

# Linear Time Series

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## Chapter 2: ARMA Models

# Outline

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  - Existence of solutions to the ARMA model
  - Generality of ARMA processes
- 2 Characterisation of the orders  $p$  and  $q$ 
  - Autocorrelation function
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- 3 Estimation, validation and predictions
  - Estimation of ARMA models
  - Validation and model choice
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## Definition of ARMA models

### Definition:

$(X_t)$  is an ARMA( $p, q$ ) if  $(X_t)$  is 2nd-order stationary and satisfies for all  $t$

$$X_t - \phi_1 X_{t-1} - \cdots - \phi_p X_{t-p} = \epsilon_t - \psi_1 \epsilon_{t-1} - \cdots - \psi_q \epsilon_{t-q}$$

where  $(\epsilon_t) \sim WN(0, \sigma^2)$ . It is said that  $(X_t)$  follows an ARMA( $p, q$ ) with mean  $\mu$  if  $(X_t - \mu)$  is an ARMA( $p, q$ ) (with mean 0).

Problem: given coefficients  $\psi_i$  and  $\phi_j$  and a WN, does a solution  $(X_t)$  of the ARMA model exist?

Existence depends on the AR and MA polynomials

$$\phi(z) = 1 - \phi_1 z - \cdots - \phi_p z^p \quad \text{and} \quad \psi(z) = 1 - \psi_1 z - \cdots - \psi_q z^q.$$

## Backward and Forward operators

### Definition:

- 1 the **backward operator**  $B$  (also noted  $L$  for lag operator) transforms  $X_t$  into  $X_{t-1}$ :

$$BX_t = X_{t-1}$$

- 2 the **forward operator**  $F$  transforms  $X_t$  into  $X_{t+1}$ :

$$FX_t = X_{t+1}$$

### Properties:

- iteration:  $B^k X_t = X_{t-k}$ ,  $F^k X_t = X_{t+k}$
- identity operator:  $I = BF = FB$

## Polynomials in $B$ or $F$

$$\begin{aligned}P(B) &= a_0 + a_1 B + \dots + a_n B^n \\ \Rightarrow P(B)X_t &= a_0 X_t + a_1 X_{t-1} + \dots + a_n X_{t-n}\end{aligned}$$

One can add, multiply polynomials in  $B$  or  $F$  as usual polynomials.

We can also consider infinite sums:

if  $\sum_{i=-\infty}^{\infty} |a_i| < \infty$ , let the **power series in  $B$**

$$a(B) = \sum_{i=-\infty}^{\infty} a_i B^i,$$

such that for any stationary process  $(X_t)$ ,

$$a(B)X_t = \sum_{i=-\infty}^{\infty} a_i X_{t-i} \text{ is stationary.}$$

► Existence

## Multiplicative inverse of a polynomial in $B$

If  $P(B)X_t = Z_t$ , can we express  $X_t$  as a function of present, past and future values of  $Z_t$ ?

$$BX_t = Z_t \Rightarrow X_t = B^{-1}Z_t = Z_{t+1}$$

Inversion of  $1 - \phi B$ : suppose  $(1 - \phi B)X_t = Z_t$

①  $|\phi| < 1$

$$(1 - \phi B)^{-1} = \sum_{i=0}^{+\infty} \phi^i B^i \Rightarrow X_t = \sum_{i=0}^{+\infty} \phi^i Z_{t-i}$$

②  $|\phi| > 1$

$$(1 - \phi B)^{-1} = - \sum_{i=1}^{+\infty} \frac{1}{\phi^i} B^i \Rightarrow X_t = - \sum_{i=1}^{+\infty} \frac{1}{\phi^i} Z_{t+i}$$

③  $|\phi| = 1$  Inversion is not possible.

## Inversion of $\phi(B) = 1 - \phi_1 B - \dots - \phi_p B^p$

Suppose that

$$\phi(z) \neq 0 \quad \text{for all complex numbers } z \text{ such that } |z| \leq 1,$$

that is, the roots of  $\phi(z)$  are outside the unit disc.

Then, there exists  $\delta > 0$  such that

$$\frac{1}{\phi(z)} = \sum_{j=0}^{\infty} c_j z^j, \quad |z| < 1 + \delta$$

where  $\sum_{j=0}^{\infty} |c_j| < \infty$ .

