# Econometrics 3

Laura Gómez Assignment 3

May 30, 2020

# Question 1

You are interested in the response of output growth  $y_{1,t}$  to fiscal policy shocks  $e_{2,t}$ . The structural model is as follows:

$$y_{1,t} = \gamma_1 y_{2,t} + \alpha_{11} y_{1,t-1} + \alpha_{12} y_{2,t-1} + e_{1,t}$$

$$y_{2,t} = \gamma_2 y_{1,t} + \alpha_{21} y_{1,t-1} + \alpha_{22} y_{2,t-1} + e_{2,t}$$
(1)

#### Question 1.a

a) Under what conditions can this structural model be transformed in structural VAR form? Write down the conditions in terms of the parameters above and explain.

The conditions that we need to transform this structural model to structural VAR form are that the parameters do not change with the time. In other words  $\gamma_1$ ,  $\alpha_{11}$ ,  $\alpha_{12}$ ,  $\gamma_2$ ,  $\alpha_{21}$  and  $\alpha_{22}$  should be the same for all periods of time in the model to estimate them. Another assumptions are  $E(e_t) = 0$ ,  $E(e_t e_t') = \Sigma_e$  and  $\Gamma$  is invertible where:

$$\Gamma = \begin{bmatrix} 1 & -\gamma_1 \\ -\gamma_2 & 1 \end{bmatrix} 
e_t = \begin{bmatrix} e_{1,t} \\ e_{2,t} \end{bmatrix} 
\Sigma_e = \begin{bmatrix} \sigma_1 & 0 \\ 0 & \sigma_2 \end{bmatrix}$$
(2)

## Question 1.b

b) Assuming the model can be written in a structural VAR form, write down the structural VAR formulation.

First we rewrite in lag terms:

$$y_{1,t} = \gamma_1 y_{2,t} + \alpha_{11} y_{1,t-1} + \alpha_{12} y_{2,t-1} + e_{1,t} = \gamma_1 y_{2,t} + \alpha_{11}(L) y_{1,t} + \alpha_{12}(L) y_{2,t} + e_{1,t}$$

$$y_{2,t} = \gamma_2 y_{1,t} + \alpha_{21} y_{1,t-1} + \alpha_{22} y_{2,t-1} + e_{2,t} = \gamma_2 y_{1,t} + \alpha_{21}(L) y_{1,t} + \alpha_{22}(L) y_{2,t} + e_{2,t}$$
(3)

Rewrite as:

$$\underbrace{\begin{bmatrix} 1 & -\gamma_1 \\ -\gamma_2 & 1 \end{bmatrix}}_{\Gamma} \underbrace{\begin{bmatrix} y_{1,t} \\ y_{2,t} \end{bmatrix}}_{Y_t} = \underbrace{\begin{bmatrix} \alpha_{11}(L) & \alpha_{12}(L) \\ \alpha_{21}(L) & \alpha_{22}(L) \end{bmatrix}}_{A(L)} \underbrace{\begin{bmatrix} y_{1,t} \\ y_{2,t} \end{bmatrix}}_{Y_t} + e_t$$

$$\Gamma Y_t = A Y_{t-1} + e_t \tag{4}$$

And the reduce form is:

$$\Gamma Y_t = A(L)Y_t + e_t$$

$$Y_t = \underbrace{\Gamma^{-1}A(L)}_{A^*(L)}Y_t + \underbrace{\Gamma^{-1}e_t}_{u_t}$$

$$Y_t = A^*(L)Y_t + u_t$$
(5)

#### Question 1.c

c) Based on b), use backward substitution to derive the expression for the IRF of output growth to a fiscal policy shock at horizons 0, 1 (i.e.  $\frac{\partial y_{1,t+h}}{\partial e_{2,t}}$ , h = 0, 1). For derive IRF we should rewrite the expression of b):

$$Y_{t} = A^{*}(L)Y_{t} + u_{t}$$

$$Y_{t} - A^{*}(L)Y_{t} = +u_{t}$$

$$(I_{2} - A^{*}(L))Y_{t} = u_{t}$$

$$Y_{t} = (I_{2} - A^{*}(L))^{-1}u_{t}$$

$$Y_{t} = (I_{2} - A^{*}(L))^{-1}\Gamma^{-1}e_{t}$$
(6)

