



# Time Series Analysis

Henrik Madsen

[hm@imm.dtu.dk](mailto:hm@imm.dtu.dk)

Informatics and Mathematical Modelling  
Technical University of Denmark  
DK-2800 Kgs. Lyngby

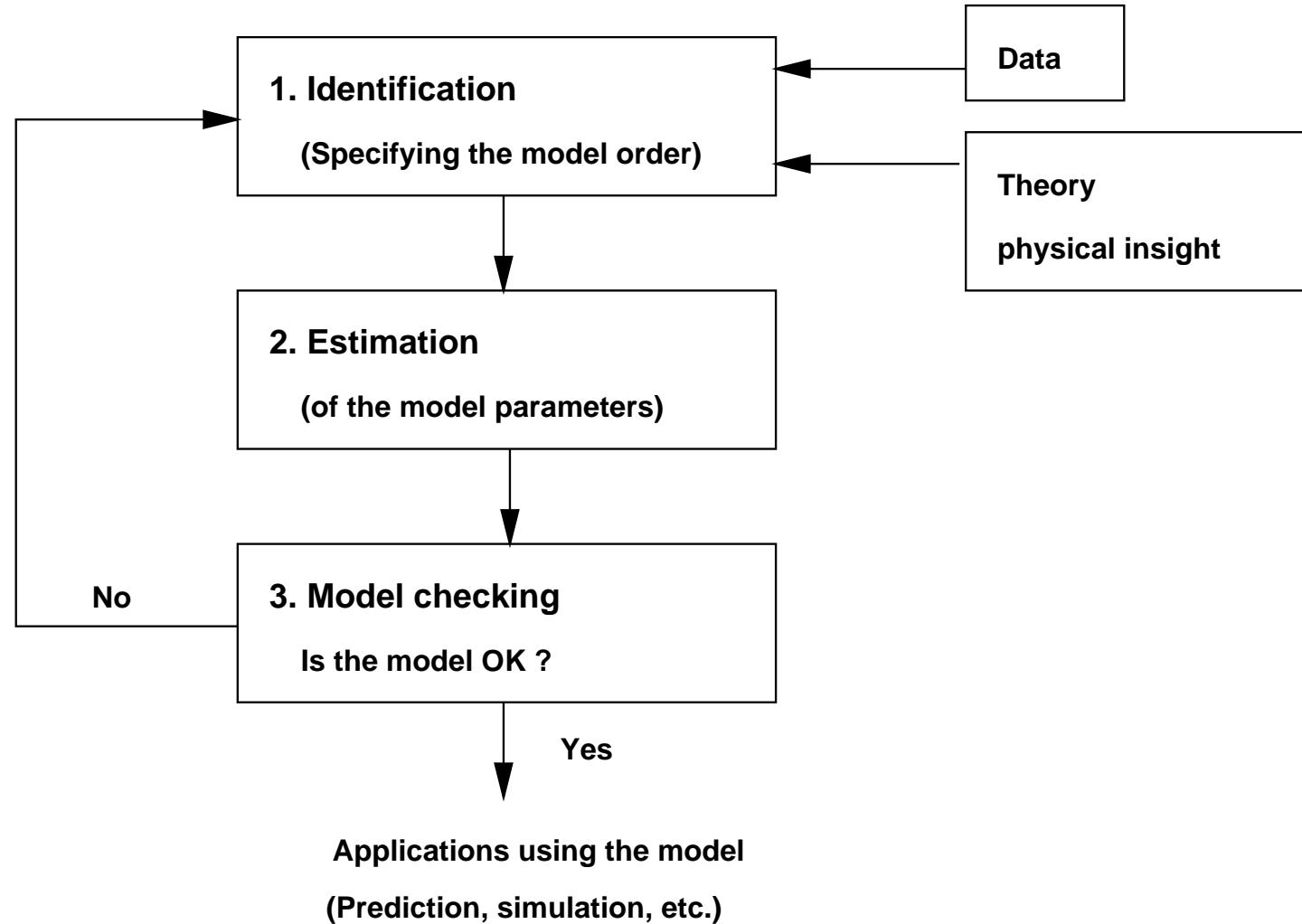


# Outline of the lecture

## Identification of univariate time series models

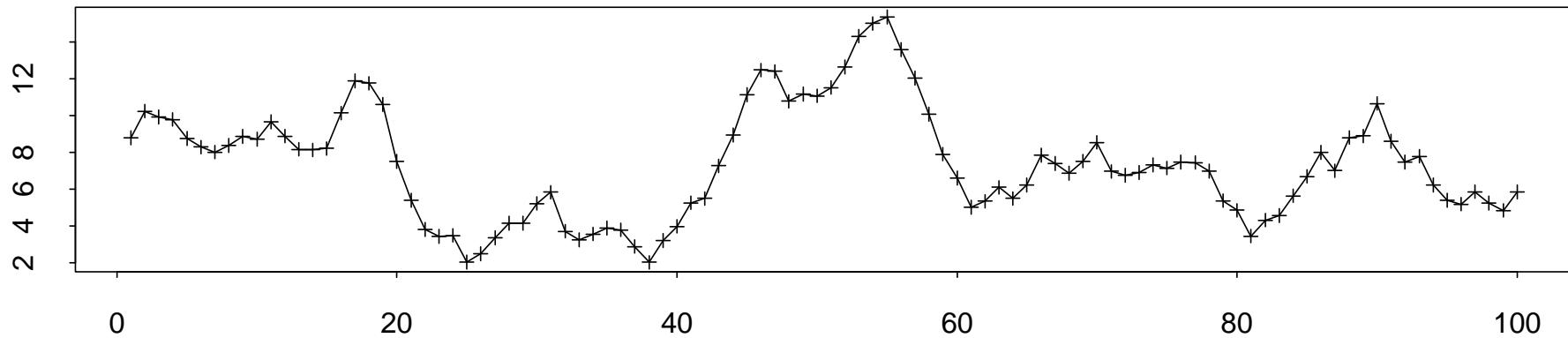
- Introduction, Sec. 6.1
- Estimation of auto-covariance and -correlation, Sec. 6.2.1  
(and the intro. to 6.2)
- Using SACF, SPACF, and SIACF for suggesting model structure, Sec. 6.3
- Estimation of model parameters, Sec. 6.4
- Examples...
- Cursory material:
  - ▶ The extended linear model class in Sec. 6.4.2  
(we'll come back to the extended model class later)

# Model building in general



# Identification of univariate time series models

- What ARIMA structure would be appropriate for the data at hand? (If any)



- Given the structure we will then consider how to estimate the parameters (next lecture)
- What do we know about ARIMA models which could help us?



## Estimation of the autocovariance function

- Estimate of  $\gamma(k)$

$$C_{YY}(k) = C(k) = \hat{\gamma}(k) = \frac{1}{N} \sum_{t=1}^{N-|k|} (Y_t - \bar{Y})(Y_{t+|k|} - \bar{Y})$$

- It is enough to consider  $k > 0$
- S-PLUS: `acf(x, type = "covariance")`



## Some properties of $C(k)$

- The estimate is a non-negative definite function (as  $\gamma(k)$ )
- The estimator is *non-central*:

$$E[C(k)] = \frac{1}{N} \sum_{t=1}^{N-|k|} \gamma(k) = \left(1 - \frac{|k|}{N}\right) \gamma(k)$$

- Asymptotically central (consistent) for fixed  $k$ :  
 $E[C(k)] \rightarrow \gamma(k)$  for  $N \rightarrow \infty$
- The estimates are autocorrelated them self (don't trust apparent correlation at high lags too much)



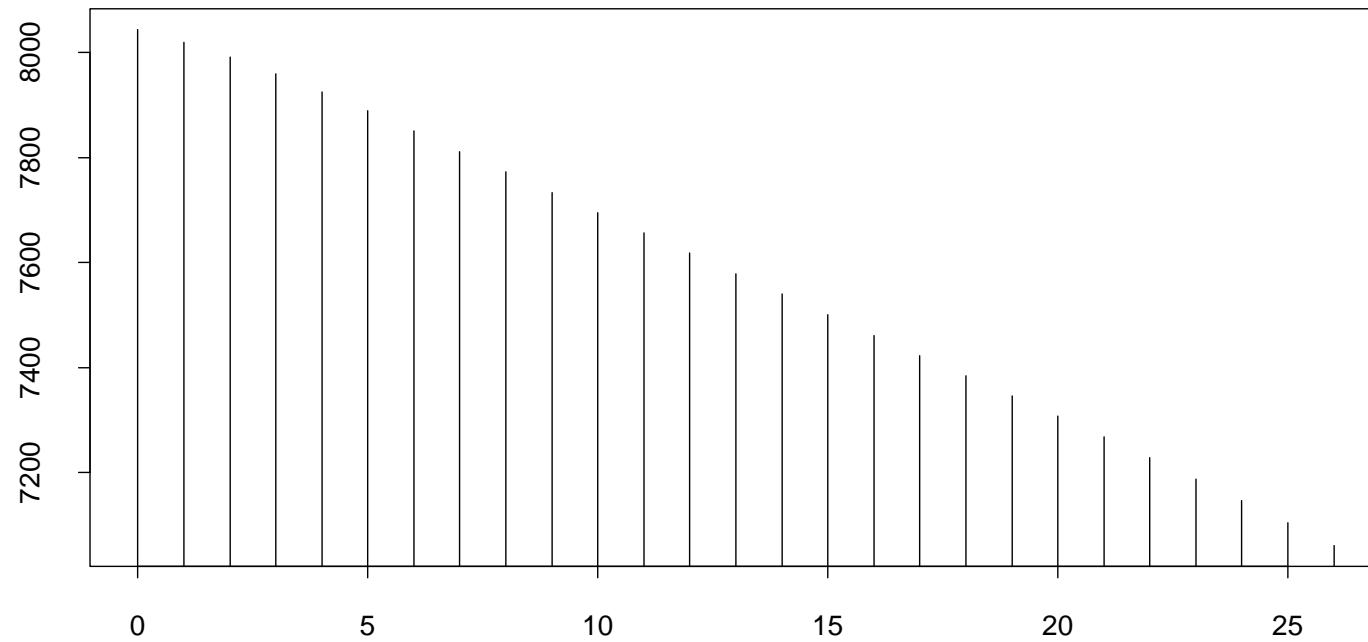
# How does $C(k)$ behave for non-stationary series?