We then have

$$\frac{1}{\phi(B)} = \sum_{j=0}^{\infty} c_j B^j.$$

## Practical derivation of the inverse

The coefficients  $c_j$  of

$$\frac{1}{\phi(B)} = \sum_{j=0}^{\infty} c_j B^j$$

can be obtained by

- Identification
- Partial fraction decomposition

**Example:**  $X_t - 0.6X_{t-1} + 0.08X_{t-2} = \epsilon_t$

$$\rightarrow X_t = \epsilon_t + \sum_{i=1}^{+\infty} [2(0.4)^i - 0.2^i] \epsilon_{t-i}$$

$$\left( \frac{1}{(1-0.4z)(1-0.2z)} = \frac{2}{1-0.4z} - \frac{1}{1-0.2z} \right)$$



## Existence of a causal solution

The ARMA( $p, q$ ) model

$$X_t - \phi_1 X_{t-1} - \cdots - \phi_p X_{t-p} = \epsilon_t - \psi_1 \epsilon_{t-1} - \cdots - \psi_q \epsilon_{t-q}$$

writes  $\phi(B)X_t = \psi(B)\epsilon_t$ . We assume that  $\text{Var}(\epsilon_t) = \sigma^2 > 0$ , and that the polynomials  $\phi(z)$  and  $\psi(z)$  have **no common root**.

### Proposition:

A **causal stationary solution** (i.e. of the form  $X_t = \sum_{j=0}^{\infty} c_j \epsilon_{t-j}$ ) exists if and only if

$$\phi(z) = 1 - \phi_1 z - \cdots - \phi_p z^p \neq 0 \quad \text{for all } z \text{ such that } |z| \leq 1.$$

We then have  $\frac{\psi(B)}{\phi(B)} = \sum_{j=0}^{\infty} c_j B^j$ .

► Proof

## Invertible solution

**Invertibility** means that  $\epsilon_t$  can be expressed as a function of  $X_t$  and its past values.

### Proposition:

The ARMA model is invertible (i.e. we have  $\epsilon_t = \sum_{j=0}^{\infty} b_j X_{t-j}$  with  $\sum_{j=0}^{\infty} |b_j| < \infty$ ) if and only if

$$\psi(z) = 1 - \psi_1 z - \dots - \psi_q z^q \neq 0 \quad \text{for all } z \text{ such that } |z| \leq 1.$$

We then have

$$\epsilon_t = X_t + \sum_{j=1}^{\infty} b_j X_{t-j}, \quad \frac{\phi(z)}{\psi(z)} = \sum_{j=0}^{\infty} b_j z^j.$$

## Common roots

If **common roots** exist in the AR and MA polynomials and **have modulus different from 1**, we get the same solutions by cancelling those roots.

For instance, the model

$$X_t - \phi X_{t-1} = \epsilon_t - \phi \epsilon_{t-1}, \quad |\phi| \neq 1$$

has the unique solution

$$X_t = \epsilon_t.$$

⚠: coefficient  $\phi$  is not identifiable in this model.

## Canonical ARMA

### Definition and property:

The series  $(X_t)$  admits a **canonical ARMA**( $p, q$ ) representation if

$$\phi(B)X_t = \psi(B)\epsilon_t$$

where  $\phi$  and  $\psi$  have no common roots, and

$$\phi(z) = 1 - \phi_1 z - \dots - \phi_p z^p \neq 0 \quad \text{for all } z \text{ such that } |z| \leq 1,$$

$$\psi(z) = 1 - \psi_1 z - \dots - \psi_q z^q \neq 0 \quad \text{for all } z \text{ such that } |z| \leq 1.$$

- $(\epsilon_t)$  is the **linear innovation** of  $(X_t)$ :

$$X_t = \underbrace{EL(X_t | X_{t-1}, \dots)}_{\text{best **linear** pred.}} + \epsilon_t, \quad \text{with} \quad \text{Cov}(\epsilon_t, X_{t-i}) = 0, \quad \text{for } i > 0$$

- The past of  $X_t$  and  $\epsilon_t$  coincide

## ARMA with mean

A stationary ARMA with mean  $\mu$  can be written as

$$\phi(B)(X_t - \mu) = \psi(B)\epsilon_t$$

or equivalently

$$\phi(B)X_t = c + \psi(B)\epsilon_t$$

since

$$\phi(B)(X_t - \mu) = \phi(B)X_t - \phi(B)\mu = \phi(B)X_t - \phi(1)\mu$$

and  $\phi(1) \neq 0$  under the stationarity condition. The intercept  $c$  is however more difficult to interpret than the mean  $\mu$ .

## Generality of ARMA models

From the Wold (1938) theorem, any stationary process admits a linear representation after subtraction of a deterministic component:

$$X_t = \epsilon_t + \sum_{j=1}^{\infty} c_j \epsilon_{t-j}, \quad \sum_{j=1}^{\infty} c_j^2 < \infty.$$

This MA( $\infty$ ) representation is not easy to use in practice.

