Assignment2

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Problem1

(a)

$$L(\theta) = p_1 p_2 \dots p_n \tag{1}$$

$$lnL(\theta) = ln(p_1) + \dots + p_n \tag{3}$$

$$(4)$$

$$l(\theta) = -\sum_{i=1}^{n} ln\pi - \sum_{i=1}^{n} ln[1 + (x_i - \theta)^2]$$
(5)

$$(6)$$

$$l(\theta) = -nln\pi - \sum_{i=1}^{n} ln[1 + (x_i - \theta)^2]$$
(7)

$$l'(\theta) = -0 - \sum_{i=1}^{n} \frac{2(\theta - x_i)}{1 + (\theta - x_i)^2}$$
(9)

$$l'(\theta) = -2\sum_{i=1}^{n} \frac{(\theta - x_i)}{1 + (\theta - x_i)}$$
(11)

$$\begin{pmatrix}
n & 1 & 1 & 1 & 1 \\
n & 1 & 1 & 1 & 1
\end{pmatrix}$$

$$l''(\theta) = -2\sum_{i=1}^{n} \frac{1 + (\theta - x_i)^2 - 2(\theta - x_i)^2}{[1 + (\theta - x_i)^2]^2}$$
(13)

$$l''(\theta) = -2\sum_{i=1}^{n} \frac{1 - (\theta - x_i)^2}{[1 + (\theta - x_i)^2]^2}$$
(14)

$$i=1$$
 (16)

$$p'(x) = -\frac{2(x-\theta)}{\pi[1+(x-\theta)^2]^2}$$
(17)

$$(18)$$

$$I(\theta) = \frac{4n}{\pi} \int_{-\infty}^{\infty} \frac{(x-\theta)^2}{[1+(x-\theta)^2]^3} dx$$
 (19)

$$(20)$$

$$I(\theta) = \frac{4n}{\pi} \int_{-\infty}^{\infty} \frac{x^2}{[1+x^2]^3} dx$$
 (21)

(00)

(2)

(8)

$$(23)$$

if we choose x = tanx, $1 + x^2 = \frac{1}{cos^2\theta}$,

$$I(\theta) = \frac{4n}{\pi} \int_{-\infty}^{\infty} \frac{\sin^2 x}{\cos^2 x} \cos^6 x d(\tan x)$$
 (24)

$$(25)$$

$$I(\theta) = \frac{4n}{\pi} \int_{-\infty}^{\infty} \frac{1}{\cos^2 x} \sin^2 x \cos^4 x dx \tag{26}$$

$$(27)$$

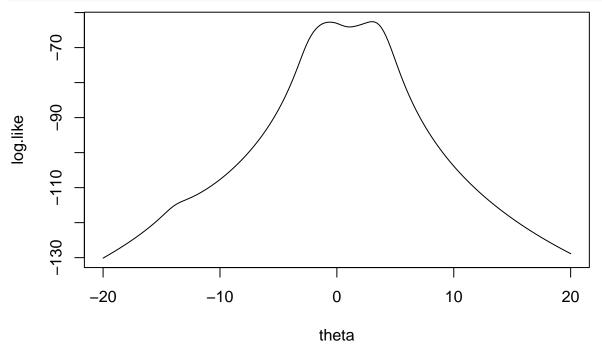
$$I(\theta) = \frac{4n}{\pi} \int_{-\frac{\pi}{2}}^{\frac{\pi}{2}} \frac{\sin^2 2x}{4} dx \tag{28}$$

$$(29)$$

$$I(\theta) = \frac{4n\pi}{\pi} \frac{\pi}{8} = n/2 \tag{30}$$

(b)

First, we graph the log-likelihood funtion $l(\theta) = -nln\pi - \sum_{i=1}^{n} ln[1 + (x_i - \theta)^2]$ as shown in the former section