$$C(k) = \frac{1}{N} \sum_{t=1}^{N-|k|} (Y_t - \bar{Y})(Y_{t+|k|} - \bar{Y})$$

# How does $C(k)$ behave for non-stationary series?

$$C(k) = \frac{1}{N} \sum_{t=1}^{N-|k|} (Y_t - \bar{Y})(Y_{t+|k|} - \bar{Y})$$

Series : arima.sim(model = list(ar = 0.9, ndiff = 1), n = 500)

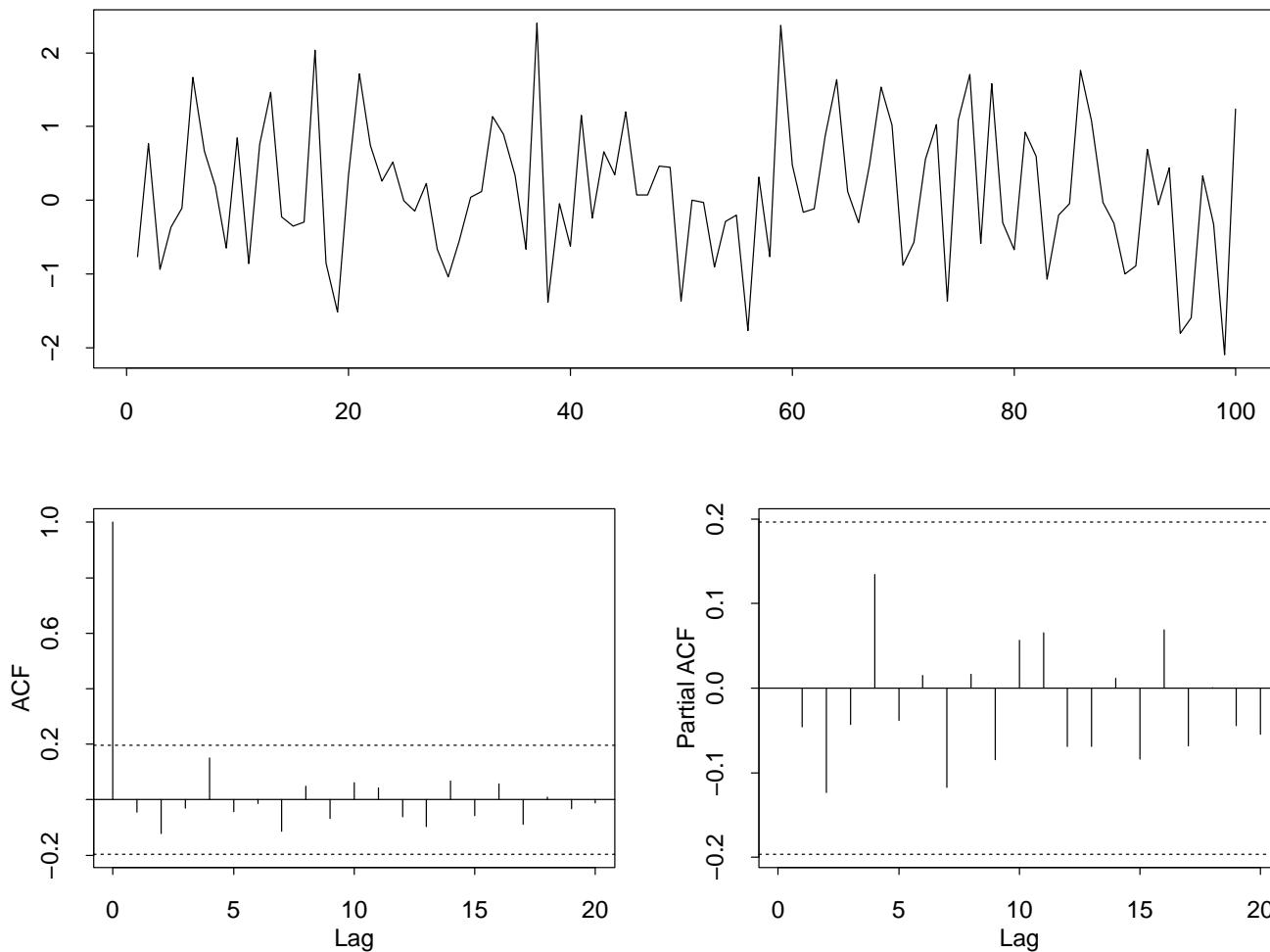




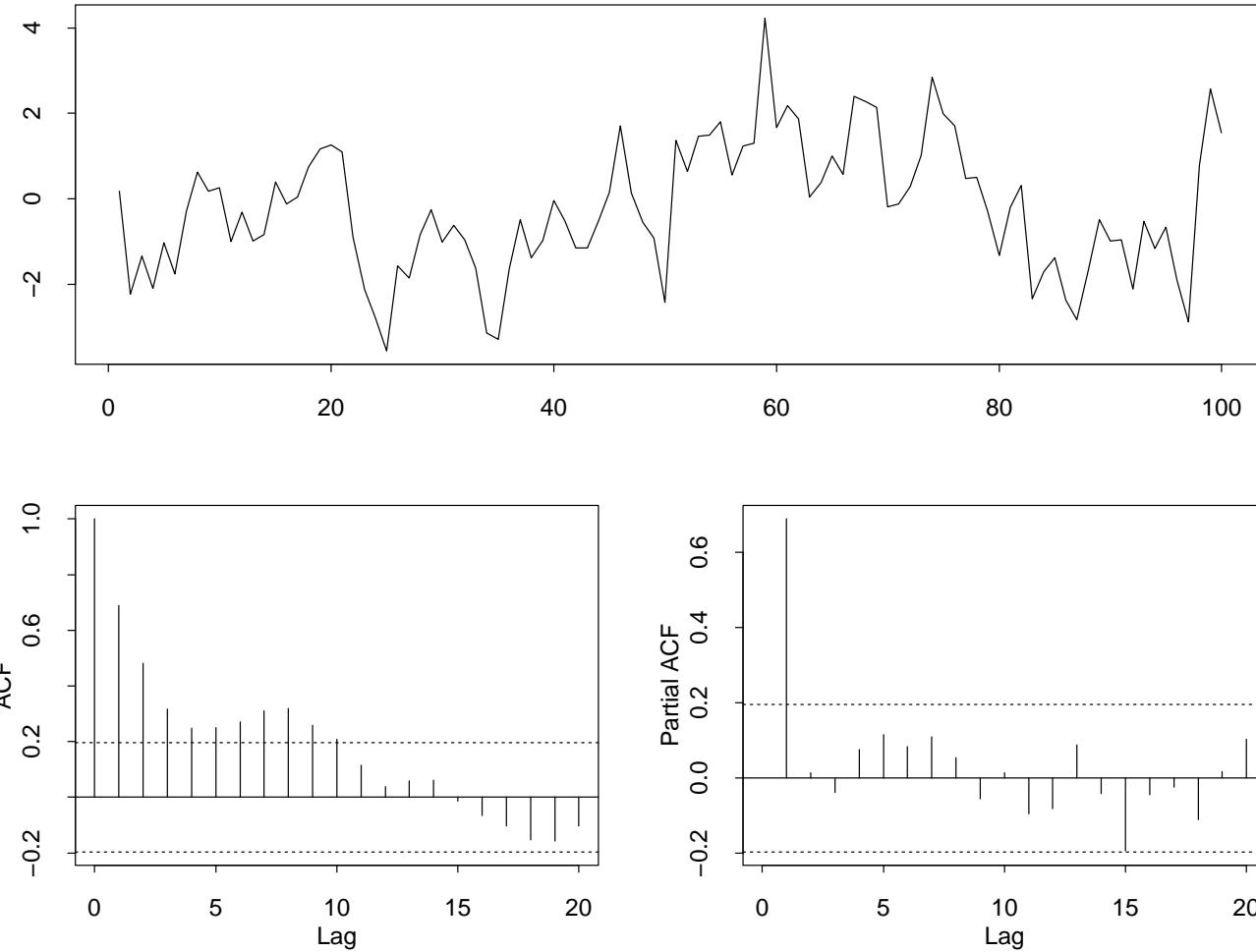
# Autocorrelation and Partial Autocorrelation

- Sample autocorrelation function (SACF):  
 $\hat{\rho}(k) = r_k = C(k)/C(0)$
  - For white noise and  $k \neq 1$  it holds that  $E[\hat{\rho}(k)] \simeq 0$  and  $V[\hat{\rho}(k)] \simeq 1/N$ , this gives the bounds  $\pm 2/\sqrt{N}$  for deciding when it is not possible to distinguish a value from zero.
  - S-PLUS: `acf(x)`
- 
- Sample partial autocorrelation function (SPACF): Use the Yule-Walker equations on  $\hat{\rho}(k)$  (exactly as for the theoretical relations)
  - It turns out that  $\pm 2/\sqrt{N}$  is also appropriate for deciding when the SPACF is zero (more in the next lecture)
  - S-PLUS: `acf(x, type="partial")`

