HW2

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$1 \quad 2.10$

$1.1 \quad (1)$

Proof that the sum of r independent geometric random variables with parameters p is a negative binomial(r,p) random variable: Suppose $X_1, \dots, X_r \stackrel{iid}{\sim} geometric(p)$, i.e. $P(X_i = x_i) = p(1-p)^{x_i-1}$. Then

$$P(\sum_{i=1}^{r} X_i = x) = \sum_{x_1 + \dots + x_r = x} P(X_1 = x_1, \dots, X_r = x_r)$$

$$= \sum_{x_1 + \dots + x_r = x} P(X_1 = x_1) \dots P(X_r = x_r)$$

$$= \sum_{x_1 + \dots + x_r = x} p^r (1 - p)^{x - r}$$

$$= C_{r-1}^{r-1} p^r (1 - p)^{x - r}$$

which is the probability distribution of negative binomial (r, p). So we can generate $r \times n$ geometric random number to obtain n negative binomial random number.

1.2 (2)

Suppose $U \sim U(0,1), X \sim \text{negative binomial}(r,p)$

$$F(x) = P(X \le x) = \begin{cases} \sum_{k=r}^{x} C_{k-1}^{r-1} p^{r} (1-p)^{k-r} & x = r, r+1, \dots \\ 0 & x = 0, \dots, r-1 \end{cases}$$

Y = x when $F(x - 1) < U \le F(x)$, then Y is the negative binomial random number.

```
inver_sampling <- function(N, F, X, seed) {
    set.seed(seed)
    U <- runif(N, 0, 1)
    Y <- rep(0, N)
    for (i in 1:N){
        k <- 1
        while(U[i] > F(k-1)) k <- k + 1
        Y[i] <- k - 1
    }
    return(Y)
}</pre>
```

$2\quad 2.11$

2.1 (1)

 $U \sim U(0,1)$. The cdf of $Beta(\frac{1}{n},1)$ is $F(x)=x^{1/n}$. Then $X=U^n$ is the $Beta(\frac{1}{n},1)$ random variable.

2.2(2)

The cdf of Beta(n,1) is $F(x)=x^n$. Then $X=U^{1/n}$ is the Beta(n,1) random variable.

2.3 (3)

The corresponding cdf is $F(x) = \int_0^x \frac{2}{\pi \sqrt{1-t^2}} dt = \frac{2}{\pi} \arcsin x$. Then $X = \sin(\frac{\pi}{2}U)$ is the target random variable.

2.4(4)

The corresponding cdf is $F(x) = \int_{-\infty}^{x} \frac{1}{\pi(1+t^2)} dt = \frac{1}{\pi} \arctan x + \frac{1}{2}$. Then $X = \tan(\pi(U - \frac{1}{2}))$ is the target random variable.

2.5 (5)

The corresponding cdf is $F(x) = \int_0^x \cos t dt = \sin x$. Then $X = \arcsin U$ is the target random variable.

2.6 (6)

The corresponding cdf is $F(x) = \int_0^x \frac{\alpha}{\eta} t^{\alpha-1} e^{-x^{\alpha}/\eta} dt = 1 - e^{-x^{\alpha}/\eta}$. Then $X = (-\eta \log(1-U))^{1/\alpha}$ is the target random variable.

3 2.13

The inverse transformation is

$$\begin{pmatrix} \alpha & = \arctan \frac{Y}{X} \\ R & = X^2 + Y^2 \end{pmatrix} \tag{1}$$

The Jacobi determinant is

$$J = \begin{vmatrix} -\sqrt{R}\sin\alpha & \frac{1}{2\sqrt{R}}\cos\alpha \\ \sqrt{R}\cos\alpha & \frac{1}{2\sqrt{R}}\sin\alpha \end{vmatrix} = -\frac{1}{2}$$
 (2)

Therefore, the joint density of (X, Y) is

 $=\frac{1}{\sqrt{2\pi}}e^{-\frac{x^2}{2}}\frac{1}{\sqrt{2\pi}}e^{-\frac{y^2}{2}}\tag{4}$

So X and Y are independent and both $\sim N(0,1)$.

