



Time Series Analysis

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Outline of the lecture

Identification of univariate time series models, cont.:

- Estimation of model parameters, Sec. 6.4 (cont.)
- Model order selection, Sec. 6.5
- Model validation, Sec. 6.6



Estimation – methods (from previous lecture)

- 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)



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 σ_ε^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 σ_ε^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 ε_N as a random component
- ε_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 (see also page 3):

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

- 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 θ and σ_ε^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 σ_ε^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. σ_ε^2 it can be shown that the ML-estimate of σ_ε^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:

$$\hat{Y}_{t|t-1} = -\phi_1 Y_{t-1} - \dots - \phi_p Y_{t-p} + 0 + \theta_1 \varepsilon_{t-1} + \dots + \theta_q \varepsilon_{t-q}$$

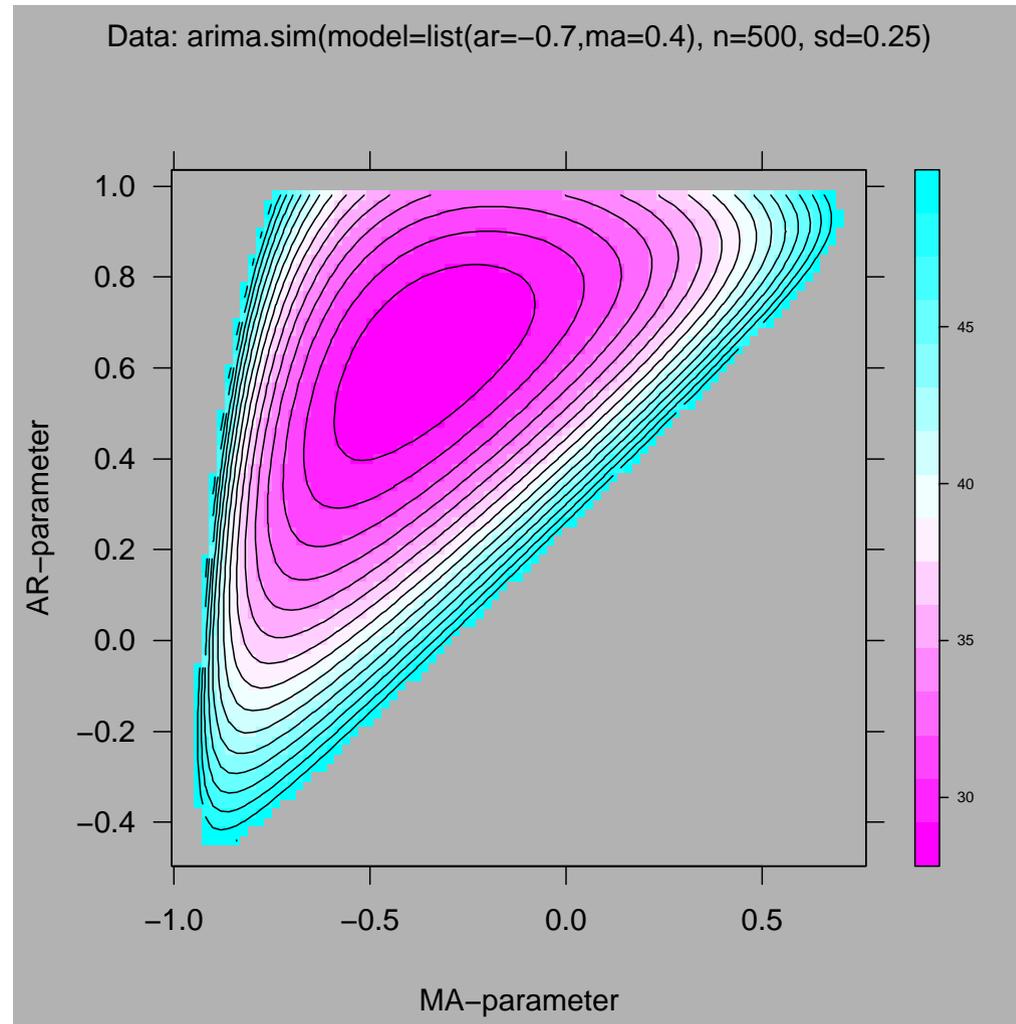
- If we use $\varepsilon_p = \varepsilon_{p-1} = \dots = \varepsilon_{p+1-q} = 0$ we can find:

$$\hat{Y}_{p+1|p} = -\phi_1 Y_p - \dots - \phi_p Y_1 + 0 + \theta_1 \varepsilon_p + \dots + \theta_q \varepsilon_{p+1-q}$$

