## Zicheng Huang, STAT243 PS4

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```
library(microbenchmark)
library(ggplot2)
```

1(a) According to the result from .Internal(inspect), we can see that only one copy of the vector 1:10 is created, the two local variables input and data are pointing to the same copy of the vector 1:10 since their addresses are the same.

```
x <- 1:10
f <- function(input){
   data <- input
    # addresses of data and input
    .Internal(inspect(data))
   .Internal(inspect(input))
   g <- function(param) return(param * data)
   return(g)
}
myFun <- f(x)

## @0x000000001852caa8 13 INTSXP g0c4 [NAM(2)] (len=10, tl=0) 1,2,3,4,5,...
## @0x0000000001852caa8 13 INTSXP g0c4 [NAM(2)] (len=10, tl=0) 1,2,3,4,5,...</pre>
```

1(b) We make 'x' a large vector with size of around 40000 bytes. In part (a) we argue that only one copy of 'x' is made, we expect to see the number of bytes that store the information in the closure to be close to the size of 'x'. However, the actual size of the serialized 'myFun' is around 80000 bytes, twice of what we expect. This difference is due to the fact that under the function frame of myFun there exist the body text of the function, the variable 'input', and the variable 'data'. Even though both 'input' and 'data' are pointing to 'x', when R saves 'myFun' what happens is that 'input' and 'data' are saved seperated as two individual objects each takes up around 40000 bytes of size. Thus, the size of 'myFun' is 80000 bytes which is twice the size of what we expected.

```
x <- 1:10000
# size of x
length(serialize(x, NULL))
## [1] 40022

f <- function(input){
   data <- input
   g <- function(param) return(param * data)
   return(g)
}
myFun <- f(x)
# size of myFun
length(serialize(myFun, NULL))
## [1] 86993</pre>
```

1(c) When 'myFun' is define by f(x), f(x) is not evaluated due to the lazy evaluation, so the user supplied argument 'data' in the function frame of 'f' does not take in the values in 'x'. When myFun(3) is called, we see that what is actually being called is g(3). When g(3) is evaluated, g will look for a variable 'data' to return (3 \* data), but since no 'data' is defined in the function frame of 'g', it will go to the enclosing environment, function frame of 'f', to look for 'data'. Due to the lazy evaluation and the fact that 'x' is removed, g cannot find 'data' in the function frame of 'f' as well. Since there exists no 'data' in the global environment, myFun(3) will return error.

```
x <- 1:10
f <- function(data){
    g <- function(param) return(param * data)
    return(g)
}
myFun <- f(x)
rm(x)
data <- 100
myFun(3)
## Error in myFun(3): object 'x' not found</pre>
```

 $\mathbf{1}(\mathbf{d})$  Since we prevented the lazy evaluation from happening using force(), when we define myFun by passing in f(x), f(x) will be evaluated in a sense that user supplied argument 'data' will take in variable of 'x'. In this way, even 'x' is removed, when myFun(3) is called, g(3) will be able to find a local vaiable 'data' in the function frame of f to return (param \* data), thus preventing the error from happening again. The resulting serialized closure has size around 40000 bytes.

```
x <- 1:10000
f <- function(data){
   force(data) # add this line to prevent the lazy evaluation from happening
   g <- function(param) return(param * data)
   return(g)
}
myFun <- f(x)
rm(x)
head(myFun(3))
## [1] 3 6 9 12 15 18
# size of modified 'myFun'
length(serialize(myFun, NULL))
## [1] 47349</pre>
```

**2(a)** According to the results from .Internal(inspect) in a plain R session, we can see that the address of the two lists, before and after modification, are the same, so no new lists are created when the change is made.

```
lst <- list(c(1, 2),c(3, 4))
.Internal(inspect(lst))
lst[[1]][1] <- 8
.Internal(inspect(lst))

## Results shown in plain R session

## @Ox000000000399b3a8 19 VECSXP gOc2 [NAM(1)] (len=2, tl=0)
## @Ox000000000399b418 14 REALSXP gOc2 [] (len=2, tl=0) 1,2
## @Ox000000000399b3e0 14 REALSXP gOc2 [] (len=2, tl=0) 3,4</pre>
```

```
## @0x00000000399b3a8 19 VECSXP g0c2 [NAM(1)] (len=2, tl=0)
## @0x00000000399b418 14 REALSXP g0c2 [] (len=2, tl=0) 8,2
## @0x000000000399b3e0 14 REALSXP g0c2 [] (len=2, tl=0) 3,4
```

**2(b)** After making a copy of the original list 'lst', we see that there is no copy-on-change going on since the address of the two lists are the same.

