Advanced Statistical Methods HW8

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Exercise 10.4

4. Verify formula (10.38) for the number of distinct bootstrap samples.

ure 10.3). This rapidly becomes impractical. The number of distinct bootstrap samples for *n* points turns out to be

$$\binom{2n-1}{n}.\tag{10.38}$$

Given n data points, we want to figure out number of distinct tuples

$$(x_1,\cdots,x_n)$$

satisfying $x_1 + \cdots + x_n = n$ and $x_i \in \{0, 1, \cdots, n\}$ for each $i = 1, \cdots, n$

It is a problem equivalent to the following: for n equally shaped balls, we are going to separate those balls into n ordered groups, with group of zero ball being allowed.

The cardinality of the first group will be equal to x_1 and the cardinality of the second group will be equal to x_2 and so on.

Hence, it is a problem of locating separation bars between those n number of balls. Since we need to separate those balls into n number of ordered groups, we need n-1 separation bars. For example, if n=8 and balls are represented by \circ and separation bars are represented by | then

the above example represents the case of

$$(x_1, x_2, \cdots, x_n) = (1, 1, 2, 0, 0, 1, 1, 2)$$

Hence, the number of distinct bootstrap samples for n points can be calculated by

$$\binom{n+(n-1)}{n} = \binom{2n-1}{n}$$

Exercise 10.5

5. A normal theory least squares model (7.28)–(7.30) yields $\hat{\beta}$ (7.32). Describe the parametric bootstrap estimates for the standard errors of the components of $\hat{\beta}$.

Linear regression, perhaps the most widely used estimation technique, is based on a version of $\hat{\mu}^{\text{MLE}}$. In the usual notation, we observe an *n*-dimensional vector $\mathbf{y} = (y_1, y_2, \dots, y_n)'$ from the linear model

$$y = X\beta + \epsilon. \tag{7.28}$$

Here X is a known $n \times p$ structure matrix, β is an unknown p-dimensional parameter vector, while the noise vector $\boldsymbol{\epsilon} = (\epsilon_1, \epsilon_2, \dots, \epsilon_n)'$ has its components uncorrelated and with constant variance σ^2 ,

$$\boldsymbol{\epsilon} \sim (\boldsymbol{0}, \sigma^2 \boldsymbol{I}), \tag{7.29}$$

where I is the $n \times n$ identity matrix. Often ϵ is assumed to be multivariate normal,

$$\epsilon \sim \mathcal{N}_n(\mathbf{0}, \sigma^2 \mathbf{I}),$$
 (7.30)

From the model above given as

$$\mathbf{y} \sim N(\mathbf{X}\beta, \sigma^2 \mathbf{I})$$

and the ordinary least squares estimates of β is derived as

$$\hat{\beta} = (\mathbf{X}^T \mathbf{X})^{-1} \mathbf{X}^T \mathbf{y}$$

Also, the estimates for variance is calculated by MSE

$$\hat{\sigma}^2 = \frac{1}{n-n} (\mathbf{y} - \mathbf{X}\hat{\beta})^T (\mathbf{y} - \mathbf{X}\hat{\beta})$$

For the parametric boostrap, we generate bootstrap sample $\mathbf{y}_1^*, \cdots, \mathbf{y}_B^*$ from

$$\mathbf{y}^* \sim N(\mathbf{X}\hat{\beta}, \hat{\sigma}^2 \mathbf{I})$$

Then, the corresponding bootstrap samples for $\hat{\beta}$ is computed by

$$\hat{\beta}^{*(1)} = (\mathbf{X}^T \mathbf{X})^{-1} \mathbf{X}^T \mathbf{y}_1^*, \cdots, \ \hat{\beta}^{*(B)} = (\mathbf{X}^T \mathbf{X})^{-1} \mathbf{X}^T \mathbf{y}_B^*$$

