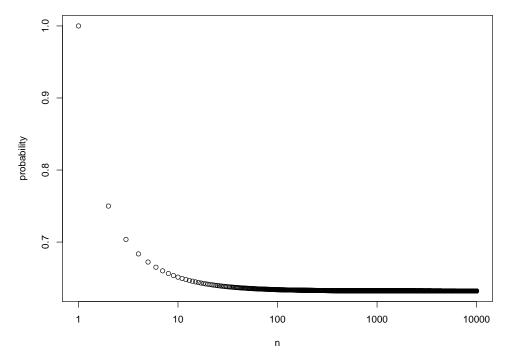
Solution to Series 5

- 1. a) We sample with replacement. The probability that given observation is not picked in every sampling iteration is 1 1/n. Each draw is independent. Hence, we apply the product rule. The solution is $(1 1/n)^n$ since we generate the samples of size n.
 - b) The solution is $1 (1 1/n)^n = 1 (1 1/100)^{100} = 1 (99/100)^{100} = 63.4\%$.
 - c) > foo <- function(n) $\{1 (1 1/n)^n\}$
 - > n <- 1:10000
 - > plot(n, foo(n), log = "x", ylab = "probability")



- d) We expect that about 2/3 of the original observations occur in the bootstrap sample, i.e. $\lim_{n\to\infty} 1 (1-1/n)^n = 1 1/e \approx 63.21\%$ and this limit is visible in the plot.
- 2. For any random variable Z, let $q_Z(\alpha)$ denote the α -quantile of Z. We have:

$$1 - \alpha = P\left(q_{\frac{\hat{\theta}^* - \hat{\theta}}{\widehat{sd}(\hat{\theta}^*)}}(\alpha/2) < \frac{\hat{\theta}^* - \hat{\theta}}{\widehat{sd}(\hat{\theta}^*)} < q_{\frac{\hat{\theta}^* - \hat{\theta}}{\widehat{sd}(\hat{\theta}^*)}}(1 - \alpha/2)\right)$$

$$\approx P\left(q_{\frac{\hat{\theta}^* - \hat{\theta}}{\widehat{sd}(\hat{\theta}^*)}}(\alpha/2) < \frac{\hat{\theta} - \theta}{\widehat{sd}(\hat{\theta})} < q_{\frac{\hat{\theta}^* - \hat{\theta}}{\widehat{sd}(\hat{\theta}^*)}}(1 - \alpha/2)\right)$$

The second equality follows from the bootstrap consistency: if the distribution of $\hat{\theta}^* - \hat{\theta}$ can be approximated by the distribution of $\hat{\theta} - \theta$ and both random variables are divided by their standard deviations, the

distribution of $\frac{\hat{\theta}^* - \hat{\theta}}{\widehat{sd}(\hat{\theta}^*)}$ can be approximated by the distribution of $\frac{\hat{\theta} - \theta}{\widehat{sd}(\hat{\theta})}$. Hence,

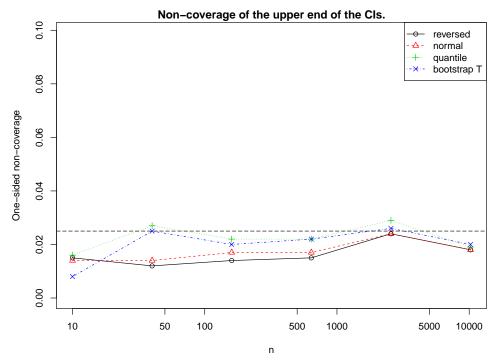
$$\begin{split} 1 - \alpha &\approx P\left(q_{\frac{\hat{\theta}^* - \hat{\theta}}{\widehat{sd}(\hat{\theta}^*)}}(\alpha/2) < \frac{\hat{\theta} - \theta}{\widehat{sd}(\hat{\theta})} < q_{\frac{\hat{\theta}^* - \hat{\theta}}{\widehat{sd}(\hat{\theta}^*)}}(1 - \alpha/2)\right) \\ &= P\left(q_{\frac{\hat{\theta}^* - \hat{\theta}}{\widehat{sd}(\hat{\theta}^*)}}(\alpha/2) \cdot \widehat{sd}(\hat{\theta}) < \hat{\theta} - \theta < q_{\frac{\hat{\theta}^* - \hat{\theta}}{\widehat{sd}(\hat{\theta}^*)}}(1 - \alpha/2) \cdot \widehat{sd}(\hat{\theta})\right) \\ &= P\left(-q_{\frac{\hat{\theta}^* - \hat{\theta}}{\widehat{sd}(\hat{\theta}^*)}}(\alpha/2) \cdot \widehat{sd}(\hat{\theta}) > \theta - \hat{\theta} > -q_{\frac{\hat{\theta}^* - \hat{\theta}}{\widehat{sd}(\hat{\theta}^*)}}(1 - \alpha/2) \cdot \widehat{sd}(\hat{\theta})\right) \\ &= P\left(-q_{\frac{\hat{\theta}^* - \hat{\theta}}{\widehat{sd}(\hat{\theta}^*)}}(1 - \alpha/2) \cdot \widehat{sd}(\hat{\theta}) < \theta - \hat{\theta} < -q_{\frac{\hat{\theta}^* - \hat{\theta}}{\widehat{sd}(\hat{\theta}^*)}}(\alpha/2) \cdot \widehat{sd}(\hat{\theta})\right) \\ &= P\left(\hat{\theta} - q_{\frac{\hat{\theta}^* - \hat{\theta}}{\widehat{sd}(\hat{\theta}^*)}}(1 - \alpha/2) \cdot \widehat{sd}(\hat{\theta}) < \theta < \hat{\theta} - q_{\frac{\hat{\theta}^* - \hat{\theta}}{\widehat{sd}(\hat{\theta}^*)}}(\alpha/2) \cdot \widehat{sd}(\hat{\theta})\right). \end{split}$$