```
lst = list(c(1, 2),c(3, 4))
lstCopy <- lst
.Internal(inspect(lst))

## @0x0000000018301a10 19 VECSXP g0c2 [NAM(2)] (len=2, tl=0)
## @0x000000001821b170 14 REALSXP g0c2 [] (len=2, tl=0) 1,2
## @0x000000001821b1a8 14 REALSXP g0c2 [] (len=2, tl=0) 3,4

.Internal(inspect(lstCopy))

## @0x0000000018301a10 19 VECSXP g0c2 [NAM(2)] (len=2, tl=0)
## @0x000000001821b170 14 REALSXP g0c2 [] (len=2, tl=0) 1,2
## @0x0000000001821b1a8 14 REALSXP g0c2 [] (len=2, tl=0) 3,4</pre>
```

After modification is made to one of the vectors in one of the lists, a copy of the entire list is made in a sense that the entire vector of addresses (or vector of pointers) that stored the two sub-vectores contained in the original list is copied. The sub-vector that is about to be modified is also copied in order to make the modification.

```
lst = list(c(1, 2),c(3, 4))
.Internal(inspect(lst))

## @0x00000001852bf10 19 VECSXP g0c2 [NAM(2)] (len=2, tl=0)

## @0x000000001852bea0 14 REALSXP g0c2 [] (len=2, tl=0) 1,2

## @0x000000001852bed8 14 REALSXP g0c2 [] (len=2, tl=0) 3,4

lst[[1]][1] <- 8
.Internal(inspect(lst))

## @0x000000018477f80 19 VECSXP g0c2 [NAM(1)] (len=2, tl=0)

## @0x0000000018477fb8 14 REALSXP g0c2 [] (len=2, tl=0) 8,2

## @0x0000000001852bed8 14 REALSXP g0c2 [NAM(2)] (len=2, tl=0) 3,4</pre>
```

**2(c)** lstC is a copy of the original list.

```
lst1 <- list(1, 2)</pre>
lst2 <- list(3, 4)</pre>
lst <- list(lst1, lst2)</pre>
lstC <- lst</pre>
.Internal(inspect(lst))
## @0x0000000181aa328 19 VECSXP g0c2 [NAM(2)] (len=2, t1=0)
##
     @0x0000000184712a0 19 VECSXP g0c2 [NAM(2)] (len=2, tl=0)
##
       @0x0000000177c5fc0 14 REALSXP g0c1 [NAM(2)] (len=1, tl=0) 1
##
       @0x0000000177c6080 14 REALSXP g0c1 [NAM(2)] (len=1, tl=0) 2
##
     @0x00000001813cd70 19 VECSXP g0c2 [NAM(2)] (len=2, t1=0)
##
       @0x0000000177c6110 14 REALSXP g0c1 [NAM(2)] (len=1, tl=0) 3
       @0x0000000177c6170 14 REALSXP g0c1 [NAM(2)] (len=1, tl=0) 4
##
```

```
.Internal(inspect(lstC))
## @0x0000000181aa328 19 VECSXP g0c2 [NAM(2)] (len=2, tl=0)
## @0x00000000184712a0 19 VECSXP g0c2 [NAM(2)] (len=2, tl=0)
## @0x00000000177c5fc0 14 REALSXP g0c1 [NAM(2)] (len=1, tl=0) 1
## @0x00000000177c6080 14 REALSXP g0c1 [NAM(2)] (len=1, tl=0) 2
## @0x000000001813cd70 19 VECSXP g0c2 [NAM(2)] (len=2, tl=0)
## @0x00000000177c6110 14 REALSXP g0c1 [NAM(2)] (len=1, tl=0) 3
## @0x00000000177c6170 14 REALSXP g0c1 [NAM(2)] (len=1, tl=0) 4
```

After adding an element to lstC, the vectors that stored the addresses of the two original sublists are copied, but the address of each of the original sublist itself is not copied. The data in the sublists stored in the two uncopied addresses are shared.