From this bootstrap sample $\hat{\beta}^*$'s , we can derive parametric bootstrap estimate for the standard errors of the components of $\hat{\beta}$ as the following :

$$\hat{\text{se}}_{boot}(\hat{\beta}_1) = \text{sd}(\hat{\beta}_1^{*(1)}, \dots, \hat{\beta}_1^{*(B)}), \dots, \hat{\text{se}}_{boot}(\hat{\beta}_p) = \text{sd}(\hat{\beta}_p^{*(1)}, \dots, \hat{\beta}_p^{*(B)})$$

where
$$sd(x_1, \dots, x_n) = \sqrt{\frac{1}{n-1} \sum_{i=1}^{n} (x_i - \overline{x})^2}$$

Exercise 10.7

7. Verify formula (10.70).

For the sample mean \bar{x} , the jackknife yields exactly the usual variance estimate (1.2), $\sum_i (x_i - \bar{x})^2 / (n(n-1))$, while the ideal bootstrap estimate $(B \to \infty)$ gives

$$\sum_{i=1}^{n} (x_i - \bar{x})^2 / n^2. \tag{10.70}$$

First, we shall show that for $\hat{\theta} = \overline{x} = \frac{1}{n} \sum_{i=1}^{n} x_i$,

$$\hat{\text{var}}_{jack}(\hat{\theta}) = \frac{\sum_{i=1}^{n} (x_i - \overline{x})^2}{n(n-1)}$$

Note that

$$var_{jack}(\hat{\theta}) = \frac{n-1}{n} \sum_{i=1}^{n} (\hat{\theta}_{(i)} - \hat{\theta}_{(\cdot)})^2$$

For $\hat{\theta} = \overline{x}$,

$$\begin{split} \hat{\theta}_{(i)} &= \frac{n\overline{x} - x_i}{n - 1} \\ \hat{\theta}_{(\cdot)} &= \frac{1}{n} \sum_{i=1}^n \hat{\theta}_{(i)} = \frac{1}{n} \sum_{i=1}^n \frac{n\overline{x} - x_i}{n - 1} = \frac{1}{n} \frac{n^2 \overline{x} - n\overline{x}}{n - 1} = \overline{x} \\ \hat{\theta}_{(i)} - \hat{\theta}_{(\cdot)} &= \frac{n\overline{x} - x_i}{n - 1} - \overline{x} = \frac{\overline{x} - x_i}{n - 1} \\ \hat{\text{vâr}}_{jack}(\hat{\theta}) &= \frac{n - 1}{n} \sum_{i=1}^n \frac{(\overline{x} - x_i)^2}{(n - 1)^2} = \frac{\sum_{i=1}^n (x_i - \overline{x})^2}{n(n - 1)} \end{split}$$

Next, we shall prove that for the original estimate $\hat{\theta} = S(P_0)$ and resampling estimate $\hat{\theta}^* = S(P)$ with resampling vector P, there is a linear approximation $S_L(P)$ of S(P) such that $S(P_{(i)}) = S_L(P_{(i)})$. Here, P is by definition a vector of nonnegative weights summing to 1 and $P_{(i)}$ is a resampling vector corresponding to i-th jackknife value $\hat{\theta}_{(i)}$, which is given by $\frac{1}{n-1}(1,1,\cdots,1,0,1,\cdots,1)$. Also, P_0 is defined as $\frac{1}{n}(1,1,\cdots,1)$. We shall define $S_L(P)$ by the following:

$$a = \hat{\theta}_{(\cdot)}$$

$$\mathbf{b} = (n-1)(\hat{\theta}_{(i)} - \hat{\theta}_{(\cdot)})$$

$$S_L(P) = a - \mathbf{b}^T P = \hat{\theta}_{(\cdot)} - (n-1) \sum_{i=1}^n p_i(\hat{\theta}_{(i)} - \hat{\theta}_{(\cdot)})$$

$$= n\hat{\theta}_{(\cdot)} - (n-1) \sum_{i=1}^n p_i \hat{\theta}_{(i)}$$

Observe that $S_L(P)$ is linear in P. Now we shall check that $S(P_{(i)}) = S_L(P_{(i)})$ for each $i = 1, \dots, n$

$$S_L(P_{(i)}) = n\hat{\theta}_{(\cdot)} - (n-1)\sum_{j \neq i} \frac{1}{n-1}\hat{\theta}_{(j)} = \sum_{j=1}^n \hat{\theta}_{(j)} - \sum_{j \neq i} \hat{\theta}_{(j)} = \hat{\theta}_{(i)} = S(P_{(i)})$$

