## MSiA 420 HW1

#### 2023-01-16

#### Question 1

See attached

#### Question 2

```
\mathbf{a}
## New names:
## Rows: 18 Columns: 5
## -- Column specification
## ------ Delimiter: "," dbl
## (2): y, x lgl (3): ...3, ...4, ...5
## i Use 'spec()' to retrieve the full column specification for this data. i
## Specify the column types or set 'show_col_types = FALSE' to quiet this message.
## * '' -> '...3'
## * '' -> '...4'
## * '' -> '...5'
##
## Call:
## lm(formula = y \sim x, data = dat)
## Residuals:
               1Q Median
                              ЗQ
## -3.8103 -1.3567 0.7347 1.5796 3.0106
## Coefficients:
              Estimate Std. Error t value Pr(>|t|)
## (Intercept) 5.42071
                         0.77963 6.953 3.25e-06 ***
## x
               0.48964
                         0.04359 11.234 5.32e-09 ***
## ---
## Signif. codes: 0 '***' 0.001 '**' 0.05 '.' 0.1 ' ' 1
## Residual standard error: 2.164 on 16 degrees of freedom
## Multiple R-squared: 0.8875, Adjusted R-squared: 0.8805
## F-statistic: 126.2 on 1 and 16 DF, p-value: 5.317e-09
```

As seen above, the coefficients  $\hat{\beta}_0$  and  $\hat{\beta}_1$  are 5.4207 and 0.4896, respectively. Thus, the initial guesses of the form  $\hat{\gamma_0} = \frac{1}{\hat{\beta}_0} = \frac{1}{5.4207} = 0.1845$ , and  $\hat{\gamma_1} = \frac{\hat{\beta}_1}{\hat{\beta}_0} = \frac{0.48964}{5.4207} = 0.0903$ 

```
\mathbf{b}
## [1] 28.13688 12.57428
## Nonlinear regression model
     model: y \sim ((p1 * x)/(p2 + x))
##
      data: dat
##
      p1
             p2
## 28.14 12.57
    residual sum-of-squares: 4.302
##
## Number of iterations to convergence: 9
## Achieved convergence tolerance: 9.921e-06
Question 3
\mathbf{a}
##
              [,1]
                         [,2]
## [1,] 15.37455 -13.73842
## [2,] -13.73842 13.92197
              [,1]
                         [,2]
## [1,] 0.5502797 0.5430248
## [2,] 0.5430248 0.6076944
## [1] 0.7418084 0.7795475
b
##
             p1
## p1 0.529951 0.5202800
## p2 0.520280 0.5822471
Although not exactly the same, the covariance matrices are very similar.
\mathbf{c}
## [1] 26.68294
## [1] 29.59083
## [1] 11.04637
## [1] 14.10219
```

2.5 %

## p1 26.71021 29.56383 ## p2 11.07887 14.06997

97.5 %

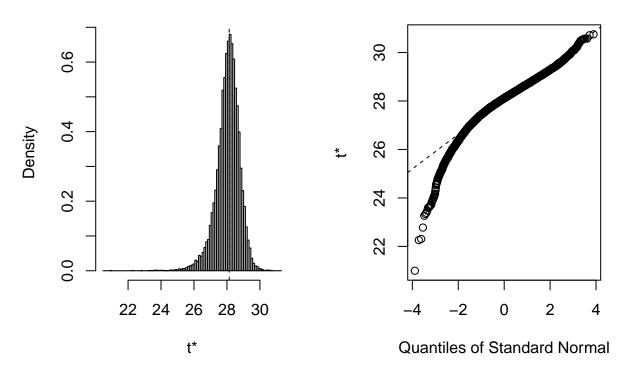
##

As seen from the above comparison, the "crude" confidence interval is similar to the output by R, but it is wider.

