STAT 428: Homework 4: Chapter 6 Monte Carlo Methods in Inference

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Please refer to the [detailed homework policy document] on Course Page for information about homework formatting, submission, and grading.

Exercise 1

An Exploration of Standard Error in Monte Carlo Estimation

Consider the following integral:

$$\int_0^1 \frac{\ln(x+1)}{\pi \sqrt{x(1-x)}} dx.$$

a. Estimate the integral using naive Monte Carlo. What is the standard error of this estimate?

```
u = runif(1000)
I0 = mean(log(u+1)/(pi*sqrt(u*(1-u))))
I0
## [1] 0.3844
```

```
x = numeric()
for(i in 1:1000){
    x[i] = mean(log(u[i]+1)/(pi*sqrt(u[i]*(1-u[i]))))
}
E0 = sd(x)/sqrt(1000)
E0
## [1] 0.02287
```

b. Let's see if we can improve the standard error. Implement Monte Carlo with antithetic sampling to estimate this integral. What is the standard error of this estimate?

```
I1 = numeric()
I11 = numeric()
I12 = numeric()

for(i in 1:1000){
    k1 <- 500
    u1 <- runif(500)
    I11[i] <- mean(log(u[i]+1)/(pi*sqrt(u[i]*(1-u[i]))))

    k2 <- 500
    u2 <- runif(500)
    I12[i] <- mean(log(1-u[i]+1)/(pi*sqrt(u[i]*(1-u[i]))))

I1[i] <- (I11[i] + I12[i])/2
}

E1 <- sd(I1)/sqrt(1000)
E1

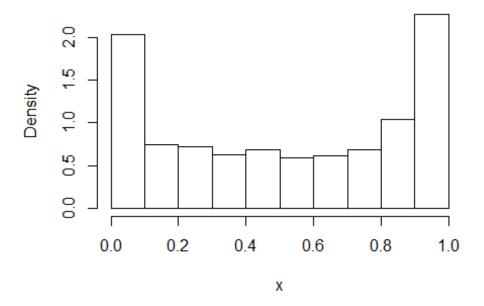
## [1] 0.01227</pre>
```

c. Would stratified sampling seem to help here? Why or why not? (Whatever you decide, you do not need to implement it).

I believe that stratified sampling will also work in reducing the standard error. It uses simaple random sampling from uniform distributions but stratify these to ensure balance over a partition into k subintervals of (0, 1)

d. $f(x) = \frac{1}{\pi\sqrt{x(1-x)}}$ for $x \in (0,1)$ is the probability density function for the Arcsine distribution. Using inverse transformation method, sample 1000 random values from the Arcsine distribution.

Histogram of x



e. Use importance sampling and the code you wrote in part d to estimate this integral. What is the standard error?

The importance function we choose is the pdf of Arcsine distribution.

```
m = 1000

gf = function(x){
    g = log(x+1)/(pi*sqrt(x*(1-x)))
    f = (x<1)*(x>0)
    g*f
}

##try our importance function
u = runif(m)
x = (sin(pi*u/2))^2
gfphi = gf(x)*(pi*sqrt(x*(1-x)))
E2 = sd(gfphi)/sqrt(m)
E2

## [1] 0.007742
```

f. Are all methods equally effective? Which method is the most efficient?

No. It seems like importance sampling is the most efficient way to reduce the error variance.

Exercise 2

Comparing MSE of estimators using MC.

Let $f(x|\theta) = t(\nu, \mu)$, the non-central t-distribution, where μ is a location parameter and ν is the degrees of freedom.

Estimate the MSE of the level k trimmed means for random samples of size 20 generated from a a non-central t-distribution with degrees of freedom 3 and mean 4 (with $\nu = 3$ and $\mu = 4$). Summarize the estimates of MSE in a table for k = 1, 2, ..., 9.

```
n = 20
K = n/2 - 1
m = 1000
mse = matrix(0, n/2, 2)
trimmed.mse = function(n, m, k){
 tmean = numeric(m)
  for(i in 1:m){
   x = sort(rt(m, 3, 4))
    tmean[i] = mean(x[(k+1):(n-k)])
  mse.est = mean(tmean^2)
  se.mse = sqrt(mean((tmean - mean(tmean))^2))/sqrt(m)
 return (c(mse.est, se.mse))
}
for(k in 1:K){
  mse[k, 1:2] = trimmed.mse(n,m,k)
round(mse, 3)
##
          [,1] [,2]
## [1,] 1.885 0.003
## [2,] 1.918 0.003
## [3,] 1.931 0.003
## [4,] 1.960 0.003
## [5,] 1.983 0.003
## [6,] 1.975 0.003
## [7,] 1.995 0.003
## [8,] 2.006 0.003
## [9,] 2.016 0.003
## [10,] 0.000 0.000
```

Exercise 3

Bayesian Statistics Suppose $X_1, ..., X_n$ are n independent and identical distributed random variables from $Exp(\theta)$, where θ is the unknown parameter. So,

$$f(x|\theta) = \theta e^{-\theta x}, \quad x \ge 0.$$

We assume the prior distribution on θ is the Gamma distribution (Gamma(3,2)).