- 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 $\varepsilon_{p+2} \dots$ 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



$S(\boldsymbol{\theta})$ 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 (page 104). We simply plug in the estimated autocorrelation function in lags 1 to p :

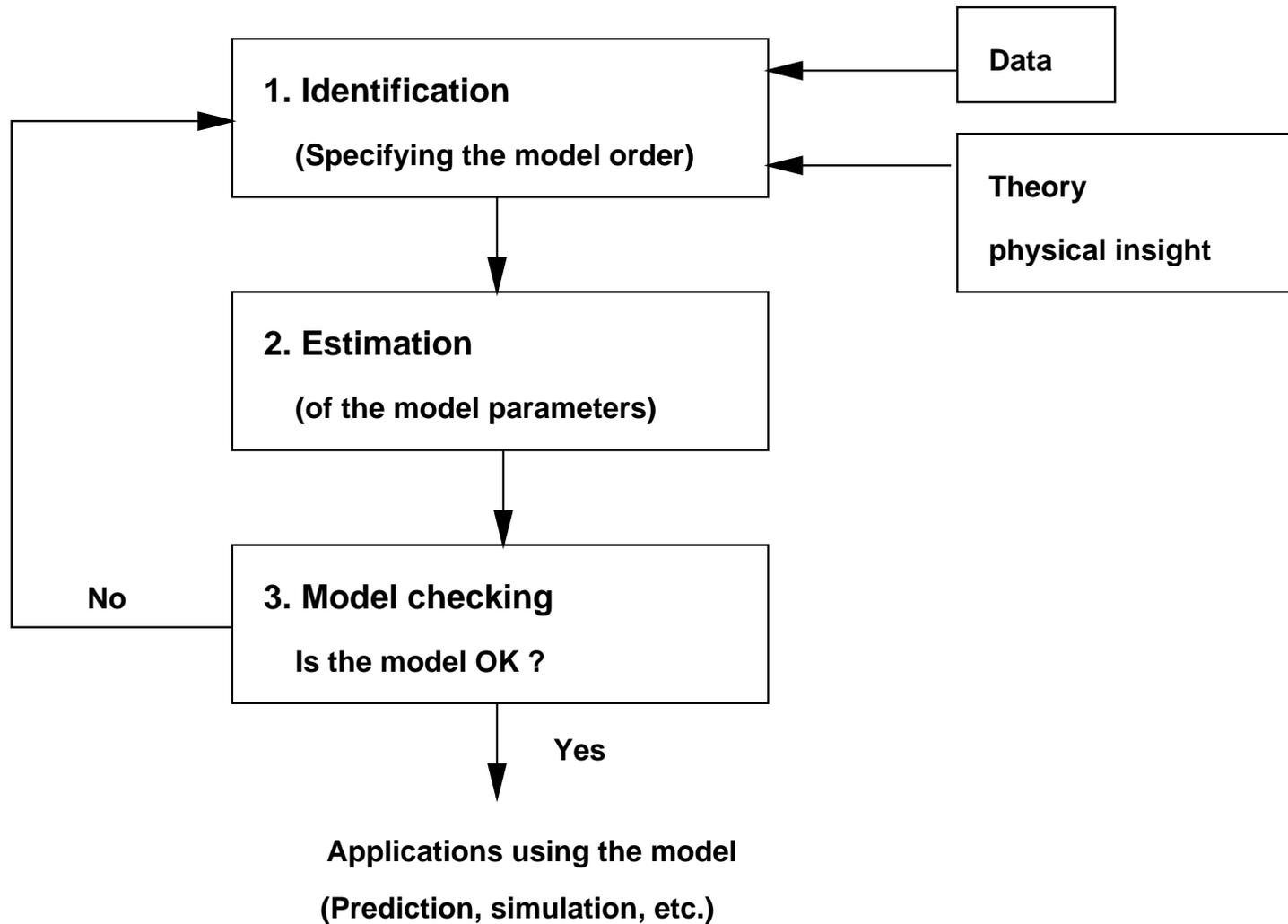
$$\begin{bmatrix} \hat{\rho}(1) \\ \hat{\rho}(2) \\ \vdots \\ \hat{\rho}(p) \end{bmatrix} = \begin{bmatrix} 1 & \hat{\rho}(1) & \cdots & \hat{\rho}(p-1) \\ \hat{\rho}(1) & 1 & \cdots & \hat{\rho}(p-2) \\ \vdots & \vdots & & \vdots \\ \hat{\rho}(p-1) & \hat{\rho}(p-2) & \cdots & 1 \end{bmatrix} \begin{bmatrix}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