Now, we should consider what ideal bootstrap estimate is. For boootstrap estimate $\hat{\theta}^* = S(P^*)$, ideal bootstrap estimate of variance of $\hat{\theta}$ can be written as $Var(S(P^*))$ where P^* denotes the bootstrap resampling

vector with the form $\frac{1}{n}(N_1, \dots, N_n)$ and $nP^* \sim Multinomial(n, P_0)$ Here, by the property of multinomial distribution, we have

$$Var(nP^*) = n \left[\text{Diag}(P_0) - P_0 P_0^T \right]$$

Now, we shall claim that

$$Var(S_L(P^*)) = \frac{n-1}{n} \hat{\text{var}}_{jack}(\hat{\theta})$$

In above, we've shown that $S_L(P^*)$ is a linear approximation of $S(P^*)$ given as $S_L(P^*) = a - \mathbf{b}^T P^*$. Observe that randomness only lies in P^* for $a - \mathbf{b}^T P^*$. Thus,

$$Var(S_L(P^*)) = Var(a - \mathbf{b}^T P^*) = Var(\mathbf{b}^T P^*) = \mathbf{b}^T Var(P^*) \mathbf{b} = \frac{1}{n^2} \mathbf{b}^T Var(nP^*) \mathbf{b}$$

$$= \frac{1}{n^2} \mathbf{b}^T n \left[\text{Diag}(P_0) - P_0 P_0^T \right] \mathbf{b}$$

$$= \frac{1}{n} \{ \mathbf{b}^T \text{Diag}(P_0) \mathbf{b} - (P_0^T \mathbf{b})^2 \}$$

$$= \frac{1}{n} \mathbf{b}^T \text{Diag}(P_0) \mathbf{b} \quad \because P_0^T \mathbf{b} = \frac{n-1}{n} \sum_{i=1}^n (\hat{\theta}_{(i)} - \hat{\theta}_{(\cdot)}) = 0$$

$$= \frac{(n-1)^2}{n^2} \sum_{i=1}^n (\hat{\theta}_{(i)} - \hat{\theta}_{(\cdot)})^2 = \frac{n-1}{n} \text{vâr}_{jack}(\hat{\theta})$$

Hence, we've shown our claim above.

Note that $\hat{\theta} = \overline{x}$ can be written as $\hat{\theta} = S(P_0)$ with $S(P) = \sum_{i=1}^n p_i x_i = \mathbf{x}^T P$ so that S(P) is linear in P. Furthermore, in this case $S_L(P)$ agrees with S(P).

$$S_L(P) = a - \mathbf{b}^T P = \hat{\theta}_{(\cdot)} - (n-1) \sum_{i=1}^n p_i (\hat{\theta}_{(i)} - \hat{\theta}_{(\cdot)}) = \overline{x} - (n-1) \sum_{i=1}^n p_i \frac{\overline{x} - x_i}{n-1} = \overline{x} - \sum_{i=1}^n p_i (\overline{x} - x_i) = \sum_{i=1}^n p_i x_i = S(P)$$

Hence, combining all the results above, the ideal bootstrap variance estimate of the sample mean \overline{x} is given as

$$Var(S(P^*)) = \frac{n-1}{n} \cdot \frac{\sum_{i=1}^{n} (x_i - \overline{x})^2}{n(n-1)} = \frac{1}{n^2} \sum_{i=1}^{n} (x_i - \overline{x})^2$$

Exercise 10.9

9. A survey in a small town showed incomes x_1, x_2, \ldots, x_m for men and y_1, y_2, \ldots, y_n for women. As an estimate of the differences,

$$\hat{\theta} = \text{median}\{x_1, x_2, \dots, x_m\} - \text{median}\{y_1, y_2, \dots, y_n\}$$

was computed.

- (a) How would you use nonparametric bootstrapping to assess the accuracy of $\hat{\theta}$?
- (b) Do you think your method makes full use of the bootstrap replications?