Then we find the MLE for θ through application of the Newton-Raphson method for the nine given starting points. By setting the the sample mean, which turns out to be 3.257778, as the starting point

```
library(elliptic)
## Attaching package: 'elliptic'
## The following objects are masked from 'package:stats':
##
       sd, sigma
## The following object is masked from 'package:base':
##
##
       is.primitive
x \leftarrow c(1.77, -0.23, 2.76, 3.80, 3.47, 56.75, -1.34, 4.24, -2.44,
       3.29, 3.71, -2.40, 4.53, -0.07, -1.05, -13.87, -2.53, -1.75
f \leftarrow function(theta) \{ -2 * sum((theta - x) / (1 + (theta - x)^2)) \}
fdash \leftarrow function(theta) \{ -2 * sum((1 - (theta - x)^2) / ((1 + (theta - x)^2))^2) \}
s.p \leftarrow c(-11, -1, 0, 1.5, 4, 4.7, 7, 8, 38)
for (i in 1:9) {
    result <- newton_raphson(s.p[i], f, fdash, maxiter=100)
  print(paste0("starting point is: ", s.p[i]))
  print(result)
## [1] "starting point is: -11"
## $root
## [1] -2.080688e+17
##
## $f.root
## [1] 1.730197e-16
##
## $iter
## [1] 55
## [1] "starting point is: -1"
## $root
## [1] -0.5914735
##
## $f.root
## [1] 3.469447e-17
##
## $iter
## [1] 4
##
## [1] "starting point is: 0"
## $root
## [1] -0.5914735
##
## $f.root
## [1] -1.873501e-16
## $iter
## [1] 5
```

```
## [1] "starting point is: 1.5"
## $root
## [1] 1.09273
## $f.root
## [1] -1.387779e-17
## $iter
## [1] 5
## [1] "starting point is: 4"
## $root
## [1] 3.021345
##
## $f.root
## [1] -6.938894e-17
##
## $iter
## [1] 4
##
## [1] "starting point is: 4.7"
## $root
## [1] -0.5914735
##
## $f.root
## [1] -1.873501e-16
## $iter
## [1] 5
## [1] "starting point is: 7"
## $root
## [1] 1.736482e+17
## $f.root
## [1] -2.073157e-16
##
## $iter
## [1] 56
## [1] "starting point is: 8"
## $root
## [1] 1.838207e+17
## $f.root
## [1] -1.95843e-16
##
## $iter
## [1] 57
## [1] "starting point is: 38"
## $root
## [1] 3.226576e+17
```

```
##
## $f.root
## [1] -1.115734e-16
##
## $iter
## [1] 54
mean(x)
## [1] 3.257778
#sample mean 3.257778
newton_raphson(3.257778, f, fdash, maxiter=100)
## $root
## [1] 3.021345
##
## $f.root
## [1] -6.938894e-17
##
## $iter
## [1] 5
```

As shown in the results, taking the sample mean 3.257778 as the starting point gives the root 3.021345 while the other trials have more obvious difference between the starting point and the root. Therefore, the sample mean is a good starting point. Better choice might be available though.

(c)