using backward substitution, first we found the matriz Gamma inverse

$$\Gamma^{-1} = \begin{bmatrix} 1 & -\gamma_1 \\ -\gamma_2 & 1 \end{bmatrix}^{-1} = \begin{bmatrix} \frac{1}{1-\gamma_1\gamma_2} & \frac{\gamma_1}{1-\gamma_1\gamma_2} \\ \frac{\gamma_2}{1-\gamma_1\gamma_2} & \frac{1}{1-\gamma_1\gamma_2} \end{bmatrix}$$
 (7)

With this we found  $A^*(L)$ 

$$A^{*}(L) = \Gamma^{-1}A(L) = \begin{bmatrix} \frac{1}{1-\gamma_{1}\gamma_{2}} & \frac{\gamma_{1}}{1-\gamma_{1}\gamma_{2}} \\ \frac{\gamma_{2}}{1-\gamma_{1}\gamma_{2}} & \frac{1}{1-\gamma_{1}\gamma_{2}} \end{bmatrix} \begin{bmatrix} \alpha_{11}(L) & \alpha_{12}(L) \\ \alpha_{21}(L) & \alpha_{22}(L) \end{bmatrix}$$

$$A^{*}(L) = \begin{bmatrix} \frac{\alpha_{11}(L) + \alpha_{21}(L)\gamma_{1}}{1-\gamma_{1}\gamma_{2}} & \frac{\alpha_{12}(L) + \gamma_{1}\alpha_{22}(L)}{1-\gamma_{1}\gamma_{2}} \\ \frac{\gamma_{2}\alpha_{11}(L) + \alpha_{21}(L)}{1-\gamma_{1}\gamma_{2}} & \frac{\gamma_{2}\alpha_{12}(L) + \alpha_{22}(L)}{1-\gamma_{1}\gamma_{2}} \end{bmatrix}$$

$$(8)$$

Using  $A^*(L)$  we can calculate  $I_2 - A^*(L)$ :

$$I_{2} - A^{*}(L) = \begin{bmatrix} 1 & 0 \\ 0 & 1 \end{bmatrix} - \begin{bmatrix} \frac{\alpha_{11}(L) + \alpha_{21}(L)\gamma_{1}}{1 - \gamma_{1}\gamma_{2}} & \frac{\alpha_{12}(L) + \gamma_{1}\alpha_{22}(L)}{1 - \gamma_{1}\gamma_{2}} \\ \frac{\gamma_{2}\alpha_{11}(L) + \alpha_{21}(L)}{1 - \gamma_{1}\gamma_{2}} & \frac{\gamma_{2}\alpha_{12}(L) + \alpha_{22}(L)}{1 - \gamma_{1}\gamma_{2}} \end{bmatrix}$$

$$I_{2} - A^{*}(L) = \begin{bmatrix} \frac{1 - \gamma_{1}\gamma_{2} - \alpha_{11}(L) - \alpha_{21}(L)\gamma_{1}}{1 - \gamma_{1}\gamma_{2}} & -\frac{\alpha_{12}(L) + \gamma_{1}\alpha_{22}(L)}{1 - \gamma_{1}\gamma_{2}} \\ -\frac{\gamma_{2}\alpha_{11}(L) + \alpha_{21}(L)}{1 - \gamma_{1}\gamma_{2}} & \frac{1 - \gamma_{1}\gamma_{2} - \gamma_{2}\alpha_{12}(L) - \alpha_{22}(L)}{1 - \gamma_{1}\gamma_{2}} \end{bmatrix}$$

$$(9)$$

The inverse of  $I_2 - A^*(L)$  is:

$$[I_2 - A^*(L)]^{-1} = \begin{bmatrix} \frac{1 - \gamma_1 \gamma_2 - \gamma_2 \alpha_{12}(L) - \alpha_{22}(L)}{\theta} & \frac{\alpha_{12}(L) + \gamma_1 \alpha_{22}(L)}{\theta} \\ \frac{\gamma_2 \alpha_{11}(L) + \alpha_{21}(L)}{\theta} & \frac{1 - \gamma_1 \gamma_2 - \alpha_{11}(L) - \alpha_{21}(L)\gamma_1}{\theta} \end{bmatrix}$$
(10)