Finite order MA models can be seen as an approximation of the Wold representation:

$$X_t = \epsilon_t + \sum_{j=1}^q c_j \epsilon_{t-j},$$

But a very large order  $q$  may be required.

A more parsimonious approximation is obtained by explicitly involving the past values of  $X_t$ .

► Wold theorem

## Recursive relation between autocorrelations

Let the ARMA be written under the canonical form:

$$X_t - \phi_1 X_{t-1} - \cdots - \phi_p X_{t-p} = \epsilon_t - \psi_1 \epsilon_{t-1} - \cdots - \psi_q \epsilon_{t-q}.$$

We have

$$\begin{aligned} \gamma(h) &= \text{Cov}(X_t, X_{t-h}) = \text{Cov}\left(\sum_{i=1}^p \phi_i X_{t-i} + \epsilon_t - \sum_{i=1}^q \psi_i \epsilon_{t-i}, X_{t-h}\right) \\ &= \sum_{i=1}^p \phi_i \text{Cov}(X_{t-i}, X_{t-h}) + \text{Cov}(\epsilon_t, X_{t-h}) - \sum_{i=1}^q \psi_i \text{Cov}(\epsilon_{t-i}, X_{t-h}) \end{aligned}$$

Thus

$$\gamma(h) = \sum_{i=1}^p \phi_i \gamma(h-i), \quad \forall h > q.$$

## Characteristic property of ARMA( $p, q$ )

### Proposition:

For a stationary and centred process  $(X_t)$ , the autocorrelations satisfy a recursive equation of order  $p$ , starting from rank  $q+1$ :

$$\rho(h) = \sum_{i=1}^p \phi_i \rho(h-i), \quad \forall h > q \text{ (with } \phi(z) \neq 0 \text{ for } |z| \leq 1)$$

**if and only if**  $X_t \sim \text{ARMA}(p, q)$ , where the  $\phi_i$ 's are the AR coefficients.

► Proof

In particular:  $\text{MA}(q) \iff \rho(h) = 0, \quad h > q$

In an ARMA, the autocorrelations decrease at exponential rate.



## Application: identification of a model

Given observations  $x_1, \dots, x_n$ , one approach to **identify the orders** is to compare the  $\hat{\rho}(h)$ 's with the  $\rho(h)$ 's of a given model.

In particular, if  $\hat{\rho}(h) \approx 0$ , for  $h > q$ , one can identify a  $\text{MA}(q)$ .

$\text{MA}(q)$ : the  $\hat{\rho}(h)$  are asymptotically Gaussian for  $h > q$ ,

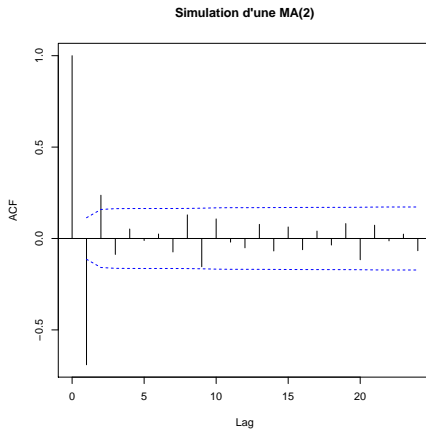
$$\sqrt{n}\hat{\rho}(h) \xrightarrow{\mathcal{L}} \mathcal{N}(0, 1 + 2\rho^2(1) + \dots + 2\rho^2(q)) \quad h > q.$$

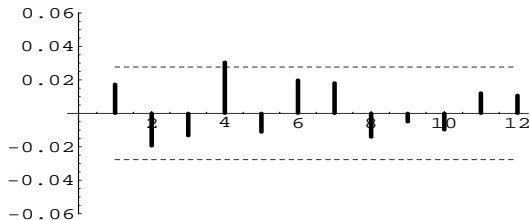
Significance bounds at 95%:

$$\pm 1.96 \sqrt{1 + 2\hat{\rho}^2(1) + \dots + 2\hat{\rho}^2(h-1)} / \sqrt{n}.$$

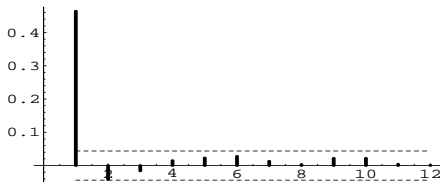
In practice, it is standard to use the smaller bounds  $\pm 1.96 / \sqrt{n}$ .

```
n<-300; epsilon<-rnorm(n)
X<-epsilon[3:n]-1.804*epsilon[2:(n-1)]+0.806*epsilon[1:(n-2)]
acf(X,ci.type="ma",main="Simulation of a MA(2)")
```





Empirical autocorrelations of a strong WN, for  $n=5000$ .  
In dotted lines, the significance bounds:  $\pm 1.96/\sqrt{n}$



Empirical autocorrelations of a MA(1), for  $\theta = 0.5$  and  $n=2000$ .  
In dotted lines, the significance bounds:  $\pm 1.96/\sqrt{n}$  (invalid here)

## Specific influence of a lagged variable

Like correlations, **partial autocorrelations** convey crucial information on the dependence structure of a stationary process. Both depend only on the second-order properties of the process.

