

Resampling methods — 2010



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1 Introduction

Homepage: <http://www.kirchkamp.de/>

Literature:

- Bradley Efron and Robert J. Tibshirani, An Introduction to the Bootstrap, Chapman & Hall, 1994.
- A. C. Davison, D. V. Hinkley, Bootstrap Methods and their Application, Cambridge University Press, 1997.

Aim of the course In this course we will discuss resampling methods like the bootstrap, jackknife, or permutation tests. Resampling methods provide solutions to a couple of interesting problems without making specific distributional assumptions.

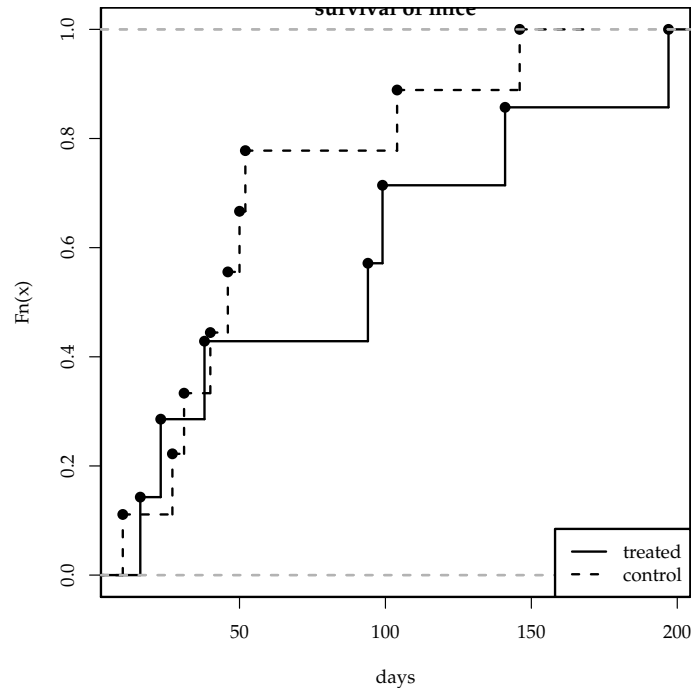
Purpose of this handout In this handout you find all slides from the lecture (in a more printer friendly version). You also find (most of) the examples in R I plan to use in the lecture.

1.1 Motivation

Let us start with a simple example: a comparison of two treatments:

```
library(bootstrap)
library(boot)

par(mar = c(4, 4, 0, 0))
plot(ecdf(mouse.t), xlim = range(mouse.c, mouse.t), verticals = TRUE,
     main = "survival of mice", xlab = "days")
lines(ecdf(mouse.c), lty = "dashed", verticals = TRUE)
legend("bottomright", c("treated", "control"), lty = c("solid",
  "dashed"), bg = "white")
```



Treated mice have a larger mean survival time:

```
| mean(mouse.t)
```

```
[1] 86.85714
```

```
| mean(mouse.c)
```

```
[1] 56.22222
```

Question: Is the difference significant?

Traditionally we estimate

$$\bar{x} = \frac{1}{n} \sum_{i=1}^n x_i \quad \hat{\sigma}_x^2 = \frac{1}{n-1} \sum_{i=1}^n (x_i - \bar{x})^2$$

and hence

$$\hat{\sigma}_{\bar{x}} = \sqrt{\frac{\hat{s}_x^2}{n}}$$

The estimated standard error of the difference between the two means \bar{x} and \bar{y} is

$$\hat{\sigma}_{\bar{x}-\bar{y}} = \sqrt{\frac{\hat{s}_x^2}{n} + \frac{\hat{s}_y^2}{m}}$$

```
| sdDiff <- sqrt(var(mouse.t)/length(mouse.t) + var(mouse.c)/length(mouse.c))
```

```
[1] 28.92647
```

Hence $t = \frac{\bar{x} - \bar{y}}{\sigma_{\bar{x} - \bar{y}}}$ is

```
| t <- (mean(mouse.t) - mean(mouse.c))/sdDiff
```

```
[1] 1.059062
```

Is this is large number?

```
| 2 * (1 - pnorm(t))
```

```
[1] 0.2895715
```

Why are we allowed to do this?

Central limit theorem

Be X_1, \dots, X_n a sequence of independently and identically (i.i.d.) distributed random variables, then with expected value $E(X)$ and variance σ_X^2 , then

$$\lim_{n \rightarrow \infty} \frac{\sum_{i=1}^n X_i - nE(X)}{\sigma \sqrt{n}} = \lim_{n \rightarrow \infty} \left(\frac{\sum_{i=1}^n X_i}{n} - E(X) \right) \cdot \frac{\sqrt{n}}{\sigma} \sim N(0, 1).$$

When do we have a problem?

- When n is small.
- When we are interested in a statistic other than the mean.

Example: Let us assume we are interested in the median:

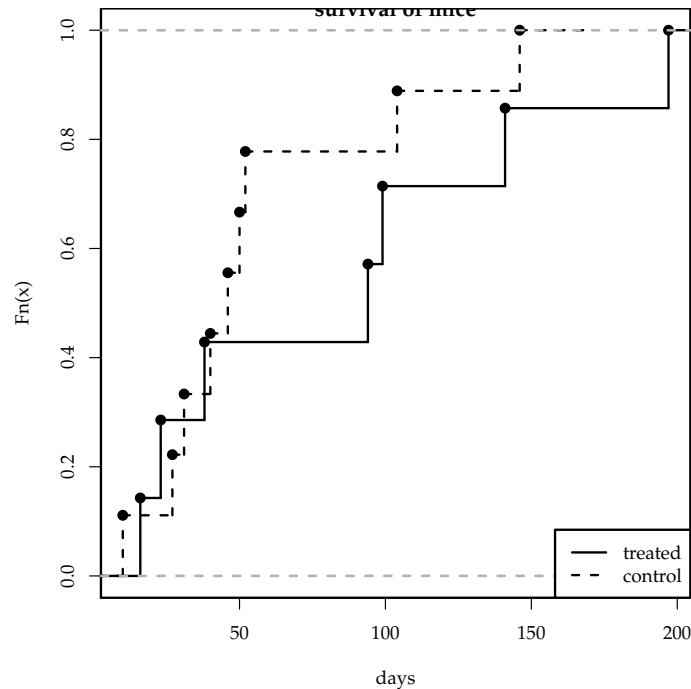
```
| mean(mouse.t) - mean(mouse.c)
```

```
[1] 30.63492
```

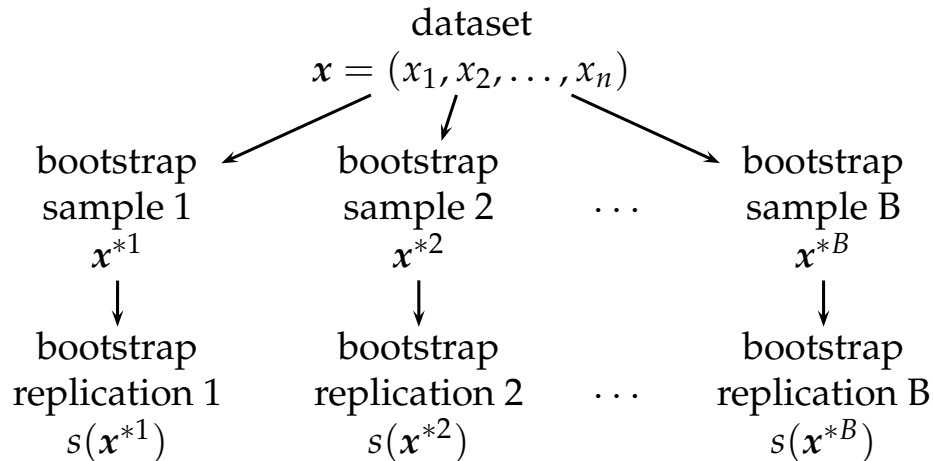
```
| median(mouse.t) - median(mouse.c)
```

```
[1] 48
```

- In this example the difference in medians is larger than the difference in means. But what is the standard error of the difference in medians?
- There are formulas for the standard error of medians, but they require specific distributions of X .
- Do we know the distribution of X ? — No, but we can estimate it!
- Above we have estimated $E(X)$ by using our sample to calculate the statistic: $\bar{x} = \widehat{E(X)}$
- Now we use our sample as an estimation of the distribution of X .



Let us assume that we are interested in the statistic s . Our estimation of the distribution of X is given by the distribution of the dataset x .



Then our estimate of the distribution of s is given by the distribution of the bootstrap replications $s(x^{*1}), s(x^{*2}), \dots, s(x^{*B})$.

Our estimate of the standard error of $s(x)$ is given by the standard deviation of the bootstrap replications $s(x^{*1}), s(x^{*2}), \dots, s(x^{*B})$.

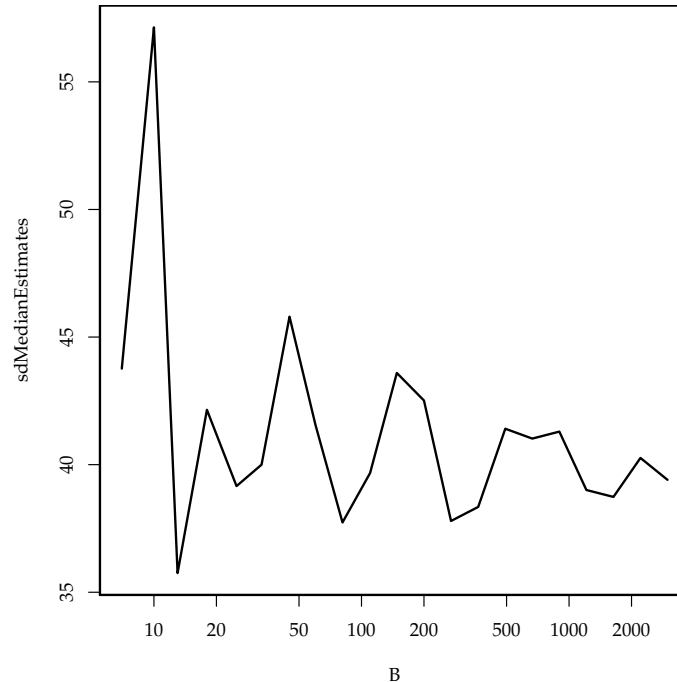
$$\hat{\sigma}_{s,BS} = \text{se}_{b=1}^B(s(x^{*b})) = \sqrt{\frac{1}{B-1} \sum_{b=1}^B \left(s(x^{*b}) - \frac{1}{B} \sum_{b=1}^B s(x^{*b}) \right)^2}$$

```
| set.seed(123)
```

```
| sdDiff2 <- sd(replicate(100, mean(sample(mouse.t, replace = TRUE)) -
|   mean(sample(mouse.c, replace = TRUE))))
| sdMedianDiff2 <- sd(replicate(100, median(sample(mouse.t,
|   replace = TRUE)) - median(sample(mouse.c, replace = TRUE))))
```

```
| sdMedian <- function(B) sd(replicate(B, median(sample(mouse.t,
|   replace = TRUE)) - median(sample(mouse.c, replace = TRUE))))
| set.seed(123)
| B <- round(exp(seq(2, 8, 0.3)))
| sdMedianEstimates <- sapply(B, sdMedian)
```

```
| par(mar = c(4, 4, 0, 0))
| plot(sdMedianEstimates ~ B, log = "x", t = "l")
```



- A number of 50 ... 200 bootstraps is usually sufficient to estimate standard errors.

→ we increase computational complexity by a factor of 50 ... 200.

→ we do not have to worry about “unusual” statistics or distributions.

We replace complicated derivations based on specific distributions with computing power.

Exercise 1.1 *The dataset `x1.csv` contains two variables, `x1` and `x2`.*

- *What is the interquartile range of `x1`?*
- *What is the standard deviation of your estimate?*
- *Draw a graph with the number of bootstrap replications on the horizontal axis and the estimated standard deviation on the vertical axis. How quick does your estimation converge?*

2 Parameters, distributions and the plug-in principle

| | | |
|--|----------------------------|--|
| universe \mathcal{U} | $U_1 \ U_2 \ \cdots \ U_N$ | \leftarrow units |
| population \mathcal{X} (measurements of \mathcal{U}) | $X_1 \ X_2 \ \cdots \ X_N$ | population distribution F |
| random sample x | x_1, x_2, \dots, x_n | empirical distribution \hat{F} |

Note that with independence \hat{F} is a sufficient statistic for F (we can use either x or \hat{F} to learn something about F).

What can we learn from x about X ?

Parameters and distributions

F is a distribution of X iff

$$F(x, \theta, \dots) \equiv P(X < x | \theta, \dots)$$

We call θ a parameter of F .

Parameters and statistics

$$\theta = t(F)$$

$t(\cdot)$ is a statistic of F .