and solve w.r.t. the ϕ 's

The function `ar` in S-PLUS or R use this approach as default



Model building





Validation of the model and extensions / reductions

- Residual analysis (Sec. 6.6.2): Is it possible to detect problems with residuals? (the 1-step prediction errors using the estimates, i.e. $\{\varepsilon_t(\hat{\theta})\}$, should be white noise)
- If the SACF or the SPACF of $\{\varepsilon_t(\hat{\theta})\}$ points towards a particular ARMA-structure we can derive how the original model should be extended (Sec. 6.5.1)
- If the model pass the residual analysis it makes sense to test null hypotheses about the parameters (Sec. 6.5.2)



Residual analysis

- Plot $\{\varepsilon_t(\hat{\theta})\}$; do the residuals look stationary?
- Tests in the autocorrelation. If $\{\varepsilon_t(\hat{\theta})\}$ is white noise then $\hat{\rho}_\varepsilon(k)$ is approximately Gaussian distributed with mean 0 and variance $1/N$.
If the model fails calculate SPACF also and see if an ARMA-structure for the residuals can be derived (Sec. 6.5.1)
- Since $\hat{\rho}_\varepsilon(k_1)$ and $\hat{\rho}_\varepsilon(k_2)$ are independent (Eq. 6.4) the test statistic $Q^2 = \sum_{k=1}^m \left(\sqrt{N} \hat{\rho}_{\varepsilon_t(\hat{\theta})}(k) \right)^2$ is approximately distributed as $\chi^2(m - n)$, where n is the number of parameters.

S-PLUS: `arima.diag('output from arima.mle')`



Residual analysis (continued)

- Test for the number of changes in sign. In a series of length N there is $N - 1$ possibilities for changes in sign. If the series is white noise (with mean zero) the probability of change is $1/2$ and the changes will be independent. Therefore the number of changes is distributed as $Bin(N - 1, 1/2)$