```
lstC[[3]] <- list(5, 6)</pre>
.Internal(inspect(lst))
## @0x0000000181aa328 19 VECSXP g0c2 [NAM(2)] (len=2, t1=0)
##
     @0x0000000184712a0 19 VECSXP g0c2 [NAM(2)] (len=2, tl=0)
       @0x0000000177c5fc0 14 REALSXP g0c1 [NAM(2)] (len=1, tl=0) 1
##
       @0x0000000177c6080 14 REALSXP g0c1 [NAM(2)] (len=1, t1=0) 2
##
     @0x00000001813cd70 19 VECSXP g0c2 [NAM(2)] (len=2, t1=0)
##
##
       @0x0000000177c6110 14 REALSXP g0c1 [NAM(2)] (len=1, tl=0) 3
       @0x0000000177c6170 14 REALSXP g0c1 [NAM(2)] (len=1, tl=0) 4
##
.Internal(inspect(lstC))
## @0x0000000194f2560 19 VECSXP g0c3 [NAM(1)] (len=3, tl=0)
##
     @0x0000000184712a0 19 VECSXP g0c2 [NAM(2)] (len=2, tl=0)
##
       @0x0000000177c5fc0 14 REALSXP g0c1 [NAM(2)] (len=1, tl=0) 1
       @0x0000000177c6080 14 REALSXP g0c1 [NAM(2)] (len=1, tl=0) 2
##
##
     @0x00000001813cd70 19 VECSXP g0c2 [NAM(2)] (len=2, t1=0)
##
       @0x0000000177c6110 14 REALSXP g0c1 [NAM(2)] (len=1, tl=0) 3
##
       @0x0000000177c6170 14 REALSXP g0c1 [NAM(2)] (len=1, tl=0) 4
##
     @0x0000000194f1cc8 19 VECSXP g0c2 [NAM(2)] (len=2, tl=0)
       @0x00000001921cf78 14 REALSXP g0c1 [NAM(2)] (len=1, tl=0) 5
##
       @0x00000001921cf48 14 REALSXP g0c1 [NAM(2)] (len=1, tl=0) 6
```

2(d) The result from object.size() show that the size of the object 'tmp' is 160 Mb, however results from gc() shows that only 80 Mb of memeory is being used. Since 'tmp' is the list which contains two elements. Both elements are 'x' so they point to the same vector of size 80 Mb. The result from object.size() shows the size of 'tmp' being 160 Mb since it contains two 'x', while gc() reports the amount of memory being used to be 80 Mb since what is using up the memory is just the object 'x' in that 'tmp' is built by 'x'.

```
gc()
## used (Mb) gc trigger (Mb) max used (Mb)
## Ncells 557375 29.8   940480 50.3   940480 50.3
## Vcells 968782 7.4   1926429 14.7 1334875 10.2

tmp <- list()
x <- rnorm(1e7)
tmp[[1]] <- x
tmp[[2]] <- x
.Internal(inspect(tmp))</pre>
```

```
## @0x0000000019506668 19 VECSXP g0c2 [NAM(1)] (len=2, t1=0)
## @0x00007ff5facb0010 14 REALSXP g0c7 [NAM(2)] (len=10000000, t1=0) -0.133067,1.57412,-0.762538,-1.92128
## @0x00007ff5facb0010 14 REALSXP g0c7 [NAM(2)] (len=10000000, t1=0) -0.133067,1.57412,-0.762538,-1.92128

object.size(tmp)
## 160000136 bytes

gc()
## used (Mb) gc trigger (Mb) max used (Mb)
## Ncells 558111 29.9 940480 50.3 940480 50.3
## Vcells 10970639 83.7 16197621 123.6 11027723 84.2
```

3 In order to increase the efficiency in this chunk of codes, we replace the three nested for loops with a matrix multiplication which will do what is essentially the same thing. In addition, we pass in the result of the calculation of 'A\*q[,z]' to a variable so that this calculation will not be repeated twice in the second for loop.