The linear influence of  $X_{t-h}$  on  $X_t$  can be measured by  $\rho(h)$   
But between these 2 variables we have:  $X_{t-h+1}, \dots, X_{t-1}$ .

$$X_{t-h} \rightarrow X_{t-h+1} \rightarrow \dots \rightarrow X_{t-1} \rightarrow X_t$$

Question: what is the **specific influence of  $X_{t-h}$  on  $X_t$** ? (adjusted for the intermediate variables)

## Partial autocorrelations of a stationary process

For  $h > 1$  let

- $\tilde{X}_t$  = linear regression of  $X_t$  on  $1, X_{t-1}, \dots, X_{t-h+1}$ .
- $\tilde{X}_{t-h}^*$  = linear regression of  $X_{t-h}$  on  $1, X_{t-1}, \dots, X_{t-h+1}$

To adjust for the influence of the intermediate variables let

$$X_t - \tilde{X}_t \quad \text{and} \quad X_{t-h} - \tilde{X}_{t-h}^*$$

### Partial autocorrelation function $r$

$$r(1) = \rho(1)$$

$$r(h) = \text{Corr}(X_t - \tilde{X}_t, X_{t-h} - \tilde{X}_{t-h}^*) := \text{Corr}(X_t, X_{t-h} \mid X_{t-1}, \dots, X_{t-h+1})$$

$r(h)$  can be interpreted as a measure of the (linear) dependence between  $X_t$  and  $X_{t-h}$  that is not conveyed by the intermediate variables

## Alternative definition of $r(h)$

### Proposition:

$r(h)$  is the coefficient of  $X_{t-h}$  in the regression of  $X_t$  on  $\{1, X_{t-1}, \dots, X_{t-h}\}$ :

$$X_t = \alpha_0 + \sum_{i=1}^{h-1} \alpha_i X_{t-i} + r(h) X_{t-h} + \epsilon_t$$

Proof: see Brockwell Davis (Corr. 5.2.1, 1991)

⚠: the other coefficients  $\alpha_i$  cannot be interpreted as  $r(i)$ 's!

## Link with the autocorrelation function

### Yule-Walker equations

$$X_t = \alpha_0 + \sum_{i=1}^h \alpha_i X_{t-i} + \epsilon_t, \quad \epsilon_t \perp 1, X_{t-1}, \dots, X_{t-h},$$

entails

$$\begin{bmatrix} \rho(1) \\ \vdots \\ \vdots \\ \vdots \\ \rho(h) \end{bmatrix} = \begin{bmatrix} 1 & \rho(1) & \dots & \rho(h-1) \\ \rho(1) & 1 & \rho(1) & \dots & \rho(h-2) \\ \vdots & \ddots & \ddots & \ddots & \vdots \\ \vdots & \ddots & \ddots & 1 & \rho(1) \\ \rho(h-1) & \rho(h-2) & \dots & \rho(1) & 1 \end{bmatrix} \begin{bmatrix} \alpha_1 \\ \vdots \\ \vdots \\ \alpha_{h-1} \\ r(h) \end{bmatrix}$$

If the matrix  $R_h$  is non-singular (generally the case), one can solve the system in  $(\alpha_1, \dots, r(h))$ . The solution is unique.

The calculation can be done fastly using Durbin-Levinson's algorithm (see BD, Chapter 5).



## Partial autocorrelation of an $AR(p)$

For a **causal**  $AR(p)$

$$X_t - \phi_1 X_{t-1} - \cdots - \phi_p X_{t-p} = \epsilon_t$$

$(\epsilon_t)$  is the innovation:

$$\text{Cov}(\epsilon_t, X_{t-i}) = 0, \quad i > 0.$$

Thus

$$r(p) = \phi_p$$

and

$$r(h) = 0, \quad \text{for } h > p.$$

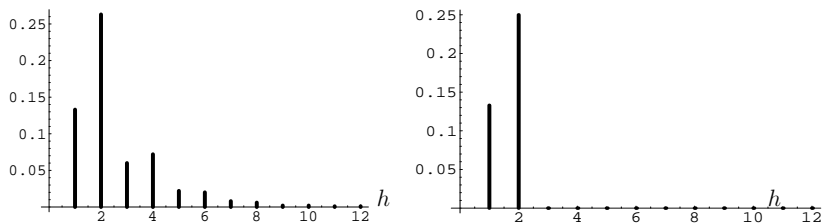


FIG. 1 – Fonction d'autocorrélation (graphe de gauche) et fonction d'autocorrélation partielle (graphe de droite) d'un AR(2) :