4 3.2

4.1 (1)

```
[2]: # random point sampling
     randompoint_samp <- function(N, h, low, high, m = 0, M, B = 1, seed = NA) { \#N_{\sqcup}
      \rightarrow is the sample size, h(x) is the target function
         Iall <- rep(0, B)</pre>
                                                                    #B is repeated_
      → times(when calculate variance)
         for(i in 1:B) {
                                                                    #M and m are the upper
      \rightarrow bound and lower bound of h(x)
              if (!is.na(seed)) {set.seed(seed[i])}
                                                                    #low and high are the
      \rightarrowupper bound and lower bound of x
              else {set.seed(i)}
              x <- runif(N, low, high)
              y <- runif(N, m, M)
              p < -sum(y < h(x))
              Iall[i] \leftarrow p / N * (M-m) * (high-low)
         I <- mean(Iall)</pre>
         Var <- mean((Iall - I)^2)</pre>
         return(list(mean = I, var = Var))
     }
     N <- 10000
     I1 <- randompoint_samp(N, exp, -1, 1, 0, exp(1))
     I1$mean
```

2.37577831807321

```
else {set.seed(i)}
    u <- runif(N, low, high)
    Iall[i] <- (high - low) * sum(h(u)) / N
}
I <- mean(Iall)
Var <- mean((Iall - I)^2)
    return(list(mean = I, var = Var))
}
I2 <- average_samp(N, exp, -1, 1)
I2$mean</pre>
```

2.35766992054428

```
[4]: # importance sampling using U(-1,1) as g(x)
     importance_samp <- function(N, g, h, Ginver, B = 1, seed = NA) { #g is the_
      →selected function, Ginver is the inverse of its cdf
         Iall <- rep(0, B)</pre>
         for (i in 1:B) {
              if (!is.na(seed)) {set.seed(seed[i])}
              else {set.seed(i)}
              u <- runif(N, 0, 1)
              x <- Ginver(u)
              Iall[i] \leftarrow sum(h(x)/g(x)) / N
         I <- mean(Iall)</pre>
         Var \leftarrow mean((Iall - I)^2)
         return(list(mean = I, var = Var))
     g <- function(x) {</pre>
         return(1/2)
     Ginver <- function(y) {</pre>
         return(2*y-1)
     I3 <- importance_samp(N, g, exp, Ginver)</pre>
     I3$mean
```

2.35766992054428

```
    I <- mean(Iall)
    Var <- mean((Iall - I)^2)
    return(list(mean = I, var = Var))

}
g1 <- function(x) {
    return(1/2*(x+1))
}
Ginver1 <- function(y) {
    return(2*sqrt(y)-1)
}
I3_1 <- importance_samp(N, g1, exp, Ginver1)
I3_1$mean</pre>
```

2.35460530650625

(I do not try this g(x) in the following analysis, but it may be better)

```
[6]: # hierachical sampling, mixed with average sampling
     hierachical_samp <- function(n, h, seg, B = 1, seed = NA) {
         m <- length(seg) - 1 #number of segments
         Iall <- rep(0, B)</pre>
         for (j in 1:B) {
              for (i in 1:m) {
                  if (!is.na(seed)) {Iall[j] <- Iall[j] + average_samp(n[i], h, u
      \rightarrowseg[i], seg[i+1], seed = seed[j])$mean}
                  else {Iall[j] <- Iall[j] + average_samp(n[i], h, seg[i], seg[i+1], u
      \rightarrowseed = c(j))$mean}
             }
         }
         I <- mean(Iall)</pre>
         Var <- mean((Iall - I)^2)</pre>
         return(list(mean = I, var = Var))
     I4 <- hierachical_samp(c(5000, 5000), exp, c(-1, 0, 1))
     I4$mean
```

2.34722253862014

4.2 (2)

random point sampling

 $Var(\hat{I}_1) = [M(b-a)]^2 p(1-p)/N$, where M = e, b = 1, a = -1 and $p = (e-e^{-1})/2e$.

$$Var(\hat{I}_1) = \frac{(2e)^2(e - e^{-1})(2e - e + e^{-1})}{N \times 2e \times 2e}$$
$$= \frac{7.254}{N}$$

In order to store only 3 digits after the decimal point, we need to keep $z_{0.975}\sqrt{Var(\hat{I}_1)} < 0.0005$, i.e. N > 111467866.

average sampling

 $Var(\hat{I}_2) = (b-a)^2 Var(h(U))/N$, where b = 1, a = -1 and $Var(h(U)) = \int_a^b [h(u) - E(h(U))]^2 du/(b-a) = 1/2 - 1/2 \times e^{-2}$.

$$Var(\hat{I}_2) = \frac{2^2(1/2 - 1/2 \times e^{-2})}{N}$$
$$= \frac{1.729}{N}$$

In order to store only 3 digits after the decimal point, we need to keep $z_{0.975}\sqrt{Var(\hat{I}_2)} < 0.0005$, i.e. N > 26568506.

