Unit root problem, solution of difference equations

Simple deterministic model, question of unit root

$$(1 - \phi_1 L)X_t = \mu$$

,

$$X_t - \phi_1 X_{t-1} = \mu$$

Solution

$$X_t = A + Bz^t$$

with unknown z and unknown A (clearly $X_{t-1} = A + Bz^{t-1}$)

$$0 = X_t - \phi_1 X_{t-1} - \mu = A + Bz^t - \phi_1 (A + Bz^{t-1}) - \mu$$

$$= A + Bz^{t} - \phi_{1}\left(A + \frac{B}{z}z^{t}\right) - \mu$$

$$=\underbrace{-\mu + A(1-\phi_1)}_{=0} + B\underbrace{\left(1-\frac{\phi_1}{z}\right)}_{=0} z^t$$

 $z = \phi_1$, $A = \frac{\mu}{1 - \phi_1}$ Elimination B by the initial condition

$$X_t = \frac{\mu}{1-\phi_t} + B\phi_1^t$$

$$X_{t_0} = rac{\mu}{1-\phi_1} + B\phi_1^{t_0} \longrightarrow B = \left[X_{t_0} - rac{\mu}{1-\phi_1}
ight]\phi_1^{-t_0}$$
 $X_t = rac{\mu}{1-\phi_1} + \left[X_{t_0} - rac{\mu}{1-\phi_1}
ight]\phi_1^{-t_0}\phi_1^t$

$$egin{align} \left(X_t - rac{\mu}{1 - \phi_1}
ight) &= \left(X_{t_0} - rac{\mu}{1 - \phi_1}
ight) \phi_1^{t - t_0} \ & rac{X_t - rac{\mu}{1 - \phi_1}}{X_{t_0} - rac{\mu}{1 - \phi_1}} &= \phi_1^{t - t_0} \ \end{aligned}$$

Qualitative difference for
$$\phi_1 < 1$$
 and $\phi_1 > 1$ (unit root means $z = \phi_1 = 1$) $\phi_1 = 0.8$: $t_1 - t_0 = 0$, $\phi_1^0 = 1$; $t - t_0 = 1$, $\phi_1^1 = 0.8$; $t - t_0 = 2$,

Unit root signature of nonstationary behaviour $z=\phi_1=1$ Take homogeneous equation $(1-\phi_1 L)X_t=0, \longrightarrow (1-L)X_t=0$ formally via the "trick" L=1/z we get characteristic equation $(1-\phi_1/z)=0$ where the unit root is achieved for $z=\phi_1=1$

Generalization for AR(p) without stochastic component:

$$X_t - \sum_{i=1}^p \phi_i X_{t-i} = 0$$

$$X_t - \phi_1 X_{t-1} - \phi_2 X_{t-2} - \phi_3 X_{t-3} \dots \phi_p X_{t-p} = 0$$

$$(1 - \phi_1 L - \phi_2 L^2 - \phi_3 L^3 \dots - \phi_p L^p) X_t = 0$$

expected form of solution

$$X_t = z^t X_0, t = 0, 1, 2 \dots$$

$$(1 - \phi_1(1/z) - \phi_2(1/z)^2 - \phi_3(1/z)^3 \dots - \phi_p(1/z)^p) z^t X_0 = 0$$

$$\left[1-\phi_1(1/z)-\phi_2(1/z)^2-\phi_3(1/z)^3\ldots-\phi_p(1/z)^p\right]z^tX_0=0$$

characteristic equation

$$z^p-\phi_1z^{p-1}-\phi_2z^{p-2}-\phi_3z^{p-3}-\ldots-\phi_p=0$$
 generally has exactly p roots

$$(z-\lambda_1)(z-\lambda_2)\dots(z-\lambda_p)=0$$

$$\prod_{k=1}^{p} (z - \lambda_k)^k = 0$$

For non-identical (also complex) roots

$$X_t = B_1 \lambda_1^t + B_2 \lambda_2^t + \ldots + B_p \lambda_p^t$$

when roots are identical for e.g. p = 2

$$X_t = B_1 \lambda_1^t + B_2 t \lambda_1^t$$

Augmented Dickey-Fuller (ADF)

unit root test

(Rozšírený DF test)

The test is widely used by practioners. In statistics, the traditional DickeyFuller test tests (1979) whether a unit root is present in an autoregressive model.