$$X_t = 1 + 0.1X_{t-1} + 0.25X_{t-2} + \epsilon_t$$

## Empirical partial autocorrelations

The **empirical partial autocorrelations**,  $\hat{r}(h)$ , are obtained by replacing the  $\rho(k)$  by  $\hat{\rho}(k)$  in the previous matrix equation

In a **strong**  $AR(p)$  (with strong WN), the asymptotic distribution of the  $\hat{r}(h)$ ,  $h > p$ , is very simple.

### Proposition:

If  $(X_t)$  is the causal stationary solution of an  $AR(p)$  model with iid WN,

$$\sqrt{n}\hat{r}(h) \xrightarrow{\mathcal{L}} \mathcal{N}(0, 1), \quad \forall h > p.$$

## Identification of pure MA and AR models

Recall that a  $\text{MA}(q)$  satisfies

$$\rho(h) = 0 \quad \text{for all } h > q,$$

and that an  $\text{AR}(p)$  satisfies

$$r(h) = 0 \quad \text{for all } h > p.$$

In practice we use  $\hat{\rho}(h)$  and  $\hat{r}(h)$ . If for instance

$$\hat{\rho}(1) \neq 0, \quad \text{and for all } h > 1, \hat{\rho}(h) \approx 0,$$

a  $\text{MA}(1)$  can be fitted.

If now,

$$\hat{r}(3) \neq 0, \quad \text{and for all } h > 4, \hat{r}(h) \approx 0,$$

an  $\text{AR}(3)$  can be fitted.

## Identification of mixed models

For (ARMA( $p, q$ ) models with  $pq \neq 0$ ), more sophisticated statistical methods can be used:

- the **corner method** (Béguin, Gouriéroux, Monfort, 1980);

▶ Corner method

- the **spectral density**;

▶ Spectral density

One can also proceed by estimating a lot of models, and by testing

- (i) the significance of the estimated parameters, and
- (ii) the independence of the residuals.

One can also use information criteria.

## Least Squares (LS) estimator

From observations  $X_1, X_2, \dots, X_n$ , one can approximate  $\epsilon_t(\theta)$ , for  $0 < t \leq n$ , by  $e_t(\theta)$  recursively defined by

$$e_t(\theta) = X_t - \sum_{i=1}^p \phi_i X_{t-i} + \sum_{i=1}^q \psi_i e_{t-i}(\theta)$$

where  $e_0(\theta) = e_{-1}(\theta) = \dots = e_{-q+1}(\theta) = X_0 = X_{-1} = \dots = X_{-p+1} = 0$ .

An (approximated) LS estimator  $\hat{\theta}_n$  is defined by

$$Q_n(\hat{\theta}_n) = \min_{\theta \in \Theta} Q_n(\theta), \quad Q_n(\theta) = \frac{1}{n} \sum_{t=1}^n e_t^2(\theta).$$

## Case of the AR(1)

Let  $(X_t)$  a causal AR(1) with a strong white noise

$$X_t = aX_{t-1} + \epsilon_t, \quad |a| < 1.$$

We have  $\theta = a$ ,  $e_1(a) = X_1$ ,  $e_t(a) = X_t - aX_{t-1}$  for  $t = 2, \dots, n$  and  $Q_n(\hat{a}) = 0$  iff

$$\hat{a} = \frac{\sum_{t=2}^n X_t X_{t-1}}{\sum_{t=2}^n X_{t-1}^2}.$$

By the ergodic theorem,  $\hat{a} \rightarrow a$  a.s. A CLT for non iid sequences shows that

$$\sqrt{n}\{\hat{a} - a\} = \frac{n^{-1/2} \sum_{t=2}^n \epsilon_t X_{t-1}}{n^{-1} \sum_{t=2}^n X_{t-1}^2} \xrightarrow{\mathcal{L}} \mathcal{N}(0, 1 - a^2).$$

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\*For other models, for instance the MA(1), the LS estimator is not explicitly defined.

## Validation

To evaluate the **goodness of fit** of an ARMA model, **2 types of tests**:

- tests on the nullity of the **parameters**  $\phi_j$  and  $\psi_j$   
[Aim: check if the model cannot be simplified].
- tests on the **residuals**: if the fit is appropriate, the residuals should resemble a **VN sequence**.  
[Aim: check if the model is rich enough].