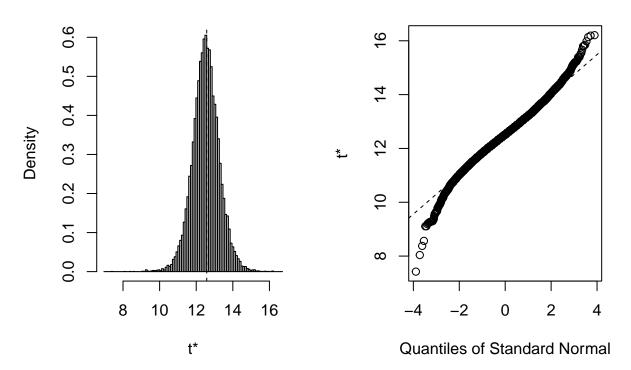
## Question 4

```
\mathbf{a}
##
## ORDINARY NONPARAMETRIC BOOTSTRAP
##
##
## Call:
## boot(data = dat, statistic = dat_fit, R = 20000, theta0 = c(0.1845,
##
       0.0903))
##
##
## Bootstrap Statistics :
       original
                      bias
                              std. error
## t1* 28.13688 -0.08372891
                               0.7143639
## t2* 12.57428 -0.05167943
                               0.7408931
```

# Histogram of t Histogram of t0



## Histogram of t Histogram of t1



```
b
## BOOTSTRAP CONFIDENCE INTERVAL CALCULATIONS
## Based on 20000 bootstrap replicates
##
## boot.ci(boot.out = dat_boot, conf = 0.95, type = "norm", index = 1)
##
## Intervals :
## Level
              Normal
         (26.82, 29.62)
## 95%
## Calculations and Intervals on Original Scale
## BOOTSTRAP CONFIDENCE INTERVAL CALCULATIONS
## Based on 20000 bootstrap replicates
##
## CALL :
## boot.ci(boot.out = dat_boot, conf = 0.95, type = "norm", index = 2)
## Intervals :
## Level
              Normal
         (11.17, 14.08)
## 95%
## Calculations and Intervals on Original Scale
```

 $\mathbf{c}$ 

```
## BOOTSTRAP CONFIDENCE INTERVAL CALCULATIONS
## Based on 20000 bootstrap replicates
## CALL:
## boot.ci(boot.out = dat_boot, conf = 0.95, type = "basic", index = 1)
## Intervals :
## Level
              Basic
## 95%
         (27.01, 29.86)
## Calculations and Intervals on Original Scale
## BOOTSTRAP CONFIDENCE INTERVAL CALCULATIONS
## Based on 20000 bootstrap replicates
## boot.ci(boot.out = dat_boot, conf = 0.95, type = "basic", index = 2)
##
## Intervals :
## Level
              Basic
         (11.14, 14.09)
## 95%
## Calculations and Intervals on Original Scale
```

For the most part, the confidence intervals in part b and part c agree. The confidence interval for  $\gamma_1$  is much closer since the histogram is pretty much completely normal. This means in part b where we use normal approximation, the confidence intervals will be much closer. For  $\gamma_0$ , we can see from part a that the histogram is slightly left skewed. Thus, the reflected confidence interval is slightly shifted to the right compared to the "crude" confidence interval. Note that the histogram for  $\gamma_0$  is still very close to normal despite skewedness, and thus the confidence intervals do not differ by much.

#### Question 5

 $\mathbf{d}$ 

```
## [1] 18.10226 20.29115

## BOOTSTRAP CONFIDENCE INTERVAL CALCULATIONS
## Based on 20000 bootstrap replicates
##
## CALL:
## boot.ci(boot.out = dat_boot2, conf = 0.95, type = c("basic"))
##
## Intervals:
## Level Basic
## 95% (18.91, 19.72)
## Calculations and Intervals on Original Scale
```

As seen above, the prediction interval for "future" Y is wider than the confidence interval. This is because of the added uncertainty from the standard errors of the prediction of the model. A prediction interval is a better representation because it not only takes into account the error of the estimated parameters, but also the error in the model that generates the prediction.

## Question 6

```
## 'log Lik.' 2.836148 (df=3)
## 'log Lik.' 1.628871 (df=3)
```

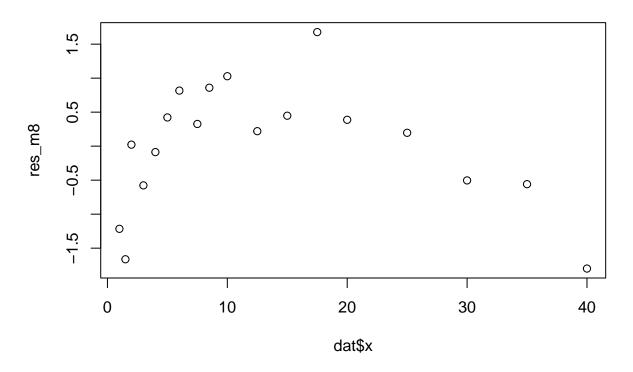
The nls model is better since it has a lower AIC.