S-PLUS: `binom.test(N-1, 'No. of changes')`

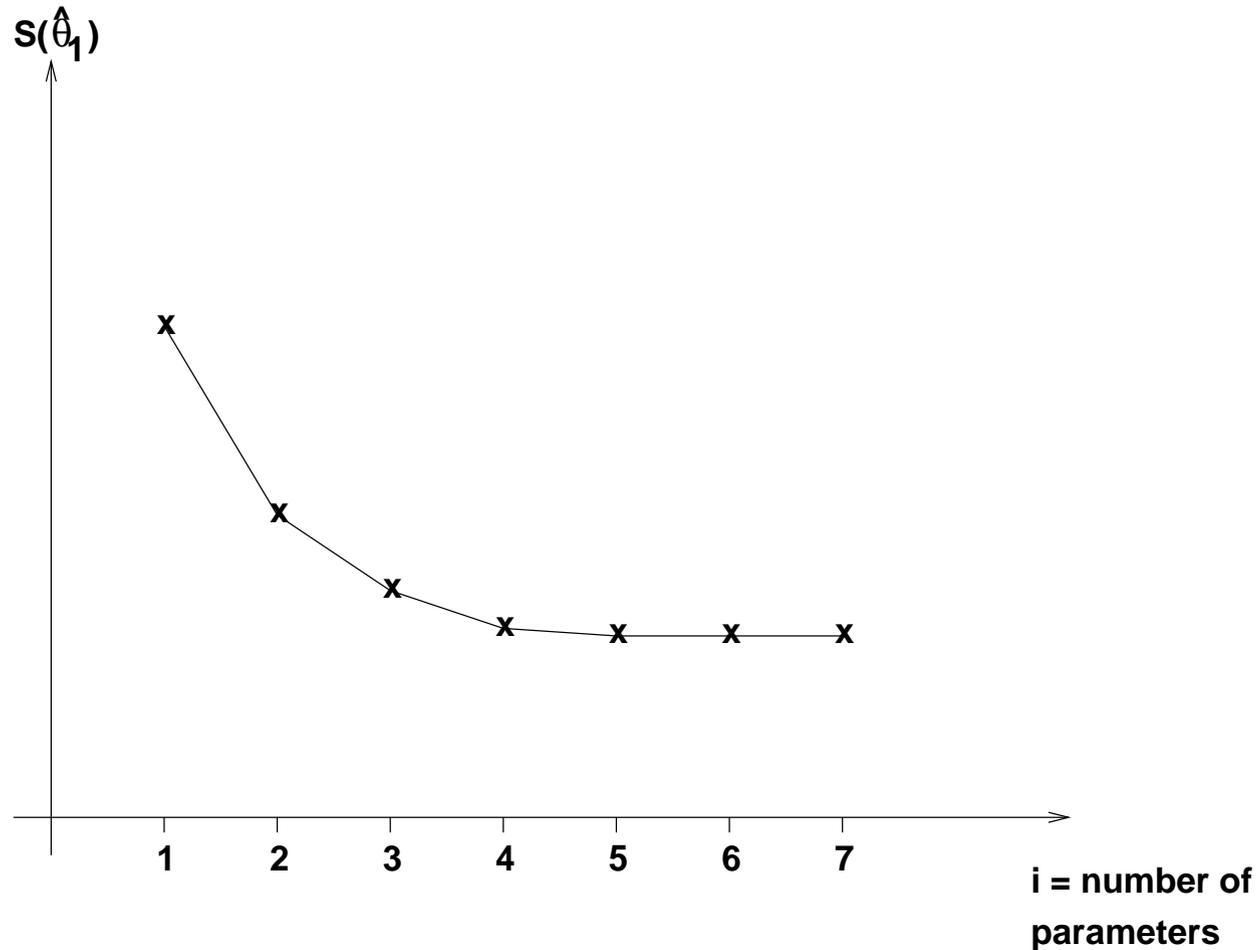
- Test in the *scaled cumulated periodogram* of the residuals is done by plotting it and adding lines at $\pm K_\alpha / \sqrt{q}$, where $q = (N - 2)/2$ for N even and $q = (N - 1)/2$ for N odd. For $1 - \alpha$ confidence limits K_α can be found in Table 6.2

S-PLUS (95% confidence interval):

```
library(MASS)
cpgram('residuals')
```



Sum of squared residuals depend on the model size



(It is assumed that the models are nested)



Test is the model

- The test essentially checks if the reduction in SSE ($S_1 - S_2$) is large enough to justify the *extra* parameters in model 2 (n_2 parameters) as compared to model 1 (n_1 parameters). The number of observations used is called N .
- If vector θ_{extra} is used to denote the extra parameters in model 2 as compared to model 1, then the test is formally:

$$H_0 : \theta_{extra} = \mathbf{0} \text{ vs. } H_0 : \theta_{extra} \neq \mathbf{0}$$

- If H_0 is true it (approximately) hold that

$$\frac{(S_1 - S_2)/(n_2 - n_1)}{S_2/(N - n_2)} \sim F(n_2 - n_1, N - n_2)$$

(The likelihood ratio test is also a possibility)



Testing one parameter for significance

$$H_0 : \theta_i = 0 \text{ against } H_1 : \theta_i \neq 0$$

- Can be done as described on the previous slide
- Alternatively we can use a t-test based on the estimate and its standard error: $\hat{\theta}_i / \sqrt{\hat{V}(\hat{\theta}_i)}$
- Under H_0 and for an $ARMA(p, q)$ -model this follows a $t(N - p - q)$ distribution (or $t(N - 1 - p - q)$ if we estimated an overall mean of the series)
- Often N is so large compared to the number of parameters that we can just use the standard normal distribution



Information criteria

- Select the model which minimize some information criterion
- Akaike's Information Criterion

$$AIC = -2 \log(L(\mathbf{Y}_N; \hat{\boldsymbol{\theta}}, \hat{\sigma}_\varepsilon^2)) + 2 n_{par}$$

- Bayesian Information Criterion

$$BIC = -2 \log(L(\mathbf{Y}_N; \hat{\boldsymbol{\theta}}, \hat{\sigma}_\varepsilon^2)) + \log N n_{par}$$

- Except for an additive constant this can also be expressed as

$$AIC = N \log \hat{\sigma}_\varepsilon^2 + 2 n_{par}$$

$$BIC = N \log \hat{\sigma}_\varepsilon^2 + \log N n_{par}$$

- BIC yields a consistent estimate of the model order



Example

A model for CO_2 ...