```
load('ps4prob3.Rda') # should have A, n, K
11 <- function(Theta, A) {</pre>
  sum.ind <- which(A==1, arr.ind=T)</pre>
 logLik <- sum(log(Theta[sum.ind])) - sum(Theta)</pre>
 return(logLik)
oneUpdate <- function(A, n, K, theta.old, thresh = 0.1) {
  theta.old1 <- theta.old
  Theta.old <- theta.old %*% t(theta.old)
 L.old <- 11(Theta.old, A)
 q \leftarrow array(0, dim = c(n, n, K))
  # compress the three nested for loops to one for loop doing what is essentially
  # the matrix multiplication
  for (i in 1:K) {
    q[ , , i] <- theta.old[, i] %*% t(theta.old[, i]) / Theta.old
 theta.new <- theta.old
  for (z in 1:K) {
    # create a variable B to prevent from calculating A*q[,,z] twice in every
    # for loop for K times
   B = A*q[,,z]
    theta.new[,z] <- rowSums(B)/sqrt(sum(B))</pre>
 Theta.new <- theta.new %*% t(theta.new)
 L.new <- ll(Theta.new, A)
  converge.check <- abs(L.new - L.old) < thresh</pre>
 theta.new <- theta.new/rowSums(theta.new)</pre>
 return(list(theta = theta.new, loglik = L.new,
              converged = converge.check))
# initialize the parameters at random starting values
temp <- matrix(runif(n*K), n, K)</pre>
theta.init <- temp/rowSums(temp)
# do single update
out <- oneUpdate(A, n, K, theta.init)</pre>
```

```
# time to execute the code
system.time(out <- oneUpdate(A, n, K, theta.init))
## user system elapsed
## 0.67 0.12 0.79</pre>
```

## 4. First Algorithm:

```
# original algorithm
PIKK <- function(x, k) {
  x[sort(runif(length(x)), index.return = TRUE)$ix[1:k]]
}</pre>
```

Use runif() to generate 2k numbers between 0 and 1. Multiply these 2k numbers with the length(x) to obtain 2k numbers ranging from 0 to the length(x). These 2k numbers now have decimal places, use round() to convert them to integers. Use unique() to filter and keep the unique numbers. This step makes sure the characteristic of without replacement. Select the first k numbers in the sequences of unique numbers we just obtain. Return these k numbers as sample without replacement.

```
modPIKK <- function(x, k) {
  x[unique(round(runif(2*k)*length(x)))[1:k]]
}</pre>
```

Now we how how efficient is the modified algorithm comparing to the original algorithm as k and n vary. Here we use  $k = \frac{n}{20}$  and  $n = \{5000, 6000, 7000, 8000, 9000, 10000\}$ .

```
# different length of x
n \leftarrow c(5000, 6000, 7000, 8000, 9000, 10000)
# time for original PIKK
timePIKK <- c()</pre>
# time for modified PIKK
timePIKKmod <- c()</pre>
# obtain time for original PIKK
for (i in n) {
 x <- rnorm(i)
 k < -i/20
 timePIKK <- c(timePIKK, mean(microbenchmark(PIKK(x,k))$time))</pre>
# obtain time for modified PIKK
for (i in n) {
 x <- rnorm(i)
 k < -i/20
 timePIKKmod <- c(timePIKKmod, mean(microbenchmark(modPIKK(x,k))$time))
# report time of original algorithm takes in nanoseconds
# with n=(5000,6000,7000,8000,9000,10000), k=n/20
timePIKK
## [1] 588527.3 629348.0 669639.0 748784.2 826496.6 916213.3
```

```
# report time of modified algorithm takes in nanoseconds # with n=(5000,6000,7000,8000,9000,10000), k=n/20 timePIKKmod ## [1] 84954.90 75910.64 81461.24 101787.37 105383.75 121514.09
```

## Second Algorithm