(a) First, by resampling m samples among x_1, \dots, x_m , we get bootstrap median $med(X)^*$. By repeating this B times, we have $med(X)^{*1}, \dots, med(X)^{*B}$. Similarly, by resampling n samples among y_1, \dots, y_n , we get bootstrap median $med(Y)^*$ and by repeating this B times, we have $med(Y)^{*1}, \dots, med(Y)^{*B}$. With large enough B, say B=200, we have B number of θ^* values derived by

$$\hat{\theta}^{*1} = med(X)^{*1} - med(Y)^{*1}, \dots, \hat{\theta}^{*B} = med(X)^{*B} - med(Y)^{*B}$$

Then we have nonparametric bootstrap estimates for standard error of $\hat{\theta}$ as

$$\hat{\text{se}}_{boot}(\hat{\theta}) = \sqrt{\frac{1}{B-1} \sum_{b=1}^{B} (\hat{\theta}^{*b} - \hat{\theta}^{*\cdot})^2} \text{ where } \hat{\theta}^{*\cdot} = \frac{1}{B} \sum_{b=1}^{B} \hat{\theta}^{*b}$$

(b) To make full use of the bootstrap replications, we can consider a parametric bootstrap. Since the survey has implemented in a small town, it is expected that m and n are not so big. Hence, it may be useful to take advantage of the power of parametric inference. For example, assume parametric model $X \sim N(\mu_x, \sigma_x^2)$ and $Y \sim N(\mu_y, \sigma_y^2)$. Then, derive MLE of $\hat{\mu}_x$, $\hat{\mu}_y$, $\hat{\sigma}_x^2$, $\hat{\sigma}_y^2$ using observed data $x_1, \cdots, x_m, y_1, \cdots, y_n$. Now, generate bootstrap sample from $X^* \sim N(\hat{\mu}_x, \hat{\sigma}_x^2)$ and $Y^* \sim N(\hat{\mu}_y, \hat{\sigma}_y^2)$. The rest of the procedure is same for the nonparametric bootstrapping above.

Exercise 11.1

1. We observe $y \sim \lambda G_{10}$ to be y = 20. Here λ is an unknown parameter while G_{10} represents a gamma random variable with 10 degrees of freedom $(y \sim G(10, \lambda))$ in the notation of Table 5.1). Apply the Neyman construction as in Figure 11.1 to find the confidence limit endpoints $\hat{\lambda}(0.025)$ and $\hat{\lambda}(0.975)$.

Our model is $Y|\lambda \sim \Gamma(10,\lambda)$ and the observed value is y=20. We will define some necessary notations.

$$f_{\lambda}(y) = \frac{1}{\Gamma(10)\lambda^{10}} y^9 \exp(-y/\lambda)$$
 : pdf of $y|\lambda$

 $q_{\alpha}(f_{\lambda}) = \alpha$ -quantile of y for given λ i.e. $P_{\lambda}(Y \leq q_{\alpha}(f_{\lambda})) = \alpha$

For given $\alpha \in (0,1)$, $q_{\alpha}(f_{\lambda})$ is a function of λ

$$I_{\lambda}(y) = I\left\{q_{\frac{\alpha}{2}}(f_{\lambda})\right\} \le y \le q_{1-\frac{\alpha}{2}}(f_{\lambda})$$

$$C(y) = \{\lambda : I_{\lambda}(y) = 1\}$$

Coverage probability: $P_{\lambda}(\lambda \in C(Y)) = P_{\lambda}(I_{\lambda}(Y) = 1) = P_{\lambda}(q_{\frac{\alpha}{2}}(f_{\lambda})) \le Y \le q_{1-\frac{\alpha}{2}}(f_{\lambda})) = 1 - \alpha$

Now, we shall claim that for given $\alpha \in (0,1)$, $q_{\alpha}(f_{\lambda})$ is an increasing function of λ . It is enough to show that $Y|\lambda_2$ is stochastically larger than $Y|\lambda_1$ whenever $\lambda_1 \leq \lambda_2$ i.e. $P_{\lambda_1}(Y \geq r) \leq P_{\lambda_2}(Y \geq r)$ for any r > 0 provided $\lambda_1 \leq \lambda_2$ Take arbitrary $0 < \lambda_1 < \lambda_2$. Write $\lambda = \lambda_1$. Then there is a constant c > 1 such that $\lambda_2 = c\lambda$. Take r > 0. Then the following holds true.