$$g(\theta) = 4\theta^2 e^{-2\theta}, \quad x \ge 0.$$

1. Write down the posterior distribution of θ , $g(\theta|X)$.

$$g(\theta|x) = \frac{L(\theta)g(\theta)}{C}$$

where,

$$L(\theta) = \prod_{n=1}^{n} f(x_i | \theta)$$

and

$$g(\theta|x) = \int \prod_{n=1}^{n} f(x_i|\theta)g(\theta)dx$$

2. Suppose n=6 and we observe that $x_1=0.4, x_2=1.1, x_3=0.2, x_4=1.6, x_5=1.4, x_6=0.9$. Estimate the posterior mean of θ based on 1000 simulated θ from its prior distribution.

```
observed_x = c(0.4, 1.1, 0.2, 1.6, 1.4, 0.9)
hx = numeric()
post.mean = numeric()
for(i in 1:length(observed x)){
theta = rgamma(n = 1000, shape = 3, rate = 2)
hx[i] = theta*exp(-theta*observed x[i])
c = mean(hx)
post.mean[i] = mean(theta*hx)/c
cbind(observed x,post.mean)
##
        observed_x post.mean
## [1,]
               0.4
                       1.466
## [2,]
               1.1
                       1.473
## [3,]
               0.2
                       1.501
## [4,]
               1.6
                       1.504
## [5,]
               1.4
                       1.477
## [6,]
               0.9
                       1.450
```

- 3. Suppose n = 6 and we observe that $x_1 = 0.4, x_2 = 1.1, x_3 = 0.2, x_4 = 1.6, x_5 = 1.4, x_6 = 0.9$.
 - a. Design an acceptance-rejection sampling algorithm to generate 1000 (accepted) samples of θ from the posterior distribution of θ . Write down your algorithm with your instrumental distribution $g(\theta)$. (Hint: for the

acceptance-rejection sampling method, the normalizing constant in the posterior distribution can be ignored.)

Given that M = 0.5, for each observed x do folowing steps 1. generate 1 sample form gamma(3, 2) as the theta 2. generate 1 sample from uniform(1) 3. calculate f(x|theta) based on the observed x and theta from 1 4. calculate g(x) based on the observed x 5, Test uf the sample from 2 less than f(x|theta) / (M*g(x)) accept theta from 1 repeat until get 1000 accepted theta

b. Implement your acceptance-rejection sampling algorithm with R code. Plot the histogram of your generated sample and compare your sample mean with your estimated posterior mean obtained in Ex.3.2.

```
observed x = c(0.4, 1.1, 0.2, 1.6, 1.4, 0.9)
fx = function(theta, x){
  theta*exp(-theta*x)*4*(theta^2)*exp(-2*theta)
AR = function(x, n){
  i = 0
  theta = numeric()
  M = 0.5
  while (i < n) {</pre>
    theta0 = rgamma(1, shape = 3, rate = 2)
    u = runif(1)
    if(M*u < fx(theta0, x)/dgamma(x, shape = 3, rate = 2)){
      i = i + 1
      theta[i] = theta0
    }
  return(theta)
x = c(0.4, 1.1, 0.2, 1.6, 1.4, 0.9)
AR(x = x, n = 6)
## [1] 0.7299 1.0390 0.4748 1.3879 1.3453 0.6148
```

Exercise 4

Do 6.1 in the book, except with n = 25 and $k = 1, 2, \dots, 10$.

```
m <- 2000
K <- 10
n <- 25
tmean <- matrix(0,m,K)
mse_est <- numeric(K)
mse_se <- numeric(K)
for(k in 1:K){
   for (i in 1:m) {
      x <- sort(reauchy(n))
      tmean[i,k] <- mean(x[(k+1):(n-k)])</pre>
```

```
}
mse_est[k] <- mean(tmean[,k]^2)
mse_se[k] <- sqrt(sum((tmean[,k] - mean(tmean[,k]))^2)) / m
}
table <- cbind(seq(1:K),round(mse_est,5),round(mse_se,5))
colnames(table) <- c("k", "Estimated MSE of level k trimmed means",
"Standard Error")
knitr::kable(table, caption = 'Estimates of MSE')
</pre>
```

Estimates of MSE

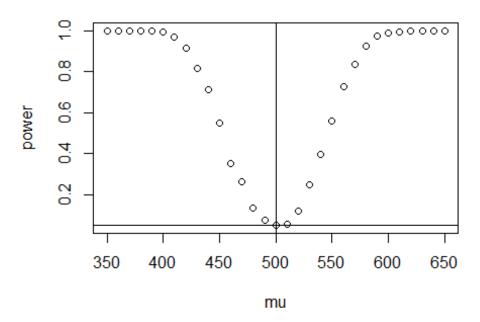
k	Estimated MSE of level k trimmed means	Standard Error
1	3.5117	0.0419
2	0.4430	0.0149
3	0.2132	0.0103
4	0.1696	0.0092
5	0.1418	0.0084
6	0.1244	0.0079
7	0.1159	0.0076
8	0.1092	0.0074
9	0.1026	0.0072
10	0.1066	0.0073

