HW4_Jingbin

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Problem 1

Given the algorithm, we compare the results with linear regression. Here is our tolerance used and the step size α .

Tolerance	Step Size	theta 0	theta 1	lm theta 0	lm theta 1
1e-07	0.01	0.9699	2.0015	0.9696	2.0016

From the above table, we found that the coefficients computed from gradient descent algorithm are close to the fitted linear regression.

```
# given the X and h below
set.seed(1256)
theta \leftarrow as.matrix(c(1, 2), nrow = 2)
X \leftarrow cbind(1, rep(1:10, 10))
h <- as.vector(X \% \% theta + rnorm(100, 0, 0.2))
# gradient descent input with theta, alpha, tolerance and m
gradient_descent <- function(theta, alpha = 0.01, tolerance = 1e-07, m = 100) {</pre>
  # set up parameters for interation
  theta0i.old <- 0
  theta1i.old <- 0
  theta0i.new <- 1
  theta1i.new <- 2
  n < 0
  # while loop stop when the absolute values less than the tolerance
  while (abs(theta0i.new - theta0i.old) > tolerance && abs(theta1i.new - theta1i.old) >
    tolerance) {
    theta0i.old <- theta0i.new
    theta1i.old <- theta1i.new
    theta.matrix <- matrix(c(theta0i.old, theta1i.old), nrow = 2)</pre>
    theta0i.new <- theta0i.old - alpha * (1/m) * sum(X %*% theta.matrix - h)
    theta1i.new <- theta1i.old - alpha * (1/m) * (t(X %*% theta.matrix - h) %*%
      X[, 2])
    n < - n + 1
  # return our results
  results <- cbind(theta_0 = theta0i.new, theta_1 = theta1i.new, iteration = n)
  return(results)
# test our function gradient descent
```

```
function.results <- gradient_descent(theta)
# lm parameter
lm(h ~ 0 + X)</pre>
```

part a

We use a step size of 1e-7, tolerace of 1e-9 and set a stopping rule for 5M. The total number of iteration is 362204. The average of theta0 is 0.9995, the average of theta1 is 1.9979.

part b

The above implementation is not a good way for gradient descent algorithm. Because it is time consuming. Although different tolence results differ, the loop takes a long run to converge.

part c

It is challenging to choose a proper parameter in our case. Maybe we should give this algorithm a smoother function to fit.

Problem 3

John Cook suggests that instead of solving for the inverse of $(X^TX)^{-1}$ we use the solve function to find $\hat{\beta}$ by solving the following system $(X^TX)\hat{\beta} = X^T\underline{y}$. However, we could also do $\hat{\beta} = (X^TX)^{-1}X^T\underline{y}$, which will take much more computation efforts. But the result will be the same.

part a

- The size of A is 112347224 bytes.
- The size of B is 1816357208 bytes.
- The time used on computation is 881.27 seconds.

```
# set up our parameters
set.seed(12456)
G \leftarrow matrix(sample(c(0, 0.5, 1), size = 16000, replace = T), ncol = 10)
R \leftarrow cor(G) # R: 10 * 10 correlation matrix of <math>G
C <- kronecker(R, diag(1600)) # C is a 16000 * 16000 block diagonal matrix
id <- sample(1:16000, size = 932, replace = F)</pre>
q <- sample(c(0, 0.5, 1), size = 15068, replace = T) # vector of length 15068
A <- C[id, -id] # matrix of dimension 932 * 15068
B <- C[-id, -id] # matrix of dimension 15068 * 15068
p <- runif(932, 0, 1)
r <- runif(15068, 0, 1)
C <- NULL #save some memory space
# the size of A and B and compute system time
object.size(A)
object.size(B)
system.time({
  y <- p + A %*% solve(B) %*% (q - r)
```

part b

The time consuming part is the inverse of the matrix. If we could come up with a method to speed up the matrix inverse.

