

How to write an efficient script

Chin Yang Shapland

MRC Integrative Epidemiology Unit

6 June 2022

Introduction

This session will cover;

- Identifying bottlenecks
- ★ The apply family
- Parallelizing a script
- ₭ High performance computing

Identifying bottlenecks: system.time()

```
system.time({
  n < -1000
  r <- numeric(n)
  for(i in 1:n) {
    x \leftarrow rnorm(n)
    r[i] \leftarrow mean(x)
})
##
      user system elapsed
               0.00
                        0.11
##
      0.11
```

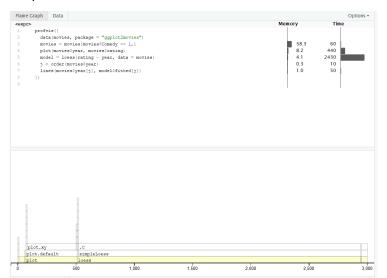
user is the time taken to excute the code, system is time for the system to open and close files, and elapsed is the stopwatch time, i.e. total time code took from start to finish.

Identifying bottlenecks: The R Profiler

```
library(profvis)
library(bench)
library(ggplot2movies)
profvis({
  data(movies, package = "ggplot2movies")
  movies = movies[movies$Comedy == 1,]
  plot(movies$year, movies$rating)
  model = loess(rating ~ year, data = movies)
  j = order(movies$year)
  lines(movies$year[j], model$fitted[j])
})
```

Output

Outputs a tab called "Profile1";



The apply family

are faster and safer than loops. We will cover

- apply()
- sapply()

but there are also others,

- vapply()
- mapply()
- rapply()
- ★ tapply()

apply()

```
str(apply)
## function (X, MARGIN, FUN, ..., simplify = TRUE)
```

X is the input, MARGIN should be row (1) or column (2), or both (c(1,2)), FUN is the function. Depending what the input is, it can return a vector, list, matrix or array.

apply() example

```
data_fun < -matrix(c(1:4,1:4,1:4,1:4),4,4)
data fun
## [,1] [,2] [,3] [.4]
## [1,] 1 1 1 1
## [2,] 2 2 2 2
## [3,] 3 3 3 3
## [4,] 4 4 4 4
apply(data_fun,2,sum)
```

[1] 10 10 10 10

apply(data_fun,2,mean)

[1] 2.5 2.5 2.5 2.5

apply() example with function

```
data_fun < -matrix(c(1:4,1:4,1:4,1:4),4,4)
data fun
```

[,1] [,2] [,3] [.4] ## [1,] 1 1 1 **##** [2,] 2 2 2 2 ## [3,] 3 3 3 3 ## [4,] 4 4 4 4

apply(data_fun, 1:2, function(x) x/2)

[.1] [.2] [.3] [.4] ## [1.] 0.5 0.5 0.5 0.5 ## [2,] 1.0 1.0 1.0 1.0 ## [3,] 1.5 1.5 1.5 1.5 ## [4,] 2.0 2.0 2.0 2.0

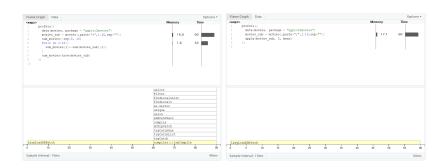
apply vs loops

```
# Using for loop
profvis({
  data(movies, package = "ggplot2movies")
  movies_sub = movies[,paste("r",1:10,sep="")]
  sum_movies<-rep(0, 10)</pre>
  for(i in 1:10){
    sum movies[i] <-sum(movies sub[,i])</pre>
  }
  sum_movies/nrow(movies_sub)
})
```

apply vs loops

```
# Using `apply()` function;
profvis({
  data(movies, package = "ggplot2movies")
  movies_sub = movies[,paste("r",1:10,sep="")]
  apply(movies_sub, 2, mean)
})
```

apply vs loops



Inputs matters

Previous examples used matrices how about data frame?

```
data df<-data.frame(1:4,1:4,1:4,1:4)
str(data df)
## 'data.frame': 4 obs. of 4 variables:
## $ X1.4 : int 1 2 3 4
## $ X1.4.1: int 1 2 3 4
## $ X1.4.2: int 1 2 3 4
## $ X1.4.3: int 1 2 3 4
apply(data_df, 2, mean)
## X1.4 X1.4.1 X1.4.2 X1.4.3
## 2.5 2.5 2.5 2.5
```