- What happens if we don't know F , but (only) \hat{F} ?

→ If F is determined by a small number of parameters (i.e. we know a lot about F , we miss only θ , e.g. we know that F is normal) then we can use ML or MM. But what if F has many other parameters?

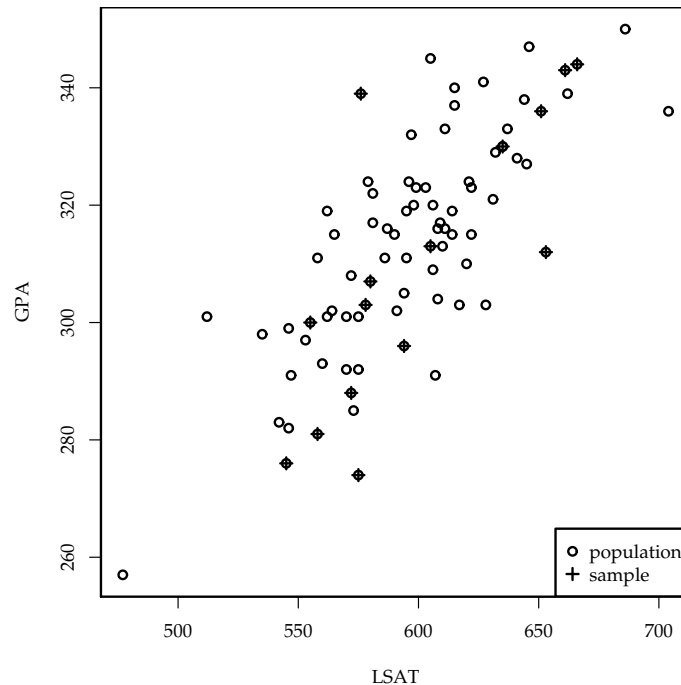
The plug-in principle

$$\hat{\theta} = t(\hat{F})$$

We (simply) apply the statistic $t(\cdot)$ to obtain an estimator for θ .

Example: the law school data

```
par(mar = c(4, 4, 0, 0))
with(law82, plot(100 * GPA ~ LSAT, ylab = "GPA"))
points(law, pch = 3)
legend("bottomright", c("population", "sample"), pch = c(1,
  3))
```



Our population are 82 law schools (dataset `law82`).

Our sample (`law`) only contains 15 observations.

We are interested in the correlation of GPA (Grade Point Average, undergraduate score) and LSAT (Law School Admission Test Score). (i.e. the θ and the $t(\cdot)$ we are talking about is the correlation.)

The true score:

```
| with(law82, cor(GPA, LSAT))
```

```
[1] 0.7599979
```

The plug-in estimate:

```
| with(law, cor(GPA, LSAT))
```

```
[1] 0.7763745
```

Convergence Why does it make sense to use a plug-in estimator?

Glivenko-Cantelli's theorem (1933)

$$\sup_x |\hat{F}_n(x) - F(x)| \xrightarrow{n \rightarrow \infty} 0 \quad \text{almost surely}$$

if $t(\hat{F})$ is continuous then

$$t(\hat{F}_n) \xrightarrow{n \rightarrow \infty} t(F) = \theta \quad \text{almost surely}$$

We can use the bootstrap to measure...

- ... the bias of the plug-in estimate
- ... the standard error of the plug-in estimate

3 Estimating standard errors

3.1 The bootstrap algorithm for standard errors

- x_1, x_2, \dots, x_n are n observed values.
- \hat{F} is the empirical distribution (with probability $1/n$ on each of x_1, x_2, \dots, x_n).
- $\mathbf{x}^* = (x_1^*, x_2^*, \dots, x_n^*)$ is a bootstrap sample, drawn from \hat{F} .
 $(\mathbf{x}^* = (x_1^*, x_2^*, \dots, x_n^*)$ is a random sample of size n drawn with replacement from the observed values x_1, x_2, \dots, x_n).
- $\hat{\theta}^* = s(\mathbf{x}^*)$ is a bootstrap replication of $\hat{\theta}$.
- $\text{se}_F(\hat{\theta})$ is the standard error of $\hat{\theta}$ (we usually can not calculate $\text{se}_F(\hat{\theta})$).
- $\text{se}_{\hat{F}}(\hat{\theta}^*)$ is the (ideal) bootstrap estimate of $\text{se}_F(\hat{\theta})$.
 $\text{se}_{\hat{F}}(\hat{\theta}^*)$ can be hard to calculate.
 We can, however, get a good approximation of $\text{se}_{\hat{F}}(\hat{\theta}^*)$.

The bootstrap algorithm for standard errors

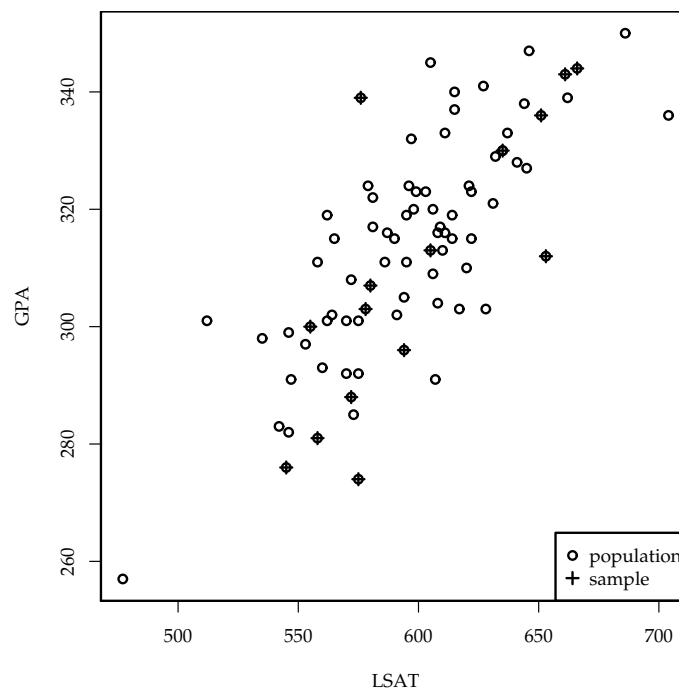
1. Draw B independent bootstrap samples $\mathbf{x}^{*1}, \mathbf{x}^{*2}, \dots, \mathbf{x}^{*B}$.
Each sample has size n and is drawn with replacement from x_1, x_2, \dots, x_n .
2. For bootstrap sample $b = 1, 2, \dots, B$ determine $\hat{\theta}^{*b} = s(\mathbf{x}^{*b})$.

$$3. \hat{\sigma}_{s,BS} = \text{se}_{b=1}^B(\hat{\theta}^{*b}) = \sqrt{\frac{1}{B-1} \sum_{b=1}^B \left(\hat{\theta}^{*b} - \frac{1}{B} \sum_{b=1}^B \hat{\theta}^{*b} \right)^2}$$

$$\lim_{B \rightarrow \infty} \hat{\sigma}_{s,BS} = \underbrace{\text{se}_{\hat{F}}(\hat{\theta}^*)}_{\text{ideal bootstrap}}$$

$$\hat{F} \rightarrow \left. \begin{array}{l} \mathbf{x}^{*1} \rightarrow s(\mathbf{x}^{*1}) = \hat{\theta}^{*1} \\ \mathbf{x}^{*2} \rightarrow s(\mathbf{x}^{*2}) = \hat{\theta}^{*2} \\ \vdots \\ \mathbf{x}^{*b} \rightarrow s(\mathbf{x}^{*b}) = \hat{\theta}^{*b} \\ \vdots \\ \mathbf{x}^{*B} \rightarrow s(\mathbf{x}^{*B}) = \hat{\theta}^{*B} \end{array} \right\} \rightarrow \text{se}_{b=1}^B(\hat{\theta}^{*b}) = \hat{\sigma}_{s,BS}$$

3.2 The law school data again



We estimated the correlation between GPA and LSAT as follows:

```
| lawCor <- with(law, cor(GPA, LSAT))
```

```
[1] 0.7763745
```

How precise is this estimate? If F is bivariate normal, then the estimated correlation coefficient $\hat{\rho}$ has a standard deviation of about

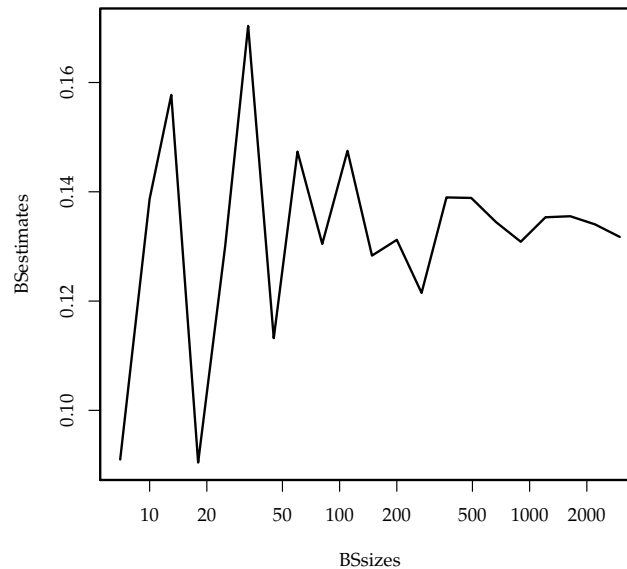
$$\hat{\sigma}_{\text{normal}} = \frac{1 - \hat{\rho}^2}{\sqrt{n - 3}} \approx 0.115$$

With a bootstrap we can live without the assumption that F is normal.

```
set.seed(123)
samplesize <- dim(law)[1]
ind <- 1:samplesize
law.boot <- replicate(200, {
  i2 <- sample(ind, replace = TRUE)
  with(law[i2, ], cor(GPA, LSAT))
})
sd(law.boot)
```

```
set.seed(123)
ind <- 1:samplesize
lawBS <- function(B) sd(replicate(B, {
  i2 <- sample(ind, replace = TRUE)
  with(law[i2, ], cor(GPA, LSAT))
}))
BSsizes <- round(exp(seq(2, 8, 0.3)))
BSestimates <- sapply(BSsizes, lawBS)
```

```
| plot(BSestimates ~ BSsizes, log = "x", t = "l")
```



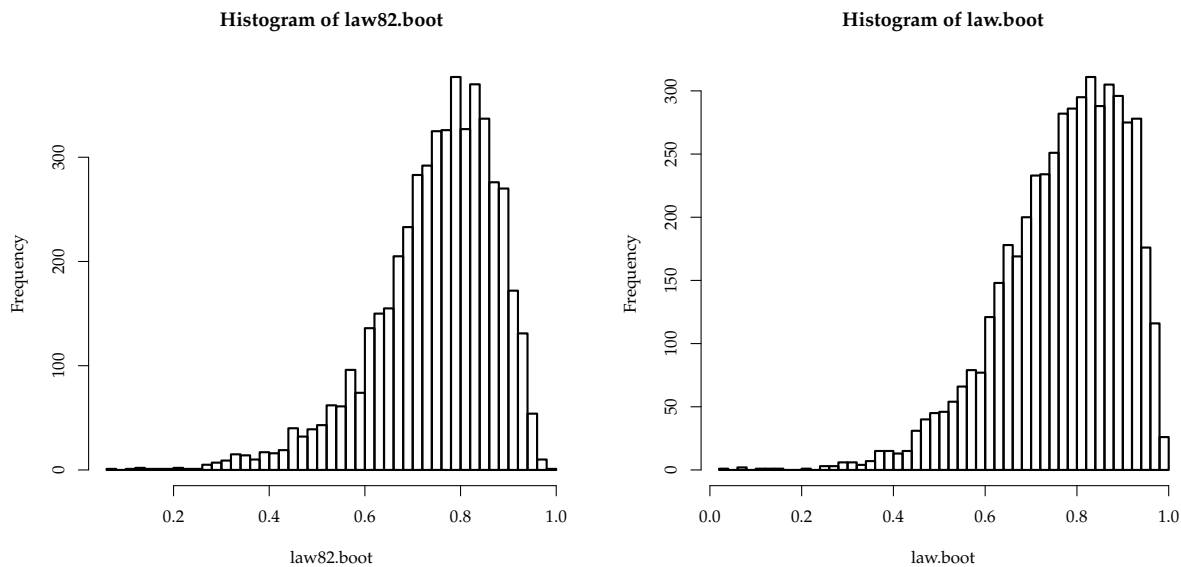
The distribution of $\hat{\theta}^*$ What is the distribution of the bootstrap replications $\hat{F}(\hat{\theta}^*)$ and how does this compare to $F(\hat{\theta}^*)$?

```

set.seed(123)
ind <- 1:dim(law)[1]
law.boot <- replicate(5000, {
  i2 <- sample(ind, size = samplesize, replace = TRUE)
  with(law[i2, ], cor(GPA, LSAT))
})
ind <- 1:dim(law82)[1]
law82.boot <- replicate(5000, {
  i2 <- sample(ind, size = samplesize, replace = TRUE)
  with(law82[i2, ], cor(GPA, LSAT))
})

```

```
| hist(law82.boot, breaks = 40) | hist(law.boot, breaks = 40)
```



Comparison of the bootstrap estimate $se_{\hat{F}}(\hat{\sigma})$ with the standard error $se_F(\hat{\sigma})$:

```
| sd(law.boot)
```

```
[1] 0.1343748
```

```
| sd(law82.boot)
```

```
[1] 0.125938
```

Exercise 3.1 Look again at the dataset `x1.csv`.