importance sampling

 $Var(\hat{I}_3) = Var(h(X)/g(X))/N$, where $X \sim g(x) = U(-1,1)$, then $Var(h(X)/g(X)) = Var(2e^x) = 2 - 2e^{-2}$.

$$Var(\hat{I}_3) = \frac{2 - 2e^{-2}}{N}$$
$$= \frac{1.729}{N}$$

In order to store only 3 digits after the decimal point, we need to keep $z_{0.975}\sqrt{Var(\hat{I}_3)} < 0.0005$, i.e. N > 26568506. hierachical sampling

 $Var(\hat{I}_4) = Var(\hat{I}_{41}) + Var(\hat{I}_{42}) = (b_1 - a_1)^2 Var(h(U_1))/(N/2) + (b_2 - a_2)^2 Var(h(U_2))/(N/2),$ where $b_1 = a_2 = 0$, $a_1 = -1$, $b_2 = 1$, $U_1 \sim U(-1,0)$, and $U_2 \sim U(0,1)$.

$$Var(\hat{I}_4) = \frac{(1 - e^{-2}) - 2(1 - e^{-1})^2}{N} + \frac{(e^2 - 1) - 2(e - 1)^2}{N}$$
$$= \frac{0.55}{N}$$

In order to store only 3 digits after the decimal point, we need to keep $z_{0.975}\sqrt{Var(\hat{I}_4)}<0.0005$, i.e. N>8451520.

4.3 (3)

```
Var hat
                               Var
                      name
                                              <dbl>
                      < chr >
                               <dbl>
                     I1
                               6.017275e-04
                                             0.0007254
A data.frame: 4 \times 3
                      I2
                               1.522177e-04
                                             0.0001729
                     Ι3
                               1.522177e-04
                                             0.0001729
                     T4
                               9.956849e-05
                                             0.0000550
```

4.4 (4)

```
[8]: I \leftarrow \exp(1) - \exp(-1)
     seed <- matrix(c(1:100), nrow = 100)</pre>
     Iall1 <- unlist(mapply(randompoint_samp, seed = seed, MoreArgs = list(N = N, h_
      \Rightarrow= exp, low = -1, high = 1, m = 0,
                                                                                   M = \exp(1), \sqcup
      \rightarrow B = 1))[1,])
     Iall2 <- unlist(mapply(average_samp, seed = seed, MoreArgs = list(N = N, h = 0
      \rightarrowexp, low = -1, high = 1, B = 1))[1,])
     Iall3 <- unlist(mapply(importance_samp, seed = seed, MoreArgs = list(N = N, g = __ )
      \rightarrowg, h = exp, Ginver = Ginver, B = 1))[1,])
     Iall4 <- unlist(mapply(hierachical_samp, seed = seed, MoreArgs = list(n = _ _
      \rightarrowc(5000, 5000), h = exp, seg = c(-1, 0, 1), B = 1))[1,])
     MAE <- data.frame(name = c("I1", "I2", "I3", "I4"), MAE = c(sum(abs(Iall1 - U))
      \hookrightarrowI)), sum(abs(Iall2 - I)),
                                                                           sum(abs(Iall3 -___
      \rightarrowI)), sum(abs(Iall4 - I))) / B)
     MAE
```

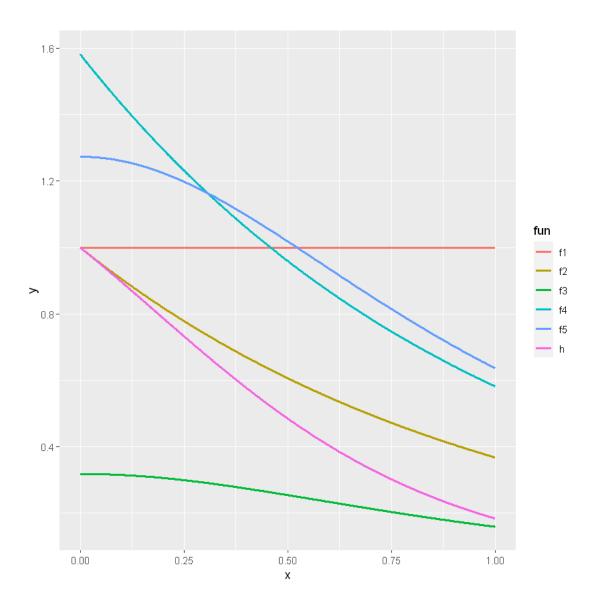
 $MAE(I_1) > MAE(I_2) = MAE(I_3) > MAE(I_4)$. Hierarchical sampling is the most accurate method.