Inputs matters

```
data df$Day<-as.factor(1:4)
str(data df)
## 'data.frame': 4 obs. of 5 variables:
## $ X1.4 : int 1 2 3 4
## $ X1.4.1: int 1 2 3 4
## $ X1.4.2: int 1 2 3 4
## $ X1.4.3: int 1 2 3 4
   $ Day : Factor w/ 4 levels "1", "2", "3", "4": 1 2 3 4
##
apply(data_df, 2, mean)
    X1.4 X1.4.1 X1.4.2 X1.4.3
                               Day
##
      NA
            NA NA NA
                               NA
```

... options

... within the function have multiple options but the most option is to use for handling missing data

```
data_miss<-matrix(c(1:4,c(1:3,NA),1:4,1:4),4,4)
data miss
## [,1] [,2] [,3] [,4]
## [1,] 1 1 1 1
## [2,] 2 2 2 2
## [3,] 3 3 3 3
## [4,] 4 NA 4 4
apply(data_miss, 2, mean)
## [1] 2.5 NA 2.5 2.5
apply(data miss, 2, mean, na.rm=T)
## [1] 2.5 2.0 2.5 2.5
```

lapply()

```
str(lapply)
## function (X, FUN, ...)
```

X is the input, FUN is the function. Only returns list object.

lapply() example

```
list fun<-list(a = 1:5, b = 10:15, c = 21:25)
list_fun
## $a
## [1] 1 2 3 4 5
##
## $b
## [1] 10 11 12 13 14 15
##
## $c
## [1] 21 22 23 24 25
```

lapply() example

```
lapply(list_fun,sum)
## $a
## [1] 15
##
## $b
## [1] 75
##
## $c
## [1] 115
```

lapply() example

```
lapply(list_fun,mean)
## $a
## [1] 3
##
## $b
## [1] 12.5
##
## $c
## [1] 23
```

lapply() example with function

```
lapply(list_fun, function(x) x/2)
## $a
## [1] 0.5 1.0 1.5 2.0 2.5
##
## $b
## [1] 5.0 5.5 6.0 6.5 7.0 7.5
##
## $c
## [1] 10.5 11.0 11.5 12.0 12.5
```

lapply() example with vectors

```
Random<-c("these", "are", "random", "example")
lapply(Random, nchar)
## [[1]]
## [1] 5
##
## [[2]]
## [1] 3
##
## [[3]]
## [1] 6
##
## [[4]]
## [1] 7
```

sapply()

This is a simplified form of lapply()

```
str(sapply)
## function (X, FUN, ..., simplify = TRUE, USE.NAMES = TRUE)
```

X is the input, FUN is the function. If simplify is changed to F then it becomes lapply() and returns only list object. Otherwise it also outputs vector, matrix, array or list.

sapply() example

```
list_fun < -list(a = 1:5, b = 10:15, c = 21:25)
list fun
## $a
## [1] 1 2 3 4 5
##
## $b
## [1] 10 11 12 13 14 15
##
```

\$c

sapply(list_fun,sum)

[1] 21 22 23 24 25

a b c ## 15 75 115

sapply() example

```
list_fun < -list(a = 1:5, b = 10:15, c = 21:25)
list fun
## $a
## [1] 1 2 3 4 5
##
## $b
## [1] 10 11 12 13 14 15
##
```

\$c

sapply(list_fun,mean) ## a b c ## 3.0 12.5 23.0

[1] 21 22 23 24 25

sapply() example

```
sapply(list fun, function(x) x/2)
## $a
## [1] 0.5 1.0 1.5 2.0 2.5
##
## $b
## [1] 5.0 5.5 6.0 6.5 7.0 7.5
##
## $c
## [1] 10.5 11.0 11.5 12.0 12.5
```

Notice that this is given as a list rather than a vector, this is because each element have different lengths

Others

vapply()

similar to sapply but offers a third argument which sets a template for the output.

mapply()

multivariate version to sapply function, but it applies functions to the argument one by one.

rapply()

Similar to lapply but you can specify the structure of the output

tapply()

applies a function to numeric data distributed across various categories.

Parallelizing a script: when

When to parallelize?

"its important to evaluate the computational efficiency of requests, and work to ensure that additional compute resources brought to bear will pay off in terms of increased work being done."

Parallelizing a script: how

When parallelizing jobs, one can:

- Use the multiple cores on a local computer through mclapply
- Use multiple processors on local (and remote) machines using makeCluster and clusterApply (Note this approach, one has to manually copy data and code to each cluster member using clusterExport. This is extra work, but sometimes gaining access to a large cluster is worth it.)