### Question 7

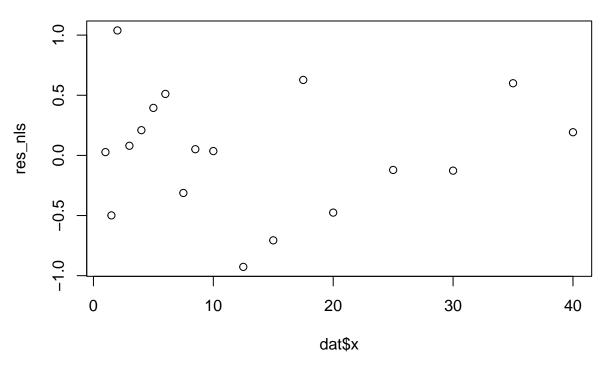
```
##
             [,1]
                       [,2]
   [1,] 1.175988 0.3116147
   [2,] 1.175988 0.3116147
   [3,] 1.175988 0.3116147
##
  [4,] 1.175988 0.3116147
## [5,] 1.175988 0.3116147
## [6,] 1.175988 0.3116147
   [7,] 1.175988 0.3116147
  [8,] 1.175988 0.3116147
  [9,] 1.175988 0.3116147
## [10,] 1.175988 0.3116147
## [11,] 1.175988 0.3116147
## [12,] 1.175988 0.3116147
## [13,] 1.175988 0.3116147
## [14,] 1.175988 0.3116147
## [15,] 1.175988 0.3116147
## [1] 1.1759878 0.3116147
```

As seen above, the better model is clearly model 2, which uses nls.

# X vs Residuals for Linear Model



## X vs Residuals for NLS



As seen from the plots above, the X vs Residuals plot for the linear model shows non-randomly distributed residuals around the 0 line. There is a clear non-linear pattern, which means that the linear model is not a good fit for the data. On the contrary, the X vs Residuals plot for the non-linear least squares model shows randomly distributed residuals around 0, which suggests there is no further relationship that we missed. Thus, the nls model is a better choice, which is in keeping with results above.