Where  $\theta = 1 - \gamma_1 \gamma_2 - \gamma_2 \alpha_{12}(L) - \alpha_{22}(L) - \alpha_{11}(L) - \alpha_{21}(L) \gamma_1 + \alpha_{11}(L) \alpha_{22}(L) - \alpha_{21}(L) \alpha_{12}(L)$ . For last we need to find  $[I_2 - A^*(L)]^{-1}\Gamma^{-1}$ , we have both matrix so:

$$[I_{2} - A^{*}(L)]^{-1}\Gamma^{-1} = \begin{bmatrix} \frac{1 - \gamma_{1}\gamma_{2} - \gamma_{2}\alpha_{12}(L) - \alpha_{22}(L)}{\theta} & \frac{\alpha_{12}(L) + \gamma_{1}\alpha_{22}(L)}{\theta} \\ \frac{\gamma_{2}\alpha_{11}(L) + \alpha_{21}(L)}{\theta} & \frac{1 - \gamma_{1}\gamma_{2} - \alpha_{11}(L) - \alpha_{21}(L)\gamma_{1}}{\theta} \end{bmatrix} \begin{bmatrix} \frac{1}{1 - \gamma_{1}\gamma_{2}} & \frac{\gamma_{1}}{1 - \gamma_{1}\gamma_{2}} \\ \frac{\gamma_{2}}{1 - \gamma_{1}\gamma_{2}} & \frac{1}{1 - \gamma_{1}\gamma_{2}} \end{bmatrix}$$

$$[I_{2} - A^{*}(L)]^{-1}\Gamma^{-1} = \begin{bmatrix} \frac{1 - \alpha_{22}(L)}{\theta} & \frac{\gamma_{1} + \alpha_{12}(L)}{\theta} \\ \frac{\gamma_{2} + \alpha_{21}(L)}{\theta} & \frac{1 - \alpha_{11}(L)}{\theta} \end{bmatrix}$$

$$(11)$$

The effect of  $e_2$  on  $y_1$  is  $\frac{\gamma_1 + \alpha_{12}(L)}{\theta}$ , the  $\frac{\partial y_{1,t+1}}{\partial e_{2,t}} = \frac{\alpha_{12}}{\theta}$  and  $\frac{\partial y_{1,t}}{\partial e_{2,t}} = \frac{\gamma_1}{\theta}$ 

#### Question 1.d

d) Are there any identification problems in estimating the structural parameters  $\gamma_1$  and  $\gamma_2$  from the structural VAR formulation? Explain, and if there are such problems, give in words a solution to these. We do have identification problem because we have eight parameters to estimate  $(\gamma_1, \alpha_{11}, \alpha_{12}, \gamma_2, \alpha_{21}, \alpha_{22}, \sigma_1 \text{ and } \sigma_2)$  but in the empirical model (the reduce form), we can estimate only seven estimates and hence there is underidentification problem in the model.

The solution for this problem implies to find a instrumental variable or make another restriction. But IV are hard to find and very unlikely to be truely exogenous. Exist others ways to make another restriction such as 1) The short-run restrictions, 2) The long-run restrictions, 3) The informal identification and 4) The sign restrictions to use the theory to make restrictions about the behaviour we expect of our model. The first one try to reduce the problem assuming an implicit ordering in the Model, with this order we can make one contemporaneous parameters (gamma one or gamma two) equal to zero, however the estimates about the error depends on the parameter equal to zero. The second one use the economic theory to impose impose a restriction on parameters but based on long-term relationships. For example, the change in monetary policy on GDP, we know that there is neutrality of money, so in the long term it must be annulled. But this condition is a non-linear constraint, so it complicates the problem and, depending on the problem, we need to identify whether or not this type of constraint can be applied. The third one model means to move the model until the impulse response function resembles what the investigator expects. This is an informal and very subjective methodology depending on what the researcher thinks. And the last one, a very similar to the third one, is the sign restriction in which the form in which it should be is not directly stated, but we must say for example that the effect of fiscal policy on GDP must be positive and the coefficients must be reduced to positive values, and with that it gives me a restriction that can be solved. For this particular case, we can use the sign restrictions to impose the effect of fiscal policy or the other alternative is to look for an instrument such as the announcement of fiscal policy.