$$P_{\lambda_{1}}(Y \geq r) = P_{\lambda}(Y \geq r) = \frac{1}{\Gamma(10)\lambda^{10}} \int_{r}^{\infty} y^{9} \exp(-y/\lambda) \, dy$$

$$P_{\lambda_{2}}(Y \geq r) = P_{c\lambda}(Y \geq r) = \frac{1}{\Gamma(10)c^{10}\lambda^{10}} \int_{r}^{\infty} y^{9} \exp(-y/c\lambda) \, dy$$

$$= \frac{1}{\Gamma(10)c^{10}\lambda^{10}} \int_{r/c}^{\infty} c^{9}z^{9} \exp(-z/\lambda) \, c \, dz \quad \because z = \frac{y}{c} , \, dz = \frac{1}{c}dy$$

$$= \frac{1}{\Gamma(10)\lambda^{10}} \int_{r/c}^{\infty} z^{9} \exp(-z/\lambda) \, dz = P_{\lambda}(Y \geq \frac{r}{c})$$

$$\Rightarrow P_{\lambda_{2}}(Y \geq r) = P_{c\lambda}(Y \geq r) = P_{\lambda}(Y \geq \frac{r}{c}) \geq P_{\lambda}(Y \geq r) = P_{\lambda_{1}}(Y \geq r)$$

$$\therefore P_{\lambda_{1}}(Y \geq r) \leq P_{\lambda_{2}}(Y \geq r)$$

We've shown that $Y|\lambda_2$ is stochastically larger than $Y|\lambda_1$ whenever $\lambda_1 \leq \lambda_2$ Hence if we denote the cdf of $y|\lambda$ as F_{λ} , then

$$1 - F_{\lambda_1}(r) \le 1 - F_{\lambda_2}(r) \; , \; F_{\lambda_1}(r) \ge F_{\lambda_2}(r) \; , \; F_{\lambda_1}^{-1}(\alpha) \le F_{\lambda_2}^{-1}(\alpha) \; , \; q_{\alpha}(f_{\lambda_1}) \le q_{\alpha}(f_{\lambda_2})$$

for given r > 0 and $\alpha \in (0,1)$ provided $\lambda_1 \leq \lambda_2$. Therefore $q_{\alpha}(f_{\lambda})$ is an increasing function of λ for fixed $\alpha \in (0,1)$. Hence $\lambda \in C(y)$ is satisfied for some closed interval $[\lambda_l(y), \lambda_u(y)]$ where

$$y = q_{1-\frac{\alpha}{2}}(f_{\lambda_l(y)})$$
 and $y = q_{\frac{\alpha}{2}}(f_{\lambda_u(y)})$

We should find such $\lambda_l(y), \lambda_u(y)$ such that

$$F_{\lambda_l(y)}(y) = 1 - \frac{2}{\alpha}$$
 and $F_{\lambda_u(y)}(y) = \frac{\alpha}{2}$

```
where F_{\lambda} is a cdf of \Gamma(10, \lambda) and y = 20 is observed.
```

```
pgamma(20, shape=10, scale = 1) # lambda=1
## [1] 0.9950046
pgamma(20, shape=10, scale = 1.5) # lambda=1.5
## [1] 0.8550948
\# Since F_{\lambda} and G(y) is decreasing in \lambda , \lambda , \lambda , \lambda
candidate=seq(from=1, to=1.5, length=1000)
p_L=0
lambda_L=0
for(i in 1:length(candidate)){
  lambda_L[i] = candidate[i]
 p_L[i] = pgamma(20, shape=10, scale = candidate[i])
i_L = which.min(abs(p_L-0.975)) # 1- 2/alpha = 0.975
lambda_L[i_L]
## [1] 1.170671
abs(p_L[i_L]-0.975) < 1e-5
## [1] TRUE
pgamma(20, shape=10, scale = 4) # lambda=4
## [1] 0.03182806
pgamma(20, shape=10, scale = 4.5) # lambda=4.5
## [1] 0.01584119
\# Since F_{\lambda} is decreasing in \lambda, \lambda with H is decreasing in \lambda.
candidate=seq(from=4, to=4.5, length=1000)
p_U=0
lambda_U=0
for(i in 1:length(candidate)){
  lambda_U[i] = candidate[i]
  p_U[i] = pgamma(20, shape=10, scale = candidate[i])
}
i_U = which.min(abs(p_U-0.025)) # 2/alpha = 0.025
lambda_U[i_U]
## [1] 4.170671
abs(p_U[i_U]-0.025) < 1e-5
```