part c

We inverse the matrix by converting the matrix by diagnosis elements, which bring down the time to 98 seconds.

```
inverse.R <- solve(R)
new.C <- kronecker(R, diag(1600))
inverse.B <- new.C[-id, -id]
# system.time({y <- p + A%*%B_inv%*%(q-r)})</pre>
```

Problem 5

part a

```
# Create a function
success.pro <- function(x = c(0, 1, 1, 0)) {
  n <- length(x)</pre>
```

```
success <- sum(x)</pre>
proportion <- success/n
return(proportion)
```

part b

```
set.seed(12345)
P4b_data <- matrix(rbinom(10, 1, prob = (31:40)/100), nrow = 10, ncol = 10, byrow = FALSE)
```

part c

```
The matrix is not random.
# First we compute the proportion of success by column
for (i in 1:10) {
  pro <- success.pro(x = P4b_data[, i])</pre>
  list.col <- c("Column: ", i, pro)</pre>
  print(list.col)
}
## [1] "Column: " "1"
                               "0.6"
## [1] "Column: " "2"
                               "0.6"
## [1] "Column: " "3"
                               "0.6"
## [1] "Column: " "4"
                              "0.6"
## [1] "Column: " "5"
                              "0.6"
## [1] "Column: " "6"
                              "0.6"
## [1] "Column: " "7"
                              "0.6"
## [1] "Column: " "8"
                              "0.6"
## [1] "Column: " "9"
                              "0.6"
## [1] "Column: " "10"
                              "0.6"
# compute the proportion of success by row
for (i in 1:10) {
  pro <- success.pro(x = P4b_data[i, ])</pre>
  list.row <- c("Row: ", i, pro)</pre>
  print(list.row)
## [1] "Row: " "1"
                        "1"
                        "1"
## [1] "Row: " "2"
## [1] "Row: " "3"
                        "1"
## [1] "Row: " "4"
                        "1"
## [1] "Row: " "5"
                        "0"
                        "0"
## [1] "Row: " "6"
## [1] "Row: " "7"
                        "0"
## [1] "Row: " "8"
                        "0"
## [1] "Row: " "9"
                        "1"
## [1] "Row: " "10"
                        "1"
```

part d

From the results, we use sapply function to create random matrix. And now the function works.

```
create.vect <- function(p = 0.5) {
   data <- rbinom(10, 1, prob = p)
   return(data)
}
prob <- c(seq(0.31, 0.4, by = 0.01))
P4b_data <- sapply(prob, create.vect)
proportion_P4b_col <- apply(P4b_data, 2, success.pro)  # by column
proportion_P4b_row <- apply(P4b_data, 1, success.pro)  # by column
proportion_P4b_col

## [1] 0.2 0.3 0.4 0.3 0.4 0.6 0.3 0.3 0.5 0.6
proportion_P4b_row

## [1] 0.7 0.3 0.5 0.5 0.3 0.1 0.8 0.4 0.1 0.2</pre>
```

[1] 0.7 0.0 0.0 0.0 0.0 0.1 0.0 0.1 0.1 0.2

Problem 6

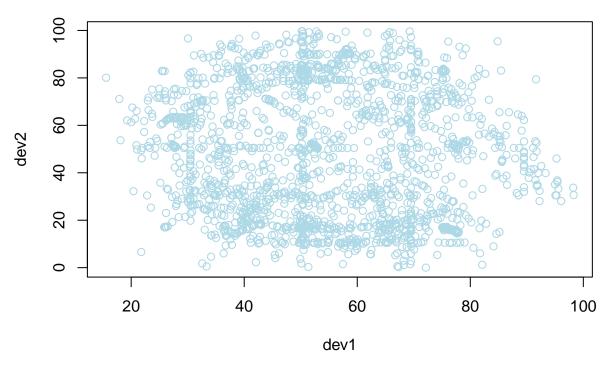
part 1

```
setwd("~/Desktop")
# Import the data from last assignment
data.device <- data.frame(readRDS("HW3_data.rds"))
data.device <- data.frame(data.device)
colnames(data.device) <- c("observer", "x", "y")
# Create a scatter plot based on x and y
scatterplot <- function(X = data.device) {
    x <- X[, 2]
    y <- X[, 3]
    plot <- return(plot(x, y, col = "lightblue"))
}</pre>
```