We define the function for the fixed-point iteration with an accuracy tolerance of 10^{-6} and the maximum iteration times to be 1000.

```
p.fixed <- function(p0,alpha,obs,tol = 1E-6,max.iter = 1000,verbose=F){
  pold <- p0
   pnew <- pold + alpha * (-2) * sum((pold - obs)/(1 + (pold - obs)^2))
   iter <- 1
  while ((abs(pnew - pold) > tol) && (iter < max.iter)){
    pold <- pnew
    pnew <- pold + alpha * (-2) * sum((pold - obs)/(1 + (pold - obs)^2))
    iter <- iter + 1
    if(verbose)
      cat("At iteration", iter, "value of p is:", pnew, "\n")
  if (abs(pnew - pold) > tol) {
    cat("Algorithm failed to converge")
    return(NULL)
  }
  else {
    cat("Algorithm converged, in :" ,iter, "iterations \n")
    return(pnew)
  }
}
x \leftarrow c(1.77, -0.23, 2.76, 3.80, 3.47, 56.75, -1.34, 4.24, -2.44,
       3.29, 3.71, -2.40, 4.53, -0.07, -1.05, -13.87, -2.53, -1.75)
```

```
p.fixed(p0 = -1, alpha = 1.00, obs = x)
## Algorithm failed to converge
## NULL
p.fixed(p0 = -1, alpha = 0.64, obs = x)
## Algorithm converged, in : 11 iterations
## [1] -0.5914734
p.fixed(p0 = -1, alpha = 0.25, obs = x)
## Algorithm converged, in : 20 iterations
## [1] -0.5914741
#alpha = 1 will not converge.
s.p \leftarrow c(-11, -1, 0, 1.5, 4, 4.7, 7, 8, 38)
alpha \leftarrow c(1, 0.64, 0.25)
i <- 1
j <- 1
for (i in 1:9){
 for (j in 1:3){
  print(paste0("p0 = ", s.p[i], " scaling alpha =", alpha[j]))
   print(p.fixed(p0 = s.p[i], alpha = alpha[j], obs = x))
 }
}
## [1] "p0 = -11 scaling alpha =1"
## Algorithm converged, in : 218 iterations
## [1] -0.591474
## [1] "p0 = -11 scaling alpha =0.64"
## Algorithm converged, in : 17 iterations
## [1] -0.5914734
## [1] "p0 = -11 scaling alpha =0.25"
## Algorithm converged, in : 28 iterations
## [1] -0.5914745
## [1] "p0 = -1 scaling alpha =1"
## Algorithm failed to convergeNULL
## [1] "p0 = -1 scaling alpha =0.64"
## Algorithm converged, in : 11 iterations
## [1] -0.5914734
## [1] "p0 = -1 scaling alpha =0.25"
## Algorithm converged, in : 20 iterations
## [1] -0.5914741
## [1] "p0 = 0 scaling alpha =1"
## Algorithm failed to convergeNULL
## [1] "p0 = 0 scaling alpha =0.64"
## Algorithm converged, in : 11 iterations
## [1] -0.5914736
## [1] "p0 = 0 scaling alpha =0.25"
## Algorithm converged, in : 20 iterations
## [1] -0.5914726
## [1] "p0 = 1.5 scaling alpha =1"
```

```
## Algorithm failed to convergeNULL
## [1] "p0 = 1.5 scaling alpha =0.64"
## Algorithm failed to convergeNULL
## [1] "p0 = 1.5 scaling alpha =0.25"
## Algorithm converged, in : 12 iterations
## [1] 3.021345
## [1] "p0 = 4 scaling alpha =1"
## Algorithm failed to convergeNULL
## [1] "p0 = 4 scaling alpha =0.64"
## Algorithm converged, in : 12 iterations
## [1] -0.5914736
## [1] "p0 = 4 scaling alpha =0.25"
## Algorithm converged, in : 8 iterations
## [1] 3.021345
## [1] "p0 = 4.7 scaling alpha =1"
## Algorithm failed to convergeNULL
## [1] "p0 = 4.7 scaling alpha =0.64"
## Algorithm converged, in : 13 iterations
## [1] -0.5914736
## [1] "p0 = 4.7 scaling alpha =0.25"
## Algorithm converged, in : 9 iterations
## [1] 3.021345
## [1] "p0 = 7 scaling alpha =1"
## Algorithm failed to convergeNULL
## [1] "p0 = 7 scaling alpha =0.64"
## Algorithm failed to convergeNULL
## [1] "p0 = 7 scaling alpha =0.25"
## Algorithm converged, in : 8 iterations
## [1] 3.021345
## [1] "p0 = 8 scaling alpha =1"
## Algorithm failed to convergeNULL
## [1] "p0 = 8 scaling alpha =0.64"
## Algorithm converged, in : 14 iterations
## [1] -0.5914736
## [1] "p0 = 8 scaling alpha =0.25"
## Algorithm converged, in : 10 iterations
## [1] 3.021345
## [1] "p0 = 38 scaling alpha =1"
## Algorithm failed to convergeNULL
## [1] "p0 = 38 scaling alpha =0.64"
## Algorithm failed to convergeNULL
## [1] "p0 = 38 scaling alpha =0.25"
## Algorithm converged, in : 93 iterations
## [1] 3.021345
```