- What is the average value of the ratio `x1/x2`.
- What is the standard deviation we would obtain with the standard formula?
- Use a bootstrap to estimate the standard deviation.
- How does the standard deviation converge? Show a graph!

```
sTest <- function(X, Y, sX, sY) {
  sim <- var((rnorm(10000, mean = X, sd = sX)/rnorm(10000,
    mean = Y, sd = sY)))
  approx <- sX^2/X^2 + sY^2/Y^2
  approx <- (sX^2/Y^2 + sY^2 * X^2/Y^4)
  c(sim, approx)
}
sTest(100, 300, 0.7, 0.2)
```

| |
|-----------------------------------|
| [1] 0.000005472840 0.000005493827 |
|-----------------------------------|

3.3 How many bootstrap samples do we need?

Coefficient of variation

$$\text{cv}_X := \frac{\sigma_X}{|\mu_X|}$$

$$\text{cv}(\widehat{\text{se}}_B) \doteq \sqrt{\text{cv}(\widehat{\text{se}}_\infty)^2 + \frac{1}{4B}(E(\Delta) + 2)}$$

- Δ is a measure for long-tailedness of the distribution of $\hat{\theta}^*$. For the normal distribution $\Delta = 0$ and usually $\Delta < 10$.
- The magnitude of $\text{cv}(\widehat{\text{se}}_\infty)$ depends on
 - \hat{F} ,
 - the statistic $s(\cdot)$,
 - and the sample size.

3.4 The parametric bootstrap

Now we assume that we know a lot about F , e.g. that F follows a (multivariate) normal distribution.

- x_1, x_2, \dots, x_n are n observed values.
- estimate the parameters of F (e.g. mean, variance, covariance in the case of a normal distribution) and obtain F_{par} .
- Apply the bootstrap algorithm for standard errors to F_{par} .

Example: The law school data again Bivariate normal random variables

If x and y have variance-covariance matrix

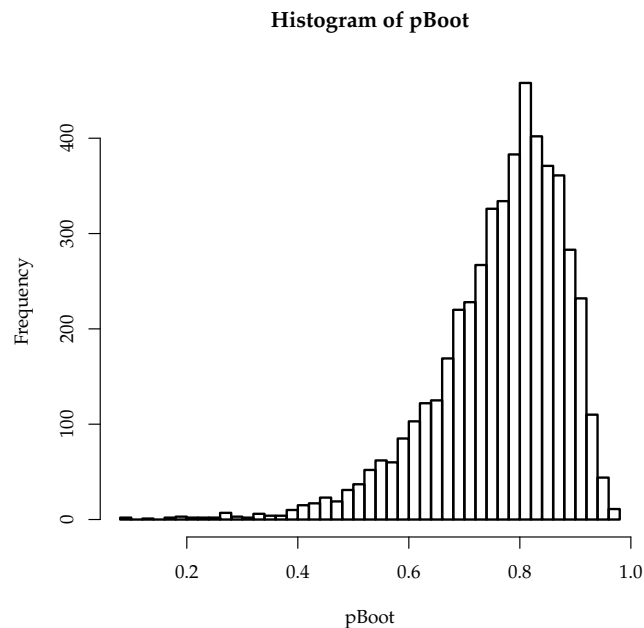
$$\Sigma_{xy} = \begin{pmatrix} \sigma_x^2 & \sigma_{xy} \\ \sigma_{xy} & \sigma_y^2 \end{pmatrix}$$

and if r_1 and r_2 are random variables that follow a standard normal distribution, then x and y can be constructed as follows:

$$\begin{aligned} x &= \mu_x + \sigma_x r_1 \\ y &= \mu_y + \sigma_y \frac{r_1 + c \cdot r_2}{\sqrt{1 + c^2}} \end{aligned}$$

with $c = \sqrt{\sigma_x^2 \sigma_y^2 / \sigma_{xy}^2 - 1}$.

```
paraBoot <- function(XY, B) {
  sigma <- cov(XY)
  mu <- rapply(XY, mean)
  c <- sqrt(prod(diag(sigma))/sigma[1, 2]^2 - 1)
  r1 <- rnorm(B)
  r2 <- rnorm(B)
  x <- mu[1] + sqrt(sigma[1, 1]) * r1
  y <- mu[2] + sqrt(sigma[2, 2]) * (r1 + c * r2)/sqrt(1 +
    c^2)
  cbind(x, y)
}
set.seed(123)
pBoot <- replicate(5000, cor(paraBoot(law, samplesize))[2,
  1])
hist(pBoot, breaks = 40)
```



And what is the parametric bootstrap estimation of the standard deviation?

```
| sd(pBoot)
```

```
[1] 0.1168820
```

Exercise 3.2 Return to exercise 3.1. What result would you obtain with a parametric bootstrap?

3.5 Eigenvalues und Eigenvectors

Let us construct a 100×4 matrix which we interpret at 100 observations of 4 different variables.

```
set.seed(123)
N <- 100
latent <- 10 * runif(N)
latent2 <- 10 * runif(N)
sd <- 1
x1 <- latent + sd * rnorm(N)
x2 <- latent + latent2 + sd * rnorm(N)
x3 <- -latent2 + sd * rnorm(N)
x4 <- latent2 - 2 * latent + sd * rnorm(N)
X <- cbind(x1, x2, x3, x4)
```

Let us assume that we do not know the underlying structure and that we use eigenvalues to learn more about the structure:

```
| lapply(eigen(cov(X)), round, 3)
```

```
$values
[1] 53.212 18.772  0.956  0.853

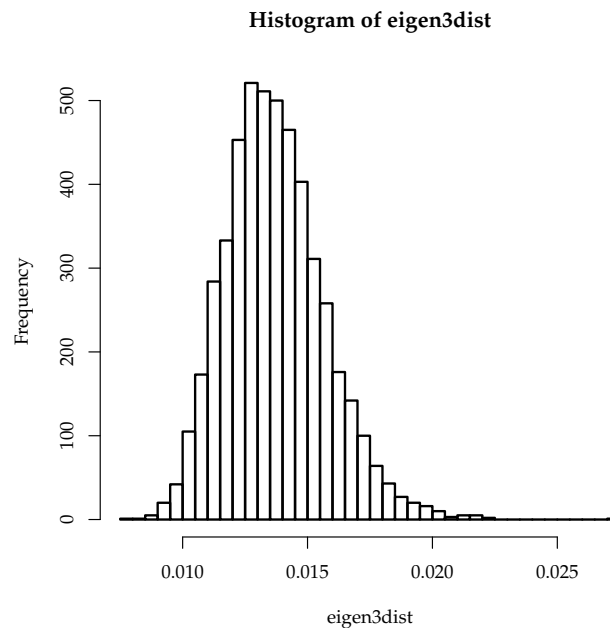
$vectors
      [,1]  [,2]  [,3]  [,4]
[1,] -0.388  0.160  0.242  0.875
[2,] -0.292  0.737  0.465 -0.393
[3,] -0.107 -0.609  0.772 -0.150
[4,]  0.867  0.245  0.360  0.241
```

We find the first two eigenvalues fairly large (indeed, we constructed the data with two main latent variables in our mind), and the last two relatively small. Does that mean that they are not significant?

If $\lambda_1, \lambda_2, \lambda_3, \lambda_4$ are the eigenvalues, then our test statistic is

$$\hat{\sigma}^* = \hat{\lambda}_3^* / \sum_{i=1}^4 \hat{\lambda}_i$$

```
| eigen3 <- function(X) {
      ee <- eigen(cov(X))["values"]
      ee[3]/sum(ee)
    }
ind <- 1:N
set.seed(123)
eigen3dist <- replicate(5000, eigen3(X[sample(ind, replace = TRUE),
]))
| hist(eigen3dist, breaks = 40)
```



Now let us compare our initial estimate...

```
| eigen3(X)
```

```
[1] 0.01295262
```

... with the mean of the bootstrap replications:

```
| mean(eigen3dist)
```

```
[1] 0.01371007
```

3.6 When bootstraps fail

Consider the following problem: X follows a uniform distribution over $[0, \theta]$. The ML estimator for θ is $\max(X_i)$. We have a sample of 20 observations. We compare:

1. Standard bootstrap of $\hat{\theta}^*$.
2. Parametric bootstrap based on a uniform distribution $[0, \hat{\theta}]$.

```

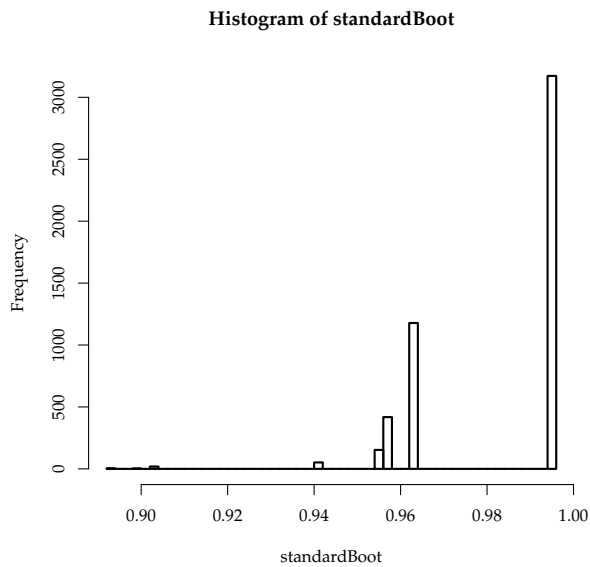
set.seed(123)
N <- 50
X <- runif(N)
thetaHat <- max(X)
standardBoot <- replicate(5000, max(sample(X, N, replace = TRUE)))
paramBoot <- replicate(5000, max(runif(N, min = 0, max = thetaHat)))

```

```

| hist(standardBoot, breaks = 40) | hist(paramBoot, breaks = 40, xlim = range(st

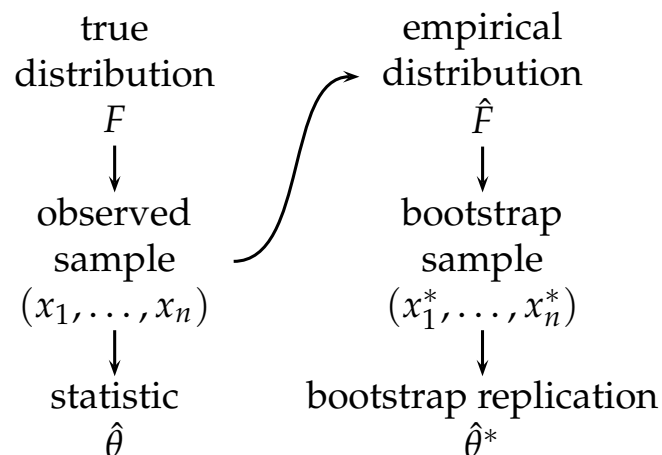
```



4 More complicated data structures

4.1 Motivation

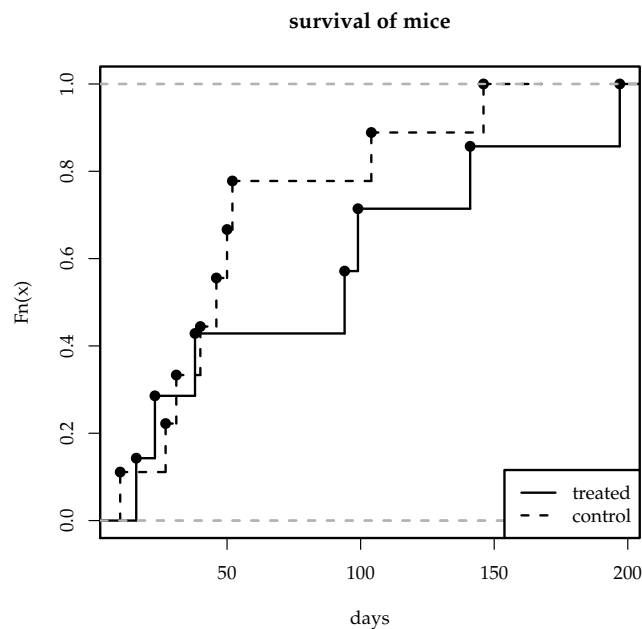
The bootstrap in one-sample problems



Crucial is the step from the observed sample (x_1, \dots, x_n) to the empirical distribution \hat{F} in the bootstrap.