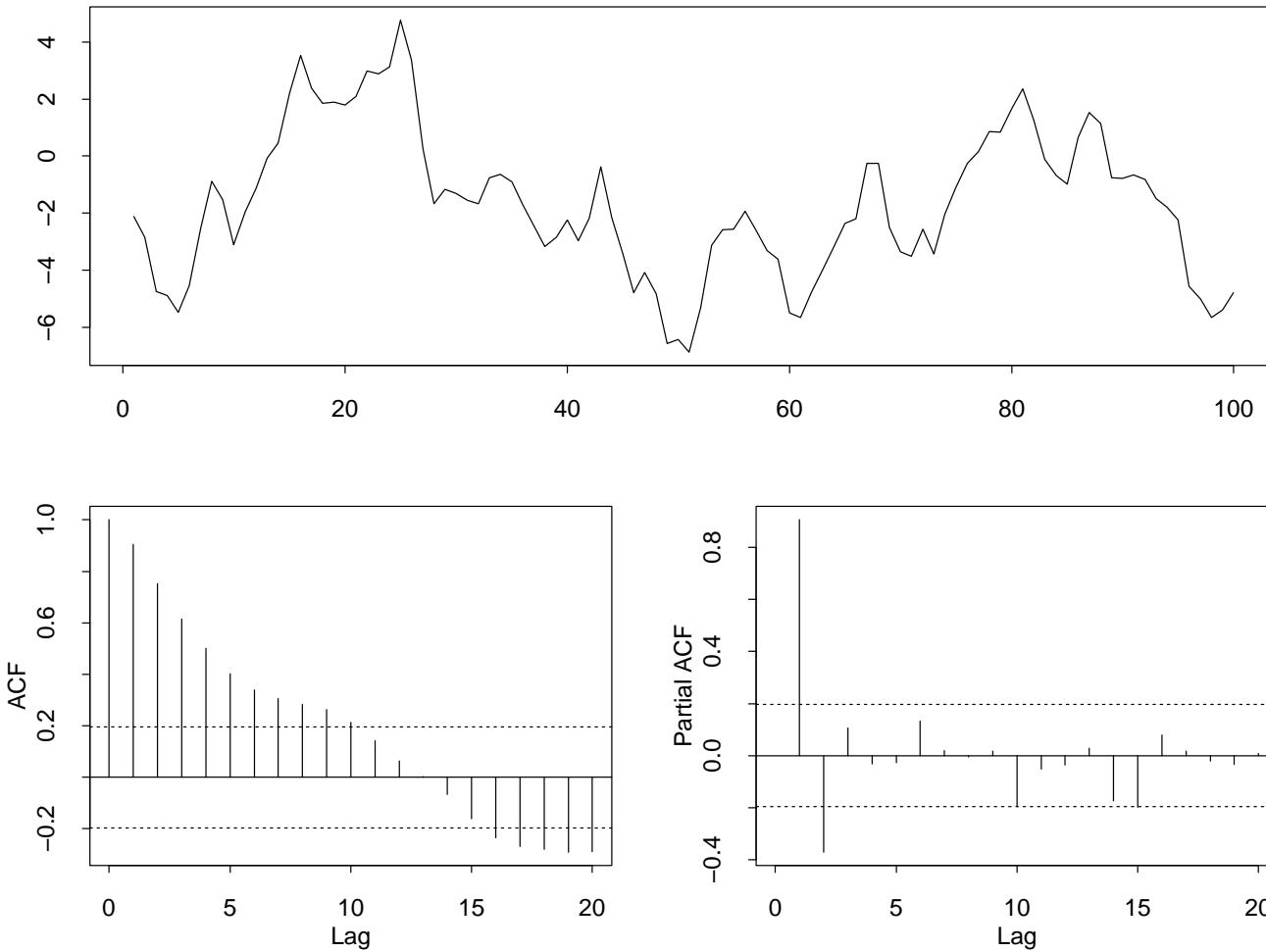
# What would be an appropriate structure?