(d)

Note that to maximize the log-likelihood is equivalent to minimise its negative value. We first use the function nlminb with the fisher score

$$I(\theta) = n/2$$

as found in part (a) to find the MLE for θ .

```
x \leftarrow c(1.77, -0.23, 2.76, 3.80, 3.47, 56.75, -1.34, 4.24, -2.44,
       3.29, 3.71, -2.40, 4.53, -0.07, -1.05, -13.87, -2.53, -1.75)
s.p \leftarrow c(-11, -1, 0, 1.5, 4, 4.7, 7, 8, 38)
i <- 1
for (i in 1:9) {
  f \leftarrow function(theta)\{length(x) * log(pi) + sum(1 + (theta - x)^2)\}
  grf \leftarrow function(theta){2 * sum((theta - x) / (1 + (theta - x)^2))}
  fs <- function(theta){matrix(length(x) / 2, nrow = 1)}</pre>
  z <- nlminb(s.p[i], f, grf,fs)}</pre>
Then we refine the estimation by running Newtion-Raphson method as we did in part (b).
  f \leftarrow function(theta) \{-2*sum((theta - x)/(1 + (theta - x)^2))\}
  fdash \leftarrow function(theta) \{-2*sum((1 - (theta - x)^2)/((1 + (theta - x)^2))^2)\}
  result <- newton_raphson(z$par, f, fdash, maxiter=1000)
  print(paste0("starting point =", s.p[i]))
## [1] "starting point =38"
  print(paste0("Fisher scoring"))
## [1] "Fisher scoring"
print(z)
## $par
## [1] 19.17511
##
## $objective
## [1] 7945.465
## $convergence
## [1] 1
##
## $iterations
## [1] 150
##
## $evaluations
## function gradient
##
        151
                  151
##
## $message
## [1] "iteration limit reached without convergence (10)"
  print(paste0("Newton-Rasphon"))
## [1] "Newton-Rasphon"
  print(result)
## $root
## [1] 2.889657e+17
##
## $f.root
## [1] -1.245823e-16
```

```
##
## $iter
## [1] 55
```

(e) From the results, we can see from the numbers of iteration that the fixed-point method is less efficient than Fisher scoring with Newton-Rasphon in this case. As for the stability, the later one performs better as well.

Problem2

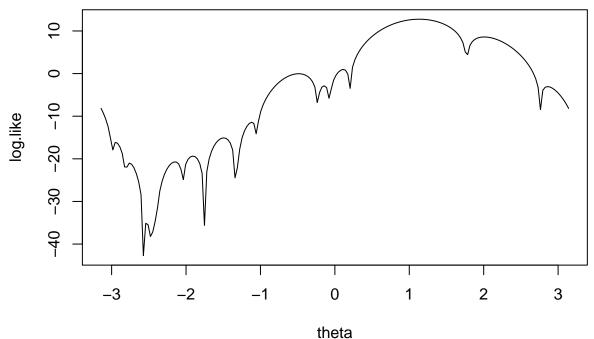
(a)

First, we find the log-likelihood function of θ

$$L(\theta) = p_1 p_2 \dots p_{19}$$
$$lnL(\theta) = lnp_1 + \dots + p_{19}$$
$$l(\theta) = \sum_{i=1}^{19} ln(\frac{1 - cos(x - \theta)}{2\pi})$$