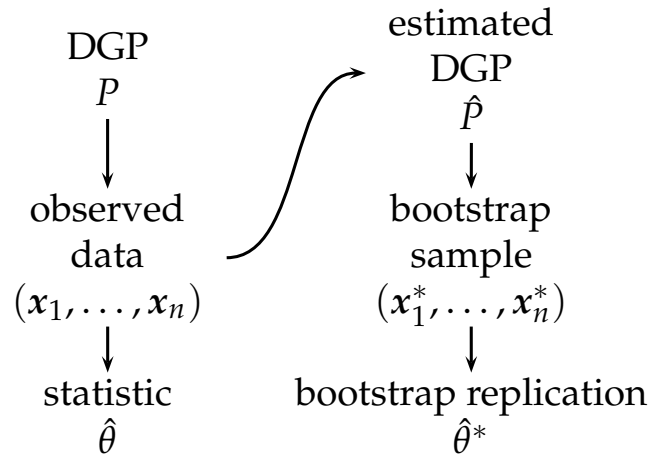
The above procedure does not work with more complicated data structure. Let us look again at the mice data:

```
plot(ecdf(mouse.t), xlim = range(mouse.c, mouse.t), verticals = TRUE,
     main = "survival of mice", xlab = "days")
lines(ecdf(mouse.c), lty = "dashed", verticals = TRUE)
legend("bottomright", c("treated", "control"), lty = c("solid",
  "dashed"), bg = "white")
```



Here we have two samples. When we move from our observations to the bootstrap we use the distribution of both samples as a starting point for the bootstrap.

The bootstrap with more general data structures Let us call the data generating process P .



4.2 Regression

In a linear model the data structure is more complicated:

$$y_i = \beta x_i + \epsilon_i$$

- “bootstrapping pairs”:
we sample with replacement from the set

$$\{(x_1, y_1), (x_2, y_2), \dots, (x_n, y_n)\}$$

- “bootstrapping residuals”:
 - We first estimate $\hat{\beta}$ from the model.
 - We then estimate the residuals $\hat{\epsilon}_i$.
 - We sample with replacement from the $\hat{\epsilon}_i$ and use the following dataset:

$$\{(x_1, \hat{\beta}x_1 + \hat{\epsilon}_{i_1}), (x_2, \hat{\beta}x_2 + \hat{\epsilon}_{i_2}), \dots, (x_n, \hat{\beta}x_n + \hat{\epsilon}_{i_n})\}$$

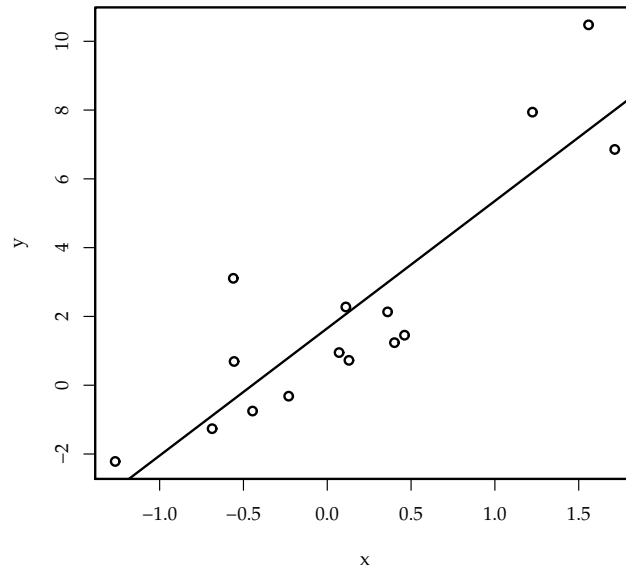
4.2.1 Bootstrapping pairs

```

set.seed(123)
N <- 15
sd <- 1.5
x <- rnorm(N)
y <- 3 * x + sd * rnorm(N)^2

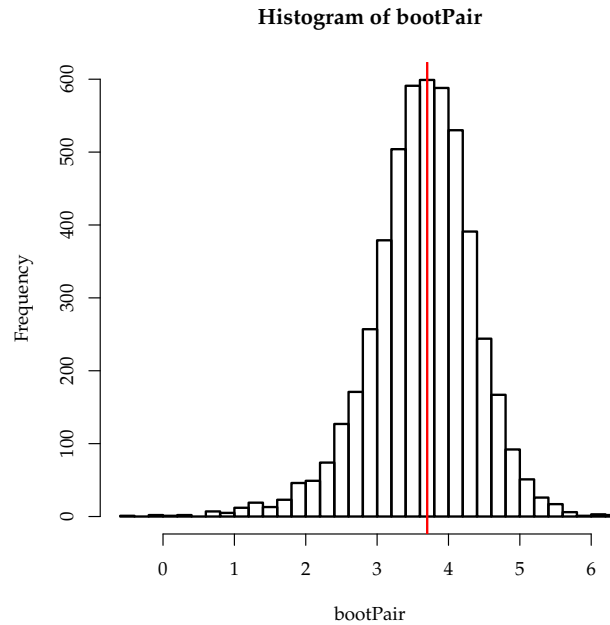
```

```
est <- lm(y ~ x)
plot(y ~ x)
abline(est)
```



```
bootPair <- replicate(5000, {
  ind <- sample(1:N, replace = TRUE)
  coef(lm(y[ind] ~ x[ind]))[2]
})
```

```
hist(bootPair, breaks = 40)
abline(v = coef(est)[2], col = "red")
```



Now we compare the bootstrapped $\hat{\sigma}_\beta$ with the one from the regression model:

```
| sd(bootPair)
```

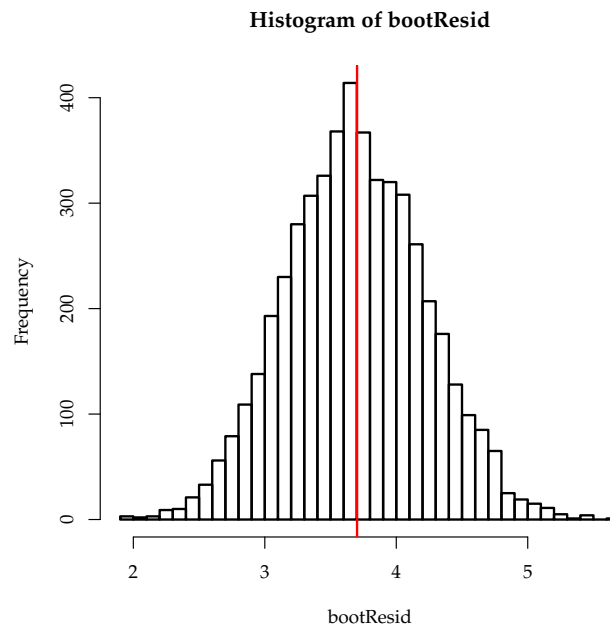
```
[1] 0.7282672
```

```
| sqrt(vcov(est)[2, 2])
```

```
[1] 0.5610004
```

4.2.2 Bootstrapping residuals

```
| rr <- residuals(est)
| beta <- coef(est)
| bootResid <- replicate(5000, {
|   epsilon <- sample(rr, size = N, replace = TRUE)
|   coef(lm((cbind(1, x) %*% beta + epsilon) ~ x))[2]
| })
| hist(bootResid, breaks = 40)
| abline(v = coef(est)[2], col = "red")
```



```
| sd(bootResid)
```

```
[1] 0.5261317
```

```
| sqrt(vcov(est)[2, 2])
```

```
[1] 0.5610004
```

Exercise 4.1 We use again the dataset `x1.csv`. Now we estimate the linear model

$$x_2 = \beta_0 + \beta_1 x_1 + u$$

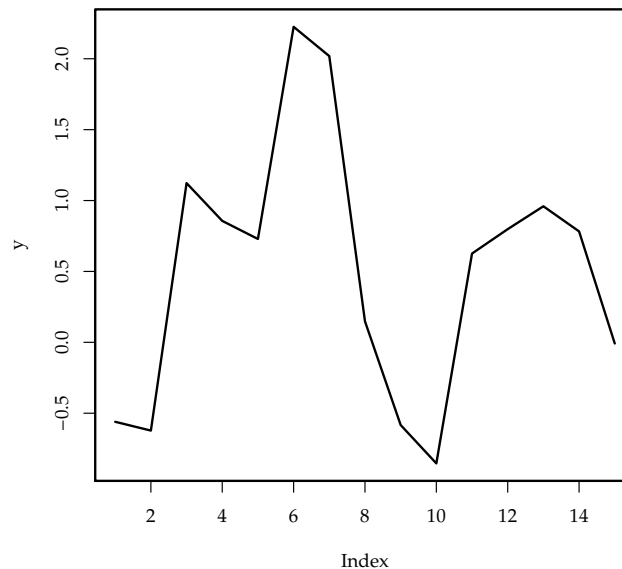
- Estimate β_0 and β_1
- What is the standard deviation of the coefficients you obtain with the standard regression command?
- What is the standard deviation you obtain when you bootstrap pairs?
- What is the standard deviation you obtain when you bootstrap residuals?

4.3 Timeseries — Example: A simple AR-1 process

Let us consider a simple AR-1 process:

$$y_t = \beta y_{t-1} + \epsilon_t$$

```
set.seed(123)
N <- 15
epsilon <- rnorm(N)
y <- epsilon
for (i in 2:N) y[i] <- y[i - 1] * 0.7 + epsilon[i]
plot(y, t = "l")
```



Here we have two options:

- “bootstrapping residuals”:
- “moving blocks”

Let us start with bootstrapping residuals:

4.3.1 Bootstrapping residuals

Let us try to estimate β :

```
| est <- lm(y[-1] ~ y[-N] - 1)
```

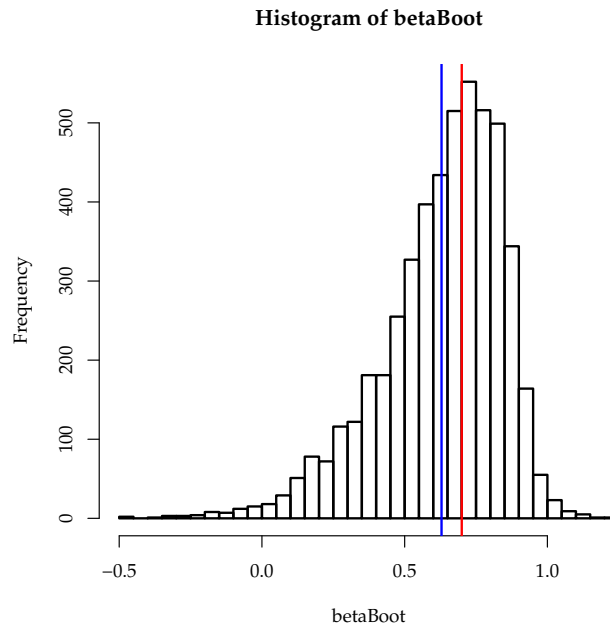
```
Call:
lm(formula = y[-1] ~ y[-N] - 1)

Coefficients:
 y[-N]
0.5994
```

To construct the bootstrap sample we will use the estimated residuals:

```
rr <- residuals(est)
beta <- coef(est)
betaBoot <- replicate(5000, {
  epsilon <- sample(rr, size = N, replace = TRUE)
  y <- epsilon
  for (i in 2:N) y[i] <- y[i - 1] * beta + epsilon[i]
  coef(lm(y[-1] ~ y[-N] - 1))
})

hist(betaBoot, breaks = 40)
abline(v = 0.7, col = "red")
abline(v = mean(betaBoot), col = "blue")
```



Now let us compare the bootstrap estimate of the standard error with the regression estimate:

```
| sd(betaBoot)
```

```
[1] 0.2142523
```

```
| sqrt(vcov(est))
```

```
          y[-N]
y[-N] 0.218549
```

Now let us assume that we (wrongly) suspect the process to be AR-2:

$$y_t = \beta_1 y_{t-1} + \beta_2 y_{t-2} + \epsilon_t$$

We estimate

```
| est <- lm(y[-c(1:2)] ~ y[-c(1, N)] + y[-c(N - 1, N)] -
           1)
```