[1] TRUE

Therefore, the confidence limit endpoints are computed as

$$\hat{\lambda}(0.025) = 1.1707$$
 and $\hat{\lambda}(0.975) = 4.1707$

Exercise 11.3

3. Suppose \widehat{G} in (11.33) was perfectly normal, say $\widehat{G} \sim \mathcal{N}(\widehat{\mu}, \widehat{\sigma}^2)$. What does $\widehat{\theta}_{BC}(\alpha)$ reduce to in this case, and why does this make intuitive sense?

$$p_0 = \#\left\{\hat{\theta}^{*b} \le \hat{\theta}\right\} / B \tag{11.31}$$

(an estimate of (11.29)), and define the bias-correction value

$$z_0 = \Phi^{-1}(p_0), \tag{11.32}$$

where Φ^{-1} is the inverse function of the standard normal cdf. The BC level- α confidence interval endpoint is defined to be

$$\hat{\theta}_{BC}[\alpha] = \hat{G}^{-1} \left[\Phi \left(2z_0 + z^{(\alpha)} \right) \right], \tag{11.33}$$

where \hat{G} is the bootstrap cdf (11.16) and $z^{(\alpha)} = \Phi^{-1}(\alpha)$ (11.25).

Suppose $\hat{G} \sim N(\hat{\mu}, \hat{\sigma}^2)$, perfectly normal with $\hat{\mu} = \frac{1}{n} \sum_{i=1}^n x_i$ and $\hat{\sigma}^2 = \frac{1}{n-1} \sum_{i=1}^{n-1} (x_i - \overline{x})^2$. Note that if $F_{\mu,\sigma}$ is a cdf of $N(\mu, \sigma^2)$ then $F_{\mu,\sigma}(x) = \Phi(\frac{x-\mu}{\sigma})$. Hence $F_{\mu,\sigma}^{-1}(\alpha) = \sigma\Phi^{-1}(\alpha) + \mu$. Why?

$$F_{\mu,\sigma}(X \le \sigma \Phi^{-1}(\alpha) + \mu) = F_{\mu,\sigma}\left(\frac{X - \mu}{\sigma} \le \Phi^{-1}(\alpha)\right) = \Phi(\Phi^{-1}(\alpha)) = \alpha \quad \forall \ \alpha \in (0,1)$$

Thus, we have

$$\hat{\theta}_{BC}(\alpha) = \hat{G}^{-1}(\Phi(2z_0 + z_{(\alpha)})) = \hat{\sigma}\Phi^{-1}(\Phi(2z_0 + z_{(\alpha)})) + \hat{\mu} = \hat{\sigma}(2z_0 + z_{(\alpha)}) + \hat{\mu}$$

 $\hat{\theta}_{BC}(\alpha)$ can be reduced to $\hat{\sigma}(2z_0+z_{(\alpha)})+\hat{\mu}$. If $z_0=0$, then $\hat{\theta}_{BC}(\alpha)=\hat{\sigma}z_{(\alpha)}+\hat{\mu}$. This is just a normal α -quantile of assumed bootstrap distribution $N(\hat{\mu},\hat{\sigma})$, which makes sense because when $z_0=0$, we can think that there is no bias in bootstrap distribution.

However, if $z_0 > 0$, then we can think that there is downward bias in bootstrap distribution so that adjusting upward is required. Hence $\hat{\theta}_{BC}(\alpha)$ is adjusted upward by $2\hat{\sigma}z_0$ from the original normal α -quantile. On the other hand, if $z_0 < 0$, then we can think there is upward bias in bootstrap distribution so that adjusting downward is necessary. Thus $\hat{\theta}_{BC}(\alpha)$ is adjusted downward by $2\hat{\sigma}|z_0|$ from the original normal α -quantile.