Then, we plot the graph for the $l(\theta)$ as required

```
theta <- seq(from = -pi, to = pi, by = 2 * pi / 199)
i <- 1
log.like <- rep(0, 200)
for (i in 1:200) {
   log.like[i] <- sum(log((1 - cos(x - theta[i])) / 2 * pi))
}
plot(theta,log.like,type = "l")</pre>
```



(b)

We solve for the value of the estimator $\hat{\theta}_{moment}$

$$E[x|\theta] = \int_{0}^{2\pi} x \frac{1 - \cos(x - \theta)}{2\pi}$$
 (31)

(32)

$$E[x|\theta] = \frac{1}{2\pi} \left[\int_0^{2\pi} x dx - \int_0^{2\pi} x \cos(x - \theta) dx \right]$$
 (33)

(34)

$$E[x|\theta] = \pi - \frac{1}{2\pi} (x \sin(x - \theta) + \cos(x - \theta))|_0^{2\pi}$$
(35)

(36)

$$E[x|\hat{\theta}_{moment}] = \pi + \sin(\hat{\theta}_{moment}) = \bar{x}$$
(37)

$$\hat{\theta}_{moment} = \arcsin(\bar{x} - \pi) \tag{39}$$

(40)

(41)

Therefore, the obeject function is $f(\theta) = \pi + sin(\hat{\theta}_{moment}) - \bar{x}$. Here we use the function uniroot mean(x)

```
## [1] 3.257778
```

```
f1<-function(theta){
   pi+sin(theta)-mean(x)
}
uniroot.1<-uniroot(f1,lower=-pi,upper=-pi/2,extendInt = "yes")$root[1]
uniroot.2<-uniroot(f1,lower=-pi/2,upper=0,extendInt = "yes")$root[1]
uniroot.3<-uniroot(f1,lower=0,upper=pi/2,extendInt = "yes")$root[1]
uniroot.4<-uniroot(f1,lower=pi/2,upper=pi,extendInt = "yes")$root[1]
theta.mom <- c(uniroot.1, uniroot.2, uniroot.3, uniroot.4)
return(theta.mom)</pre>
```

[1] -3.2580418 -3.2580413 0.1164476 3.0251450

Roots are found as -3.2580418, -3.2580413, 0.1164476 and 3.0251450. ##(c) To use the Newton-Raphson method, similar to the earlier process, we will need the object function of θ , which is

$$f(\theta) = \sum_{i=1}^{19} \frac{\sin(\theta - x_i)}{1 - \cos(\theta - x_i)}$$