Call:

```
lm(formula = y[-c(1:2)] ~ y[-c(1, N)] + y[-c(N - 1, N)] - 1)
```

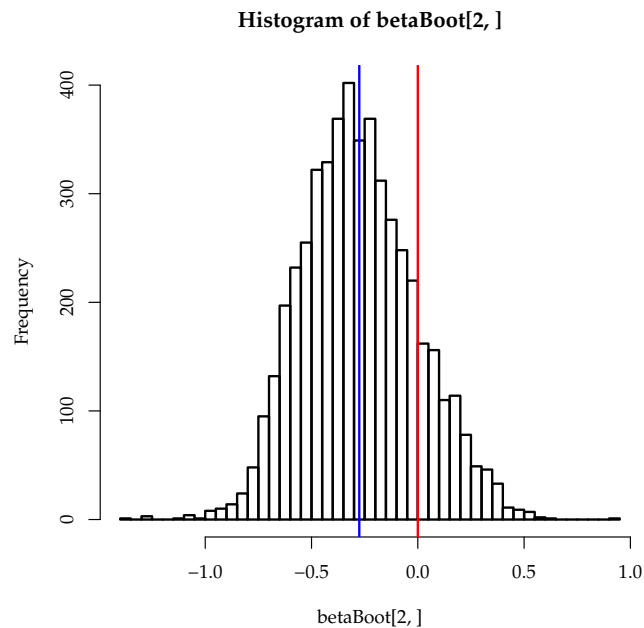
Coefficients:

```
  y[-c(1, N)]  y[-c(N - 1, N)]
    0.8086         -0.3587
```

What can we say about the accuracy of $\hat{\beta}_2$?

```
| rr <- residuals(est)
  beta <- coef(est)
  betaBoot <- replicate(5000, {
    epsilon <- sample(rr, N, replace = TRUE)
    y <- epsilon
    for (i in 3:N) y[i] <- y[i - 1] * beta[1] + y[i -
      2] * beta[2] + epsilon[i]
    coef(lm(y[-c(1:2)] ~ y[-c(1, N)] + y[-c(N - 1, N)] -
      1))
  })
```

```
| hist(betaBoot[2, ], breaks = 40)
  abline(v = 0, col = "red")
  abline(v = mean(betaBoot[2, ]), col = "blue")
```



Let us also here compare the estimated standard deviations of β_2 :

```
| sd(betaBoot[2, ])
```

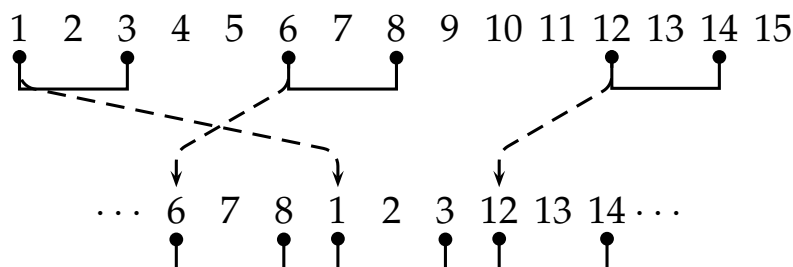
```
[1] 0.2672242
```

```
| sqrt(vcov(est)[2, 2])
```

```
[1] 0.287147
```

4.3.2 Moving blocks

The above regression model assumes that we know a lot about the data generating process. Here we just assume that a short block of data already contains a characteristic pattern.



```

blockLen <- 3
blockNum <- N/blockLen
betaBlock <- replicate(5000, {
  start <- sample(1:(N - blockLen + 1), size = blockNum,
    replace = TRUE)
  blockedIndices <- c(sapply(start, function(x) seq(x,
    x + blockLen - 1)))
  yy <- y[blockedIndices]
  coef(lm(yy[-1] ~ yy[-N] - 1))
})

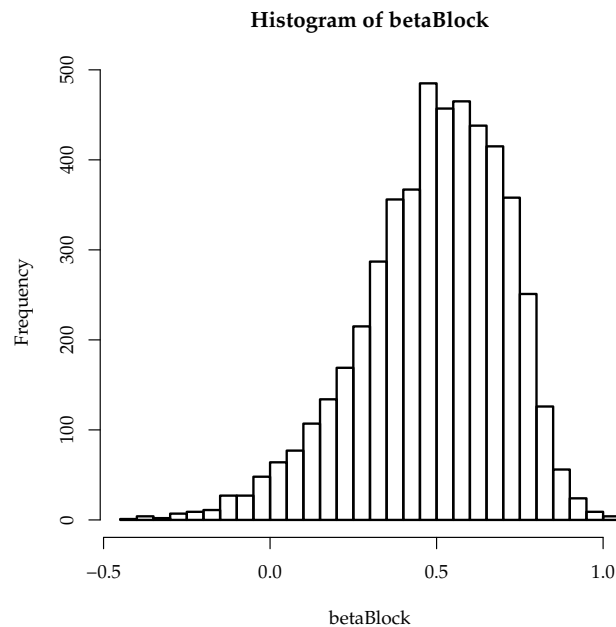
```

```

hist(betaBlock, breaks = 40)
sd(betaBlock)

```

```
[1] 0.2174306
```



Exercise 4.2 The dataset `x2.csv` contains a sequence of values for x and y . You estimate the following model:

$$y_t = \beta_0 + \beta_1 y_{t-1} + \beta_2 x_t$$

- What estimates do you obtain with OLS?
- What are the standard deviations of your estimates if you use the traditional method?

- Use the bootstrapping residuals methods to obtain standard deviations.
- Use moving blocks to obtain standard deviations.
- How does the standard deviation depend on the block size? Provide a plot with the block size on the horizontal axis and the standard deviations on the vertical axis.

5 Bias

$$\text{Bias}_F = E_F[s(\mathbf{x})] - t(F)$$

as above we replace F by \hat{F} to obtain the bootstrap estimate of the bias:

$$\text{Bias}_{\hat{F}} = E_{\hat{F}}[s(\mathbf{x}^*)] - t(\hat{F})$$

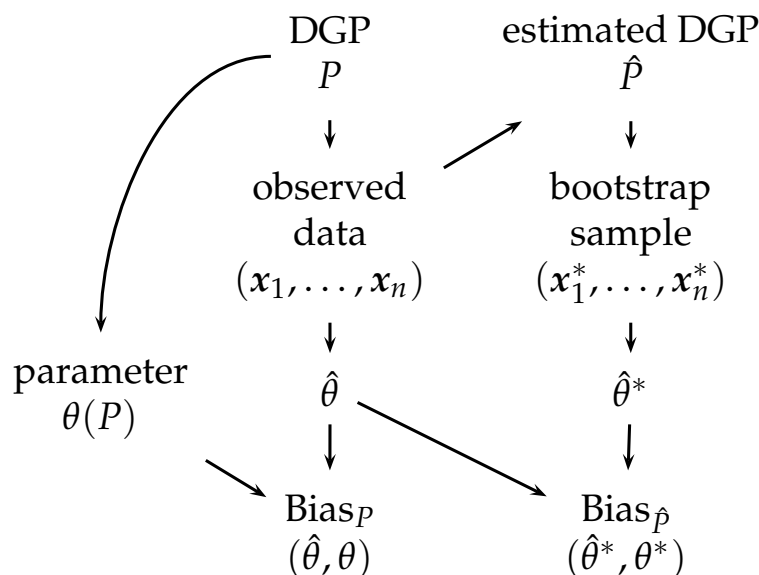
The bootstrap estimate of the bias

1. Draw B independent bootstrap samples $\mathbf{x}^{*1}, \mathbf{x}^{*2}, \dots, \mathbf{x}^{*B}$.

Each sample has size n and is drawn with replacement from x_1, x_2, \dots, x_n .

2. For bootstrap sample $b = 1, 2, \dots, B$ determine $\hat{\theta}^{*b} = s(\mathbf{x}^{*b})$.

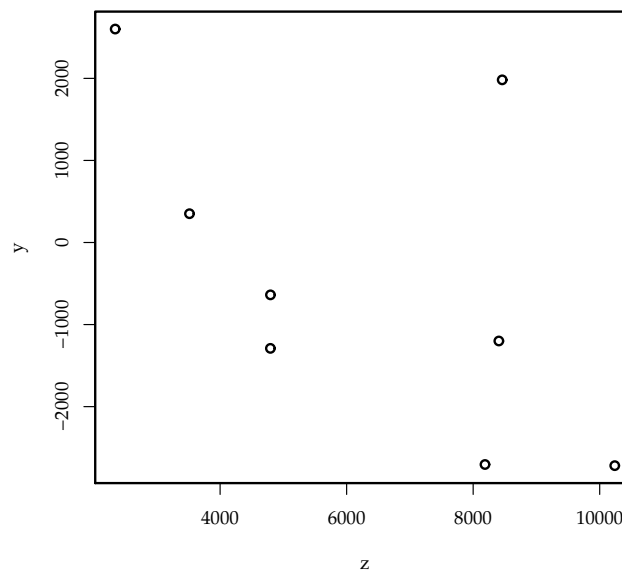
3. $\text{Bias}_{\hat{F}} \approx \widehat{\text{Bias}}_B = \frac{1}{n} \sum_{b=1}^B \hat{\theta}^{*b} - s(\mathbf{x})$



Example: The patch data We are interested in

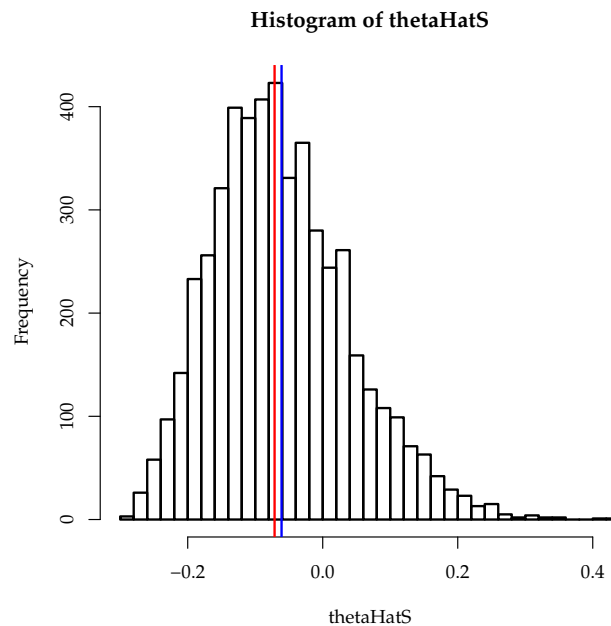
$$\frac{E(\text{new patch}) - E(\text{old patch})}{E(\text{old patch}) - E(\text{placebo patch})} = \frac{E(y)}{E(z)} \approx \frac{\bar{y}}{\bar{x}} = \frac{-452.25}{6342.375} \approx -0.0713$$

```
data(patch, package = "bootstrap")
plot(y ~ z, data = patch)
```



```
N <- dim(patch)[1]
set.seed(123)
thetaHatS <- replicate(5000, {
  ind <- sample(1:N, replace = TRUE)
  with(patch[ind, ], mean(y)/mean(z))
})
```

```
hist(thetaHatS, breaks = 40)
thetaHat <- with(patch, mean(y)/mean(z))
abline(v = thetaHat, col = "red")
abline(v = mean(thetaHatS), col = "blue")
```



Estimated bias and standard deviation

```
| bias <- mean(thetaHatS) - thetaHat
```

```
[1] 0.01032944
```

```
| sd(thetaHatS)
```

```
[1] 0.1017147
```

```
| bias/sd(thetaHatS)
```

```
[1] 0.1015531
```

5.1 A better estimate for the bias

A problem in

$$\text{Bias}_{\hat{f}} \approx \widehat{\text{Bias}}_B = \frac{1}{n} \sum_{b=1}^B \hat{\theta}^{*b} - s(\mathbf{x})$$

is that in $\frac{1}{n} \sum_{b=1}^B \hat{\theta}^{*b}$ some observations are sampled more frequently than others, while in $s(\mathbf{x})$ all observations in \mathbf{x} have the same weight. We should correct for the different weights.

- We can express the sampling process as a list of “sampled” indices.
- Alternative we can express the sampling process as frequencies:

$$P_i^* = \frac{1}{n} \cdot \#\{x_j^* = x_i\} \quad \text{for } i = 1, \dots, n$$

Hence, we define a resampling vector

$$\mathbf{P}^* = (P_1^*, \dots, P_n^*) \quad \text{with } \sum_{i=1}^n P_i^* = 1$$

In our case:

$$\hat{\theta}^* = \frac{\bar{y}^*}{\bar{z}^*} = \frac{\sum_{i=1}^n P_i^* y_i}{\sum_{i=1}^n P_i^* z_i}$$

More generally, for $\hat{\theta} = t(F)$ we write

$$\hat{\theta}^* = T(\mathbf{P}^*)$$

Let $\mathbf{P}^0 = \left(\frac{1}{n}, \frac{1}{n}, \dots, \frac{1}{n}\right)$ be a vector of length n where all elements are $1/n$, i.e. all elements of \mathbf{x} have the same weight.

$$T(\mathbf{P}^0) = \hat{\theta} = t(\hat{F})$$

Above we defined

$$\widehat{\text{Bias}}_B = \frac{1}{n} \sum_{b=1}^B \hat{\theta}^{*b} - s(\mathbf{x}) = \frac{1}{n} \sum_{b=1}^B T(\mathbf{P}_b^*) - T(\mathbf{P}^0)$$

Now we define

$$\bar{\mathbf{P}}^* = \frac{1}{B} \sum_{b=1}^B \mathbf{P}_b^*$$

Then a better estimate for the bias is

$$\overline{\text{Bias}}_B = \frac{1}{n} \sum_{b=1}^B T(\mathbf{P}_b^*) - T(\bar{\mathbf{P}}^*)$$

5.2 The jackknife

With the above “better” estimate we still have to calculate a large number of replications. Furthermore, all these replications are random. Is there a more systematic way?

