimate of the  $R^2$  is an element in the returned object from summary.
- One strategy: create an R list of summary objects

mysummaries <- lapply (myregressions, FUN= summary)

• Getting the  $R^2$  out of each one of those requires some tedious grabbing, such as

```
myrsq <- lapply(mysummaries, FUN = function(mr) {mr$
    r.square})
myrsq[1:5]</pre>
```

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Step 3. ...

[[1]] [1] 0.03758218		
[[2]] [1] 0.03746384		
[[3]] [1] 0.02569663		
[[4]] [1] 0.03390325		
[[5]] [1] 0.04059477		

```
myrsq <- unlist(myrsq)
str(myrsq)</pre>
```

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 Bootstrapping

Step 3. ...

num [1:100] 0.0376 0.0375 0.0257 0.0339 0.0406  $\ldots$ 

Iterators

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# Sapply will do that in one shot

- sapply is the "simplified apply", it attempts to convert a list into a vector or matrix.
- snoop through the regressions, grab the  $R^2$ .

```
myrsq <- sapply(mysummaries, FUN = function(mr) {mr$
    r.square})
mean(myrsq)</pre>
```

[1] 0.04510022

sd(myrsq)

[1] 0.01280801

median (myrsq)

[1] 0.04424352

## Everybody Still Loves Histograms



# Example: Balance in Logistic Regression

- Two years ago, I wondered (while auditing the categorical class), "what if we run a logistic regression comparing men and women and there are not very many men?"
- Write functions to
  - manufacture data
  - analyze data
  - summarize & plot data

Create Output Data: Need to convert real numbers to 0's and 1's

```
\eta "eta" is input, the proclivity to "vote democratic"
```

```
simLogit <- function(myeta){
    mypi <- exp(myeta) / (1 + exp(myeta)) ## SAME AS 1/(1+
        exp(-myeta))
    myunif <- runif(length(myeta))
    y <- ifelse(myunif < mypi, 1, 0)
}</pre>
```

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Example Use: Creates 1000 Observations

 $\begin{array}{l} N <- \ 1000 \\ A <- \ -1 \\ B <- \ 0.3 \\ x <- \ 1 \ + \ 10 \ * \ rnorm (N) \\ myeta <- \ A \ + \ B \ * \ x \\ y \ <- \ simLogit (myeta) \end{array}$ 

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### Illustration of Simulated Data



Iterators

### The Fitted Line from glm



Iterators

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### We are Interested in the Difference Between Two Groups



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### Now Automate That Process

- Manufacture data
- Run Regression
- Return row of estimates

```
simUnbalanced <- function(iter=0, parm){
    A <- parm$A; B<- parm$B; C<- parm$C; PrFem <- parm$PrFem
    sex <- ifelse(runif(N) < PrFem,0,1)
    myeta <- A + B * x + C * sex
    sex <- factor(sex, levels = c(0,1), labels = c("M","F"))
    y <- simLogit(myeta)
    myglm2 <- glm( y ~ x + sex, family = binomial)
    myglm2sum <- coef(summary(myglm2))
    est <- myglm2sum[3,]
}</pre>
```

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## Use sapply to run 1000 Regressions

Note: I'm combining the sapply result, along with "p", for record-keeping

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## Now Plan to Draw Some Figures

```
createFigs <- function(result){</pre>
  dat <- result [[1]]
 C <- result$parm$C
  PrFem <- result $parm $PrFem
 mybeta <- dat [1,]
  hrow1 <- hist(mybeta, breaks=50, plot=F)</pre>
  mybreaks <- hrow1$breaks
  breakMember <- cut(dat[1,], mybreaks)</pre>
  mypval < - dat[4]
  mysignif <- if else ((mypval < 0.05), 1, 0)
  df <- data.frame(mybeta, mypval, mysignif, breakMember)
  propsig <- by(df$mysignif, INDICES = list(df$breakMember)</pre>
      , mean, simplify = TRUE)
  mytrat < - dat[3]
  mycounts <- hrow1$counts
                                           (日) (同) (日) (日)
```

Now Plan to Draw Some Figures ...

```
plot(dat[1,], dat[4,], xlab = "beta estimate", ylab = "
    estimated p", cex = 0.7, main = paste("True Beta=",C,
        "Prop. Fem.=", PrFem))
gc <- c("gray98","gray70","gray50","gray40")
cut(propsig, breaks=c(-1,0.1,0.5,0.9,1.1))
catpropsig <- cut(propsig, breaks = c(-1,0.1,0.5,0.9,1.1))
, ordered = T, labels = c("0","lth","mth","1"))
barplot(hrow1$density, col = gc[as.numeric(catpropsig)],
        names = hrow1$mids)</pre>
```

For Balanced Data

For Balanced Data ...