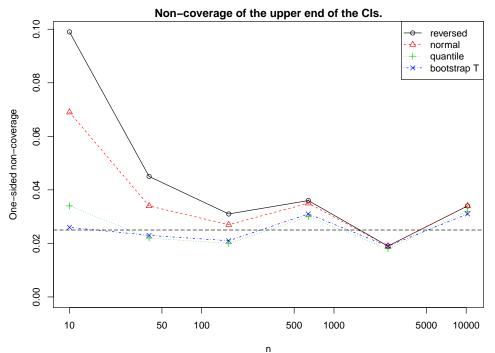
```
3. a) Approximate the true value
      > set.seed(3)
      > (true.par <- mean(rgamma(100000000, shape = 2, rate = 2), trim = 0.1))</pre>
       [1] 0.9103737
   b) > set.seed(1)
      > sample40 <- rgamma(n = 40, shape = 2, rate = 2)
      > mean(sample40, trim = 0.1)
       [1] 0.8824166
   c) > require("boot")
      > tm \leftarrow function(x, ind) \{mean(x[ind], trim = 0.1)\}
      > tm_var <- function(x, ind) {</pre>
         # trimmed mean
         t1 \leftarrow tm(x, ind)
         # bootstrap variance of the trimmed mean (required for the bootstrap T CI)
          t2 \leftarrow var(boot(data = x[ind], statistic = tm, R = 50)$t)
         return(c(t1, t2))
      > res.boot <- boot(data = sample40, statistic = tm_var, R = 10000,
                         sim = "ordinary")
      > boot.ci(res.boot, conf = 0.95, type = c("basic", "norm", "perc", "stud"),
                var.t0 = var(res.boot\$t[, 1]))
      BOOTSTRAP CONFIDENCE INTERVAL CALCULATIONS
      Based on 10000 bootstrap replicates
      boot.ci(boot.out = res.boot, conf = 0.95, type = c("basic", "norm",
           "perc", "stud"), var.t0 = var(res.boot$t[, 1]))
      Intervals:
      Level
                                      Basic
                  Normal
             (0.7022, 1.0569) (0.6945, 1.0492)
      Level
               Studentized
                                     Percentile
             (0.6995, 1.0792) (0.7156, 1.0704)
      Calculations and Intervals on Original Scale
   d) We first define two functions before running the simulation.
      > ##' Checks if a confidence interval contains the true parameter (separately
      > ##' for the lower and the upper end)
      > ##' @param ci: Output of the function boot.ci which contains CIs
      > ##' @param ty: Type of confidence interval
```

```
> ##' @param true.par: True parameter
> ##'
> ##' @return Vector with two elements where first one corresponds to the lower
> ##'
             end and the second to the upper end of the confidence interval.
> ##'
             If the CI is [CI_1, CI_u], the first element is 1 if theta < CI_1
> ##'
             and 0 otherwise. The second element is 1 if theta > CI_u and 0
> ##'
              otherwise.
> check_ci <- function(ci, ty, true.par) {</pre>
   # Get confidence interval of type ty from object ci
   lower.upper <- switch (ty,</pre>
     "norm" = ci[["normal"]][2:3],
     "perc" = ci[["percent"]][4:5],
     "basic" = ci[["basic"]][4:5],
     "stud" = ci[["student"]][4:5]
   res <- if (true.par < lower.upper[1]) {
     } else if (true.par > lower.upper[2]) {
       c(0, 1)
     } else {
       c(0, 0)
   names(res) <- c("lower", "upper")</pre>
   return(res)
> ##' Runs one simulation run, i.e. creates new data set, calculates bootstrap
> ##' CIs, and checks if true parameter is contained.
> ##'
> ##' @param n: Size of sample
> ##' @param true.par: True parameter
> ##' @param R: Number of bootstrap replicates
> ##' @param type: Type of bootstrap CIs, see function boot.ci
> ##'
> ##' @return A vector containing the result of the function check_ci for each
             of the confidence intervals
> do_sim <- function(n, true.par, R = 1000,</pre>
                    type = c("basic", "norm", "perc", "stud")) {
   # Generate the data
   x \leftarrow rgamma(n = n, shape = 2, rate = 2)
   # Constract the CIs for the trimmed mean
   res.boot <- boot(data = x, statistic = tm_var, R = R, sim = "ordinary",