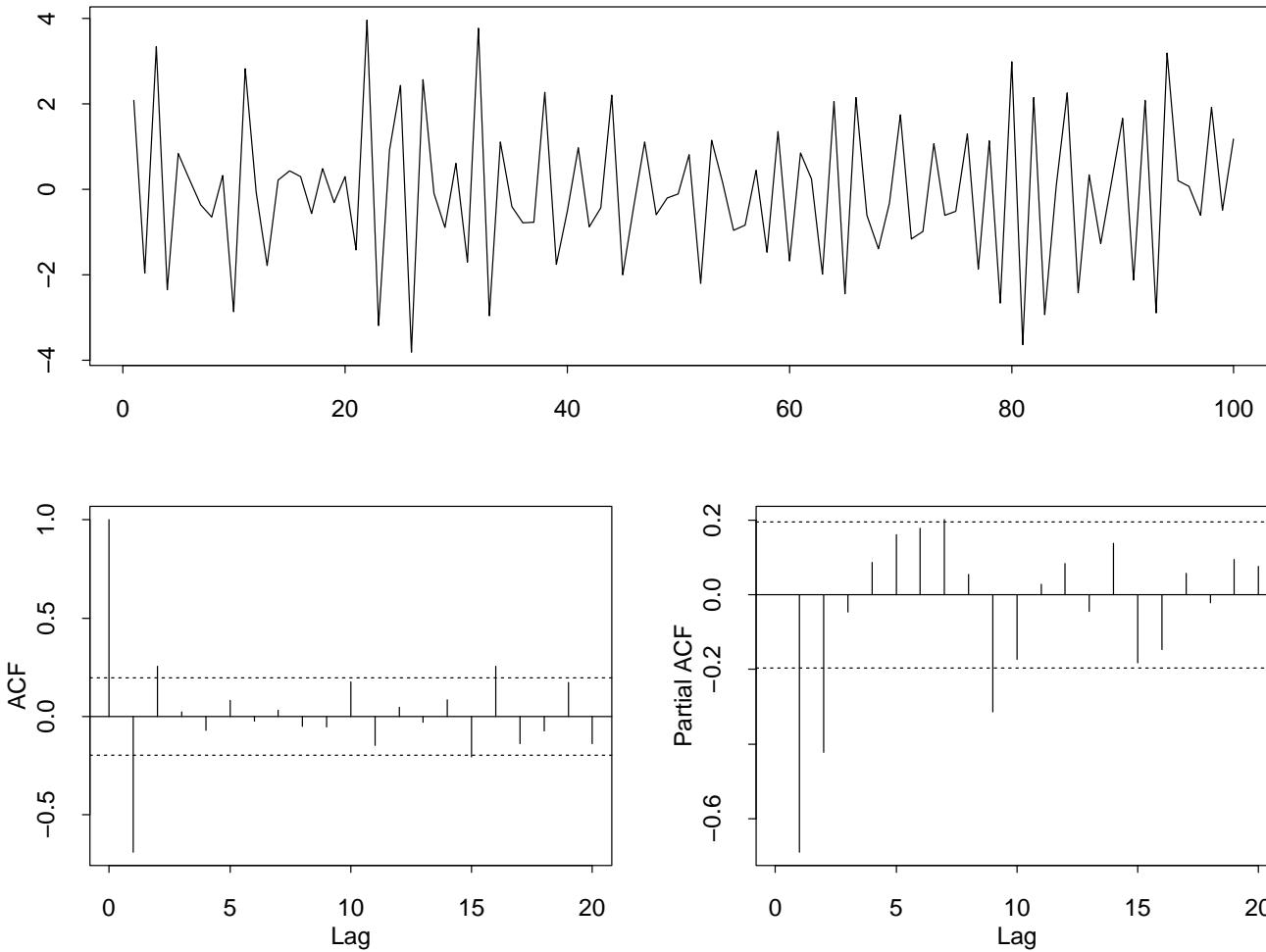
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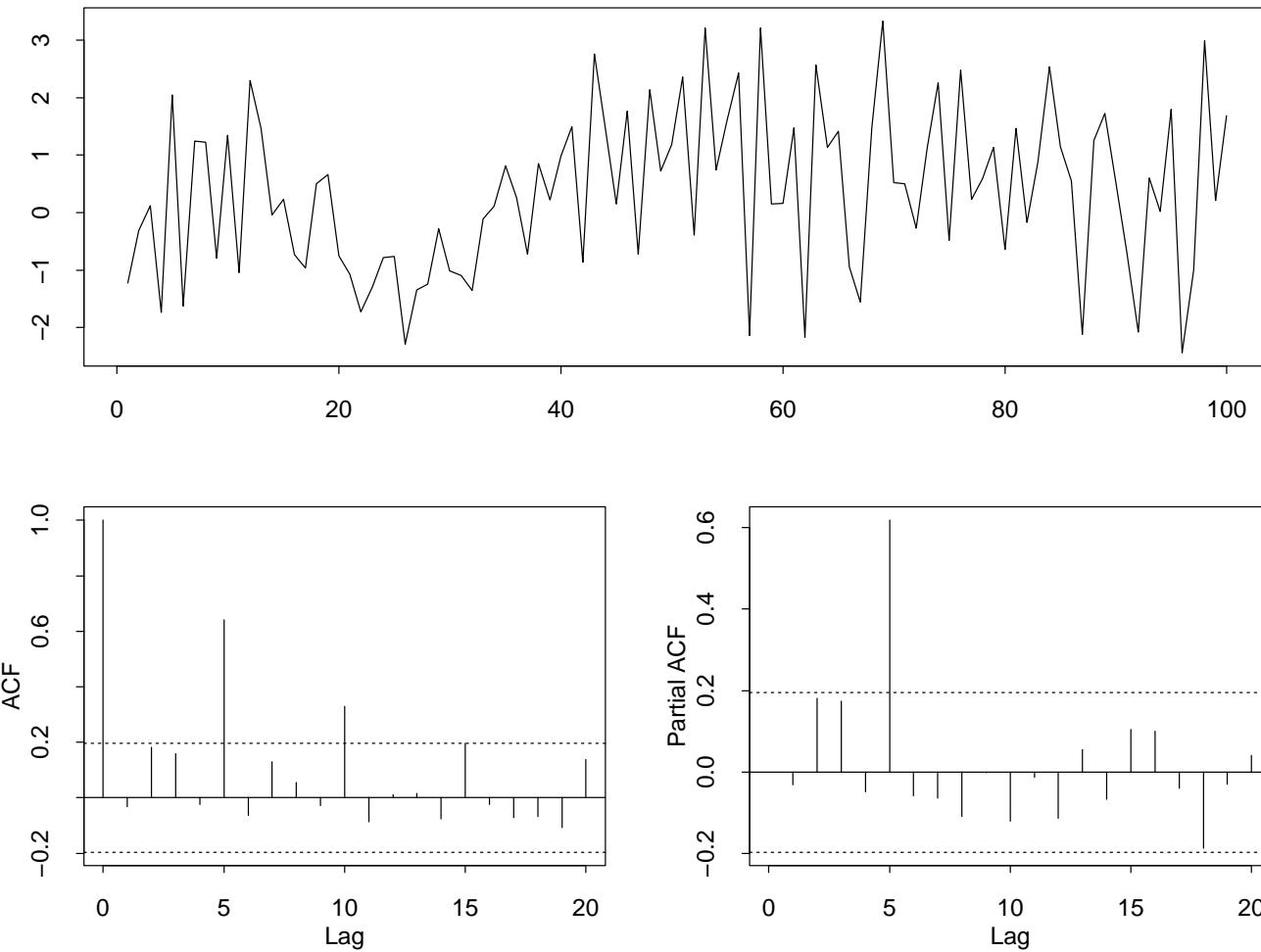
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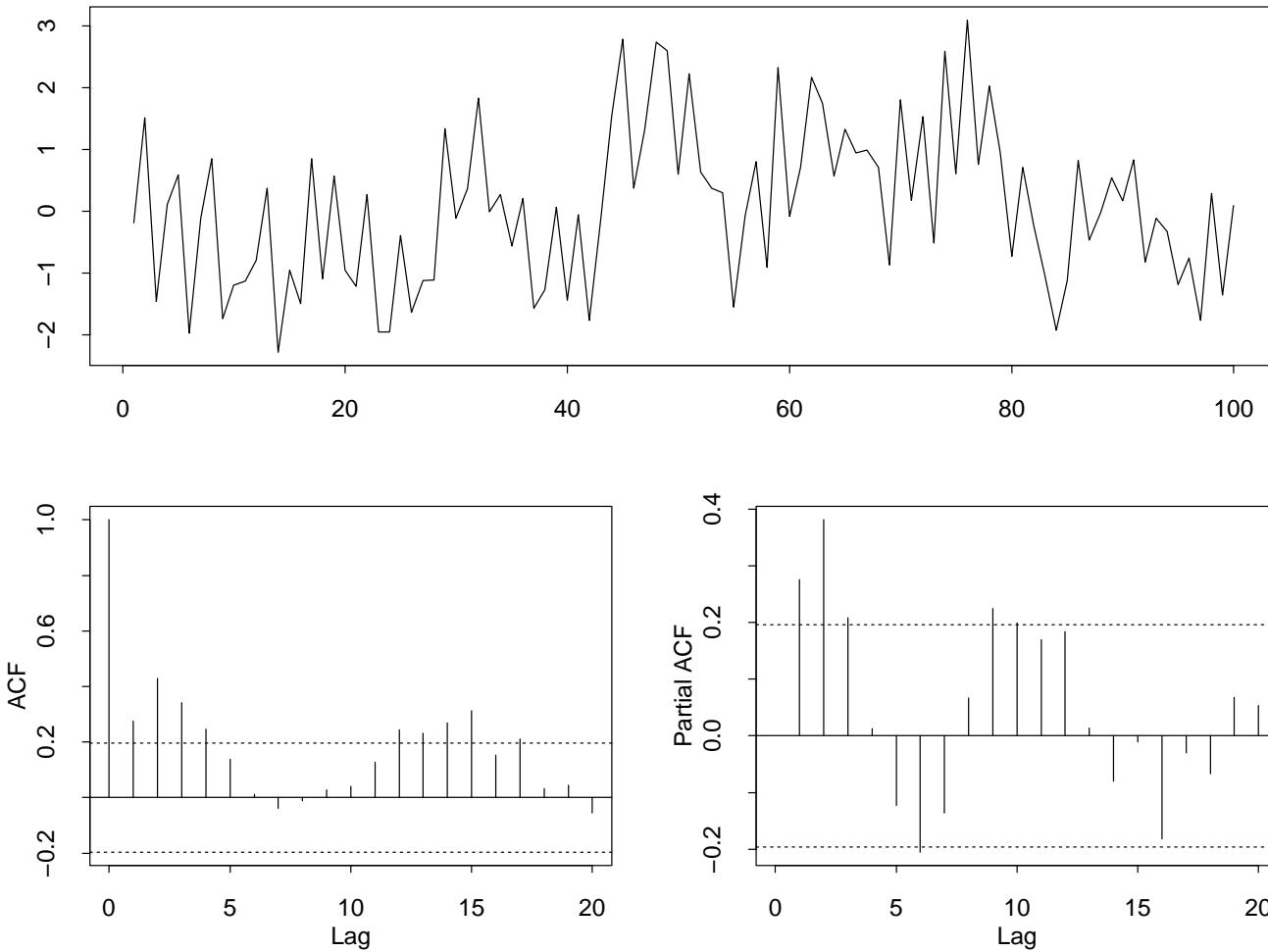
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