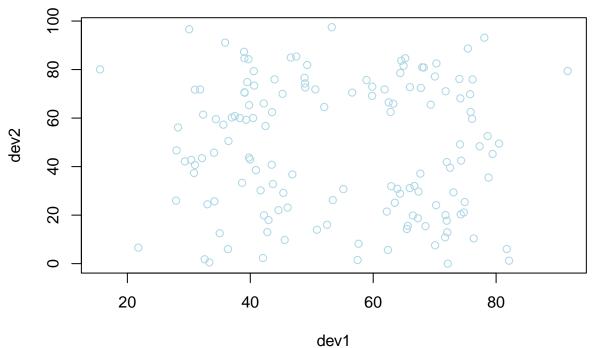
part 2

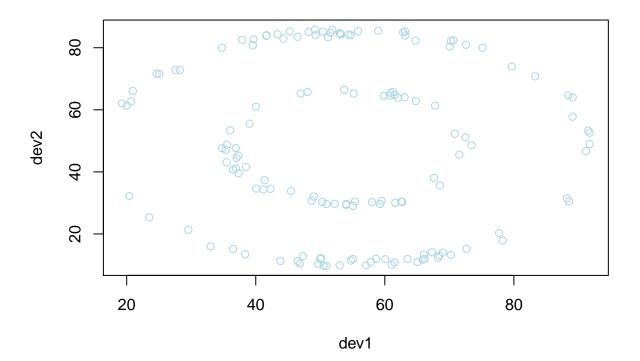
```
# the function make the plot
scatter.plot <- function(data, title, xlab, ylab) {
   plot <- plot(data$x, data$y, main = title, xlab = xlab, ylab = ylab, col = "light blue")
   return(plot)
}
scatter.plot(data.device, "device measurements", "dev1", "dev2")</pre>
```