The jackknife leaves out only one observation from n observations:

$$\mathbf{x} = (x_1, x_2, \dots, x_n)$$

$$\mathbf{x}_{(i)} = (x_1, x_2, \dots, x_{i-1}, x_{i+1}, \dots, x_n)$$

For each jackknife replication we calculate

$$\hat{\theta}_{(i)} = s(\mathbf{x}_{(i)})$$

Then

$$\widehat{\text{Bias}}_{\text{jack}} = (n - 1) \left(\frac{1}{n} \sum_{i=1}^n \hat{\theta}_{(i)} - \hat{\theta} \right)$$

```
| jacks <- sapply(1:N, function(i) with(patch[-i, ], mean(y)/mean(z)))
```

```
[1] -0.05711856 -0.12849970 -0.02145610 -0.13245033 -0.05067038
[6] -0.08404803 -0.06486298 -0.02219698
```

```
| biasJack <- (N - 1) * (mean(jacks) - thetaHat)
```

```
[1] 0.008002488
```

The jackknife works well if $t(\cdot)$ is a smooth statistic. (e.g. the median is not smooth)

There is also a jackknife estimate of the standard error:

$$\widehat{\text{se}}_{\text{jack}} = \sqrt{\frac{n-1}{n} \sum_{i=1}^n \left(\hat{\theta}_{(i)} - \frac{1}{n} \sum_{i=1}^n \hat{\theta}_{(i)} \right)^2} = \frac{n-1}{\sqrt{n}} \text{se}(\hat{\theta}_{(i)})$$

```
| sd(jacks) * (N - 1)/sqrt(N)
```

```
[1] 0.1055278
```

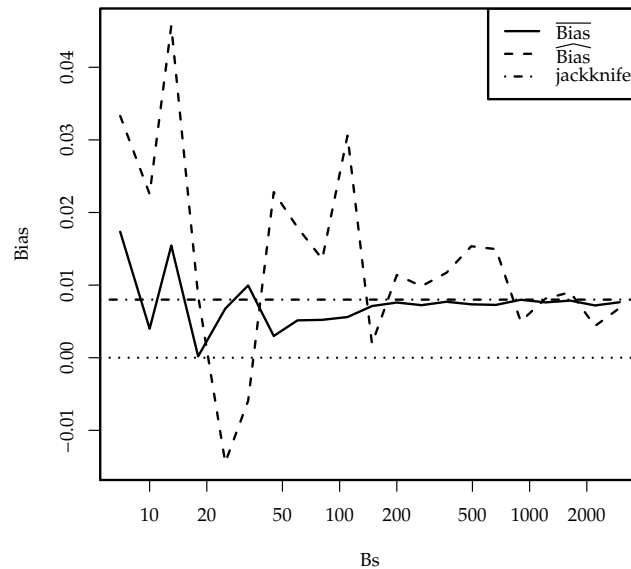
```
| sd(thetaHatS)
```

```
[1] 0.1017147
```

5.3 Coverage

```
set.seed(123)
T <- function(P) with(patch, (y %**% P)/(z %**% P))
weights <- matrix(0, N, N)
Pzero <- rep(1, N)/N
diag(weights) <- 1
Pvec <- function() apply(weights[sample(1:N, replace = TRUE),
], 2, sum)/N
estimate <- function(B) {
  Pbar <- rep(0, N)
  thetaStar <- mean(replicate(B, {
    PP <- Pvec()
    Pbar <- Pbar + PP
    T(PP)
  }))
  c(biasBar = thetaStar - T(Pbar/sum(Pbar)), biasHat = thetaStar -
    T(Pzero))
}
Bs <- round(exp(seq(2, 8, 0.3)))
estimates <- sapply(Bs, estimate)

plot(estimates["biasHat", ] ~ Bs, log = "x", t = "l",
      lty = "dashed", ylab = "Bias")
lines(estimates["biasBar", ] ~ Bs, t = "l")
abline(h = 0, lty = "dotted")
abline(h = biasJack, lty = 4)
legend("topright", c(expression(bar(Bias)), expression(widehat(Bias)),
  "jackknife"), lty = c(1, 2, 4))
```



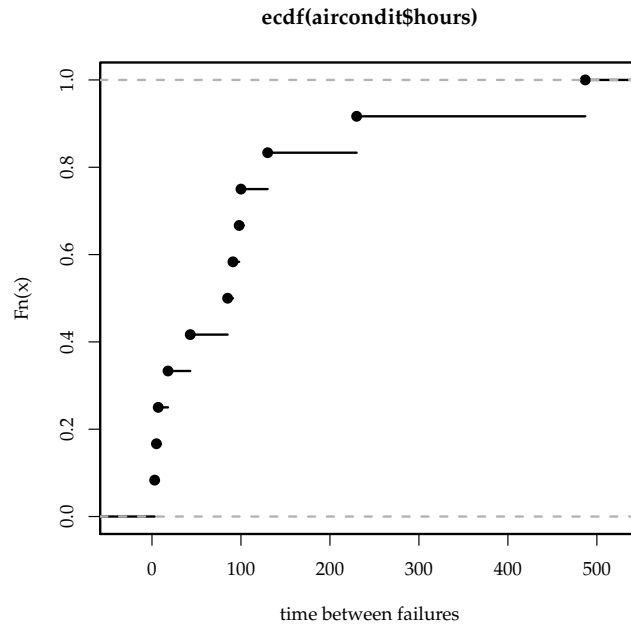
Exercise 5.1 1. You suspect that your estimate for the coefficient of y_{t-1} in exercise 4.2 might be biased.

- What is the standard bootstrap estimate of the bias?
 - What is the better estimate?
 - What is the jackknife estimate?
2. Now you use a median regression to estimate the relationship in exercise 4.2, i.e. you minimise the sum of absolute values of residuals, and not, as in OLS, the sum of squared residuals.
- What coefficients do you estimate?
 - What is the standard bootstrap estimate of the bias?
 - What is the better estimate?
 - What is the jackknife estimate?

6 Confidence intervals

The `aircondit` data: Failure times of air-conditions.

```
data(aircondit, package = "boot")
plot(ecdf(aircondit$hours), xlab = "time between failures")
```



```
| airMean <- mean(aircondit$hours)
```

The distribution does not look normal.

6.1 The exact approach:

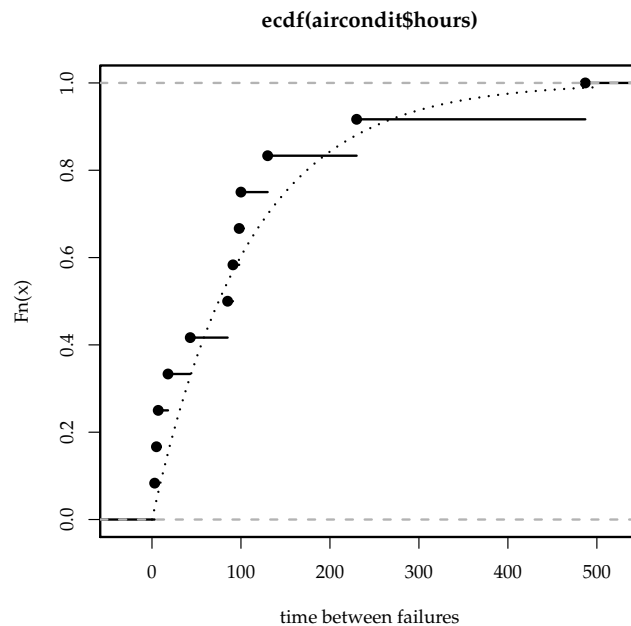
- Works even for small samples
- requires knowledge of the “true” distribution
- may require some serious transformations

If failures times follow a Poisson distribution then the times between failure follow an exponential distribution.

$$f(x, \lambda) = \begin{cases} \lambda e^{-\lambda x} & x \geq 0 \\ 0 & x < 0 \end{cases}$$

To find λ we can use ML: $\hat{\lambda} = n / \sum x_i = 1/\bar{x}$.

```
with(aircondit, {
  n <- length(hours)
  lambda <- length(hours)/sum(hours)
})
plot(ecdf(aircondit$hours), xlab = "time between failures")
with(list(x = seq(0, 500, 10)), lines(1 - exp(-lambda *
  x) ~ x, lty = "dotted"))
```



With $\lambda = 0.00925$ and $\bar{x} = 108$ we can now try to work out the distribution of the mean of an exponentially distributed random variable.

$$CI_{1/\lambda} = \left[\frac{1}{\bar{\lambda}} \frac{2n}{\chi_{2n,1-\alpha/2}^2}; \frac{1}{\bar{\lambda}} \frac{2n}{\chi_{2n,\alpha/2}^2} \right]$$

```
alpha <- 5/100
sapply(c(1 - alpha/2, alpha/2), function(alpha) 2 * n / (lambda *
  qchisq(alpha, 2 * n)))
```

```
[1] 65.89765 209.17415
```

6.2 The normal approximation:

- Asymptotically the mean follows a normal distribution
- Hence: We can do this when the sample is “large”

6.2.1 ...based on parametric estimates of σ

```
confint(lm(hours ~ 1, data = aircondit))
```

```
                2.5 %    97.5 %
(Intercept) 21.52561 194.6411
```

6.2.2 ... based on bootstrap estimates of σ

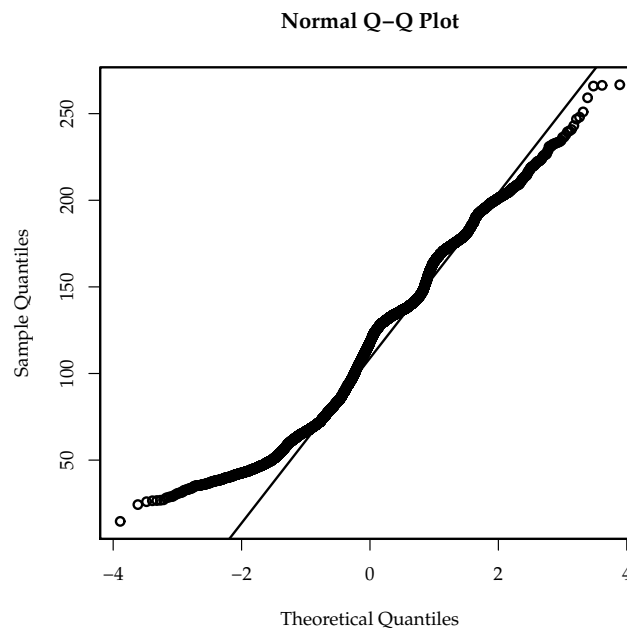
```
set.seed(123)
xx <- with(aircondit, replicate(5000, c(mean = mean(sample(hours,
  replace = TRUE)), sd = sd(sample(hours, replace = TRUE)))))

mean(xx["mean", ]) + c(1, -1) * qnorm(alpha/2) * sd(xx["mean",
  ])
```

```
[1] 33.62068 182.08342
```

Can we use a normal approximation for the distribution of the mean?

```
qqnorm(xx)
qqline(xx)
```



6.3 The bootstrap- t interval

Idea: If θ is the mean of a normally distributed random variable, then we have

$$\frac{\hat{\theta} - \theta}{\hat{\sigma}_{\hat{\theta}}} \sim t_{n-1} \quad \rightarrow \text{CI} = \left[\hat{\theta} - t_{n-1}^{(1-\alpha/2)} \hat{\sigma}_{\hat{\theta}}, \hat{\theta} - t_{n-1}^{(\alpha/2)} \hat{\sigma}_{\hat{\theta}} \right]$$