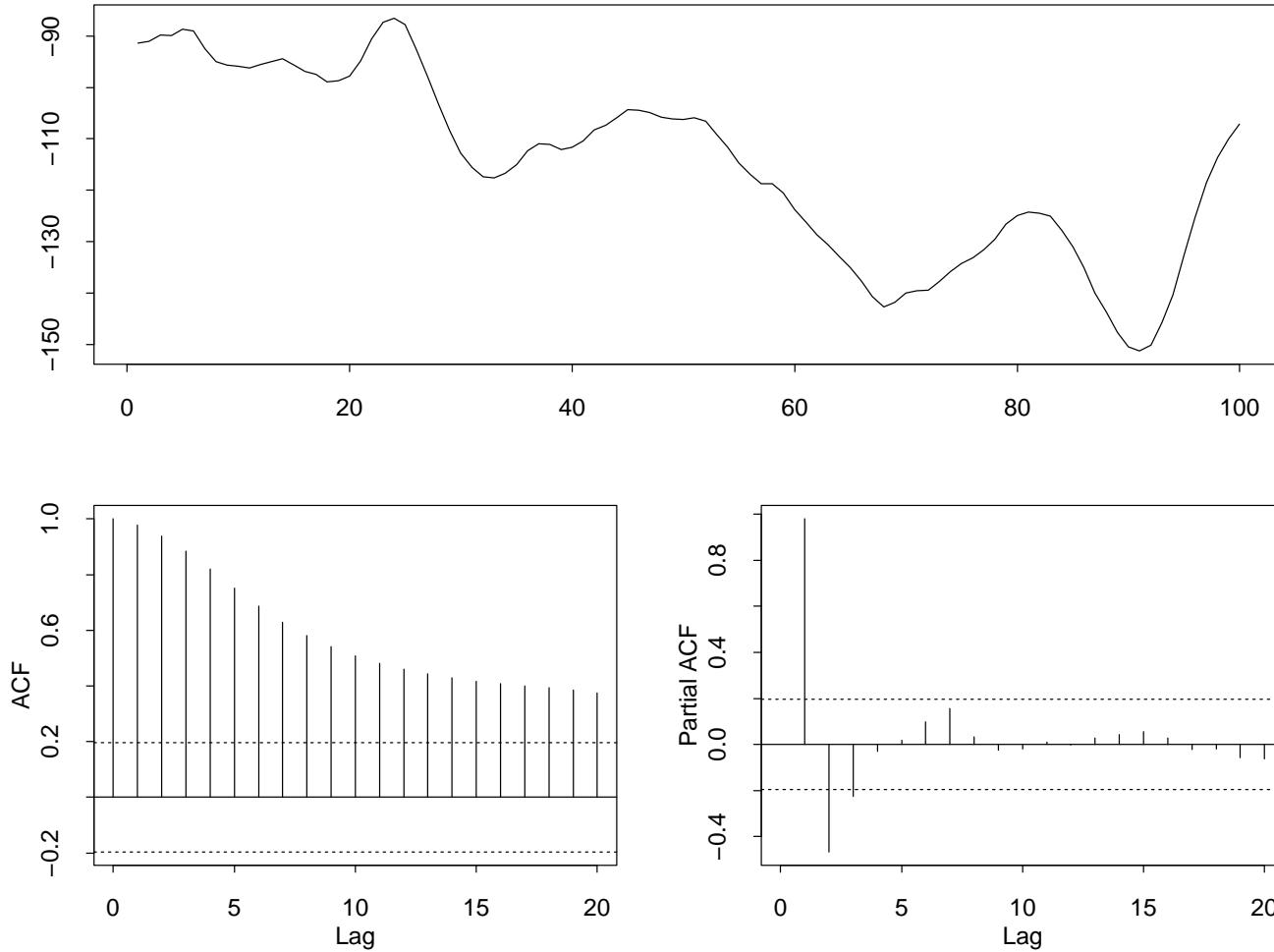
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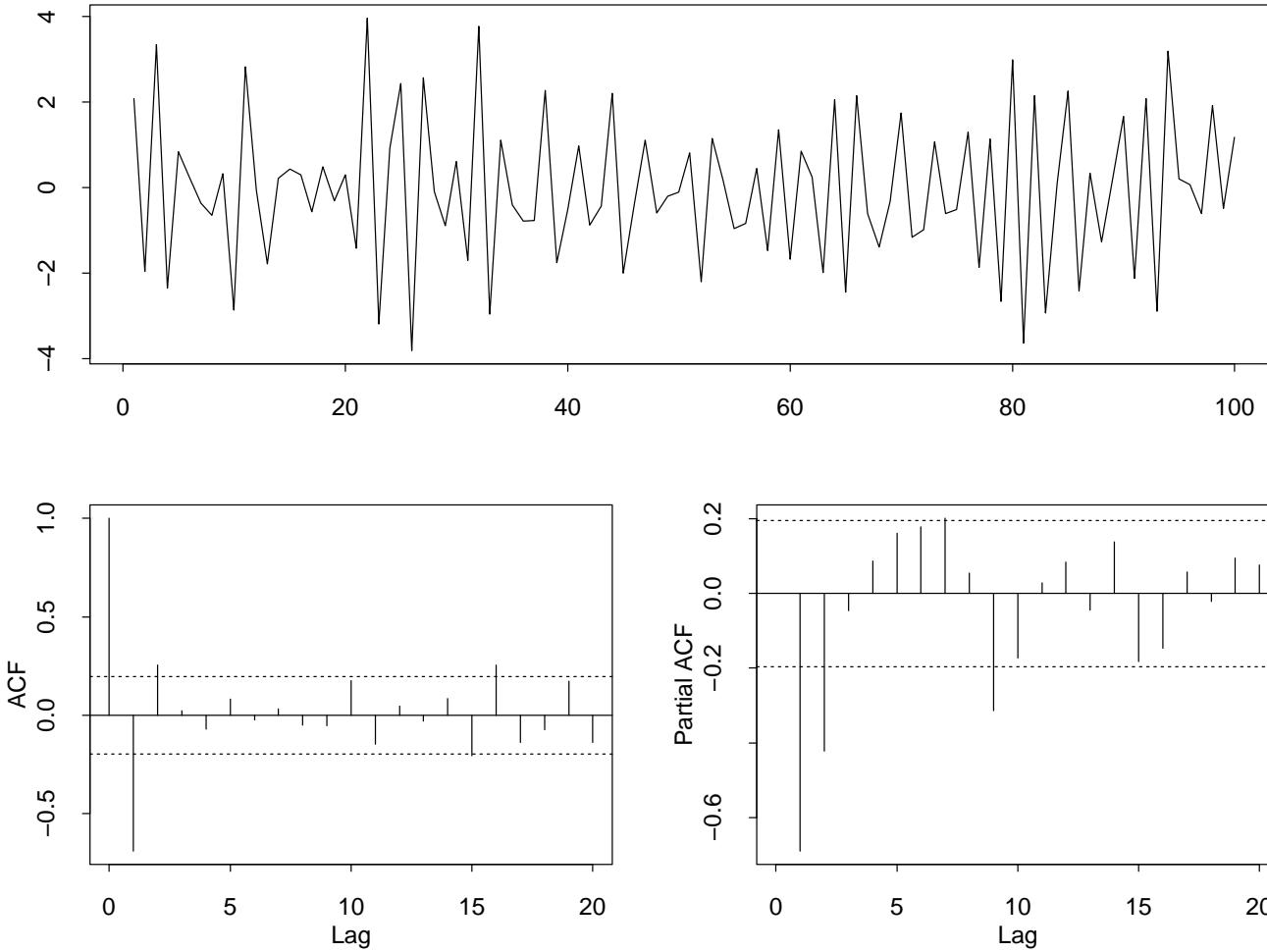
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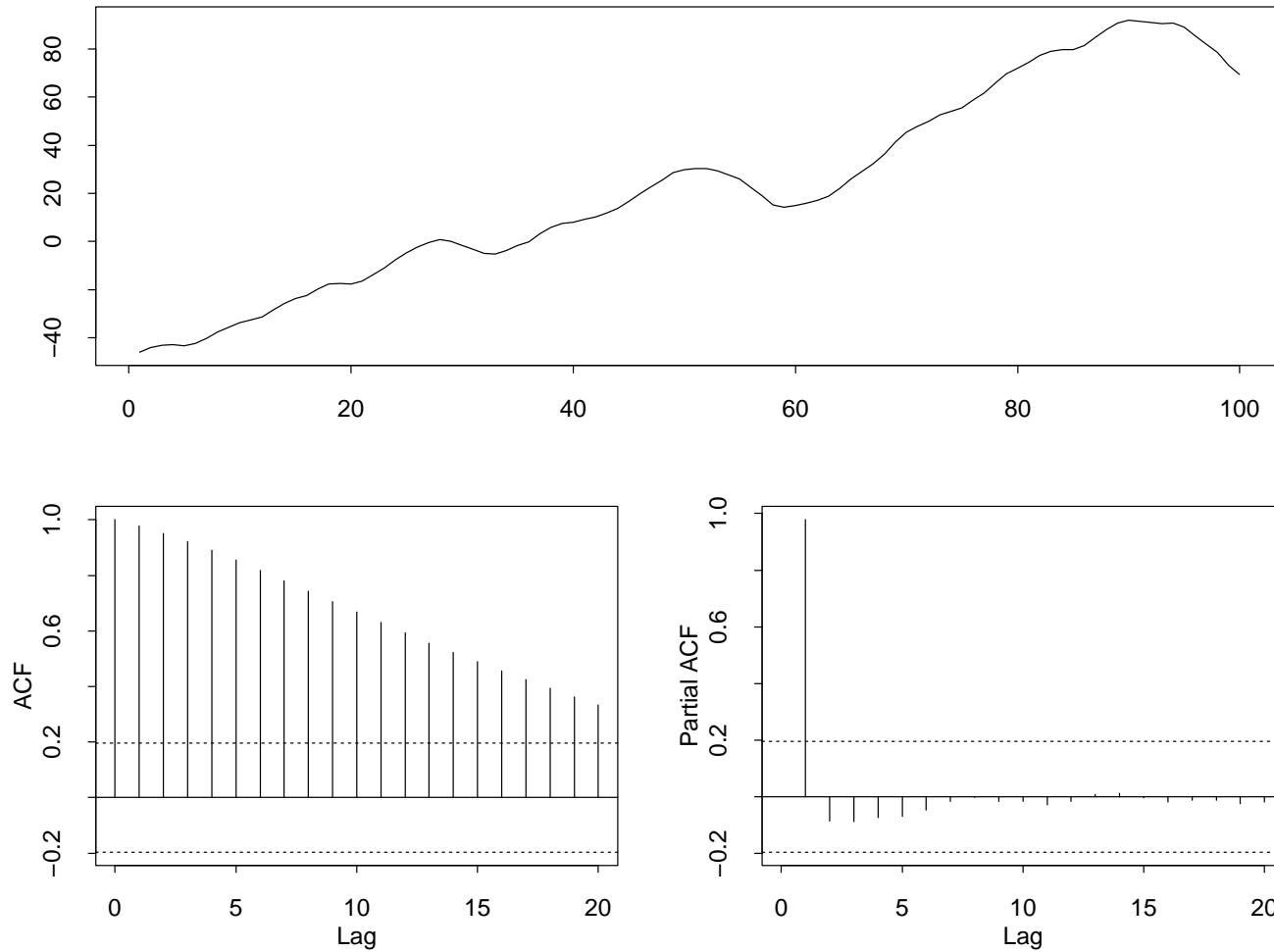
# What would be an appropriate structure?