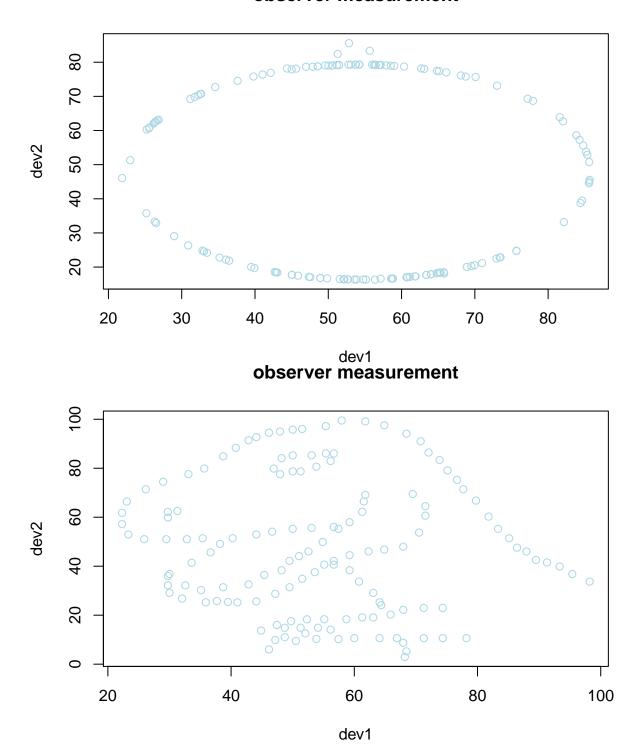
device measurements

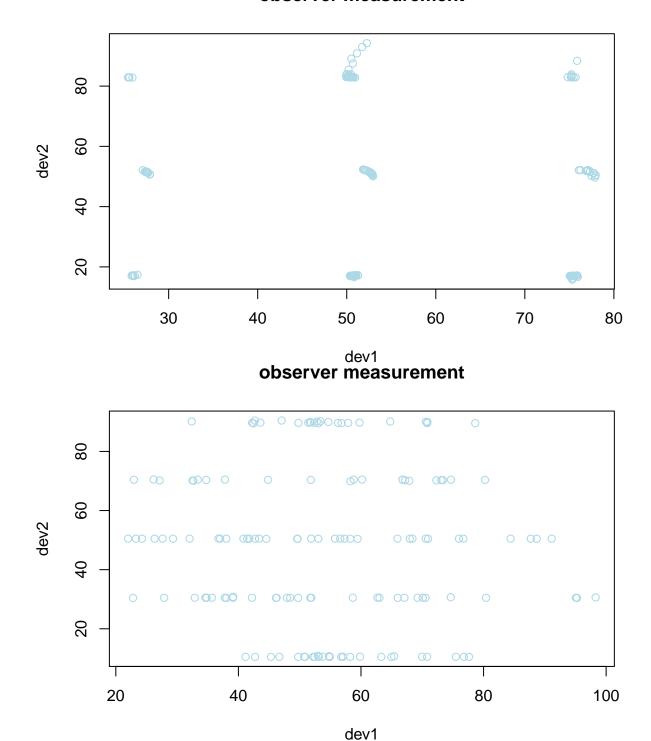


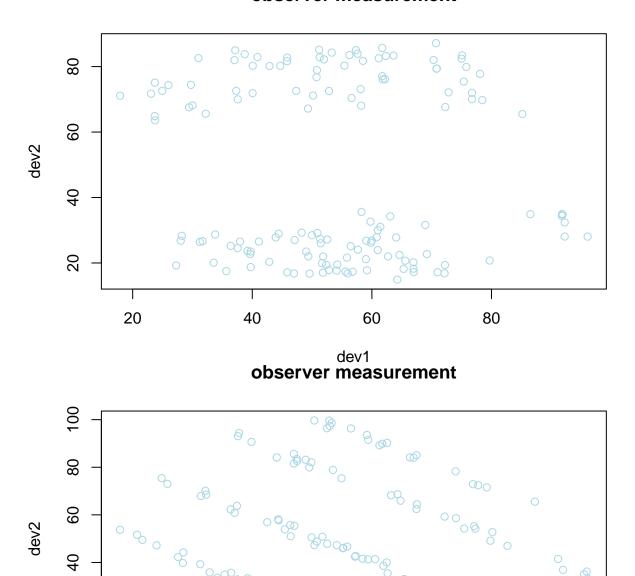
NULL



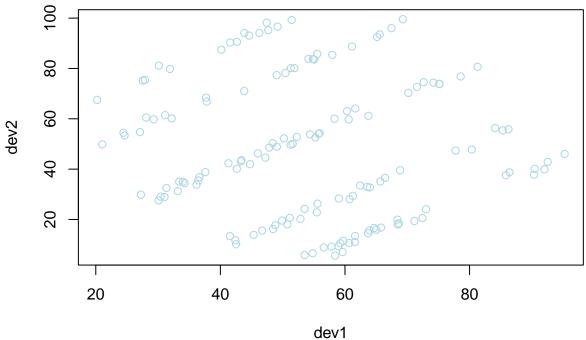


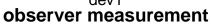


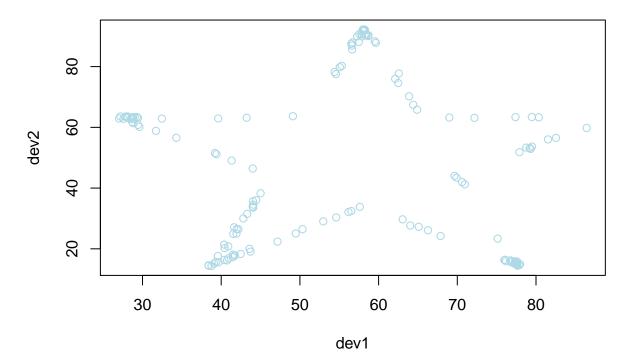


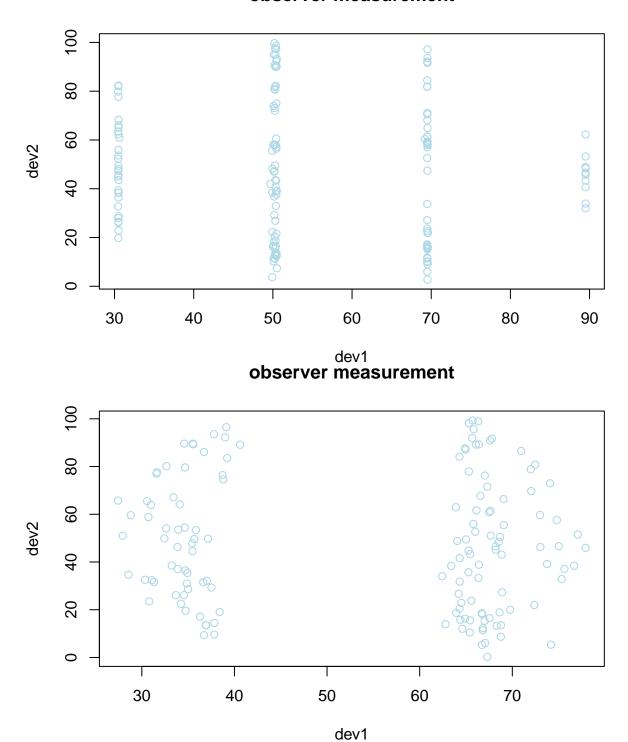


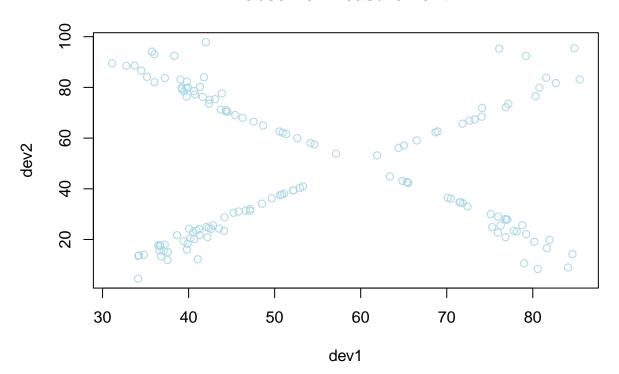
dev1











```
## [[1]]
## NULL
##
## [[2]]
## NULL
##
## [[3]]
## NULL
## [[4]]
## NULL
##
## [[5]]
## NULL
## [[6]]
## NULL
##
## [[7]]
## NULL
##
## [[8]]
## NULL
## [[9]]
## NULL
##
## [[10]]
## NULL
##
```

```
## [[11]]
## NULL
## [[12]]
## NULL
##
## [[13]]
## NULL
```

Part a

```
download("http://www.farinspace.com/wp-content/uploads/us_cities_and_states.zip",
  dest = "us_cities_states.unzip")
unzip("us_cities_states.unzip", exdir = "D:/VT/Rstudio/directory")
states <- fread(input = "D:/VT/Rstudio/directory/us_cities_and_states/states.sql",
  skip = 23, sep = "'", sep2 = ",", header = F, select = c(2, 4))
cities <- fread(input = "D:/VT/Rstudio/directory/us_cities_and_states/cities_extended.sql",
  skip = 26, sep = "'", sep2 = ",", header = F, select = c(2, 4, 6, 8, 10, 12))
colnames(cities) <- c("City", "State_Code", "Zip", "Latitude", "Longitude", "County")</pre>
```

part b

We show the first five frequency of cities for states.

```
city.count <- table(cities$State_Code)
head(city.count)
##
## AK AL AR AZ CA CO
## 273 838 709 532 2651 659</pre>
```