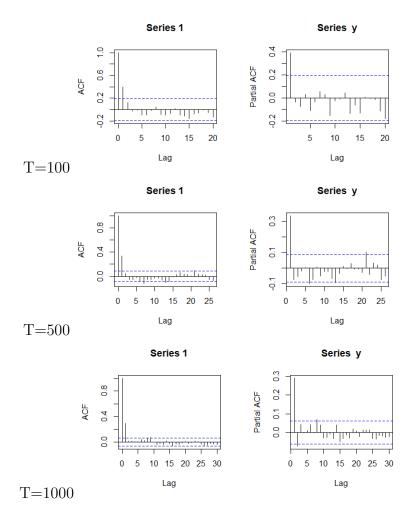
# Question 2

a) Let the true model be  $y_t = \alpha + u_t$ , where  $u_t = \epsilon_t + \beta \epsilon_{t1}$  with  $\beta = 0.7$ , and  $\epsilon_t$  i.i.d.N (0, 1), and  $\alpha = 0$ . Let T be the sample size. This is a model unknown the econometrician, who is interested in testing whether  $E(y_t) = 0$ , and knows there may be some autocorrelation in the errors.

## Question 2.a

Simulate one sample from this model letting T = 100, 500, 1000. Plot the ACF and PACF of this model, and explain whether they make sense.

After simulate the moving average model (MA1), we can see that as we increase the T from 100 to 1000, the interval of confidence shrinks. Therefore, it makes possible more accurate results. Also, as the model is MA(1), the autocorrelation function has just one significant lag. It makes sense with the model and the observation we added to the simulation.



#### Question 2.b

Simulate S = 500 replications from this model for T = 100, 500, 1000 with burn-in B = 100 observations each time (so you should have 500 datasets, each with T observations). Calculate for each simulation the t-statistic for the null hypothesis that  $\alpha = 0$   $t = \overline{y}/\hat{\sigma y}$ , where  $\overline{y} = PT$  t=1  $y_t$ , and  $\hat{\sigma}_{y}$  is the Newey-West HAC estimator of the standard error of  $\overline{y}$  using the quadratic spectral kernel. Calculate the fraction of times in S = 500 simulations that this test rejects, for each sample size, and significance level 5% (using standard normal critical values), and compare it the nominal level 5%. Are they close or not to this nominal 5% level? If yes, explain why.

After the simulation S=500 for T=100, 500 and 1000 was done, it was calculated for each simulation the t-statistic  $t = \overline{y}/\sigma_y$  where  $\sigma_y$  was estimated using Newey-West HAC estimator. After that, the fraction rejection time was calculated. As a result for T=100, 500 and 1000 the fraction rejection was of 0.082, 0.072 and 0.06, respectively. It can be concluded that the rejection rate get closer to 5% as we increase the sample size because as the sample increase, it resembles to a normal distribution.

#### Question 2.c

Let the true model now by

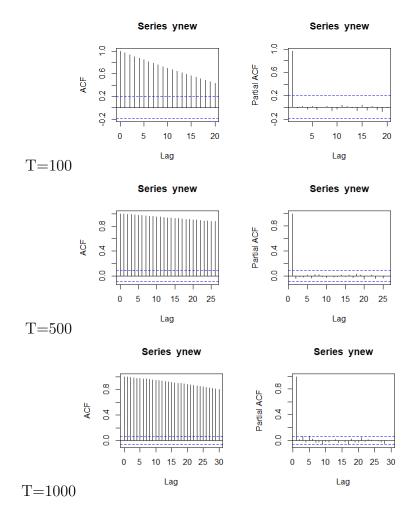
$$y_t = \alpha + 0.7y_{t1} + u_t$$
, where  $u_{1-t}$  i.i.d.N (0, 1), and  $\alpha = 0$ 

Let T be The sample size. This is again a model unknown the econometrician, who is interested in testing whether E(yt)=0, and knows there may be some autocorrelation in the errors. Simulate one sample from this model letting T = 100, 500, 1000, with B = 100 burn-in observations (so compute the model forward with T + B observations, and discard the first B observations) Explain why we need to do this "burn-in".