# For Unbalanced Data

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# For Unbalanced Data ...



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# Final Cautionary Note

- At some point, this approach will start to "bog down" under the weight of memory usage and CPU delays
- I'd suggest re-designing so that we separately create the data frames and run all of the analysis on each separately
- That would allow us to 1) stay within memory limits and 2) parallelize the work across separate cores or computers (see the R parallel package).

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- I never used mapply for the first 5 years of using R
- Now I see need for it at least once per month
- The documentation may be difficult to understand, but once you appreciate the beauty of it, you will like it.

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# When is mapply needed

- You have several vectors or lists of the same length
- You want to take the first element from each and do something.
- Then take the second element from each and do something
- Don't write nested "for" loops, as users are often tempted to do.

Example of mapply usage in rockchalk package

- rockchalk has many functions that are doing the same thing over and over for subsets of data.
- Run the examples for the addLines() function, you should see it integrates plotSlopes() and plotPlane() by transferring information.

# Example use of mapply in rockchalk 1.8

- dataSplits is a collection of data frames. We want to do the plot for each with the correct colors, which are stored in linesFrom variables col and lty.
- The small function drawLine accepts 3 arguments, one from data, one from col, one from lty.

```
if (!missing(linesFrom)) {
   dataSplits <- split(linesFrom$newdata, f = linesFrom
        $newdata[[linesFrom$call[["modx"]]]])
   drawLine <- function(nd, mycol, mylty) {
        lines(trans3d(nd[[plotx1]], nd[[plotx2]], nd$fit,
            pmat=res), col = mycol, lwd = lflwd, lty =
            mylty)
   }
   mapply(drawLine, dataSplits, linesFrom$col,
        linesFrom$lty)</pre>
```

# Example use of mapply in rockchalk 1.8 ...

• Note we are free to name the variables inside drawLine however we want. That help keep our minds clear about whether we are talking about just one color or a vector of colors.

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Bootstrapping: Some "Do it Yourself" Work Is Required

- Many R functions require users to write little functions that do little things.
- In many cases (like lapply or apply), look for FUN as an argument.
- Sometimes no builtin-exists. useR must create!

# boot Function Requires a Special Function "statistic"

```
library (boot)
?boot
```

```
Bootstrap Resampling
Description:
Generate 'R' bootstrap replicates of a statistic applied to
   data.
Both parametric and nonparametric resampling are possible.
    . . .
boot(data, statistic, R, sim = ''ordinary'', stype = ''i'',
strata=rep(1, n), L = NULL, m = 0, weights = NULL,
ran.gen=function(d, p) d, mle = NULL, simple = FALSE, ...)
statistic: A function which when applied to data returns a
    vector
containing the statistic(s) of interest...
```

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# Bootstrap: Background Explanation

- Bootstrap: draw samples repeatedly and re-estimate  $\boldsymbol{\theta}$
- $\bullet$  Resulting values approximate a sampling distribution  $\theta$
- The "boot" package asks for a data frame and a special function "statistic". statistic must
  - accept a data frame as the first argument
  - accept an "index vector" as the second argument

# Don't Panic: This is Confusing to Everybody

Example usage

boot(data, statistic = yourFunction, R = 1000)

- boot will iterate 1000 times, and yourFunction will provide the statistic of interest.
- You write yourFunction to make required calculation.
- boot will tell yourFunction which lines to use in the data frame, *over-and-over*.

# The Median of a Poisson Distribution

• Suppose you have a sample from a Poisson Process:

```
samp <- rpois(20, lambda=3)</pre>
```

• And you calculate the median:

median(samp)

[1] 2.5

• How confident are you in that estimate of the median?

# Bootstrap Your Median

```
• Here is yourFunction:
```

```
calcMed <- function(dat, ind){
    median(dat[ind])
}</pre>
```

- dat[ind] has the effect of "pulling" rows that match "ind" from "dat"
- The boot function will send 1000 "case indexes" to your function.

```
library(boot) bpois <\!\!- boot(samp, calcMed, R = 1000) bpois
```

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Bootstrap Your Median ...

```
ORDINARY NONPARAMETRIC BOOTSTRAP
Call:
boot(data = samp, statistic = calcMed, R = 1000)
Bootstrap Statistics :
original bias std. error
t1* 2.5 0.076 0.6173371
```

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Let's plot that



## Why Do They Do It That Way?

- Your instinct is to do this the "simple" way
  - (Just) "Manually" draw new random samples of rows from a data frame.
  - But: Creating 1000s of "new" re-sampled data sets would "waste" (exhaust?) memory
  - Would be especially slow if separate data sets have to be copied between systems.
- More efficient to keep 1 data frame, but 1000's of vectors of row numbers.