## Example of data from an $MA(2)$ -process

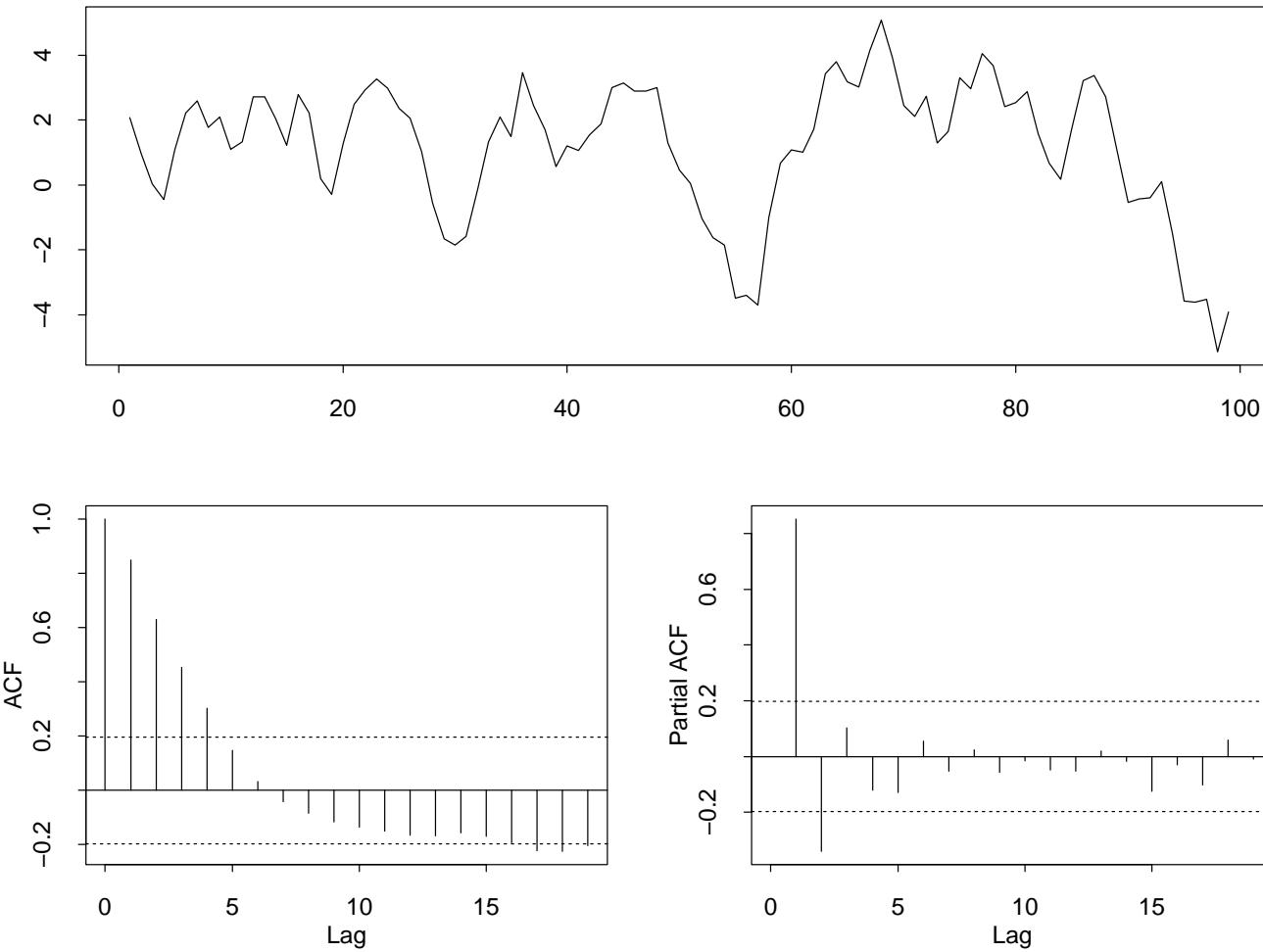


## Example of data from a non-stationary process





## Same series; analysing $\nabla Y_t = (1 - B)Y_t = Y_t - Y_{t-1}$





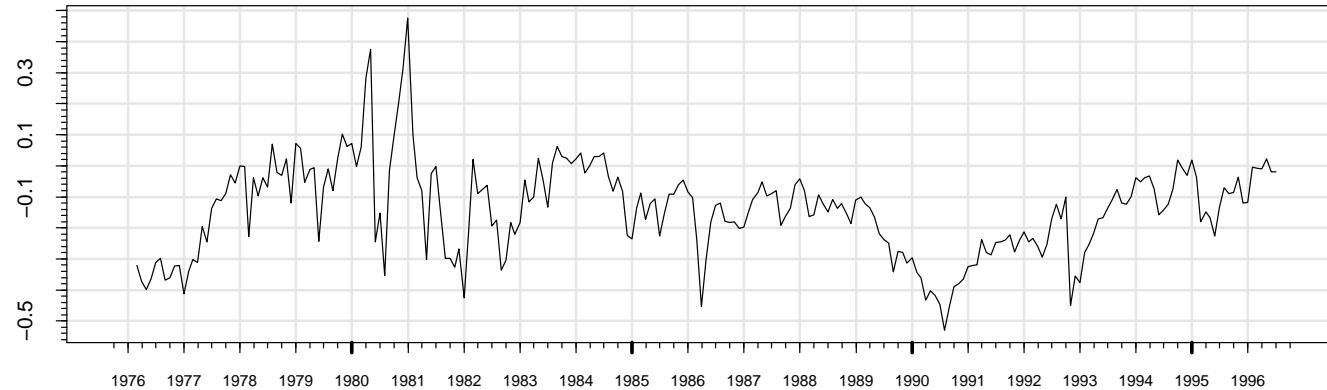
## Identification of the order of differencing

- Select the order of differencing  $d$  as the first order for which the autocorrelation decreases sufficiently fast towards 0
- In practice  $d$  is 0, 1, or maybe 2
- Sometimes a periodic difference is required, e.g.  $Y_t - Y_{t-12}$
- Remember to consider the practical application . . . it may be that the system is stationary, but you measured over a too short period

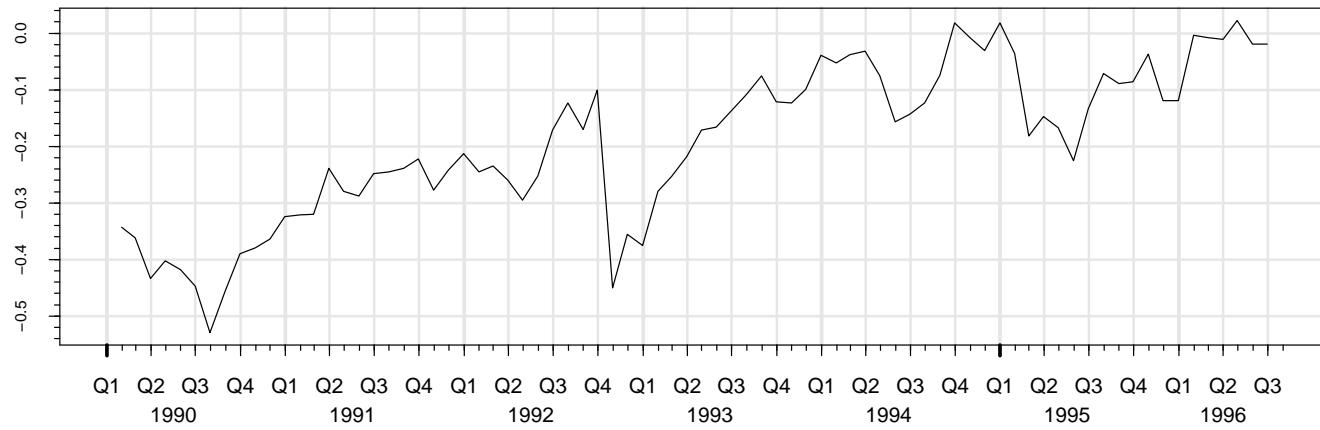


# Stationarity vs. length of measuring period

US/CA 30 day interest rate differential



US/CA 30 day interest rate differential





## Identification of the ARMA-part

Characteristics for the autocorrelation functions:

	ACF $\rho(k)$	PACF $\phi_{kk}$
$AR(p)$	Damped exponential and/or sine functions	$\phi_{kk} = 0$ for $k > p$
$MA(q)$	$\rho(k) = 0$ for $k > q$	Dominated by damped exponential and or/sine functions
$ARMA(p, q)$	Damped exponential and/or sine functions after lag $q - p$	Dominated by damped exponential and/or sine functions after lag $p - q$

The IACF is similar to the PACF; see the book page 133



## Behaviour of the SACF $\hat{\rho}(k)$ (based on $N$ obs.)

- If the process is white noise then

$$\pm 2\sqrt{\frac{1}{N}}$$

is an approximate 95% confidence interval for the SACF for lags different from 0

- If the process is a  $MA(q)$ -process then

$$\pm 2\sqrt{\frac{1 + 2(\hat{\rho}^2(1) + \dots + \hat{\rho}^2(q))}{N}}$$

is an approximate 95% confidence interval for the SACF for lags larger than  $q$



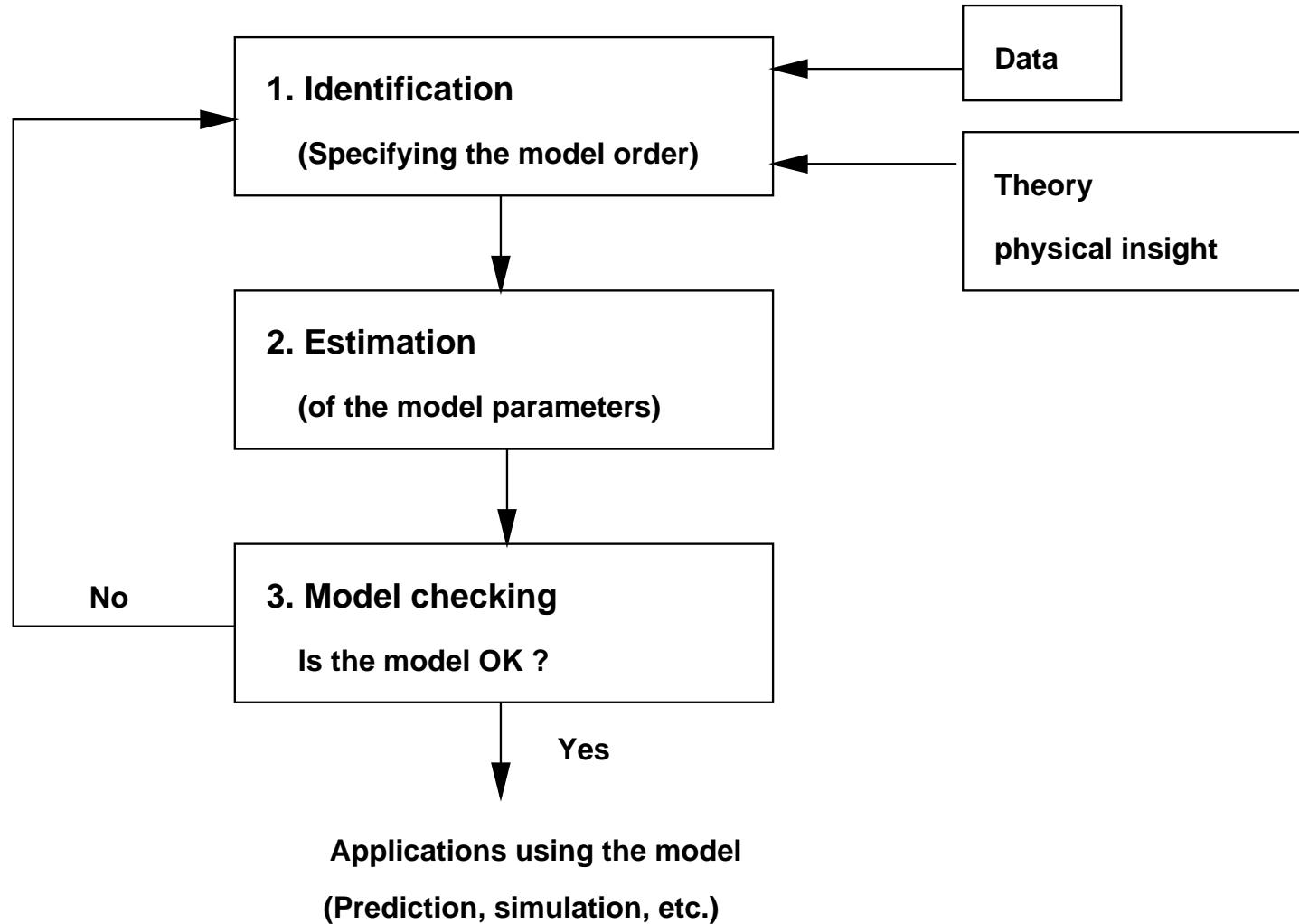
## Behaviour of the SPACF $\hat{\phi}_{kk}$ (based on $N$ obs.)