### **Appendix**

This section is to be used for including your R code. The following lines of code will take care of it. Please make sure to comment your code appropriately - in particular, demarcating codes belonging to different questions. Among other things, it will be easier for you to debug your own code.

```
labs = knitr::all_labels()
labs = labs[!labs %in% c("setup", "getlabels", "allcode")]
```

```
dat <- read csv("HW1 data.csv")</pre>
dat <- dat %>%
 select(y,x)
m1 \leftarrow lm(y\sim x, data = dat)
summary(m1)
est <- c(0.1845, 0.0903)
x_i <- dat$x
y_i <- dat$y
fn <- function(p) {</pre>
   yhat \leftarrow (p[1] * x_i) / (p[2] + x_i)
   sum((y_i-yhat)^2)
}
out_nlm <- nlm(fn, p = est, hessian=TRUE)
theta_nlm <- out_nlm$estimate #parameter estimates</pre>
theta_nlm
out_nls <- nls(y^{(p1 * x)} / (p2 + x)), data = dat, start = list(p1 = 0.1845, p2 = 0.0903))
out_nls
mse <- out_nlm$minimum / (length(y_i) - length(theta_nlm))</pre>
info_mat <- out_nlm$hessian / 2 / mse</pre>
cov_theta <- solve(info_mat)</pre>
se <- sqrt(diag(cov_theta))</pre>
info_mat
cov_theta
vcov(out_nls)
p1_left \leftarrow theta_nlm[1] - 1.96 * se[1]
p1_right <- theta_nlm[1] + 1.96 * se[1]
p2_left \leftarrow theta_nlm[2] - 1.96 * se[2]
p2_right \leftarrow theta_nlm[2] + 1.96 * se[2]
p1_left
p1_right
p2_left
p2_right
confint.default(out nls)
```

```
dat_fit <- function(Z, i, theta0){</pre>
  Zboot <- Z[i, ]</pre>
 x <- Zboot[[2]]</pre>
 y <- Zboot[[1]]
 fn <- function(p){</pre>
    yhat \leftarrow (p[1] * x) / (p[2] + x)
    sum((y - yhat)^2)
 out <- nlm(fn, p = theta0)
 theta <- out$estimate
}
dat_boot \leftarrow boot(dat, dat_fit, R = 20000, theta0 = c(0.1845, 0.0903))
dat_boot
plot(dat_boot, index = 1)
title(main = "Histogram of t0")
plot(dat_boot, index = 2)
title(main = "Histogram of t1")
# gamma O
boot.ci(dat_boot, conf = 0.95, type = "norm", index = 1)
#gamma 1
boot.ci(dat_boot, conf = 0.95, type = "norm", index = 2)
# gamma O
boot.ci(dat_boot, conf = 0.95, type = "basic", index = 1)
#gamma 1
boot.ci(dat_boot, conf = 0.95, type = "basic", index = 2)
# Bootstrap
dat_fit2 <- function(Z, i, theta0, x_pred){</pre>
 Zboot <- Z[i, ]</pre>
 x <- Zboot[[2]]
 y <- Zboot[[1]]
 fn <- function(p){</pre>
    yhat \leftarrow (p[1] * x) / (p[2] + x)
    sum((y - yhat)^2)
 out <- nlm(fn, p = theta0)
 theta <- out$estimate
 y_pred <- (theta[1] * x_pred) / (theta[2] + x_pred)</pre>
dat_boot2 \leftarrow boot(dat, dat_fit2, R = 20000, theta0 = c(0.1845, 0.0903), x_pred = 27)
Yhat0<-dat_boot2$t0
Yhatboot<-dat_boot2$t
SEY<-sqrt(var(Yhatboot)+mse)</pre>
c(Yhat0-qnorm(.975)*SEY, Yhat0+qnorm(.975)*SEY)
# Predicted
boot.ci(dat_boot2, conf = 0.95,type=c("basic"))
m6 \leftarrow lm(y\sim sqrt(x), data = dat)
n <- nrow(dat)</pre>
log_lik_m6 <- logLik(m6)</pre>
```

```
aic_m6 \leftarrow -2 * log_lik_m6 / n + 2 * 2 / n
log_lik_nls <- logLik(out_nls)</pre>
aic_nls <- -2 * log_lik_nls / n + 2 * 2 / n
aic_m6
aic_nls
Nrep <- 15 # set to 1 because we are only doing one pass
K <- nrow(dat) # set to the number of observations in the dataset
n.models = 2 #number of different models to fit and compare
FitFun1 <- function(x, p) p[1]+p[2]*sqrt(x)</pre>
FitFun2 \leftarrow function(x, p) (p[1] * x) / (p[2] + x)
n <- nrow(dat)</pre>
yhat <- matrix(0,n,n.models)</pre>
MSE <- matrix(0,Nrep,n.models)</pre>
for (j in 1:Nrep) {
  cv_idx <- cvfolds(n, K)</pre>
  for (k in cv_idx) {
    out <- lm(y~sqrt(x), data=dat[-k,])</pre>
    yhat[k,1] <- as.numeric(predict(out, dat[k,]))</pre>
    out <- nls(y \sim FitFun2(x,p), data=dat[-k,], start=list(p=c(0.1845, 0.0903)))
    yhat[k,2] <- as.numeric(predict(out, dat[k,]))</pre>
  }
  y <- dat$y
 MSE[j,] \leftarrow apply(yhat,2,function(x) sum((y-x)^2))/(n-1)
MSE
MSEAve <- apply (MSE,2,mean); MSEAve #averaged mean square CV error
m8 \leftarrow lm(y\sim sqrt(x), data = dat)
pred_m8 <- predict(m8, dat)</pre>
res_m8 <- dat$y - pred_m8
pred_nls <- predict(out_nls, dat)</pre>
res_nls <- dat$y - pred_nls</pre>
plot(dat$x, res_m8, main = "X vs Residuals for Linear Model")
plot(dat$x, res_nls, main = "X vs Residuals for NLS")
```