The choice between several appropriate models can be founded on the minimisation of **information criteria** (to find a balance between the accuracy of the fit and the simplicity of the model).



## Tests on the parameters

Goal: compare 2 formulations  $\text{ARMA}(p, q)$  and  $\text{ARMA}(p', q')$ .

Assume that one model embeds the other one: for instance  $p' \leq p$  and  $q' \leq q$

i)  $p' = p - 1$ ,  $q' = q$ : testing the significance of  $\phi_p$ .

**Student test**: At the level 0.05, the  $\text{ARMA}(p - 1, q)$  is accepted if

$$\frac{|\hat{\phi}_p|}{[\hat{V}(\hat{\phi}_p)]^{1/2}} < 1.96$$

ii)  $p' = p$ ,  $q' = q - 1$ : idem for testing the significance of  $\psi_q$ .

iii)  $p' = p - 1$ ,  $q' = q - 1$ : impossible.

One cannot reduce simultaneously the AR and MA orders (non uniqueness of the ARMA representation).

## Tests on the residuals

The residuals  $\hat{\epsilon}_t$ ,  $t = 1, \dots, n$  are obtained from

$$\hat{\epsilon}_t = \frac{\hat{\Phi}(B)}{\hat{\Psi}(B)} X_t, \quad \hat{\Phi}(B) = 1 - \hat{\phi}_1 B - \dots - \hat{\phi}_p B^p, \hat{\Psi}(B) = 1 - \hat{\psi}_1 B - \dots - \hat{\psi}_q B^q$$

and  $X_t = 0, t \leq 0$ .

- 1 **Graph of the residuals:** outliers, or deviations of the variance from a constant, are sometimes clearly indicated. Absence of correlations is more difficult to identify.
- 2 **Empirical autocorrelation function of the  $\hat{\epsilon}_t$ 's:** the sample autocorrelations of the  $\hat{\epsilon}_t$ 's (not the  $\epsilon_t$ 's) are, for  $n$  large, approximately iid with distribution  $\mathcal{N}(0, 1/n)$ .

Because each  $\hat{\epsilon}_t$  is a function of the observations, it is not an iid sequence: the asymptotic distribution is not exactly the same as in the iid case (except for large lags).

For large  $n$ , the sample autocorrelation of order  $h$  of the residuals is (approximately) distributed as

$$\hat{\rho}_{\hat{\epsilon}}(h) \sim \mathcal{N}(0, \nu)$$

where

- $\nu < 1/n$  for small  $h$ ,
- $\nu \approx 1/n$  for large  $h$ .

The variance  $\nu$  can sometimes be explicitly computed.

## Portmanteau test (Box-Pierce (1970))

Instead of checking that each  $\hat{\rho}_{\hat{e}}(h)$  is between the bounds  $\pm 1.96v^{1/2}$ , one can consider a **global statistic** depending on all sample autocorrelations.

$$Q = n \sum_{h=1}^H \hat{\rho}_{\hat{e}}^2(h).$$

The asymptotic distribution of  $Q$  est approximately a  $\chi^2$  with  $H - p - q$  degrees of freedom (**under the assumption of iid noise**).  
The model is rejected at level  $\alpha \in (0, 1)$  if

$$Q > \chi_{1-\alpha}^2(H - p - q).$$

## Modified Portmanteau test (Ljung-Box (1978))

For finite, even large,  $n$  the law of  $Q$  is far from the asymptotic distribution. Hence a modified statistic:

$$Q' = n(n+2) \sum_{h=1}^H \frac{1}{n-h} \hat{\rho}_{\hat{\epsilon}}^2(h).$$

$H$  must be chosen large enough but not too large (generally between 15 and 30). When the fit is rejected, individual autocorrelations can be inspected to modify the model.

**Drawback:** lack of power. Even poor fits pass the test. Other tests are aimed at selecting the best model.

## Model choice using information criteria

**Idea:** prevent against over-parametrisation by introducing a cost for the introduction of any additional parameter.

The approach introduced by *Akaike (1969)* is based on the Kullback-Leibler distance between the estimated and true models.

Several estimators of the information were proposed:

- ①  $AIC(p,q) = \log \hat{\sigma}^2 + \frac{2(p+q)}{n};$
- ②  $BIC(p,q) = \log \hat{\sigma}^2 + (p+q) \frac{\log n}{n};$
- ③  $\phi(p,q) = \log \hat{\sigma}^2 + c(p+q) \frac{\log \log n}{n},$  with  $c > 2$

where

$$\hat{\sigma}^2 = \frac{1}{n} \sum_{t=1}^n \hat{\epsilon}_t^2.$$

The model will be selected by **minimizing the estimated information criterion**.