In time series models in econometrics (the application of statistical methods to economics), a **unit root** is a feature of processes that evolve through time that can cause **problems** in statistical inference if it is not adequately dealt with. Such a process is **non-stationary**.

(If the other roots of the characteristic equation lie inside the unit circle that is, have a modulus (absolute value) less than one then the first difference of the process will be stationary). Other words: series of the first difference (previously considered integrated of order one) turned to be stationary.

R: test available in **adf.test() tseries()** [package for Time series analysis and computational financel

The testing procedure is applied to the model

$$\Delta X_t = \underbrace{\alpha}_{\text{drift}} + \underbrace{\delta t}_{\text{lin.trend}} + \underbrace{\varrho}_{\text{par. tested}} X_{t-1} + \sum_{j=1}^p \psi_j \Delta X_{t-j}$$

 $\Delta X_{t-j} = X_{t-j} - X_{t-j-1}$, $\Delta X_t = X_t - X_{t-1}$ number of lags of autoregression; default: $p = (\text{length of d.s.} - 1)^{1/3}$; (example: length of data series = 65, p = 4); Swert criterion: $12 \times (\text{length of d.s.}/100)^{1/4}$ Alternative form:

$$X_t = \alpha + \delta t + \underbrace{(1+\varrho)}_{\phi_1 \text{ in former notation}} X_{t-1} + \sum_{j=1}^p \psi_j \Delta X_{t-j}$$

 $\sum_{j=1}^p \psi_j \Delta X_{t-j}$ the autoregression terms absorbing dynamic structure

<u>H₀</u>: $\varrho = 0$ for given model **has a unit root**, process is I(1), i.e. the data needs to be differenced to make it stationary;

 $\underline{H_1}$: $\varrho < 1$ i.e. the data is **trend stationary**, process is I(0) and needs to be analyzed by means of using a trend in the regression model instead of differencing the data;

The unit root test is carried out under H_0 : $\varrho=0$ against the alternative hypothesis of $\varrho<0$.

<u>Note</u> 1: *Caution*: Sometimes if data is exponentially trending then you might need to take the log of the data first before differencing it.

Note 2: $\varrho = 0$ means $X_t \sim X_{t-1}$ (i.e. random walk process)

test statistic [one-sided left tail test]:

$$ADF_{\tau} = \frac{\hat{\varrho}}{SE(\hat{\varrho})}$$

(remind that
$$t_{ ext{score}} = rac{\hat{arrho}}{SE(\hat{arrho})}$$
)

 $SE(\hat{\varrho})$ is the usual standard error estimate critical value: $t_{1-\alpha}^*(n)$

If the test statistic is <u>less than</u> the **critical value** $t_{1-\alpha}^*(n)$, then the H₀ ($\varrho = 0$) is rejected and **no unit root is present**.

Example: A model that includes a constant and a time trend is estimated using sample of n=50 observations and yields the ADF $_{\tau}$ statistic of -4.57. This is more negative than the tabulated $t^*=-3.50$, so at the 95 per cent level the null hypothesis of a unit root will be rejected.

R implementation

```
> x=arima.sim(n = 1000, list(ma = c(-0.2, 0.2)),sd=1)
> adf.test(x)
```

Augmented Dickey-Fuller Test

data: x

Dickey-Fuller = -9.8277, Lag order = 9, p-value = 0.01 alternative hypothesis: stationary

Warning message:

In adf.test(x) : p-value smaller than printed p-value

[Note: rejection of null hypothesis that data has a unit root]

Augmented Dickey-Fuller Test

data: x

Dickey-Fuller = -0.038, Lag order = 9, p-value = 0.99 alternative hypothesis: stationary

Warning message:

In adf.test(x) : p-value greater than printed p-value

KPSS test

Kwiatkowski, Phillips, Schmidt, Shin (1992)

observed time series x_t is expressed as the sum $x_t = T_t + RW_t + u_t$:

- deterministic trend T_t
- pure random walk RW_t (i.e. I(1)), $RW_t = RW_{t-1} + \epsilon_t$
- stationary error term u_t (i.e. I(0))

$$\underline{\mathsf{H}_0\colon\thinspace \sigma_\epsilon^2=\mathsf{0}}$$

observable time series is stationary around a deterministic trend, x_t is I(0)

$$H_1$$
: $\sigma_{\epsilon}^2 > 0$

```
> s<-rnorm(1000)
> kpss.test(s)
>
> KPSS Test for Level Stationarity
> data: s
> KPSS Level = 0.0429, Truncation lag parameter = 7,
> p-value = 0.1
```

in R:

... stationarity cannot be rejected here ...

Second order Stochastic Difference Equation and Business Cycles

Samuelson (1939) applied only second order difference equations to explain the business cycles:

Investment I_t is dependent on the changes in income rather then the level of income

$$I_t = b_1(Y_{t-1} - Y_{t-2})$$

Consumption C_t depends on the past income by the first order nonhomogeneous equation

$$C_t = b_2 Y_{t-1} + b_3$$

Production/consumption balance with error term

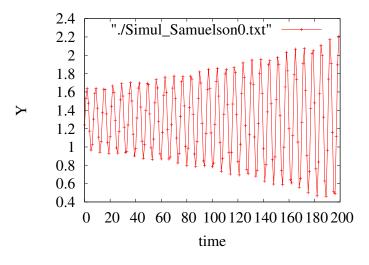
$$Y_t = C_t + I_t + \epsilon_t$$

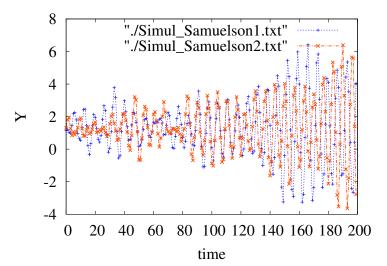


Consequence - accelerator principle

$$Y_t = b_1(Y_{t-1} - Y_{t-2}) + b_2Y_{t-1} + b_3 + \epsilon_t$$

simulation in R





Solution of the homogeneous (non-stochastic) Samuelson's system

$$Y_t = z^t Y_0$$

$$z =, b_3 = 0$$

$$Y_t = b_1(Y_{t-1} - Y_{t-2}) + b_2 Y_{t-1}$$

$$z^{t}Y_{0} = b_{1}(z^{t-1}Y_{0} - z^{t-2}Y_{0}) + b_{2}Y_{0}z^{t-1}$$

$$1 = b_1 \left(\frac{1}{z} - \frac{1}{z^2} \right) + b_2 \frac{1}{z}$$

lag operator formally produces $\sim 1/z$

$$b_1\left(\frac{1}{z}-\frac{1}{z^2}\right)+b_2\frac{1}{z}-1=0$$

$$\frac{b_1 + b_2}{7} - \frac{b_1}{7^2} - 1 = 0$$

$$z^2 - (b_1 + b_2)z + b_1 = 0$$

$$z_{1,2} = \frac{b_1 + b_2 \pm \sqrt{(b_1 + b_2)^2 - 4b_1}}{2}$$

$$b_1 = 1.01, b_2 = 0.2, b_3 = 1.04$$

$$z_{1,2} = \frac{1.21 \pm \sqrt{1.21^2 - 4.04}}{2} = 0.605 \pm i0.8024$$

$$|z_{1,2}| = 1.00492$$
, $\arccos(0.605/|z_{1,2}|) = 0.1465(2\pi)$

$$Y_t = Y_0 z^t = (1.00492)^t \left[\cos(\pm (2\pi)0.147t) + i \sin(\pm (2\pi)0.147t) \right]$$

=
$$(1.00492)^t \exp(i2\pi 0.147t) = (1.00492)^t \exp\left(i2\pi \frac{t}{6.8}\right)$$