The simulation was done for an AR(1) model with T = 100, 500, 1000, B = 100 and we added a burn-in equal to T+B and then the first observation of B was deleted. The burn-in method is a good way to find a good starting point when the one we have is bias. This is because, unlike MA(1), AR(1) is not always a stationary model. Therefore if we include initial values the estimates would be bias. Therefore, using burn-in we guarantee that the model will be stationary and unbiased.

## Question 2.d

Based on c), plot the ACF and PACF of this model, and explain whether they make sense



The plots of PACF for each sample show a significant lag of one because as we know the model is a AR(1). The plots of ACF for each sample shows several significant lags (that increase with the sample size) decaying. It represent an AR(1) that is increasing with the sample size. Finally, we can notice that the confidence values get shrink as the sample size increase. It allows us more precision in the final results.

## Question 2.e

Simulate S = 500 replications from this model for T = 100, 500, 1000 with burn-in B = 100 observations each time (so you should have 500 datasets, each with T observations). Calculate for each simulation the t-statistic for the null hypothesis that  $\alpha = 0$   $t = \bar{y}/\hat{\sigma y}$ , where  $\bar{y} = PT$  t=1  $y_t$ , and  $\hat{\sigma}_{y_l}$  is the Newey-West HAC estimator of the standard error of  $\bar{y}$  using the quadratic spectral kernel. Calculate the fraction of times in S = 500 simulations that this test rejects, for each sample size, and significance level 5% (using standard normal critical values), and compare it the nominal level 5%. Are they close or not to this nominal 5% level? If yes, explain why. If not, explain and investigate what is the source of distortions, keeping in mind that the simulated (also called)

"empirical" rejection frequency should be close to the nominal level as the sample size increases, because asymptotically, the t test is normally distributed.

After the simulation S=500 for T=100, 500 and 1000 was done, it was calculated for each simulation the t-statistic  $t = \overline{y}/\sigma_y$  where  $\sigma_y$  was estimated using Newey-West HAC estimator. After that, the fraction rejection time was calculated. As a result for T=100, 500 and 1000 the fraction rejection was of 0.168, 0.088 and 0.08, respectively. It can be concluded that the rejection rate decays as we increase the sample size. However, in comparison with MA(1), the rejection rate do not get close to nominal 5% level even for higher sample sizes. One reason of this is because the Newey-West HAC estimator could be biased therefore the rejection come out to be more than 5%. In addition, the AR(1) model is much more sensitive for this bias than MA process therefore the difference between the rejection rates as the sample size increase.

