R profiling

Stat 580: Statistical Computing

• Theme: Black - White

Printable version

References

- "Advanced R", by Hadley Wickham.
- "Writing R Extensions", by R Core Team.

Performance

- R is not a fast language
 - easy for you (to do data analysis and statistics)
 - not easy for the computer
 - improve speed of your code:
 - profiling
 - R's C interface
 - Rcpp

Why is R slow?

- R as both language and implementation of that language
 - R-language: defines what R code means and how it should work
 - implementation: reads R code and computes a result.
 - the most popular implementation is the one from r-project.org. (GNU-R)
- Both language and implementation (GNU-R) limitations can be the reasons of slowness.
 - We will look into a few reasons from the language performance perspectives.
- Poorly written R code

Microbenchmarking

- For investigation of (language and implementation) performance, we need measurement of performance of a very small piece of code
 - take microseconds or naoseconds
 - for deeper understanding
 - not very practical for optimizing your code

Microbenchmarking

```
library(microbenchmark)

x <- runif(100)
microbenchmark(sqrt(x), x^0.5)</pre>
```

- microbenchmark()
 - more accurate replacement of system.time(replicate(1000, expr))
 expression
 - randomized ordering (the default)
 - warm-up iterations performed before the actual benchmark

```
Unit: nanoseconds

expr min lq mean median uq max neval

sqrt(x) 839 890.0 1865.58 979.5 2666.5 35456 100

x^0.5 4837 4956.5 5497.08 5070.5 5176.5 14552 100
```

- sqrt() is a lot faster! Why?
- Should I replace every occurrence of x^0.5 by sqrt(x)?
 - What is the unit?
 - set argument unit=eps:

```
Unit: evaluations per second
expr min lq mean median uq max neval
sqrt(x) 217627.86 1281293.6 1272596.7 1456671.2 1579778.8 1683501.7 100
x^0.5 71797.82 253132.6 245315.7 258431.5 265287.3 270270.3 100
```

Extreme dynamism

- R is extremely dynamic, you can modify almost anything
 - after it is created
 - outside of the local environment (e.g., <<-)
 - to completely different object (atomic vector to function)

```
o x <- c(1,2); x <- function(){return(c(3,4))}</pre>
```

■ more...

```
x <- 0L

for (i in 1:1e6) {

 x <- x + 1

}
```

- R doesn't know that x is always an integer.
- R has to look for the right + method in every iteration.

Extreme dynamism

- Both good and bad:
 - Good (to us): flexibility, minimal upfront planning
 - Recall: in C, we have to plan ahead by declaring the variables as the right types.
 - Bad (to the computer): difficulty to predict exactly what will happen with a given function call
 - difficult for an interpreter or compiler to make an optimization
 - if an interpreter can't predict what's going to happen, it has to consider many options.

Name lookup with mutable environments

- Scoping rules allow the same name to be used for different objects.
- R uses lexical (static) scoping (vs. dynamic scoping):
 - if an "unknown" variable is used in a function, its value are searched for in the environment in which the function was defined.

```
a = 1
f <- function(){
  return(a)
}

g <- function(){
  a = 2
  f()
}</pre>
```

```
a <- 1
f <- function() {
    g <- function() {
        print(a)
        assign("a", 2, envir = parent.frame())
        print(a)
        a <- 3
        print(a)
    }
    print(a)
    }
    print(a)
    g()
    print(a)
}</pre>
```

Name lookup with mutable environments

- Name lookup has to be done each time (partly due to extreme dynamism)
- Even worse, almost every operation is a lexically scoped function call.

```
■ e.g. +, -
```

- e.g. (, {
- Since these functions are in global environment, R has to look through every environment in the search path.
- R asks: Who knows someone wouldn't change + to -?
 - extreme dynamism: R has to worry that these functions are defined (in environments) on the search path

```
f <- function(x){
  (2 * x) ^ 2
}</pre>
```

```
random env <- function(parent = globalenv()) {</pre>
 letter list <- setNames(as.list(runif(26)), LETTERS)</pre>
  list2env(letter list, envir = new.env(parent = parent))
set env <- function(f, e) {</pre>
  environment(f) <- e
 f
f2 <- set env(f, random env())</pre>
f3 <- set env(f, random env(environment(f2)))</pre>
f4 <- set env(f, random env(environment(f3)))</pre>
microbenchmark(
 f(1),
 f2(1),
 f3(1),
 f4(1),
 times = 10000
```

Lazy evaluation overhead

R uses lazy evaluation:

function arguments are evaluated lazily - evaluated if they're actually used.

```
f <- function(x) {
1+1
}
f(cat("hello!\n"))</pre>
```

R uses promise object to contain the expression and environment needed to compute the result.