AIC is by far the **most widely used**. But it often leads to over-parametrisation.

Only the BIC and  $\phi$  criteria lead to **consistent** estimators of  $p$  and  $q$ . But for an  $\text{AR}(\infty)$  model, only the AIC criterion leads to an asymptotically efficient estimator.

**Remark:** sometimes, the criteria do not select the same orders.

## Theoretical predictions of an ARMA

Let an  $\text{ARMA}(p, q)$

$$X_t - \sum_{i=1}^p \phi_i X_{t-i} = \epsilon_t - \sum_{i=1}^q \psi_i \epsilon_{t-i},$$

assumed to be **causal and invertible**.

The **linear optimal prediction at horizon 1** of  $X_t$  is

$$\hat{X}_{t|t-1} = \sum_{i=1}^p \phi_i X_{t-i} - \sum_{i=1}^q \psi_i \epsilon_{t-i}$$

because  $\epsilon_t = X_t - \hat{X}_{t|t-1} \perp \mathcal{H}_X(t-1)$  and  $\epsilon_{t-i} = \frac{\phi(B)}{\psi(B)} X_{t-i} \in \mathcal{H}_X(t-1)$ .<sup>†</sup>

► Complements and predictions at horizon  $h$

---

<sup>†</sup>( $\mathcal{H}_X(t-1)$ ): Hilbert space generated by the linear combinations of  $1, X_{t-1}, X_{t-2}, \dots$ )



## Prediction using an estimated ARMA model

Having estimated the ARMA coefficients from  $X_1, \dots, X_n$  with  $n \leq T$ , using initial values (for instance 0) for  $X_0, \dots, X_{1-p}, \tilde{\epsilon}_0, \dots, \tilde{\epsilon}_{1-q}$  we compute

$$\tilde{\epsilon}_t = X_t - \sum_{i=1}^p \hat{\phi}_i X_{t-i} + \sum_{i=1}^q \hat{\psi}_i \tilde{\epsilon}_{t-i}$$

for  $t = 1, \dots, T$  which allows to predict  $X_{T+1}$  by

$$\hat{X}_{T+1|T} = \sum_{i=1}^p \hat{\phi}_i X_{T+1-i} - \sum_{i=1}^q \hat{\psi}_i \tilde{\epsilon}_{T+1-i}.$$

End of chapter 2 😊 !

## Proof of the necessary and sufficient existence condition for a causal solution

SC: If the condition is satisfied, then there exists  $\delta > 0$  such that

$$\frac{1}{\phi(z)} = \sum_{j=0}^{\infty} d_j z^j, \quad |z| < 1 + \delta$$

where  $\sum_{j=0}^{\infty} |d_j| < \infty$ . Since  $\psi(B)\epsilon_t$  is stationary, multiply by  $\phi^{-1}(B)$  the equality  $\phi(B)X_t = \psi(B)\epsilon_t$ , to get

$$X_t = \phi^{-1}(B)\psi(B)\epsilon_t = \sum_{j=0}^{\infty} c_j \epsilon_{t-j}.$$

## Proof of the necessary and sufficient existence condition for a causal solution

NC: Let  $X_t$  a causal solution of the form  $X_t = c(B)\epsilon_t$ , then

$$\psi(B)\epsilon_t = \phi(B)X_t = \phi(B)c(B)\epsilon_t := \sum_{i=0}^{\infty} e_i \epsilon_{t-i}.$$

Multiply each side by  $\epsilon_{t-h}$ ,  $h \geq 0$ , and take the expectation, to get  $e_0 = 1$ ,  $e_i = -\psi_i$  for  $i = 1, \dots, q$  and  $e_i = 0$  for  $i > q$ .

Thus

$$\psi(z) = \phi(z)c(z), \quad |z| \leq 1.$$

Since  $\psi(z)$  and  $\phi(z)$  have no common root and  $|c(z)| < \infty$  for  $|z| \leq 1$ , we cannot have  $\phi(z) = 0$  for  $|z| \leq 1$ .

◀ Return

## Wold decomposition (1938)

$\mathcal{H}_X(t-1)$  : linear past of  $X_t$ , i.e. the closed subset of  $L^2$  generated by  $1, X_{t-1}, X_{t-2}, \dots$

$(X_t)$  is called deterministic if

$$X_t \in \mathcal{H}_X(-\infty) = \bigcap_{s=-\infty}^{\infty} \mathcal{H}_X(s)$$

Examples:  $X_t = m$  ( $m$  constant), or  $X_t = X$  ( $X$  r.r.v. in  $L^2$ ).