Exercise 5

Do exercise 6.2 from the book.

```
library(Hmisc) #for errbar
alpha = 0.05
mu <- c(seq(350, 650, 10)) #alternative H
n <- 20
sigma <- 100
m < -1000
mu0 <- 500
M <- length(mu)</pre>
power <- numeric(M)</pre>
for (i in 1 : M) {
  pvalues <- replicate(m, expr = {x <- rnorm(n, mean = mu[i], sd =</pre>
sigma)
      ttest <- t.test(x, alternative = "two.sided", mu = mu0)</pre>
      ttest$p.value})
  power[i] <- mean(pvalues <= alpha)</pre>
}
plot(mu, power)
```

```
abline(v = mu0, lty = 1)
abline(h = alpha, lty = 1)
```



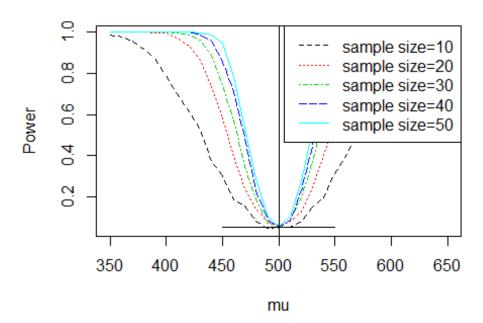
Exercise 6

Do exercise 6.3 from the book.

```
n <- seq(10,50,10) #sample size
mu <- c(seq(350, 650, 10))
m <- 1000
M <- length(mu)</pre>
N <- length(n)
power <- matrix(0,M,N)</pre>
for(j in 1:N){
  for (i in 1:M){
    mu1 <- mu[i]
    pvalues <- replicate(m, expr = {</pre>
      #simulate under alternative mu1
      x \leftarrow rnorm(n[j], mean = mu1, sd = 100)
      ttest <- t.test(x,</pre>
      alternative = "two.sided", mu = 500)
      ttest$p.value })
    power[i, j] <- mean(pvalues <= .05)</pre>
  }
}
plot(mu, power[,1], type="1", lty = 2, col = 1, main="Power curve",
xlab="mu", ylab="Power")
```

```
lines(mu, power[,2], lty = 3, col = 2)
lines(mu, power[,3], lty = 4, col = 3)
lines(mu, power[,4], lty = 5, col = 4)
lines(mu, power[,5], lty = 1, col = 5)
legend("topright", c("sample size=10","sample size=20","sample
size=30","sample size=40","sample size=50"), lty =
c(2,3,4,5,1),col=c(1,2,3,4,5))
abline(v = 500, lty = 1)
lines(c(450,550),c(0.05,0.05))
```

Power curve



Exercise 7

Do exercise 6.5 from the book.

```
alpha <- 0.05
m <- 1000
n <- 20
qt <- qt(1-alpha/2, df = n-1)
LCL <- replicate(m, expr = {
    x <- rchisq(n, 2)
    return(mean(x) - qt * sd(x) / sqrt(n))
})
UCL <- replicate(m, expr = {
    x <- rchisq(n, 2)
    return(mean(x) + qt * sd(x) / sqrt(n))
})
mean((LCL < 2) * (UCL > 2))
```

Exercise 8

Do exercise 6.8 from the book. Use 15 as small sample size, 50 as medium sample size, and 250 as large sample size.

```
alpha = 0.055
mu1 <- mu2 <- 0
sigma1 <- 1
sigma2 <- 1.5
sample_num <- c(15, 50, 250)
m < -1000
tests_F <- numeric(3)</pre>
tests CF <- numeric(3)
testF <- function(x, y) {
  f_test <- var.test(x, y, alternative = "two.sided", conf.level = 1-</pre>
alpha)
  return(f test$p.value < alpha)</pre>
}
tests5 = function(X, Y) {
  outx <- sum(X > max(Y)) + sum(X < min(Y))
  outy \leftarrow sum(Y > max(X)) + sum(Y \leftarrow min(X))
  return(as.integer(max(c(outx, outy)) > 5))
}
for (i in 1 : 3) {
  n <- sample num[i]</pre>
  power F <- mean(replicate(m, expr = {</pre>
  x <- rnorm(n, mu1, sigma1)</pre>
  y <- rnorm(n, mu2, sigma2)</pre>
  testF(x, y)
}))
  power CF <- mean(replicate(m, expr = {</pre>
  x <- rnorm(n, mu1, sigma1)</pre>
  y <- rnorm(n, mu2, sigma2)</pre>
  tests5(x, y)
}))
  tests_F[i] <- power_F
  tests_CF[i] <- power_CF
}
(table = rbind(tests_F, tests_CF))
              [,1] [,2] [,3]
## tests_F 0.321 0.800 1.000
## tests_CF 0.290 0.655 0.973
```