Exercise 11.5

5. Suppose $\hat{\theta} \sim \text{Poisson}(\theta)$ is observed to equal 16. Without employing simulation, compute the 95% central BCa interval for θ . (You can use the good approximation $z_0 = a = 1/(6\hat{\theta}^{1/2})$.)

The *BCa method* ("bias-corrected and accelerated") takes its level- α confidence limit to be

$$\hat{\theta}_{BCa}[\alpha] = \hat{G}^{-1} \left[\Phi \left(z_0 + \frac{z_0 + z^{(\alpha)}}{1 - a(z_0 + z^{(\alpha)})} \right) \right]. \tag{11.39}$$

 $\hat{\theta} \sim Poi(\theta)$ and $\hat{\theta}=16$ is observed. By the hint , we shall plug in $z_0=a=\frac{1}{6\hat{\theta}^{1/2}}=\frac{1}{24}$ on formula 11.39 with $\alpha=0.025$ and 0.975 .

For bootstrap distribution \hat{G} , from $\hat{\theta} \sim Poi(\theta)$, we expect $\hat{\theta}^* \sim Poi(\hat{\theta})$. Thus, without employing simulation, we can plug in $\hat{G} = \operatorname{cdf}$ of $Poi(\hat{\theta})$

```
theta.hat = 16
z0 = 1/(6*sqrt(theta.hat))
a = 1/(6*sqrt(theta.hat))
alpha = 0.05
L = pnorm(z0+ (z0+qnorm(alpha/2)) / (1-a*(z0+qnorm(alpha/2))))
U = pnorm(z0+ (z0+qnorm(1-alpha/2)) / (1-a*(z0+qnorm(1-alpha/2))))
BCa = c(qpois(L, lambda=theta.hat) ,qpois(U, lambda=theta.hat))
BCa
```

[1] 9 26

Thus, 95% central BCa interval for θ is (9, 26)

Exercise 11.6

6. Use the R program bcajack (available with its help file from efron.web.stanford. edu under "Talks") to find BCa confidence limits for the student score eigenratio statistic as in Figure 10.2.

First, bring R program bcajack from efron.web.stanford.edu - 2018 : Supplement files for bcajack Next, load the student score data.

```
data=read.table("https://web.stanford.edu/~hastie/CASI_files/DATA/student_score.txt", header=T)
```

To take advantage of bcajack function created by Efron, we need to create a function EigenRatio which calculates the ratio $\frac{\text{largest eigenvalue}}{\text{sum of all eigenvalues}}$ of correlation matrix for input dataset.

```
EigenRatio<-function(X){
  R = cor(X)
  result = max(eigen(R, only.values=TRUE)$values) / sum(eigen(R, only.values=TRUE)$values)
  return(result)
}</pre>
```

Then, we can use bcajack function with EigenRatio

```
library(matlib) # To read a `len` function in `bcajack` , we need this library
set.seed(123)
bca_result = bcajack(x = data , B = 1000 , func = EigenRatio , alpha= c(0.025, 0.975), m=20, catj = 0)
```

```
## {1}{2}
```

```
bca_result
```

```
## bcajack(x = data, B = 1000, func = EigenRatio, m = 20, alpha = c(0.025,
       0.975), catj = 0)
##
##
## $lims
##
         bcalims jacksd standard
## 0.025
           0.516
                    NaN
                            0.542 0.028
## 0.975
           0.820
                    NaN
                            0.843 0.978
##
## $stats
##
        thet sdboot
                       z0
                               a sdjack
```

```
## est 0.693 0.077 -0.07 0.051 0.081
## jsd 0.000 0.003 NaN 0.000 0.000
##
## $B.mean
## [1] 1000.0000000 0.6873815
```

From the result above, for $\theta =$ eigenratio , BCa confidence limit endpoints $\hat{\theta}_{BCa}(0.025)$ and $\hat{\theta}_{BCa}(0.975)$ are given as (0.516, 0.820)