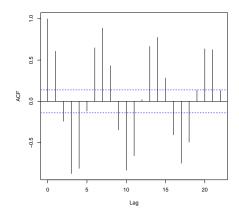


X1.3751803863379

Qualitative check of the period

x 1	0	066	046	1000	201

- 1.02464804231704
 - 1.30411746998738
 - 1.58308761594452
- 1.63837737060561
- 1.42351812632883
- 1.10769578854622
- x 8 0.9425585965488
- - 1.06172315539237
- 1.37270083551048 10
- 11 1.62862762402139
- 12 1.62421158120029
- 13 1.36038211299075
- 14 1.04560865970652
- x15 0.931200544124225



Comparison Samuelson's accelerator ($b_3 = 0$) with ARIMA(p,d,.)

$$Y_t = b_1(Y_{t-1} - Y_{t-2}) + b_2Y_{t-1}$$

 $\left(1 - \sum_{i=1}^{p} \phi_i L^i\right) (1 - L)^d Y_t = 0$

Try
$$p = 1, d = 1$$

$$(1 - \phi_1 L)(1 - L)Y_t = (1 -$$

$$(1 - \phi_1 L)(1 - L)Y_t = (1 - \phi_1 L)(Y_t - Y_{t-1})$$

$$=Y_t-Y_{t-1}-\phi_1(LY_t-LY_{t-1})$$

$$= Y_{t} - Y_{t-1} - \phi_{1}(Y_{t-1} - Y_{t-2})$$

Try
$$p = 2, d = 1$$

$$(1 - \phi_1 - \phi_2 L)(1 - L)Y_t = (1 - L - L\phi_1 + L^2\phi_1 + L^2\phi_2 + L^3\phi_2)$$

$$= Y_t - Y_{t-1} - \phi_1 Y_{t-1} + \phi_1 Y_{t-2} - \phi_2 Y_{t-2} + \phi_2 Y_{t-3})$$
Try $p = 2, d = 0$

$$(1 - \phi_1 L - \phi_2 L^2) Y_t = Y_t - \phi_1 Y_{t-1} - \phi_2 Y_{t-2}$$
$$Y_t - \phi_1 Y_{t-1} - \phi_2 Y_{t-2} = 0$$
$$Y_t = \phi_1 Y_{t-1} + \phi_2 Y_{t-2}$$

$$Y_t = b_1(Y_{t-1} - Y_{t-2}) + b_2Y_{t-1} = \underbrace{(b_1 + b_2)}_{\phi_1}Y_{t-1} + \underbrace{(-b_1)}_{\phi_2}Y_{t-2}$$

<u>Conclusion</u>: Model is transformable to ARIMA(2,0,q) i.e. d = 0 model

Box Jenkins metodology

- 1. identification of model, mostly AR(1)
 - making sure that model is stationary
 - using plots of ACF, and PACF
- 2. estimation of the model parameters: say $x_t = 0.68x_{t-1} + \epsilon_t$, $\hat{\sigma} = 11.24$
 - using the maximum likelihood
 - non-linear least-squares
- 3. diagnostics of the model: prediction abilities, Ljung-Box test Model identification by means of autocorrelation function:

	AR(p)	MA(q)	ARMA(p,q)
ρ_k	k_0 does not exist	$k_0 = q$	k_0 does not exist
ACF	ρ_{k} is U shaped		ρ_k becomes
			U-shaped after first q
ρ_{kk}	$k_0 = p$	k ₀ does not exist	k ₀ does not exist
$ ho_{kk}$ PACF		ρ_{kk} is limited	ρ_{kk} becomes
		by U-shaped curve	U-shaped after the firs

Table: U denotes the curve of geometrically or sin-like decaying amplitudes.

- Rule 1 If the series has positive autocorrelations out to a high number of lags, then it probably needs a higher order of differencing.
- Rule 2 If the lag-1 autocorrelation is zero or negative, or the autocorrelations are all small a patternless, then the series does not need a higher order of differencing. If the lag-1 autocorrelation is -0.5 or more negative, the series may be overdifferenced. Beware of overdifferencing.
- Rule 3 The optimal order of differencing is then often the order at which the standard deviation is lowest.
- Rule 4 A model with no orders of differencing assumes that the original series is stationary.
- Rule 5 A model with no orders of differencing normally includes a constant term
- Rule 6 If the part of PACF of the differenciated series displays a sharp cutoff and/or the lag 1 autocorrelation is positive, i.e. the