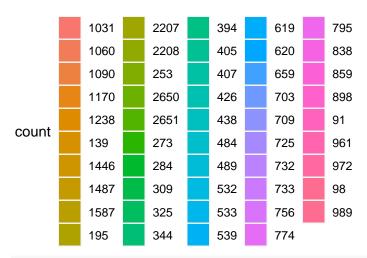
part c

```
getCount <- function(state_name, letter) {
  temp <- unlist(strsplit(tolower(state_name), split = ""))
  count <- 0
  for (i in 1:length(temp)) {
    if (temp[i] == letter) {
      count <- count + 1
      }
  }
  return(count)
}
letter.count <- data.frame(matrix(NA, nrow = 51, ncol = 26))
letter.count.results <- for (i in 1:51) {
  letter.count[i, ] <- sapply(1:26, function(n) {
      getCount(states$V2[i], letters[n])</pre>
```

```
})
}
```

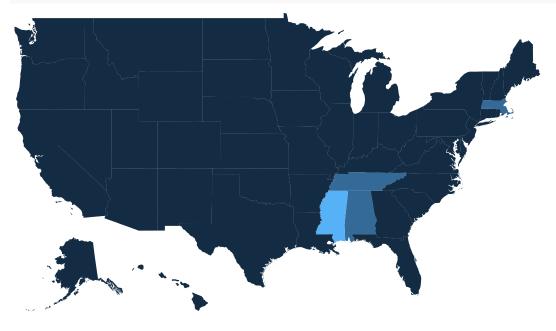
part d

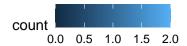




```
state.letter <- data.frame(state = tolower(states$V2), rowSums(letter.count > 3))
colnames(state.letter) <- c("state", "count")
p2 <- ggplot(state.letter, aes(map_id = state)) + geom_map(aes(fill = count), map = fifty_states) +
    expand_limits(x = fifty_states$long, y = fifty_states$lat) + coord_map() + scale_x_continuous(breaks)</pre>
```

```
scale_y_continuous(breaks = NULL) + labs(x = "", y = "") + theme(legend.position = "bottom",
  panel.background = element_blank())
p2
```





part a

The re-sampling result of each iteration is the same. So we have to correct the code. The author mistype the "Boot" as "Boot_times". "logapple08" and "logrm08" are not the variable names of "df08". When sampling, there is no need to use argument: "replace=TRUE".

```
set.seed(19941028)
# AAPL prices
apple08 <- getSymbols("AAPL", auto.assign = FALSE, from = "2008-1-1", to = "2008-12-31")[,
6]
# market proxy
rm08 <- getSymbols("^ixic", auto.assign = FALSE, from = "2008-1-1", to = "2008-12-31")[,
6]
# log returns of AAPL and market
logapple08 <- na.omit(ROC(apple08) * 100)
logrm08 <- na.omit(ROC(rm08) * 100)
# OLS for beta estimation
beta_AAPL_08 <- summary(lm(logapple08 ~ logrm08))$coefficients[2, 1]
# create df from AAPL returns and market returns
df08 <- cbind(logapple08, logrm08)
colnames(df08) <- c("logapple08", "logrm08") #Add one line here</pre>
```

```
Boot.times <- 1000
sd.boot <- rep(0, Boot.times)
for (i in 1:Boot.times) {
  bootdata <- df08[sample(nrow(df08), size = 251, replace = TRUE), ]
  sd.boot[i] <- coef(summary(lm(logapple08 ~ logrm08, data = bootdata)))[2, 2]
}</pre>
```