# **Appendix**

```
install.packages("lmtest")
2 library(lmtest)
  install.packages("forecast")
  library(forecast)
  install.packages("tseries")
6 library(tseries)
  install.packages("dplyr")
  library(dplyr)
  library (sandwich)
11
  # Solution 2(a) -----
12
13
  #Sample 100#
14 | Ti = 100
| 15 | beta = 0.7
  alpha = 0
17
  e = matrix(0,nrow = Ti,ncol = 1)
18
  e[1,1] = rnorm(1, mean=0, sd = 1)
19
  for (i in 2:Ti){
    e[i,1] = rnorm(1, mean = 0, sd = 1)
22
23
  plot(e)
24
25
  u = matrix(0, nrow = Ti, ncol = 1)
  u[1,1] = e[1,1]
27
  for (i in 2:Ti){
   u[i,1] = e[i,1] + beta*e[i-1,1] + rnorm(1,sd = 1)
29
30 }
31
32 y = matrix(0, nrow = Ti, ncol = 1)
  y = alpha + u
33
34
35
  acf(y)
36
  pacf(y)
37
38
  #Sample 500#
39 | Ti = 500
40 | beta = 0.7
41 alpha = 0
42
  e = matrix(0,nrow = Ti,ncol = 1)
|44| e[1,1] = rnorm(1, mean=0, sd = 1)
```

```
45 for (i in 2:Ti){
   e[i,1] = rnorm(1, mean = 0, sd = 1)
46
47
48
49 plot(e)
51 u = matrix(0, nrow = Ti, ncol = 1)
52 u[1,1] = e[1,1]
53 for (i in 2:Ti){
    u[i,1] = e[i,1] + beta*e[i-1,1] + rnorm(1,sd = 1)
54
55 }
56
   y = matrix(0, nrow = Ti, ncol = 1)
57
58 y = alpha + u
59
60 acf(y)
61 pacf(y)
63 #Sample 1000
64 | Ti = 1000
65 | beta = 0.7
66 alpha = 0
68 = \text{matrix}(0, \text{nrow} = \text{Ti}, \text{ncol} = 1)
69 | e[1,1] = rnorm(1, mean=0, sd = 1)
70 for (i in 2:Ti){
    e[i,1] = rnorm(1, mean = 0, sd = 1)
71
72 }
73
74 plot(e)
75
76 \mid u = matrix(0, nrow = Ti, ncol = 1)
|u[1,1]| = e[1,1]
78 for (i in 2:Ti){
    u[i,1] = e[i,1] + beta*e[i-1,1] + rnorm(1,sd = 1)
79
80 }
81
82 y = matrix(0, nrow = Ti, ncol = 1)
y = alpha + u
84
   acf(y)
85
86 pacf(y)
88 # Solution 2(b) -----
89 #Sample=100
90 rm(list = ls())
   set.seed (4534938432)
92
93 S = 500
94 B = 100
95 Ti = 100
96 | beta = 0.7
97 alpha = 0
98 e = matrix(0, nrow = Ti, ncol = S)
99 for (i in 1:S){
    e[1,i] = rnorm(1, mean = 0, sd = 1)
100
     for (j in 2:Ti){
101
       e[j,i] = rnorm(1, mean = 0, sd = 1)
102
103
104 }
105
   u = matrix(0, nrow = Ti, ncol = S)
106 for (i in 1:S){
    u[1,i] = e[1,i]
107
108
   for (j in 2:Ti){
       u[j,i] = e[j,i] + beta*e[j-1,i]
109
110
111 }
112 y = matrix(0, nrow = Ti, ncol = S)
```

```
113 y = alpha + u
114
115
   ymean = matrix(0, nrow = 1, ncol = S)
   for (i in 1:S){
116
117
    ymean[,i] = mean(y[,i])
118 }
   sigmahat = matrix(0, nrow = 1, ncol = S)
119
120 for (i in 1:S){
    sigmahat[,i] = sqrt(kernHAC(lm(y[,i] ~ 1), bw = bwNeweyWest, prewhite = 0))
121
122 }
123 tstat = matrix(0, nrow = 1, ncol = S)
124 for (i in 1:S){
125
     tstat[,i] = ymean[,i]/sigmahat[,i]
126 }
127 reject = matrix(0, nrow = 1, ncol = S)
128 for (i in 1:S){
    reject[,i] = isTRUE(tstat[,i]>1.96) | isTRUE(tstat[,i]<(-1.96))
129
130
131 table (reject)
132 hist(tstat)
133 plot(density(tstat))
134
   #Sample 500#
135
136
137 S = 500
138 B = 100
139 Ti = 500
140 | beta = 0.7
141 alpha = 0
142 e = matrix(0, nrow = Ti, ncol = S)
143 for (i in 1:S){
    e[1,i] = rnorm(1, mean = 0, sd = 1)