- series appears slightly "underdifferenced" then consider adding one or more AR terms to the model
- Rule 7 The autocorrelation function ACF of the differenced series displays a sharp cutoff and or the lag-1 autocorrelation function is negative .
- Rule 8 It is possible for an AR term and MA term to cancel each other's effects, so if a mixed AR-MA model seems to fit data also try a model with one fewer AR term.
- Rule 9 If there is a unit root in the AR part of the model the sum of AR coefficients is almost exactly 1
- Rule 10 If the long-term forecasts appear erratic or unstable, there may be a unit root.
- Rule 11 If the series has a strong and consistent seasonal pattern, then you should use an order of seasonal differencing.

Autoregression of simulated time series of the price formation process

Model of random process: p_0 , p_1 , p_2 , p_3 , ..., p_t , ...

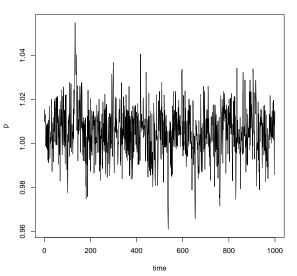
$$\mathsf{if}\, \pmb{p}_t > \pmb{p}_{\mathrm{fun}}\, \left\{ \pmb{\tilde{p}}_{t+1} = \pmb{p}_t \, \times 0.99 \right\} \\ \mathsf{else} \quad \ \pmb{\tilde{p}}_{t+1} = \pmb{p}_t$$

$$\begin{array}{c} \text{if } \tilde{\pmb{p}}_{t+1} < \pmb{p}_{\text{fun}} \ \left\{ \underline{\pmb{p}}_{t+1} = \tilde{\pmb{p}}_{t+1} \times 1.01 \right\} \\ \text{else} \quad \underline{\pmb{p}}_{t+1} = \tilde{\pmb{p}}_{t+1} \end{array}$$

$$p_{t+1} = \underline{p}_{t+1} \exp [N(0, 0.01)]$$

```
R project: simulation of price dynamics
pfun<-1
                  # "fundamental" price
p<-(pfun+0.01)
                  # initial price setting
                     auxilliary initial vector setting for pl
vp<-p
Ns<-1000
                  # number of steps
time=c(0:Ns) # vector of times
for(i in 1:Ns){
if(p>pfun) p<-(p * 0.99) # deterministic reduction
if(p<pfun) p<-(p * 1.01) # deterministic increase
p<-(p*exp(rnorm(1,mean=0,sd=0.01))) #randomization
vp<-cbind(vp,p)</pre>
vp<-as.vector(vp)</pre>
                                     ◆□▶ ◆□▶ ◆□▶ ◆□▶ □ ◆○○○
```

```
pdf("tp_simul.pdf")
plot(time,vp, type="l", pch=21, xlab="time", ylab="p")
dev.off()
```



Testing of simulated price series for stationarity

adf.test(vp)

Augmented Dickey-Fuller Test

data: vp

Dickey-Fuller = -9.0344, Lag order = 9, p-value = 0.01

alternative hypothesis: stationary

Warning message:

In adf.test(vp) : p-value smaller than printed p-value

conclusion: reject unit root

Selection of the "best" arima model

based on the Aikake information criterion

The Akaike information criterion is a measure of the **relative**. **goodness of fit** of a statistical model. It is grounded in the concept of information entropy, in effect offering a relative measure of the information lost when a given model is used to describe reality. In the general case, the AIC is $AIC = 2k - 2\ln(L)$, where k is the number of parameters in the statistical model, and L is the maximized value of the likelihood function for the estimated model. To apply AIC in practice, we start with a set of candidate models, and then find the models corresponding AIC values. There will almost always be information lost due to using one of the candidate models to represent the "true" model. We wish to select, from among *n* candidate models, the model that minimizes the estimated **information loss**. Denote the AIC values of the candidate models by AIC_1 , AIC_2 , AIC_3 , ..., AIC_n . Let AIC_{\min} be the minimum of those values. Then $e^{(AIC_{\min}-AIC_i)/2}$ can be interpreted as the relative probability that the i-th model minimizes the (estimated) information loss.

```
Fit an ARIMA model to a univariate time series.
Usage
arima(x. order = c(0. 0. 0).
     seasonal = list(order = c(0, 0, 0), period = NA),
     optim.control = list(), kappa = 1e6)
 # Example: R code for selection of the "best" model
 # Aikake's information criterion, select among 10 candidate models
 a1=arima(vp,order=c(0,0,0))
 a2=arima(vp,order=c(0,0,1))
 a3=arima(vp,order=c(0,1,0))
 a4=arima(vp,order=c(1,0,0))
 a5=arima(vp,order=c(1,1,0))
 a6=arima(vp,order=c(0,1,1))
 a7=arima(vp,order=c(1,0,1))
 a8=arima(vp,order=c(1,1,1))
 a8=arima(vp,order=c(0,0,2))
 a9=arima(vp,order=c(0,2,0))
 a10=arima(vp,order=c(2,0,0))
 veca=as.vector(cbind(a1$aic,a2$aic,a3$aic,a4$aic,
              a5$aic,a6$aic,a7$aic,a8$aic,a9$aic,a10$aic))
```

4 D > 4 B > 4 B > 4 B > 9 Q P

which.min(veca)

```
Model Outputs:
> which.min(veca)
[1] 4
Check the best:
> a4
Call:
arima(x = vp, order = c(1, 0, 0))
Coefficients:
         ar1 intercept
      0.3941 1.0042
s.e. 0.0290 0.0005
sigma<sup>2</sup> estimated as 0.0001101: log likelihood = 3140.99,
aic = -6275.98
!!! intercept indicates fundamental price p_fun=1
```

```
### Check another one = not the best
> a3
Call:
arima(x = vp, order = c(0, 1, 0))
sigma^2 estimated as 0.0001581:
log likelihood = 2957.18,
aic = -5912.36
```

Forecast by Arima

```
predict.Arima {stats} R Documentation
Forecast from ARIMA fits
Description: Forecast from models fitted by arima.
predict(object, n.ahead = 1, newxreg = NULL,
        se.fit = TRUE, ...)
Arguments:
object The result of an arima fit.
n.ahead
         The number of steps ahead for
          which prediction is required.
```

```
Make the prediction:
> a4=arima(x=vp, order=c(1,0,0))
> predict(a4,n.ahead=6)
\$pred
Time Series:
Start = 1002
End = 1007
Frequency = 1
[1] 1.005989 1.004901 1.004472 1.004303 1.004236
1.004210
\$se
Time Series:
Start = 1002
End = 1007
Frequency = 1
[1] 0.01049478 0.01128027 0.01139739 0.01141547
0.01141828 0.01141871
```

```
Testing of forecast
  vp1=vp[1:990]
> a4=arima(vp1,order=c(1,0,0))
> vp2=vp[991:1000]
> pre=predict(a4,n.ahead=10)
Time Series:
Start = 991
End = 1000
Frequency = 1
 [1] 1.002855 1.003630 1.003936 1.004057 1.004104 1.004123 1.004131 1.004133
 [9] 1.004135 1.004135
\$se
Time Series:
Start = 991
End = 1000
Frequency = 1
 [1] 0.01048887 0.01127568 0.01139327 0.01141146 0.01141429 0.01141473
 [7] 0.01141480 0.01141481 0.01141481 0.01141481
> vp2
 [1] 1.0091107 0.9884819 0.9970753 1.0119451 0.9983571 1.0193618 1.0258751
 [8] 1.0078596 1.0182782 1.0172228
```

vpred=pre\\$pred[1:10]

> vp2-vpred

- [1] 0.006256168 -0.015148120 -0.006860733 0.007888411 -0.005747281
- [6] 0.015238707 0.021744591 0.003726142 0.014143558 0.013087774

compare errors with a mean standard deviation

(vp2-vpred)/mean(se1)

- [1] 0.5533614 -1.3398593 -0.6068355 0.6977342 -0.5083501 1.3478718
- [7] 1.9233207 0.3295793 1.2510053 1.1576206