part b

```
# import the sensory data
set.seed(19941028)
library(data.table)
url1 <- "http://www2.isye.gatech.edu/~jeffwu/wuhamadabook/data/Sensory.dat"
sensory_data_raw <- fread(url1, fill = TRUE, data.table = FALSE)</pre>
saveRDS(sensory_data_raw, "sensory_data_raw.RDS")
sensory_data_raw <- readRDS("sensory_data_raw.RDS")</pre>
# Redo the data cleaning
sensory_data_raw_dl <- sensory_data_raw[-1, -2]</pre>
# Then we convert the value from string to numeric
library(stringr)
sensory data raw dl nu <- as.numeric(unlist(str extract all(sensory data raw dl,
  "[.-9]+")))
# We know that 1-10 are number of item, not true value in our table, so we delete
sensory_data_raw_value <- sensory_data_raw_dl_nu[-c(1, 17, 33, 49, 65, 81, 97, 113,
 129, 145)]
## Then we reconstruct data
sensory_data_tidy_br <- data.frame(item = sort(rep(1:10, 15)), operator = rep(c(1,
  1, 1, 2, 2, 2, 3, 3, 3, 4, 4, 4, 5, 5, 5), 10), values = sensory_data_raw_value)
sensory <- sensory_data_tidy_br[, -1]</pre>
colnames(sensory) <- c("operator", "y")</pre>
sd.boot <- rep(0, Boot.times)</pre>
# Bootstrap using the sensory data
time1 <- system.time(for (i in 1:Boot.times) {</pre>
  bootdata <- sensory[sample(nrow(sensory), size = 100, replace = TRUE), ]</pre>
  sd.boot[i] <- coef(summary(lm(y ~ operator, data = bootdata)))[2, 2]</pre>
})
time1
##
      user system elapsed
    0.862 0.013 0.877
##
```

parb c

```
cores <- max(1, detectCores() - 1)
# Create a cluster via makeCluster
cl <- makeCluster(cores)
registerDoParallel(cl)
# Parallelize
Boot_times <- 1000
sd.boot <- rep(0, Boot_times)</pre>
```

```
set.seed(777)
clusterExport(cl, list("sensory", "Boot_times", "sd.boot"))
# Bootstrap using the sensory data
time2 <- system.time(foreach(i = 1:Boot_times) %dopar% {
    # nonparametric bootstrap
    bootdata <- sensory[sample(nrow(sensory), size = 100, replace = TRUE),]
    sd.boot[i] <- coef(summary(lm(y ~ operator, data = bootdata)))[2, 2]
})
# time2<-system.time(foreach(i=1:Boot_times)%dopar%{ nonparametric bootstrap
# bootdata<-sensory[sample(nrow(sensory), size = 100, replace = TRUE),]
# sd.boot[i]<- coef(summary(lm(y~operator, data = bootdata)))[2,2] })
time2
stopCluster(cl)</pre>
```

part a

First build up the newton method algorithm, and create a vector from -1 to 1. Then apply the function and calculate the system time it takes.

```
fun <- function(x = 1) {
  f <- 3^x - \sin(x) + \cos(5 * x)
  return(f)
dr \leftarrow function(x = 1) {
  d \leftarrow (3^x) * \log(3, base = \exp(1)) - \cos(x) - 5 * \sin(5 * x)
}
newton.method <- function(m = 1) {</pre>
  tolerance <- 0.5
  while (abs(fx(m)) > tolerance) {
    new.point \leftarrow m - (fun(m)/dr(m))
    m <- new.point
  }
  return(m)
}
vector \leftarrow seq(-1, 1, by = 0.002)
# Using on of the apply functions and find the roots system.time(result <-
# lapply(vec, newton))
```

part b

We use 8 workers to accelerate computation speed.

```
cores <- 8
cl <- makeCluster(cores)
clusterExport(cl, list("fx", "dr"))
vecctor <- seq(-1, 1, by = 0.001)
# Using on of the apply functions and find the roots
system.time(result <- parSapply(cl, vector, newton.method))
stopCluster(cl)</pre>
```