- If the process is a  $AR(p)$ -process then

$$\pm 2\sqrt{\frac{1}{N}}$$

is an approximate 95% confidence interval for the SPACF for lags larger than  $p$

# Model building in general

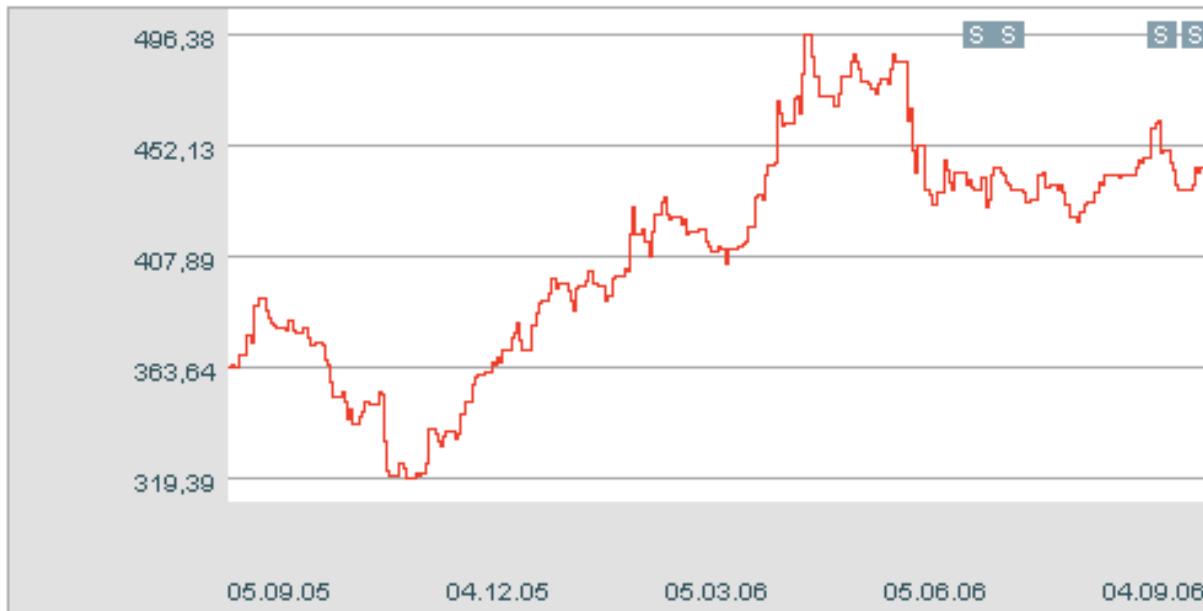




# Estimation

- We have an appropriate model structure  $AR(p)$ ,  $MA(q)$ ,  $ARMA(p, q)$ ,  $ARIMA(p, d, q)$  with  $p$ ,  $d$ , and  $q$  known
- **Task:** Based on the observations find appropriate values of the parameters
- The book describes many methods:
  - ▶ Moment estimates
  - ▶ LS-estimates
  - ▶ Prediction error estimates
    - Conditioned
    - Unconditioned
  - ▶ ML-estimates
    - Conditioned
    - Unconditioned (exact)

## Example



Using the autocorrelation functions we agreed that

$$\hat{y}_{t+1|t} = a_1 y_t + a_2 y_{t-1}$$

and we would select  $a_1$  and  $a_2$  so that the sum of the squared prediction errors got so small as possible when using the model on the data at hand

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$$\hat{y}_{t+1|t} = a_1 y_t + a_2 y_{t-1}$$

and we would select  $a_1$  and  $a_2$  so that the sum of the squared prediction errors got so small as possible when using the model on the data at hand

To comply with the notation of the book we will write the 1-step forecasts as  $\hat{y}_{t+1|t} = -\phi_1 y_t - \phi_2 y_{t-1}$



## The errors given the parameters ( $\phi_1$ and $\phi_2$ )

- Observations:  $y_1, y_2, \dots, y_N$
- Errors:  $e_{t+1|t} = y_{t+1} - \hat{y}_{t+1|t} = y_{t+1} - (-\phi_1 y_t - \phi_2 y_{t-1})$



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$$e_{3|2} = y_3 + \phi_1 y_2 + \phi_2 y_1$$

$$e_{4|3} = y_4 + \phi_1 y_3 + \phi_2 y_2$$

$$e_{5|4} = y_5 + \phi_1 y_4 + \phi_2 y_3$$

$$\vdots$$

$$e_{N|N-1} = y_N + \phi_1 y_{N-1} + \phi_2 y_{N-2}$$



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$$\vdots$$

$$e_{N|N-1} = y_N + \phi_1 y_{N-1} + \phi_2 y_{N-2}$$

$$\begin{bmatrix} y_3 \\ \vdots \\ y_N \end{bmatrix} = \begin{bmatrix} -y_2 & -y_1 \\ \vdots & \vdots \\ -y_{N-1} & -y_{N-2} \end{bmatrix} \begin{bmatrix} \phi_1 \\ \phi_2 \end{bmatrix} + \begin{bmatrix} e_{3|2} \\ \vdots \\ e_{N|N-1} \end{bmatrix}$$



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⋮

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Or just:

$$\mathbf{Y} = \mathbf{X}\boldsymbol{\theta} + \boldsymbol{\varepsilon}$$



## Solution

To minimize the sum of the squared 1-step prediction errors  $\varepsilon^T \varepsilon$  we use the result for the General Linear Model from Chapter 3:

$$\hat{\boldsymbol{\theta}} = (\mathbf{X}^T \mathbf{X})^{-1} \mathbf{X}^T \mathbf{Y}$$

With

$$\mathbf{X} = \begin{bmatrix} -y_2 & -y_1 \\ \vdots & \vdots \\ -y_{N-1} & -y_{N-2} \end{bmatrix} \quad \text{and} \quad \mathbf{Y} = \begin{bmatrix} y_3 \\ \vdots \\ y_N \end{bmatrix}$$

- The method is called the LS-estimator for dynamical systems
- The method is also in the class of prediction error methods since it minimize the sum of the squared 1-step prediction errors



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- The method is called the LS-estimator for dynamical systems
- The method is also in the class of prediction error methods since it minimize the sum of the squared 1-step prediction errors
- **How does it generalize to AR(p)-models?**



## Small illustrative example using S-PLUS

```
> obs
[1] -3.51 -3.81 -1.85 -2.02 -1.91 -0.88
> N <- length(obs); Y <- obs[3:N]
> Y
[1] -1.85 -2.02 -1.91 -0.88
> X <- cbind(-obs[2:(N-1)], -obs[1:(N-2)])
> X
     [,1] [,2]
[1,] 3.81 3.51
[2,] 1.85 3.81
[3,] 2.02 1.85
[4,] 1.91 2.02
> solve(t(X) %*% X, t(X) %*% Y) # Estimates
     [,1]
[1,] -0.1474288
[2,] -0.4476040
```



# Maximum likelihood estimates

- $ARMA(p, q)$ -process:

$$Y_t + \phi_1 Y_{t-1} + \cdots + \phi_p Y_{t-p} = \varepsilon_t + \theta_1 \varepsilon_{t-1} + \cdots + \theta_q \varepsilon_{t-q}$$

- Notation:

$$\boldsymbol{\theta}^T = (\phi_1, \dots, \phi_p, \theta_1, \dots, \theta_q)$$

$$\mathbf{Y}_t^T = (Y_t, Y_{t-1}, \dots, Y_1)$$

- The Likelihood function is the joint probability distribution function for all observations for given values of  $\boldsymbol{\theta}$  and  $\sigma_\varepsilon^2$ :

$$L(\mathbf{Y}_N; \boldsymbol{\theta}, \sigma_\varepsilon^2) = f(\mathbf{Y}_N | \boldsymbol{\theta}, \sigma_\varepsilon^2)$$

- Given the observations  $\mathbf{Y}_N$  we estimate  $\boldsymbol{\theta}$  and  $\sigma_\varepsilon^2$  as the values for which the likelihood is maximized.