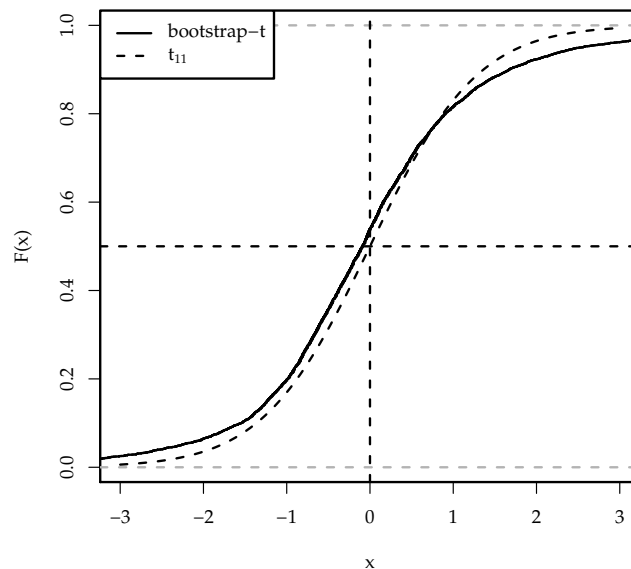
If θ is something else, we can bootstrap our own distribution of \hat{t} :

For each bootstrap sample b we calculate approximate pivots:

$$Z^*(b) = \frac{\hat{\theta}^*(b) - \hat{\theta}}{\hat{\sigma}_{\hat{\theta}}^*(b)}$$

The quantiles of the distribution of Z^* are the quantiles of \hat{t} .

```
N <- dim(aircondit)[1]
thetaHat <- mean(aircondit[["hours"]])
bT <- (xx["mean", ] - thetaHat)/(xx["sd", ]/sqrt(N))
plot(function(x) pt(x, N - 1), -3, 3, lty = 2, ylab = "F(x)")
plot(ecdf(bT), xlim = c(-1.5, 1.5), add = TRUE)
abline(v = 0, h = 0.5, lty = 2)
legend("topleft", c("bootstrap-t", expression(t[11])),
      lty = c(1, 2), bg = "white")
```



(note that we need a lot of bootstrap samples to obtain reliable quantiles)
 Now let us compare the quantiles of the original t -distribution with the bootstrap- t :

```
| quantile(bT, c(alpha/2, 1 - alpha/2))
```

| | |
|-----------|----------|
| 2.5% | 97.5% |
| -3.011323 | 3.608497 |

```
| qt(c(alpha/2, 1 - alpha/2), N - 1)
```

```
[1] -2.200985  2.200985
```

$$\text{CI} = \left[\hat{\theta} - \hat{t}_{n-1}^{(1-\alpha/2)} \hat{\sigma}_{\hat{\theta}}, \hat{\theta} - \hat{t}_{n-1}^{(\alpha/2)} \hat{\sigma}_{\hat{\theta}} \right]$$

```
| mean(xx["mean", ]) + quantile(bT, c(alpha/2, 1 - alpha/2)) *
  sd(xx["mean", ])
```

```
      2.5%      97.5%
-6.198339 244.519703
```

6.3.1 Two problems with bootstrap-*t*

- Above we did use a standard formula to obtain $\hat{\sigma}_{\hat{\theta}}^*(b)$. What, if such a formula does not exist? We can, of course, bootstrap $\hat{\sigma}_{\hat{\theta}}^*(b)$, but this requires to run a bootstrap for each of our bootstrap samples.
- The bootstrap-*t* procedure is not invariant with respect to transformations of the variable we are studying.

6.4 Percentile intervals

$$\text{CI}_{\%} = \left[\hat{\theta}^{*(\alpha/2)}, \hat{\theta}^{*(1-\alpha/2)} \right]$$

```
| quantile(xx["mean", ], c(alpha/2, 1 - alpha/2))
```

```
      2.5%      97.5%
46.8250 193.5896
```

Why is this an “interesting” confidence interval?

Percentile interval lemma

[Efron and Tibshirani] Suppose the transformation $\hat{\phi} = m(\hat{\theta})$ perfectly normalises the distribution of $\hat{\theta}$:

$$\hat{\phi} \sim N(\phi, c^2)$$

Then

$$\text{CI}_{\%} = \left[\hat{\theta}^{*(\alpha/2)}, \hat{\theta}^{*(1-\alpha/2)} \right] = \left[m^{-1} \left(\hat{\phi} - c \cdot z^{1-\alpha/2} \right), m^{-1} \left(\hat{\phi} - c \cdot z^{\alpha/2} \right) \right]$$

What happens if the intervals are asymmetric? Note that the bootstrap- t and the percentile interval come to opposite conclusions. If θ_b^* has a long left tail, then the percentile interval will deviate from the mean more to the left than to the right.

The bootstrap- t will do the opposite.

6.5 BC_a -intervals

Remember: Percentile intervals:

$$CI_{\alpha\%} = \left[\hat{\theta}^{*(\alpha/2)}, \hat{\theta}^{*(1-\alpha/2)} \right]$$

Now we adjust the $\alpha/2$ and $1 - \alpha/2$ levels:

$$BC_a = \left[\hat{\theta}^{*(\alpha_1)}, \hat{\theta}^{*(\alpha_2)} \right]$$

with

$$\alpha_1 = \Phi \left(\hat{z}_0 + \frac{\hat{z}_0 + z^{(\alpha/2)}}{1 - \hat{a}(\hat{z}_0 + z^{(\alpha/2)})} \right)$$

$$\alpha_2 = \Phi \left(\hat{z}_0 + \frac{\hat{z}_0 + z^{(1-\alpha/2)}}{1 - \hat{a}(\hat{z}_0 + z^{(1-\alpha/2)})} \right)$$

- $\hat{z}_0 = \Phi^{-1} \left(\frac{\#\{\hat{\theta}^*(b) < \hat{\theta}\}}{B} \right)$: bias correction
- $\hat{a} = \frac{\sum_{i=1}^n (\hat{\theta}_{(\cdot)} - \hat{\theta}_{(i)})^3}{6(\sum_{i=1}^n (\hat{\theta}_{(\cdot)} - \hat{\theta}_{(i)})^2)^{3/2}}$: acceleration

Note first that if $\hat{z}_0 = 0$ and $\hat{a} = 0$ then $\alpha_1 = \Phi(z^{(\alpha/2)}) = \alpha/2$ and $\alpha_2 = \Phi(z^{(1-\alpha/2)}) = 1 - \alpha/2$.

\hat{z}_0 shifts the interval to the left or to the right. \hat{a} makes the interval wider or smaller.

- BC_a is transformation respecting
- BC_a converges faster than the standard and the percentile method.

```
jack <- function(x, FUN) sapply(1:length(x), function(i) FUN(x[-i]))
u <- jack(aircondit$hours, mean)
acc <- sum((mean(u) - u)^3)/(6 * sum((mean(u) - u)^2)^(3/2))
z0 <- qnorm(mean(xx["mean", ] < thetaHat))
alphaI <- function(alpha) {
  zA <- z0 + qnorm(alpha)
```

```

    pnorm(z0 + zA/(1 - acc * zA))
  }
  quantile(xx["mean", ], c(alphaI(alpha/2), alphaI(1 -
    alpha/2)))

```

```

6.904699% 99.60277%
57.18049 231.38074

```

Since people need BC_a frequently, the function is provided in the library *bootstrap*.

```

bcanon(aircondit$hours, 5000, mean, alpha = c(alpha/2,
  1 - alpha/2))

```

```

$confpnts
  alpha bca point
[1,] 0.025    58.25
[2,] 0.975   230.75

$z0
[1] 0.1226301

$acc
[1] 0.09379807

$u
[1] 117.63636 117.45455 117.27273 116.27273 114.00000 110.18182
[7] 109.63636 109.00000 108.81818 106.09091  97.00000  73.63636

$call
bcanon(x = aircondit$hours, nboot = 5000, theta = mean, alpha = c(alpha/2,
  1 - alpha/2))

```

6.6 Comparison

| | BC_a | standard | percentile | bootstrap- t |
|------------------------------|--------|--------------|--------------|----------------|
| transformation respecting | + | | + | |
| speed | $1/n$ | $1/\sqrt{n}$ | $1/\sqrt{n}$ | $1/n$ |

Exercise 6.1 *The file x3.csv contains a vector of numbers.*

1. *Determine the mean of these numbers.*
2. *You assume that these numbers follow a χ^2 distribution with 1 degree of freedom. What is the distribution of the mean of these numbers?*
3. *Based in the χ^2 distribution, determine a 95% confidence interval for the mean.*
4. *Assume that the mean follows a normal distribution. Determine the confidence interval on the basis of the parametric estimate of the standard deviation.*
5. *Assume that the mean follows a normal distribution. Determine the confidence interval on the basis of the bootstrap estimate of the standard deviation.*
6. *Determine the bootstrap-t confidence interval.*
7. *Determine the percentile confidence interval.*
8. *Determine the BC_a confidence interval.*

7 Hypothesis Testing

7.1 Permutation tests

Null hypotheses Idea: we have two samples, \mathbf{x} and \mathbf{y} , drawn from perhaps the same, perhaps different distributions:

$$\begin{aligned} F &\rightarrow \mathbf{x} = (x_1, x_2, \dots, x_n) \\ G &\rightarrow \mathbf{y} = (y_1, y_2, \dots, y_m) \end{aligned}$$

Our null hypothesis is

$$H_0 : F = G$$

Let us assume in this example that we do a one-sided test, i.e. our alternative hypothesis would assume that values of \mathbf{x} are larger on average than those of \mathbf{y} . Let us rephrase the problem in terms of a test statistic:

$$\hat{\theta} = \bar{x} - \bar{y}$$

What can we conclude from observing that $\hat{\theta}$ is large?

Let us assume that we expect $\hat{\theta}$ to be small as long as H_0 holds.

We define the achieved significance level as

$$\text{ASL} = \Pr\{\hat{\theta}_{H_0}^* \geq \hat{\theta}\}$$

If we know that $F = N(\mu_x, \sigma^2)$ and $G = N(\mu_y, \sigma^2)$ our null hypothesis H_0 is equivalent to $\mu_x = \mu_y$.

If H_0 holds, $\hat{\theta} \sim N\left(0, \sigma^2 \cdot \left(\frac{1}{n} + \frac{1}{m}\right)\right)$

Then

$$\begin{aligned} \text{ASL} &= \Pr\{\hat{\theta}_{H_0}^* \geq \hat{\theta}\} \\ &= \Pr\left\{Z \geq \frac{\hat{\theta}}{\sigma\sqrt{1/n + 1/m}}\right\} \\ &= \Phi\left(-\frac{\hat{\theta}}{\sigma\sqrt{1/n + 1/m}}\right) \end{aligned}$$

In a normal world $\hat{\theta}$ follows a “known” distribution.

We can bootstrap the distribution of $\hat{\theta}$ under H_0 :

- We join the two samples (x, y) and split it randomly into two subsamples x' and y' of size n and m , respectively (i.e. we forget the original association to F and G).

The two-sample permutation test statistic

1. We choose B independent subsamples x' and y' taken (without replacement) from the joint sample (x, y) .
2. We calculate $\hat{\theta}^*(b)$ for each sample.
3. $\widehat{\text{ASL}}_{\text{perm}} = \frac{1}{B} \cdot \#\{\hat{\theta}^*(b) > \hat{\theta}\}$

Ideally, we would take $B = \binom{N}{n}$ many subsamples to calculate the exact ASL:

$$\text{ASL}_{\text{perm}} = \#\{\hat{\theta}^*(b) > \hat{\theta}\} / \binom{N}{n}$$

| `t.test(mouse.t, mouse.c, alternative = "greater", var.equal = TRUE)`

Two Sample t-test

```
data: mouse.t and mouse.c
t = 1.1214, df = 14, p-value = 0.1405
alternative hypothesis: true difference in means is greater than 0
95 percent confidence interval:
 -17.48178      Inf
sample estimates:
mean of x mean of y
 86.85714  56.22222
```

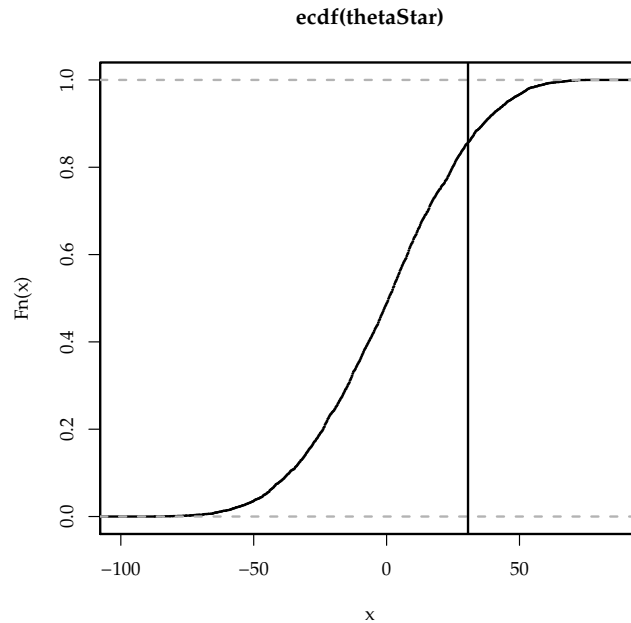
```
| thetaHat <- mean(mouse.t) - mean(mouse.c)
```

```
[1] 30.63492
```

```
| xy <- c(mouse.c, mouse.t)
| n <- length(mouse.c)
| N <- length(xy)
| thetaStar <- replicate(5000, {
|   indx <- sample(1:N, n)
|   mean(xy[indx]) - mean(xy[-indx])
| })
```

```
| plot(ecdf(thetaStar), do.points = FALSE)
| abline(v = thetaHat)
| mean(thetaStar > thetaHat)
```

```
[1] 0.1428
```