                    parallel = "multicore", ncpus = 20)
   res.ci <- boot.ci(res.boot, conf = 0.95, type = type,
                     var.t0 = var(res.boot\$t[, 1]))
   # Check if CIs contain true.par
   res <- vector(mode = "integer", length = 0)
   for (ty in type) {
     res <- c(res, check_ci(ci = res.ci, ty = ty, true.par = true.par))
     names(res)[(length(res) - 1):length(res)] <-</pre>
       paste(c(ty, ty), c("lower", "upper"), sep = "_")
   # Alternatively, one could use a function of the apply family, e.g. sapply.
   return(res)
```

```
> ### Run simulation
> ##########################
> set.seed(22)
> require("boot")
> sample.size <- c(10, 40, 160, 640, 2560, 10240)
> n.sim <- 1000
> type <- c("basic", "norm", "perc", "stud")</pre>
> # The object RES is used to store the results, i.e. each row corresponds
> # to non-coverage rate for the lower and upper end of the confidence intervals,
> # i.e. the percentage of times that theta < CI_{-}1 and the percentage of times
> # that theta > CI_u if the CI is denoted by (CI_1, CI_u). The last column
> # corresponds to the number of observations.
> RES <- matrix(NA, nrow = length(sample.size), ncol = length(type) * 2 + 1)
> colnames(RES) <- c(paste(rep(type, each = 2),
                          rep(c("lower", "upper"), times = length(type)),
                          sep = "_"), "n")
> for (j in 1:length(sample.size)) {
   n <- sample.size[j]</pre>
   # The object res.sim is used to store the results, i.e. each row corresponds
   \# to the output of the function do_sim. This means that each row contains 0
   # and 1 encoding whether the true parameter lied in the CI or outside. See
   # function check_ci.
   res.sim <- matrix(NA, nrow = n.sim, ncol = length(type) * 2)
   for (i in 1:n.sim) {
     # Calculate CIs and check if true.par is contained
     res.sim[i, ] <- do_sim(n = n, true.par = true.par, type = type, R = 2000)
   # Calculate the upper and lower non-coverage rate
   RES[j, ] \leftarrow c(apply(res.sim, 2, mean), n)
Note that the above code runs in parallel on 20 cores. We chose larger values for some parameters
then you were ask to do on the exercise sheet, i.e. R = 2000, n.sim = 1000, and sample.size =
c(10, 40, 160, 640, 2560, 10240).
The plots have the same limits on the y-axis and we use log-scale for the x-axis.
> y.lim <- max(RES[, -ncol(RES)])</pre>
> # Plot of lower non-coverage
> plot(basic_lower ~ n, data = RES, col = 1, pch = 1, ylim = c(0, y.lim),
      log = "x", ylab = "One-sided non-coverage",
      main = "Non-coverage of the upper end of the CIs.")
> points(norm_lower ~ n, data = RES, col = 2, pch = 2, xlog = TRUE)
> points(perc_lower ~ n, data = RES, col = 3, pch = 3, xlog = TRUE)
> points(stud_lower ~ n, data = RES, col = 4, pch = 4, xlog = TRUE)
> lines(basic_lower ~ n, data = RES, col = 1, lty = 1, xlog = TRUE)
> lines(norm_lower ~ n, data = RES, col = 2, lty = 2, xlog = TRUE)
> lines(perc_lower ~ n, data = RES, col = 3, lty = 3, xlog = TRUE)
> lines(stud_lower ~ n, data = RES, col = 4, lty = 4, xlog = TRUE)
> abline(h = 0.025, 1ty = 5)
> legend("topright", legend = c("reversed", "normal", "quantile", "bootstrap T"),
```

pch = 1:4, lty = 1:4, col = 1:4)





In this setting, the reversed bootstrap CI and the normal approximation CI are biased in the sense that they estimate a too small upper end of the CI for small sample sizes. There are only small differences between the CIs for large sample sizes.