- overhead of creating such objects
- recall that gcc compiler (with -Wall) will warn you if you have unused variables

```
f0 <- function() NULL
f1 <- function(a = 1) NULL
f2 <- function(a = 1, b = 1) NULL
f3 <- function(a = 1, b = 2, c = 3) NULL
f4 <- function(a = 1, b = 2, c = 4, d = 4) NULL
f5 <- function(a = 1, b = 2, c = 4, d = 4, e = 5) NULL
microbenchmark(f0(), f1(), f2(), f3(), f4(), f5(), times = 10000)</pre>
```

Profiling your code

- Before you get your hands dirty to optimize your code, think about:
 - which part of the code is the bottleneck
 - Come up with solutions to optimize
 - which level: R (e.g., vectorization), lower level (e.g. C, Fortran)?
 - algorithm (approximations are sometimes allowed.)
 - Time investment vs. efficiency gain
 - choose a solution
 - do nothing

Profiling for speed

- To detect which part of the code is the bottleneck, we use profiler.
- R uses sampling profiler:
 - it checks which function is being used at fixed time intervals (e.g. every 20 msecs)
 - stochastic
 - variability in results due to the intervals between sampling
 - Rprof() is the profiler that comes with R
- Limitation: R profiling does not extend to C code, primitive functions or byte code compiled code.

```
library(lineprof) # for the use of pause
f <- function() {</pre>
 pause(0.1)
 g()
 h()
g <- function() {</pre>
 pause(0.1)
 h()
h <- function() {</pre>
 pause(0.1)
tmp <- tempfile()</pre>
Rprof(tmp, interval = 0.1)
f()
Rprof(NULL)
summaryRprof(tmp)
unlink(tmp)
```

Using lineprof package

```
install.packages("devtools")
devtools::install_github("hadley/lineprof")
```

- the fundamental unit of analysis in lineprof() is a line of code
- note that a line of code can contain multiple function calls
- but it's easier to understand the context

```
1 <- lineprof(f())

library(shiny)
shine(1)</pre>
```

Improving speed

Simple strategies to improve speed of R code without resorting to lower level implementation:

- Vectorization
- Avoid growing an object
- Byte code compilation
- Parallelization

Vectorization

- use vector (matrix, array) rather than scalar as "your working object"
 - to make use of the vectorized function
 - the loops in a vectorized function are written in low-level language instead of R
 - you should find the existing R function that is implemented in low-level language and most closely applies to your problem.
- replace loops by vector/matrix operations (find a matrix algebra version)
- use rowSums(), colSums(), rowMeans(), and colMeans() instead of apply (if possible)
- vectorized subsetting (e.g. x[is.na(x)] <- 0)
- Other useful vectorized functions: cut(), findInverval(), cumsum(), diff()

Avoid growing an object

- Functions like c(), append(), cbind(), rbind() or paste() can be used to build a bigger object on top of an old object.
 - R must first allocate space for the new object and then copy the old one to the new location.
 - hurts the performance, especially using in a loop.

```
x <- NULL
for (i in (1:10)){
x <- c(x, i)
}</pre>
```

```
random_string <- function() {
   paste(sample(letters, 50, replace = TRUE), collapse = "")
}
strings100 <- replicate(100, random_string())

collapse <- function(xs) {
   out <- ""
   for (x in xs) {
      out <- paste0(out, x)
   }
   out
}

microbenchmark(
   collapse(strings100),
   paste(strings100, collapse = "")
)</pre>
```

Bytecode compilation

- You can use bytecode compiler to improve the speed of the R code.
 - compile your R code into bytecode (more optimized because it is closer to the machine instructions)
 - fast and easy
 - considerable performance gain (usually 5-10%) (for the amount of effort you spend)

```
# old R version of lapply
la1 <- function(X, FUN, ...) {</pre>
   FUN <- match.fun(FUN)</pre>
    if (!is.list(X))
   X <- as.list(X)</pre>
   rval <- vector("list", length(X))</pre>
   for(i in seq(along = X))
   rval[i] <- list(FUN(X[[i]], ...))
    names(rval) <- names(X) # keep `names' !</pre>
    return(rval)
la1c <- compiler::cmpfun(la1)</pre>
x \leftarrow list(1:10, letters, c(F, T), NULL)
microbenchmark(
 la1(x, is.null),
 la1c(x, is.null),
 lapply(x, is.null)
```

Parallelization via foreach package

- It is easy to implement a type of parallel computing called "embarassingly parallel" in R
 - can be divided into small problems that can be solved independently
 - does not require exchange of information between workers
- You can use foreach package to implement a for loop using parallel computing
 - requires a "parallel backend" to "define the workers"
 - doParallel package provides the parallel backend

```
library(doParallel)

cl <- makeCluster(3)
registerDoParallel(cl)
foreach(i=1:10) %dopar% {
   max(svd(matrix(rnorm(100000), nr=100))$d)
}
stopCluster(cl)</pre>
```