7.2 Bootstrap tests

The two-sample bootstrap test statistic

1. We choose B independent subsamples x' and y' taken (with replacement) from the joint sample (x, y) .
2. We calculate $\hat{\theta}^*(b)$ for each sample.
3. $\widehat{ASL}_{BS} = \frac{1}{B} \cdot \#\{\hat{\theta}^*(b) > \hat{\theta}\}$

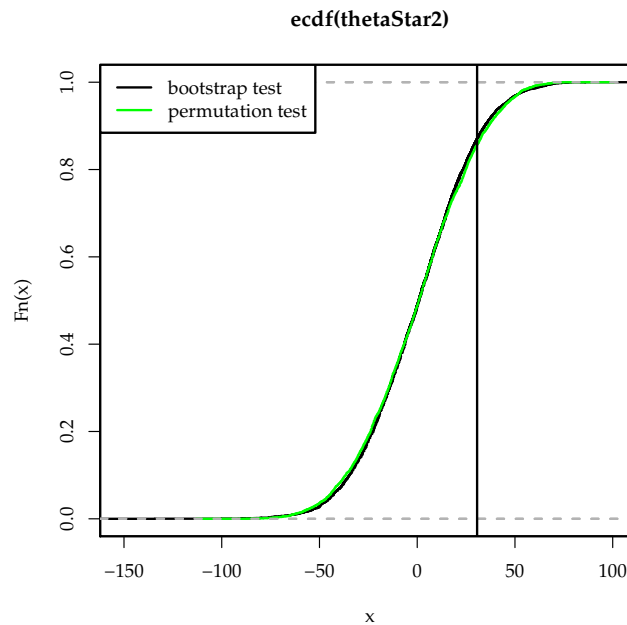
(The main difference to permutation tests is that here we sample with replacement, and not without.)

```
xy <- c(mouse.c, mouse.t)
n <- length(mouse.c)
m <- length(mouse.t)
N <- length(xy)
thetaStar2 <- replicate(5000, mean(sample(xy, n, replace = TRUE)) -
  mean(sample(xy, m, replace = TRUE)))
```

```
plot(ecdf(thetaStar2), do.points = FALSE)
lines(ecdf(thetaStar), do.points = FALSE, col = "green")
legend("topleft", c("bootstrap test", "permutation test"),
```

```
col = c("black", "green"), lty = 1, bg = "white")
abline(v = thetaHat)
mean(thetaStar > thetaHat)
```

```
[1] 0.1428
```

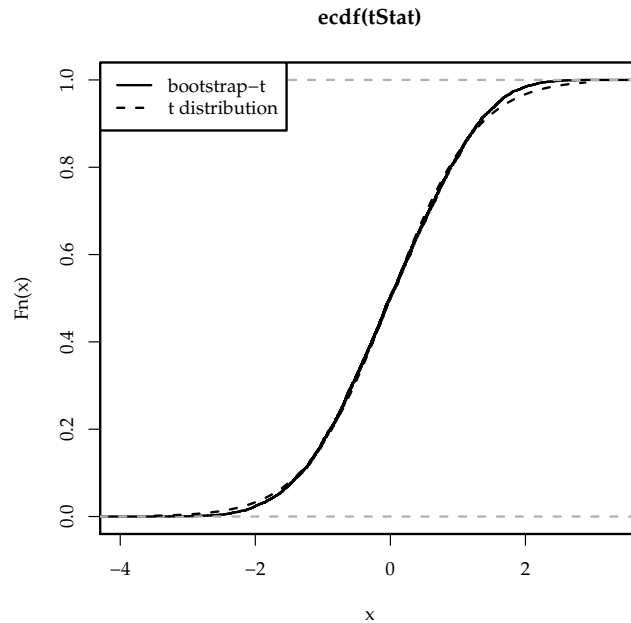


7.3 Bootstrap- t tests

Let us first consider H_0 that the means and the standard deviations in both samples are the same:

```
tStat <- replicate(5000, {
  xx <- sample(xy, n, replace = TRUE)
  yy = sample(xy, m, replace = TRUE)
  (mean(xx) - mean(yy))/(sd(c(xx, yy) * sqrt(1/n +
    1/m)))
})
```

```
plot(ecdf(tStat))
plot(function(x) pt(x, df = n + m - 2), add = TRUE, lty = "dashed",
  min(tStat), max(tStat))
legend("topleft", c("bootstrap-t", "t distribution"),
  lty = c(1, 2), bg = "white")
```



```
| ASLt <- mean(tStat > thetaHat/(sd(xy) * sqrt(1/n + 1/m)))
```

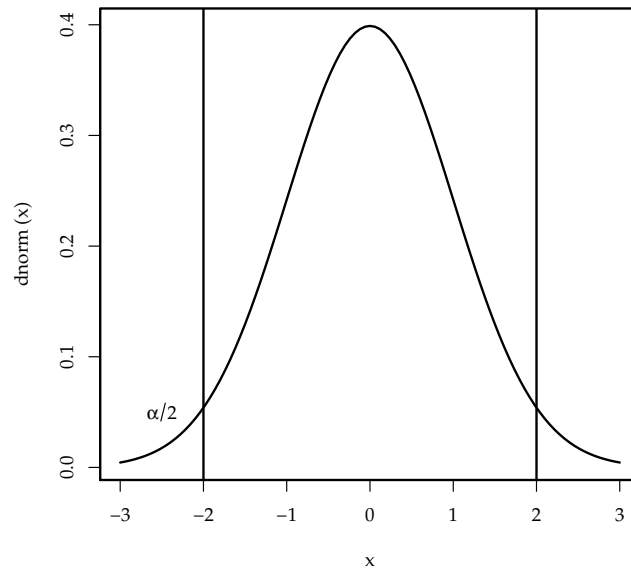
```
[1] 0.1418
```

Exercise 7.1 *At the beginning of this section we assumed identical standard deviations in both samples. Perform the same test without making this assumption.*

7.4 BC_a -tests

Above we have seen that the BC_a method can help us to determine confidence intervals more efficiently. Can we use this method for tests of hypotheses, too? Let us assume that all we have is a machine that calculates confidence intervals, but we have to calculate an ASL.

Idea:



We have to work out a level α such that the upper or lower end of the CI coincides with the value of θ under H_0 .

Let us assume that we are interested in the lower end of the CI (i.e. H_1 states that $\theta > 0$).

Remember that

$$\text{BC}_a = [\hat{\theta}^{*(\alpha_1)}, \hat{\theta}^{*(\alpha_2)}]$$

with

$$\alpha_1 = \Phi \left(\hat{z}_0 + \frac{\hat{z}_0 + z^{(\alpha/2)}}{1 - \hat{a}(\hat{z}_0 + z^{(\alpha/2)})} \right)$$

In our permutation test we observe α_1 . We can use the above formula to work out $\alpha/2$:

$$\begin{aligned} \Phi^{-1}(\alpha_1) &= \hat{z}_0 + \frac{\hat{z}_0 + z^{(\alpha/2)}}{1 - \hat{a}(\hat{z}_0 + z^{(\alpha/2)})} \\ \underbrace{\Phi^{-1}(\alpha_1) - \hat{z}_0}_M &= \frac{\hat{z}_0 + z^{(\alpha/2)}}{1 - \underbrace{\hat{a}(\hat{z}_0 + z^{(\alpha/2)})}_Q} \end{aligned}$$

$$\begin{aligned}
 M &= \frac{Q}{1 - \hat{a}Q} \\
 M \cdot (1 - \hat{a}Q) &= Q \\
 M - \hat{a}QM &= Q \\
 M &= Q \cdot (1 + \hat{a}M) \\
 \frac{M}{1 + \hat{a}M} &= Q = \hat{z}_0 + z^{(\alpha/2)} \\
 \frac{M}{1 + \hat{a}M} - \hat{z}_0 &= z^{(\alpha/2)} \\
 \Phi \left(\frac{M}{1 + \hat{a}M} - \hat{z}_0 \right) &= \alpha/2 \\
 \Phi \left(\frac{\Phi^{-1}(\alpha_1) - \hat{z}_0}{1 + \hat{a}(\Phi^{-1}(\alpha_1) - \hat{z}_0)} - \hat{z}_0 \right) &= \text{ASL}_{\text{BC}_a}
 \end{aligned}$$

We can still use the same formula for \hat{z}_0 :

- $\hat{z}_0 = \Phi^{-1} \left(\frac{\#\{\hat{\theta}^*(b) < \hat{\theta}\}}{B} \right)$: bias correction

With a two sample problem the formula for \hat{a} is slightly different:

We define $U_{x,i} = (n - 1)(\hat{\theta}_{x,(\cdot)} - \hat{\theta}_{x,(i)})$ and $U_{y,i} = (m - 1)(\hat{\theta}_{y,(\cdot)} - \hat{\theta}_{y,(i)})$

- $\hat{a} = \frac{1}{6} \frac{\frac{1}{n^3} \sum_{i=1}^n U_{x,i}^3 + \frac{1}{m^3} \sum_{i=1}^m U_{y,i}^3}{\left(\frac{1}{n^2} \sum_{i=1}^n U_{x,i}^2 + \frac{1}{m^2} \sum_{i=1}^m U_{y,i}^2 \right)^{3/2}}$: acceleration

```

set.seed(123)
FUN <- function(x, y) mean(x) - mean(y)
thetaStar <- replicate(5000, FUN(sample(mouse.t, replace = TRUE),
  sample(mouse.c, replace = TRUE)))
thetaHat <- FUN(mouse.t, mouse.c)
z0 <- qnorm(mean(thetaStar < thetaHat))
jack2 <- function(x, y) sapply(1:length(x), function(i) FUN(x[-i],
  y))
U <- function(x, y) {
  u <- jack2(x, y)
  (length(u) - 1) * (mean(u) - u)
}
Ux <- U(mouse.t, mouse.c)

```

```

Uy <- U(mouse.c, mouse.t)
n <- length(mouse.t)
m <- length(mouse.c)
ahat <- (sum(Ux^3)/n^3 + sum(Uy^3)/m^3)/(sum(Ux^2)/n^2 +
        sum(Uy^2)/m^2)^(3/2)/6
a1 <- mean(thetaStar < 0)
ASLBCa <- pnorm((qnorm(a1) - z0)/(1 + ahat * (qnorm(a1) -
        z0)) - z0)

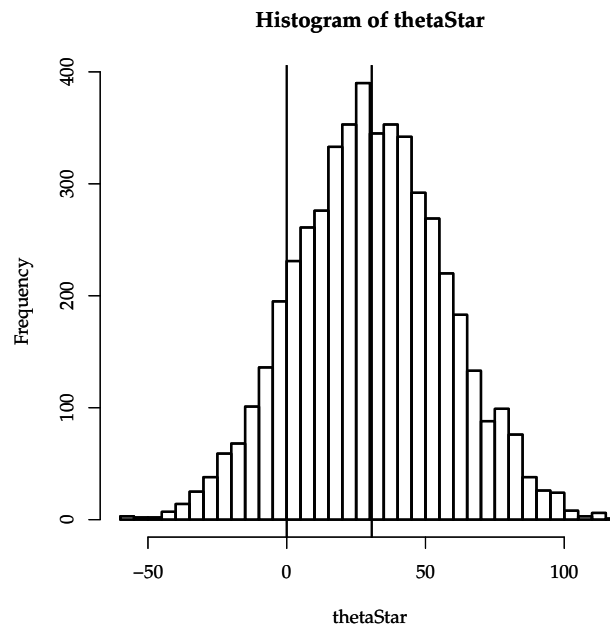
```

We obtain $z_0 = 0.018$, $\hat{a} = 0.0272$, $a_1 = 0.13$, and $ASL_{BC_a} = 0.115$.

```

plot(hist(thetaStar, breaks = 40))
abline(v = c(0, thetaHat))

```



Exercise 7.2 The dataset `x4.csv` contains data on x and y . You estimate a linear relationship

$$y = \beta_0 + \beta_1 x + \epsilon$$

Your null hypothesis is that $\beta_1 = 0$. The alternative hypothesis is that $\beta_1 \neq 0$.

1. What is the ASL of a parametric test?
2. What is the ASL of a permutation test?
3. What is the ASL of a bootstrap- t test?
4. What is the ASL of a BC_a test?