Parallel package: mclapply

```
library(parallel)
library(MASS)
starts \leftarrow rep(100, 40)
fx <- function(nstart) kmeans(Boston, 4, nstart=nstart)</pre>
numCores <- detectCores()</pre>
numCores
## [1] 8
system.time(
  results <- lapply(starts, fx)
##
      user system elapsed
    1.06 0.03 1.11
##
```

Parallel package: mclapply

```
system.time(
  results <- mclapply(starts, fx, mc.cores = numCores)
)

## user system elapsed
## 0.801 0.178 0.367</pre>
```

foreach package: %do% and %dopar%

```
for (i in 1:3) {
   print(sqrt(i))
}
## [1] 1
## [1] 1.414214
## [1] 1.732051
```

```
In foreach package %do% evaluates the expression sequentially,
library(foreach)
foreach (i=1:3, .combine=c) %do% {
   sqrt(i)
}
## [1] 1.000000 1.414214 1.732051
```

foreach package with doParallel

while $\dopar\%$ evaluates it in parallel, but $\dopar\%$ has to work with \dopar allel package

```
library(foreach)
library(doParallel)
```

Use multicore, set to the number of our cores

```
registerDoParallel(numCores)
foreach (i=1:3, .combine=c) %dopar% { sqrt(i) }
## [1] 1.000000 1.414214 1.732051
```

%do%

Using the iris data set to do a parallel bootstrap from the doParallel vignette

```
x \leftarrow iris[which(iris[,5] != "setosa"), c(1,5)]
trials <- 10000
system.time({
  r <- foreach(icount(trials), .combine=rbind) %do% {
    ind <- sample(100, 100, replace=TRUE)</pre>
    result1 \leftarrow glm(x[ind,2]~x[ind,1],
                    family=binomial(logit))
    coefficients(result1)
})
##
      user system elapsed
     20.04 0.00 20.06
##
```

%dopar%

```
system.time({
  r <- foreach(icount(trials), .combine=rbind) %dopar% {
    ind <- sample(100, 100, replace=TRUE)</pre>
    result1 \leftarrow glm(x[ind,2]~x[ind,1],
                    family=binomial(logit))
    coefficients(result1)
})
##
      user system elapsed
      2.98 0.47 6.47
##
```

Outputs

To simplify output, foreach has the .combine parameter that can simplify return values

```
# Default: Return a list
foreach (i=1:3) %dopar% { sqrt(i) }
## [[1]]
## [1] 1
##
## [[2]]
## [1] 1.414214
##
## [[3]]
## [1] 1.732051
```

Outputs

```
# Return a vector
foreach (i=1:3, .combine=c) %dopar% { sqrt(i) }
## [1] 1.000000 1.414214 1.732051
# Return a data frame
foreach (i=1:3, .combine=rbind) %dopar% { sqrt(i) }
                Γ.17
##
## result.1 1.000000
## result.2 1.414214
## result.3 1.732051
```

Further tips parallel computing

★ To clear up and stop parallel use;

```
stopImplicitCluster()
```

★ To set the cluster seed if you want reproducible results, don't use use set.seed(), use clusterSetRNGStream() instead

High performance computing

When all else fails running your script on a high powered server:

```
BlueCrystal Phase 4
```

```
#!/bin/bash
#SBATCH -- job-name=job_name
#SBATCH --output=job_name
#SBATCH --error=job name
#SBATCH --nodes=1
#SBATCH --ntasks-per-node=16
\#SBATCH --time=2:0:0
module load languages/r/4.1.0
R CMD BATCH /your dir/your Rscript.R
```

Check your shell script!

Use **cat -v** command within BlueCrystal, as sometime when you write your shell script(.sh) in RStudio or notepad it leaves a bunch of none printable characters like this;

```
#!/bin/bash^M
^ M
#SBATCH -- job-name=job_name^M
#SBATCH --output=job_name^M
#SBATCH --error=job name^M
\#SBATCH --nodes=1^M
#SBATCH --ntasks-per-node=16^M
#SBATCH --time=2:0:0^M
^ M
module load languages/r/4.1.0^M
^M
Rscript /your dir/your Rscript.R^M
```

Solution: Check your shell script!

To remove these characters, use a text editor in BlueCrystal, such nano, emacs or vi (very user friendly editors).

References

Help from HPC https://www.bristol.ac.uk/acrc/high-performance-computing/hpc-documentation-support-and-training/

Top 5 tips for efficient performance by Colin Gillespie https://csgillespie.github.io/efficientR/performance.html#top-5-tips-for-efficient-performance

Quick Intro to Parallel Computing in R by Matt Jones https://nceas.github.io/oss-lessons/parallel-computing-in-r/parallel-computing-in-r.html

Join "code-clinic"! Email Louise Millard