, and its first derivative

$$f'(\theta) = \sum_{i=1}^{19} \frac{1}{\cos(\theta - x_i) - 1}$$

```
library(elliptic)
```

```
x \leftarrow c(3.91, 4.85, 2.28, 4.06, 3.70, 4.04, 5.46, 3.53, 2.28, 1.96,
       2.53, 3.88, 2.22, 3.47, 4.82, 2.46, 2.99, 2.54, 0.52)
f<-function(theta){</pre>
  sum( sin(theta-x)/(1-cos(theta-x)) )
fdash<-function(theta){</pre>
  sum( 1 / (cos(theta-x)-1))
}
new.raphson <- function (initial, f, fdash, maxiter, give = TRUE, tol = .Machine$double.eps)
{
  old.guess <- initial
  for (i in seq_len(maxiter)) {
    new.guess <- old.guess - f(old.guess)/fdash(old.guess)</pre>
    jj <- f(new.guess)</pre>
    if (is.na(jj) | is.infinite(jj)) {
      break
    }
    if (near.match(new.guess, old.guess) | abs(jj) < tol) {</pre>
      if (give) {
        return(list(root = new.guess, f.root = jj, iter = i))
      else {
        return(new.guess)
    old.guess <- new.guess
  return("did not converge")
for (i in 1:length(theta.mom) ){
print(new.raphson(theta.mom[i],f,fdash,maxiter=1000))
}
## $root
## [1] -3.112471
##
## $f.root
## [1] 1.765255e-14
##
## $iter
## [1] 8
##
## $root
## [1] -3.112471
##
## $f.root
## [1] 1.765255e-14
##
## $iter
## [1] 8
##
```

```
## [1] "did not converge"
## $root
## [1] 3.170715
##
## $f.root
## [1] -3.719247e-15
## $iter
## [1] 8
Applying the values of \hat{\theta}_{moment} found in part (b), the results -3.112471, -3.114472, 3.170715 were obtained for
each value. Note that if start at 0.1164476, it won't converge. \#\#(d)
newton_raphson(-2.7,f,fdash,maxiter=100)
## $root
## [1] -2.668857
##
## $f.root
## [1] -2.926999e-13
##
## $iter
## [1] 5
newton_raphson(2.7,f,fdash,maxiter=100)
## $root
## [1] 2.848415
##
## $f.root
## [1] -3.008704e-14
##
## $iter
## [1] 6
(e)
N_R <- function (initial, f, fdash, maxiter, give = TRUE, tol = .Machine$double.eps)
  old.guess <- initial
  for (i in seq_len(maxiter)) {
    new.guess <- old.guess - f(old.guess)/fdash(old.guess)</pre>
    jj <- f(new.guess)</pre>
    if (is.na(jj) | is.infinite(jj)) {
      break
    if (near.match(new.guess, old.guess) | abs(jj) < tol) {</pre>
      if (give) {
        return(list(root = new.guess, f.root = jj, iter = i))
      }
      else {
        return(new.guess)
      }
    }
```

```
old.guess <- new.guess
  }
  return(list(root = "Failed to Converge", f.root = jj, iter = i))
s.p <- seq(from = -pi, to = pi, by = 2*pi/199)
newton_raphson(s.p[1],f,fdash,maxiter=1000)
## $root
## [1] -3.112471
##
## $f.root
## [1] 1.765255e-14
## $iter
## [1] 5
newton_raphson(s.p[2],f,fdash,maxiter=1000)
## $root
## [1] -3.112471
##
## $f.root
## [1] 1.765255e-14
##
## $iter
## [1] 4
newton_raphson(s.p[3],f,fdash,maxiter=1000)
## $root
## [1] -3.112471
##
## $f.root
## [1] 1.765255e-14
## $iter
## [1] 5
out <- data.frame(</pre>
 start.point <- s.p[1:200],
 root <- rep(0,200)
names(out) <- c("start.point","root")</pre>
for(i in 1:200) {
  result <- N_R(s.p[i],f,fdash,maxiter=1000)
  out[i,2] <- result$root</pre>
}
out$root <-as.factor(out$root)</pre>
#There are 18 levels
i <- 1
for (i in 1:18){
```

```
subgrp <- data.frame(
    start.point <- rep(0,length(which(out$root == levels(out$root)[i]))),
    root <- rep(0,length(which(out$root == levels(out$root)[i])))
)
names(subgrp) <- c("start.point","root")
subgrp$start.point <- out[which(out$root == levels(out$root)[i]),1]
subgrp$root <- out[which(out$root == levels(out$root)[i]),2]
assign(paste0("root.",i), subgrp)
}</pre>
```

Problem3

(a)

Since it is given that

with $N_0 = 2$, we have

$$\begin{split} N_t &= \frac{KN_0}{N_0 + (K - N_0)exp(-rt)} \\ N_t &= \frac{2K}{2 + (K - 2)exp(-rt)} \\ exp(-rt) &= \frac{2(K - N_t)}{N_t(K - 2)} \\ rt &= ln \frac{N_t(K - 2)}{2(K - N_t)} \\ r &= ln \frac{N_t(K - 2)}{2(K - N_t)} \frac{1}{t} \end{split}$$