5 3.3

5.1 (1)

plot all the functions in $x \in [0,1]$

```
[9]: library(ggplot2)
x1 <- seq(length = 1000, from = 0, to = 1)
h <- function(x) {
    return(exp(-x)/(1+x^2))
}</pre>
```



5.2(2)

The domain of f_2 and f_3 is larger than (0,1). When sampling a number outside the domain (0,1), we need to drop it.

```
[11]: F1inver <- function(y) {</pre>
          return(y)
      F2inver <- function(y) {
          return(-log(1-y))
      F3inver <- function(y) {
          return(tan(pi*(y-1/2)))
      }
      F4inver <- function(y) {
          return(-log(1-y*(1-exp(-1))))
      F5inver <- function(y) {
          return(tan(pi/4*y))
      }
      set.seed(1)
      I1 <- importance_samp(N, f1, h, F1inver, B = 100)</pre>
      I2 \leftarrow importance\_samp2(N, f2, h, F2inver, 0, 1, B = 100)
      I3 <- importance_samp2(N, f3, h, F3inver, 0, 1, B = 100)
      I4 <- importance_samp(N, f4, h, F4inver, B = 100)
      I5 <- importance_samp(N, f5, h, F5inver, B = 100)</pre>
      result <- data.frame(name = c("I1", "I2", "I3", "I4", "I5"),
                            mean = c(I1\$mean, I2\$mean, I3\$mean, I4\$mean, I5\$mean),
                            var = c(I1$var, I2$var, I3$var, I4$var, I5$var))
      result
```

	$_{\mathrm{name}}$	mean	var
A data.frame: 5×3	<chr $>$	<dbl $>$	<dbl></dbl>
	I1	0.5244436	4.849249e-06
	I2	0.5242430	1.510319 e-05
	I3	0.5246926	6.861769 e-05
	I4	0.5246850	8.236157e-07
	I5	0.5246004	1.625485 e - 06

5.3(3)

From the plot in (1), the shape of $f_1(x)$ is very different from h(x), so it may have larger variance. Compared with $f_4(x)$ and $f_5(x)$, the domain of f_2 and f_3 is larger than (0,1) and we drop many sample points, which can explain their larger variace. Actually, if we transform f_2 and f_3 to (0,1), they are the same as f_4 and f_5 .

5.4(4)

```
[12]: n <- rep(1000, 10)
    seg <- seq(from = 0, to = 1, length = 11)
    I6 <- hierachical_samp(n, h, seg, B = 100)
    I6$mean
    I6$var</pre>
```

0.524789484149965

4.88366826222391e-07

 $Var(I_6)$ is smaller than the above.

5.5 (5)

```
[13]: #hierachical sampling another version
hierachical_samp2 <- function(m, h, B = 1) {
        Iall <- rep(0, B)
        for (i in 1:B) {
            set.seed(i)
            u <- runif(m, 0, 1)
            n <- c(1:m)
            Iall[i] <- sum(h((n - 1 + u) / m)) / m
        }
        I <- mean(Iall)
        Var <- mean((Iall - I)^2)
        return(list(mean = I, var = Var))
    }
    I7 <- hierachical_samp2(N, h, B = 100)
    I7$mean
    I7$var</pre>
```

0.524797107715719

5.48934860538895e-14

 $Var(I_7)$ is smaller than the above. Hierarchical sampling further reduces the variance.

6 3.9

4.90653697166515

0.00342042897686122

```
[15]: # antithetic sampling
      set.seed(2)
      antithetic_sampling <- function(N, m, h, low, high, B = 1) {
           Iall <- rep(0, B)</pre>
           for (j in 1:B) {
               set.seed(j)
               u <- matrix(0, nrow = N, ncol = m)</pre>
               for (i in 1:m) {
                    u[,i] <- runif(N, low[i], high[i])</pre>
               Iall[j] \leftarrow prod(high - low) * (sum(apply(u, 1, h)) + 
       \rightarrowsum(apply(low+high-u, 1, h))) / (2 * N)
           I <- mean(Iall)</pre>
           Var <- mean((Iall - I)^2)</pre>
           return(list(mean = I, var = Var))
      }
      low < -c(0, 0)
      high <- c(1, 1)
      h <- function(x) {
          return(exp((sum(x))^2))
      I <- antithetic_sampling(N, 2, h, low, high, B = 100)</pre>
      I$mean
      I$var
```

4.89894833859011

0.00107924962051474

Antithetic sampling can reduce variance.