```
# original algorithm
FYKD <- function(x, k) {
    n <- length(x)
    for(i in 1:n) {
        j = sample(i:n, 1)
        tmp <- x[i]
        x[i] <- x[j]
        x[j] <- tmp
    }
    return(x[1:k])
}</pre>
```

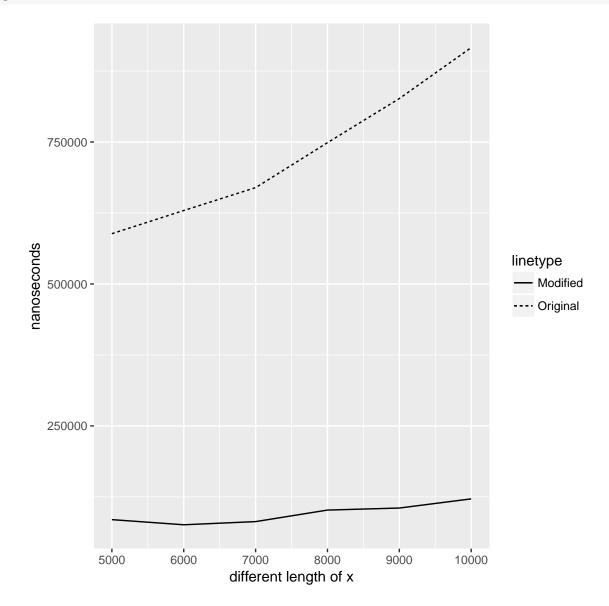
Modify FYKD by only swapping the first k numbers in x, instead of swapping all numbers in x. Then select the first k numbers.

```
# modified FYKD
modFYKD <- function(x, k) {</pre>
 n <- length(x)</pre>
  # instead of 1:n, use 1:k
 for(i in 1:k) {
    j = sample(i:n, 1)
    tmp <- x[i]</pre>
    x[i] \leftarrow x[j]
    x[j] <- tmp
 return(x[1:k])
# time for original FYKD
timeFYKD <- c()</pre>
# time for modified FYKD
timeFYKDmod <- c()</pre>
# obtain time for original FYKD
for (i in n) {
 x <- rnorm(i)
 k < -i/20
 timeFYKD <- c(timeFYKD, mean(microbenchmark(FYKD(x,k))$time))</pre>
# obtain time for modified FYKD
for (i in n) {
 x <- rnorm(i)
 k < -i/20
 timeFYKDmod <- c(timeFYKDmod, mean(microbenchmark(modFYKD(x,k))$time))</pre>
```

```
# report time of original FYKD algorithm takes in nanoseconds
# with n=(5000,6000,7000,8000,9000,10000), k=n/20
timeFYKD
## [1] 31415549 37840587 44039258 56390590 61210690 70917921
\# report time of modified FYKD algorithm takes in nanoseconds
# with n=(5000,6000,7000,8000,9000,10000), k=n/20
timeFYKDmod
## [1] 1752497 2114348 2778473 3167778 3783631 4893548
# create a dataframe collecting all information
df <- data.frame(xLength = n, PIKKtime = timePIKK,</pre>
                modPIKKtime = timePIKKmod, FYKDtime = timeFYKD,
                modFYKDtime = timeFYKDmod)
# all time are in nanoseconds
df
##
   xLength PIKKtime modPIKKtime FYKDtime modFYKDtime
## 1 5000 588527.3
                      84954.90 31415549
                                             1752497
## 2
     6000 629348.0 75910.64 37840587
                                             2114348
## 3
     7000 669639.0 81461.24 44039258
                                            2778473
## 4
     8000 748784.2 101787.37 56390590
                                             3167778
     9000 826496.6 105383.75 61210690
## 5
                                             3783631
## 6 10000 916213.3 121514.09 70917921 4893548
```

Plot the comparison of the PIKK algorithm as follows. Line on top is original algorithm, line on bottom is the modified algorithm. The size of k is always  $\frac{n}{20}$ .

```
# plot comparison for first algorithm
g1 <- ggplot(df) +
   geom_line(aes(x = xLength, y = PIKKtime, lty = "Original")) +
   geom_line(aes(x = xLength, y = modPIKKtime, lty = "Modified")) +
   xlab("different length of x") +
   ylab("nanoseconds")
g1</pre>
```



Plot the comparison of the FYKD algorithm as follows. Line on top is original algorithm, line on bottom is the modified algorithm. The size of k is always  $\frac{n}{20}$ .

```
# plot comparison for second algorithm
g2 <- ggplot(df) +
   geom_line(aes(x = xLength, y = FYKDtime, lty = "Original")) +
   geom_line(aes(x = xLength, y = modFYKDtime, lty = "Modified")) +
   xlab("different length of x") +
   ylab("nanoseconds")
g2</pre>
```