144
145
     for (j in 2:Ti){
       e[j,i] = rnorm(1, mean = 0, sd = 1)
146
147
148 }
   u = matrix(0, nrow = Ti, ncol = S)
149
150 for (i in 1:S){
    u[1,i] = e[1,i]
151
152
     for (j in 2:Ti){
153
       u[j,i] = e[j,i] + beta*e[j-1,i]
154
155 }
|y| = matrix(0, nrow = Ti, ncol = S)
|y| = alpha + u
158
159
   ymean = matrix(0, nrow = 1, ncol = S)
160 for (i in 1:S){
    ymean[,i] = mean(y[,i])
161
162 }
163
   sigmahat = matrix(0, nrow = 1, ncol = S)
164
   for (i in 1:S){
    sigmahat[,i] = sqrt(kernHAC(lm(y[,i] ~ 1), bw = bwNeweyWest, prewhite = 0))
165
166 }
   tstat = matrix(0, nrow = 1, ncol = S)
167
   for (i in 1:S){
168
    tstat[,i] = ymean[,i]/sigmahat[,i]
169
170 }
171 reject = matrix(0, nrow = 1, ncol = S)
172 for (i in 1:S){
    reject[,i] = isTRUE(tstat[,i]>1.96) | isTRUE(tstat[,i]< (-1.96))
173
174 }
175 table (reject)
176 hist(tstat)
177 plot (density (tstat))
179 #Sample 1000#
180
```

```
181
182 S = 500
183 B = 100
184 Ti = 1000
185 | beta = 0.7
186 alpha = 0
   e = matrix(0, nrow = Ti, ncol = S)
187
   for (i in 1:S){
188
    e[1,i] = rnorm(1, mean = 0, sd = 1)
189
     for (j in 2:Ti){
190
       e[j,i] = rnorm(1, mean = 0, sd = 1)
191
192
193
   u = matrix(0, nrow = Ti, ncol = S)
194
195 for (i in 1:S){
    u[1,i] = e[1,i]
196
    for (j in 2:Ti){
197
198
      u[j,i] = e[j,i] + beta*e[j-1,i]
199
200 }
201 y = matrix(0, nrow = Ti, ncol = S)
202 y = alpha + u
204 ymean = matrix(0, nrow = 1, ncol = S)
205 for (i in 1:S){
   ymean[,i] = mean(y[,i])
206
207
   sigmahat = matrix(0, nrow = 1, ncol = S)
208
   for (i in 1:S){
209
   sigmahat[,i] = sqrt(kernHAC(lm(y[,i] ~ 1), bw = bwNeweyWest, prewhite = 0))
211
   tstat = matrix(0, nrow = 1, ncol = S)
212
   for (i in 1:S){
213
   tstat[,i] = ymean[,i]/sigmahat[,i]
214
215 }
216 reject = matrix(0, nrow = 1, ncol = S)
   for (i in 1:S){
   reject[,i] = isTRUE(tstat[,i]>1.96) | isTRUE(tstat[,i]< (-1.96))
218
219 }
220
221
   table(reject)
222
   hist(tstat)
   plot(density(tstat))
223
225
226
228 # Solution 2(c, d) ------
230
231
   # If T (sample size) = 100
   rm(list = ls())
232
233 set.seed(600000)
235
236 Ti = 100
237 B = 100
238 | beta = 0.7
239 alpha = 0
240 burnin = Ti+B
   y = matrix(0, nrow = burnin, ncol = 1)
242
243 for (i in 2:burnin) {
244
   y[i,1] = alpha + y[i-1,1] + rnorm(1, mean = 0, sd = 1)
245 }
246
247 ynew = tail(y,-B)
248
```

```
249
250 acf (ynew)
251
  pacf (ynew)
254 # If T (sample size) = 500
255 Ti = 500
256 B = 100
257 beta = 0.7
258 alpha = 0
259 burnin = Ti+B
260
  y = matrix(0, nrow = burnin, ncol = 1)
261
262 for (i in 2:burnin){
   y[i,1] = alpha + y[i-1,1] + rnorm(1, mean = 0, sd = 1)
263
264 F
265
266
  ynew = tail(y, -B)
267
268 acf (ynew)
269 pacf (ynew)
270
271
  272 # sample 1000
273 Ti = 1000
274 B = 100
275 beta = 0.7
276 alpha = 0
277 burnin = Ti+B
279 y = matrix(0, nrow = burnin, ncol = 1)
280 for (i in 2:burnin){
   y[i,1] = alpha + y[i-1,1] + rnorm(1, mean = 0, sd = 1)
281