## The likelihood function for $ARMA(p, q)$ -models

- The random variable  $Y_N | \mathbf{Y}_{N-1}$  only contains  $\varepsilon_N$  as a random component
- $\varepsilon_N$  is a white noise process at time  $N$  and does therefore not depend on anything
- We therefore know that the random variables  $Y_N | \mathbf{Y}_{N-1}$  and  $\mathbf{Y}_{N-1}$  are independent, hence:

$$f(\mathbf{Y}_N | \boldsymbol{\theta}, \sigma_\varepsilon^2) = f(Y_N | \mathbf{Y}_{N-1}, \boldsymbol{\theta}, \sigma_\varepsilon^2) f(\mathbf{Y}_{N-1} | \boldsymbol{\theta}, \sigma_\varepsilon^2)$$

- Repeating these arguments:

$$L(\mathbf{Y}_N; \boldsymbol{\theta}, \sigma_\varepsilon^2) = \left( \prod_{t=p+1}^N f(Y_t | \mathbf{Y}_{t-1}, \boldsymbol{\theta}, \sigma_\varepsilon^2) \right) f(\mathbf{Y}_p | \boldsymbol{\theta}, \sigma_\varepsilon^2)$$



## The conditional likelihood function

- Evaluation of  $f(\mathbf{Y}_p | \boldsymbol{\theta}, \sigma_\varepsilon^2)$  requires special attention
- It turns out that the estimates obtained using the *conditional likelihood function*:

$$L(\mathbf{Y}_N; \boldsymbol{\theta}, \sigma_\varepsilon^2) = \prod_{t=p+1}^N f(Y_t | \mathbf{Y}_{t-1}, \boldsymbol{\theta}, \sigma_\varepsilon^2)$$

results in the same estimates as the *exact likelihood function* when many observations are available

- For small samples there can be some difference
- Software:
  - ▶ The S-PLUS function `arima.mle` calculate conditional estimates
  - ▶ The R function `arima` calculate exact estimates



## Evaluating the conditional likelihood function

- **Task:** Find the conditional densities given specified values of the parameters  $\theta$  and  $\sigma_\varepsilon^2$
- The mean of the random variable  $Y_t | \mathbf{Y}_{t-1}$  is the the 1-step forecast  $\hat{Y}_{t|t-1}$
- The prediction error  $\varepsilon_t = Y_t - \hat{Y}_{t|t-1}$  has variance  $\sigma_\varepsilon^2$
- We assume that the process is Gaussian:

$$f(Y_t | \mathbf{Y}_{t-1}, \theta, \sigma_\varepsilon^2) = \frac{1}{\sigma_\varepsilon \sqrt{2\pi}} e^{-(Y_t - \hat{Y}_{t|t-1}(\theta))^2 / 2\sigma_\varepsilon^2}$$

- And therefore:

$$L(\mathbf{Y}_N; \theta, \sigma_\varepsilon^2) = (\sigma_\varepsilon^2 2\pi)^{-\frac{N-p}{2}} \exp \left( -\frac{1}{2\sigma_\varepsilon^2} \sum_{t=p+1}^N \varepsilon_t^2(\theta) \right)$$



## ML-estimates

- The (conditional) ML-estimate  $\hat{\theta}$  is a prediction error estimate since it is obtained by minimizing

$$S(\boldsymbol{\theta}) = \sum_{t=p+1}^N \varepsilon_t^2(\boldsymbol{\theta})$$

- By differentiating w.r.t.  $\sigma_\varepsilon^2$  it can be shown that the ML-estimate of  $\sigma_\varepsilon^2$  is

$$\hat{\sigma}_\varepsilon^2 = S(\hat{\boldsymbol{\theta}})/(N - p)$$

- The estimate  $\hat{\theta}$  is asymptotically “good” and the variance-covariance matrix is approximately  $2\sigma_\varepsilon^2 \mathbf{H}^{-1}$  where  $\mathbf{H}$  contains the 2nd order partial derivatives of  $S(\boldsymbol{\theta})$  at the minimum



## Finding the ML-estimates using the PE-method

- 1-step predictions:

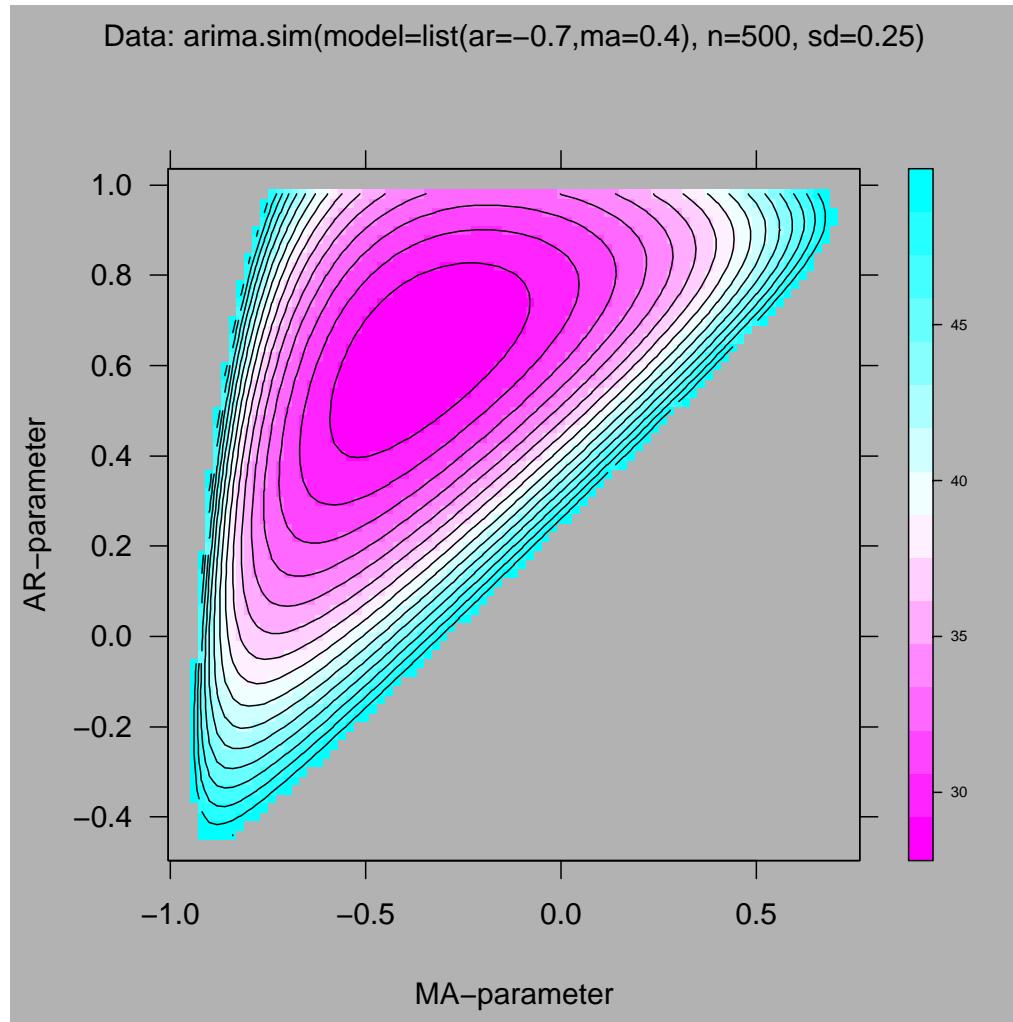
$$\hat{Y}_{t|t-1} = -\phi_1 Y_{t-1} - \cdots - \phi_p Y_{t-p} + \theta_1 \varepsilon_{t-1} + \cdots + \theta_q \varepsilon_{t-q}$$

- If we use  $\varepsilon_p = \varepsilon_{p-1} = \cdots = \varepsilon_{p+1-q} = 0$  we can find:

$$\hat{Y}_{p+1|p} = -\phi_1 Y_p - \cdots - \phi_p Y_1 + \theta_1 \varepsilon_p + \cdots + \theta_q \varepsilon_{p+1-q}$$

- Which will give us  $\varepsilon_{p+1} = Y_{p+1} - \hat{Y}_{p+1|p}$  and we can then calculate  $\hat{Y}_{p+2|p+1}$  and  $\varepsilon_{p+2}$  ... and so on until we have all the 1-step prediction errors we need.
- We use numerical optimization to find the parameters which minimize the sum of squared prediction errors

$S(\theta)$  for  $(1 + 0.7B)Y_t = (1 - 0.4B)\varepsilon_t$  with  $\sigma_\varepsilon^2 = 0.25^2$





## Moment estimates

- Given the model structure: Find formulas for the theoretical autocorrelation or autocovariance as function of the parameters in the model
- Estimate, e.g. calculate the SACF
- Solve the equations by using the lowest lags necessary
- **Complicated!**
- **General properties of the estimator unknown!**



## Moment estimates for $AR(p)$ -processes

In this case moment estimates are simple to find due to the Yule-Walker equations. We simply plug in the estimated autocorrelation function in lags 1 to  $p$ :

$$\begin{bmatrix} \hat{\rho}(1) \\ \hat{\rho}(2) \\ \vdots \\ \hat{\rho}(p) \end{bmatrix} = \begin{bmatrix} 1 & \hat{\rho}(1) & \cdots & \hat{\rho}(p-1) \\ \hat{\rho}(1) & 1 & \cdots & \hat{\rho}(p-2) \\ \vdots & \vdots & & \vdots \\ \hat{\rho}(p-1) & \hat{\rho}(p-2) & \cdots & 1 \end{bmatrix} \begin{bmatrix} -\phi_1 \\ -\phi_2 \\ \vdots \\ -\phi_p \end{bmatrix}$$

and solve w.r.t. the  $\phi$ 's

The function `ar` in S-PLUS does this