Now we rename the unknown parameter K that represents the population carrying capacity of the environment as θ_1 , and the unknown growth rate r as θ_2 . It has been shown in (a) that the error

$$f(N_t, r) = N_t - \frac{2K}{2 + (K - 2)exp(-rt)}$$

The Nonlinear Lesat Squares function is called to minimize the squared error as required. The results show that with the initial value (K, r) = (1200, 0.1716788), the optimization will be reached with (K, r) = (1049.4069185, 0.1182685)

```
beetles <- data.frame(
   days = c(0, 8, 28, 41, 63, 69, 97, 117, 135, 154),
   beetles = c(2, 47, 192, 256, 768, 896, 1120, 896, 1184, 1024))

K <- 1200
r.t <- log((beetles$beetles*(K-2))/(K - beetles$beetles)*2)
r.series <- r.t / beetles$days
mean(r.series[2:10])

## [1] 0.1716788

#theta1 <- N
#theta2 <- r</pre>
```

pop.mod <- $nls(beetles \sim N*2/(2+(N-2)*exp((-r)*days)), start=list(N = 1200, r = 0.1716788), data=beetles,$

#beetles ~ theta1*2/(2+(theta1-2)exp(-1*theta2*days))

```
## 677146.3 : 1200.0000000
                                0.1716788
## 110309.7 : 988.5766717
                              0.1367234
## 76986.61 : 1024.1437803
                                0.1224583
## 73503.58 : 1045.8468846
                                0.1189023
## 73421.83 : 1048.9806817
                                0.1183764
## 73419.7 : 1049.3958171 0.1182715
## 73419.7 : 1049.4053152
                               0.1182689
## 73419.7 : 1049.4069185
                               0.1182685
#Gaussian-Newton
i <- 1
j <- 1
t.day<- as.numeric(beetles$days)</pre>
n.b <- as.numeric(beetles$beetles)</pre>
theta <- matrix(c(1200, 0.1716788), nrow = 2)
for (j in 1:10) {
  error.matrix0 <- c()
  for (i in 1:10){
    t <- t.day[i]
    n \leftarrow n.b[i]
    K <- theta[1,1]</pre>
    r <- theta[2,1]
    error.matrix <- matrix( n - (K*2)/(2+(K-2)*exp(-r*t)), nrow = 1)
    error.matrix0 <- rbind(error.matrix0, t(error.matrix))</pre>
  }
  z.matrix <- error.matrix0</pre>
  sse <- sum(z.matrix[1:10,1]^2)</pre>
  i <- 1
  theta.note <- c("K", "r")
  gradient.matrix0 <- c()</pre>
  for (i in 1:10){
    t <- t.day[i]
    n \leftarrow n.b[i]
    K <- theta[1,1]</pre>
    r <- theta[2,1]
    dev <- numericDeriv(quote((K*2)/(2+(K-2)*exp(-r*t))),theta.note)</pre>
    gradient.matrix <- matrix(attributes(dev)$gradient, nrow = 2)</pre>
    gradient.matrix0 <- rbind(gradient.matrix0, t(gradient.matrix))</pre>
  }
  a.matrix <- gradient.matrix0
  print(theta)
  print(sse)
  #Updating Matrix
  update <- solve(t(a.matrix) %*% a.matrix) %*% t(a.matrix) %*% z.matrix
  theta <- theta + update
}
                 [,1]
## [1,] 1200.000000
```

[2,]

0.1716788

```
## [1] 110309.7
##
                 [,1]
## [1,] 1024.1437801
## [2,]
        0.1224583
## [1] 76986.61
                 [,1]
## [1,] 1045.8468842
## [2,] 0.1189023
## [1] 73503.58
##
                 [,1]
## [1,] 1048.9806817
## [2,]
        0.1183764
## [1] 73421.83
                 [,1]
## [1,] 1049.3390366
## [2,]
        0.1182867
## [1] 73419.76
                 [,1]
## [1,] 1049.3958170
## [2,] 0.1182715
## [1] 73419.7
                 [,1]
## [1,] 1049.4053151
## [2,]
        0.1182689
## [1] 73419.7
                 [,1]
## [1,] 1049.4069183
## [2,]
         0.1182685
## [1] 73419.7
                 [,1]
## [1,] 1049.4071888
## [2,]
        0.1182684
## [1] 73419.7
(b)
\#x \leftarrow as.numeric(1:5)
K \leftarrow seq(1049, 1200, by = (1200-1049)/29)
r \leftarrow seq(0.1182, 0.1716, by = (0.1716 - 0.1182)/29)
n.b <- as.numeric(beetles$beetles)</pre>
t.d <- as.numeric(beetles$days)</pre>
sse <- function(K,r,n,t){</pre>
 i <- 1
```