282 }
283
284 ynew = tail(y,-B)
285
286
287 acf (ynew)
288 pacf (ynew)
289
  # Solution 2(e) ------
291
292 rm(list = ls())
293 set.seed(600000000)
294
295
  # sample 100
296 S = 500
297 B = 100
298 Ti = 100
299 beta = 0.7
300
  alpha = 0
  burnin = Ti+B
301
303 y = matrix(0, nrow = burnin, ncol = S)
  for (i in 1:S){
304
   for (j in 2:burnin){
305
      y[j,i] = alpha + beta*y[j-1,i] + rnorm(1, mean = 0, sd = 1)
306
307
308 }
309
  ynew = tail(y, -B)
310
311
312 ymean = matrix(0, nrow = 1, ncol = S)
313 for (i in 1:S){
314
    ymean[,i] = mean(y_new[,i])
315 }
316 sigmahat = matrix(0, nrow = 1, ncol = S)
```

```
317 for (i in 1:S){
     sigmahat[,i] = sqrt(kernHAC(lm(y_new[,i] ~ 1), bw = bwNeweyWest, prewhite = 0))
318
319
   tstat = matrix(0, nrow = 1, ncol = S)
320
   for (i in 1:S){
321
     tstat[,i] = ymean[,i]/sigmahat[,i]
322
323 }
324 reject = matrix(0, nrow = 1, ncol = S)
325
   for (i in 1:S){
    reject[,i] = isTRUE(tstat[,i]>1.96) | isTRUE(tstat[,i]< (-1.96))
326
327 }
   table(reject)
328
   hist(tstat)
329
330
   plot(density(tstat))
331
   332
333
334
   # sample 500
335 S = 500
336 B = 100
337 Ti = 500
338 | beta = 0.7
339
   alpha = 0
   burnin = Ti+B
340
342 y = matrix(0, nrow = burnin, ncol = S)
343
   for (i in 1:S){
344
    for (j in 2:burnin){
       y[j,i] = alpha + beta*y[j-1,i] + rnorm(1, mean = 0, sd = 1)
345
346
347
   }
348
349
   ynew = tail(y, -B)
350
351 ymean = matrix(0, nrow = 1, ncol = S)
   for (i in 1:S){
352
353
    ymean[,i] = mean(y_new[,i])
354 }
355 sigmahat = matrix(0, nrow = 1, ncol = S)
356
   for (i in 1:S){
357
    sigmahat[,i] = sqrt(kernHAC(lm(y_new[,i] ~ 1), bw = bwNeweyWest, prewhite = 0))
358
   tstat = matrix(0, nrow = 1, ncol = S)
359
   for (i in 1:S){
360
    tstat[,i] = ymean[,i]/sigmahat[,i]
361
362 }
   reject = matrix(0, nrow = 1, ncol = S)
363
   for (i in 1:S){
364
    reject[,i] = isTRUE(tstat[,i]>1.96) | isTRUE(tstat[,i]< (-1.96))
365
366 }
367
   table(reject)
368
   hist(tstat)
   plot(density(tstat))
369
370
   *****************
371
372
373 # If T (sample size) = 1000
374 S = 500
375 B = 100
376 Ti = 1000
   beta = 0.7
377
378 alpha = 0
379 burnin = Ti+B
381 y = matrix(0, nrow = burnin, ncol = S)
382 for (i in 1:S){
   for (j in 2:burnin){
383
      y[j,i] = alpha + beta*y[j-1,i] + rnorm(1, mean = 0, sd = 1)
384
```

```
385
386
   }
387
   ynew = tail(y,-B)
388
389
390 ymean = matrix(0, nrow = 1, ncol = S)
391 for (i in 1:S){
392
    ymean[,i] = mean(y_new[,i])
393 }
394 sigmahat = matrix(0, nrow = 1, ncol = S)
395 for (i in 1:S){
    sigmahat[,i] = sqrt(kernHAC(lm(y_new[,i] ~ 1), bw = bwNeweyWest, prewhite = 0))
396
397
   tstat = matrix(0, nrow = 1, ncol = S)
398
399 for (i in 1:S){
   tstat[,i] = ymean[,i]/sigmahat[,i]
400
401 }
   reject = matrix(0, nrow = 1, ncol = S)
402
403 for (i in 1:S){
   reject[,i] = isTRUE(tstat[,i]>1.96) | isTRUE(tstat[,i]< (-1.96))
404
405 }
406 table (reject)
407 hist(tstat)
408 plot(density(tstat))
```