[1] 677146.3

error.sq <- 0

[1,] 988.5766717 ## [2,] 0.1367234

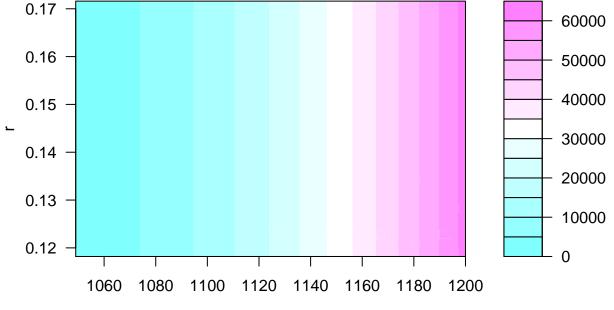
##

[,1]

```
for (i in 1:10){
  error.sq <- (n[i] - (K*2)/(2+(K-2)*exp(-r*t[i])))^2
  error.sq<- error.sq + error.sq
}
  return(error.sq)
}

z <- matrix(rep(0,900),nrow = 30)
j <- 1 #for K
k <- 1 #for r
for (j in 1:30){
  for(k in 1:30){
    z[j,k] <- sse(K[j],r[k],n.b,t.d)
}
}

filled.contour(K,r,z, xlab = 'K', ylab = 'r')</pre>
```



shown in the contour plot, the sum of squared errors decrease while (K, r) approach (1049.4069185, 0.1182685) from the initial values (K, r) = (1200, 0.1716788).

K

(c)

```
mlogl3 <- function(theta, N, days) {
    K <- theta[1]
    r <- theta[2]
    sigma <- theta[3]
    t <- days
    #mu <- log( f(...))
    mu <- log((K*2)/(2+(K-2)*exp(-r*t)))</pre>
```

```
- sum(dnorm(log(N), mu, sigma, log = TRUE))
sqrt(var(log(beetles$beetles)))
## [1] 2.031806
theta.start \leftarrow c(1200, 0.17,2.03)
out <- nlm(mlogl3, theta.start, N=beetles$beetles,days=beetles$days,hessian = TRUE)
## $minimum
## [1] 9.790127
##
## $estimate
## [1] 820.3815695  0.1926394  0.6440836
##
## $gradient
## [1] 1.138721e-08 2.900435e-05 -2.714273e-06
##
## $hessian
##
                 [,1]
                              [,2]
                                             [,3]
## [1,] 2.389705e-05 0.05442387 -3.048933e-06
## [2,] 5.442387e-02 373.52578204 -5.767475e-02
## [3,] -3.048933e-06 -0.05767475 4.817365e+01
## $code
## [1] 2
##
## $iterations
## [1] 38
theta.hat <- out$estimate</pre>
\#K = 820.3811422 , r = 0.1926394, sigma = 0.6440836
theta.hat
## [1] 820.3815695  0.1926394  0.6440836
hes <- out$hessian
hes
##
                 [,1]
                              [,2]
                                             [,3]
## [1,] 2.389705e-05 0.05442387 -3.048933e-06
## [2,] 5.442387e-02 373.52578204 -5.767475e-02
## [3,] -3.048933e-06 -0.05767475 4.817365e+01
var.matrix <- solve(hes)</pre>
# 6.262790e+04, 4.006745e-03, 2.075824e-02
diag(var.matrix)
```

[1] 6.262790e+04 4.006745e-03 2.075824e-02