### Wold decomposition Theorem (proof in BD p 187)

If  $(X_t)$  is a 2nd-order stationary process, then

$$X_t = \epsilon_t + \sum_{i=1}^{\infty} c_i \epsilon_{t-i} + V_t, \quad \sum_{i=1}^{\infty} c_i^2 < \infty,$$

where  $(\epsilon_t) \sim WN(0, \sigma^2)$ ,  $\epsilon_t \in \mathcal{H}_X(t)$ , and  $(V_t)$  is deterministic with  $EV_t \epsilon_s = 0 \quad \forall t, s$ .

## Characteristic property of an ARMA( $p, q$ )

It remains to show that

$$\gamma_X(h) - \sum_{i=1}^p \phi_i \gamma_X(h-i) = 0, \quad \forall h > q \quad \Rightarrow \quad Y_t := X_t - \sum_{i=1}^p \phi_i X_{t-i} \sim \text{MA}(q).$$

The linear innovation of  $Y_t$  is a WN defined by  $\epsilon_t = Y_t - E(Y_t | \mathcal{H}_Y(t-1))$ . The space  $\mathcal{H}_Y(t-1)$  is the direct sum of  $\mathcal{H}_Y(t-q-1)$  and of the Hilbert space  $\mathcal{H}(\epsilon_{t-1}, \dots, \epsilon_{t-q})$ , generated by  $\epsilon_{t-1}, \dots, \epsilon_{t-q}$ . The roots of  $\phi$  being outside the unit circle, we have  $\mathcal{H}_Y(t) = \mathcal{H}_X(t)$ . Since  $EY_t = 0$  and  $EY_t X_{t-h} = 0 \quad \forall h > q$ , we have

$$Y_t \perp \mathcal{H}_X(t-q-1) = \mathcal{H}_Y(t-q-1).$$

Therefore,

$$Y_t = \epsilon_t + E(Y_t | \mathcal{H}(\epsilon_{t-1}, \dots, \epsilon_{t-q})) = \epsilon_t + \sum_{i=1}^q \theta_i \epsilon_{t-i},$$

with  $E(Y_t - \theta_i \epsilon_{t-i}) \epsilon_{t-i} = 0$ , that is  $\theta_i = EY_t \epsilon_{t-i} / E\epsilon_1^2$ .

## Proof of the spectral density for a linear transform

We have

$$\gamma_X(h) = \sum_{j,\ell} a_j a_\ell \gamma_Y(\ell - j + h),$$

thus

$$\begin{aligned} f_X(\lambda) &= \frac{1}{2\pi} \sum_h \sum_{j,\ell} a_j a_\ell \gamma_Y(\ell - j + h) e^{-i\lambda(\ell - j + h)} e^{i\lambda\ell} e^{-i\lambda j} \\ &= \frac{1}{2\pi} \sum_j a_j e^{i\lambda j} \sum_\ell a_\ell e^{-i\lambda\ell} \sum_k \gamma_Y(k) e^{-i\lambda k} \\ &= \left| \sum_j a_j e^{i\lambda j} \right|^2 f_Y(\lambda). \end{aligned}$$

◀ Return

## Corner method

The orders  $p$  and  $q$  of an  $\text{ARMA}(p, q)$  are characterized by the autocorrelation function:

$$\rho(h) = \sum_{i=1}^p \phi_i \rho(h-i), \quad \forall h > q.$$

Let  $D(i, j)$  the  $j \times j$  Toeplitz matrix

$$D(i, j) = \begin{pmatrix} \rho(i) & \rho(i-1) & \cdots & \rho(i-j+1) \\ \rho(i+1) & \ddots & & \vdots \\ \vdots & & \ddots & \rho(i-1) \\ \rho(i+j-1) & \cdots & \rho(i+1) & \rho(i) \end{pmatrix}$$

and

$$\nabla(i, j) = \det D(i, j).$$

## Characterisation of the orders of an ARMA

The  $\nabla(i, j)$  can be obtained from the recursion on  $j$

$$\nabla(i, j)^2 = \nabla(i+1, j)\nabla(i-1, j) + \nabla(i, j+1)\nabla(i, j-1),$$

and by setting  $\nabla(i, 0) = 1$ ,  $\nabla(i, 1) = \rho(|i|)$

### Proposition:

The orders  $p$  and  $q$  are **minimal** (i.e.  $(X_t)$  does not admit an  $\text{ARMA}(p', q')$  representation with  $p' < p$  or  $q' < q$ ) if and only if

$$\left\{ \begin{array}{l} \nabla(i, j) = 0 \quad \text{for all } i > q \text{ and all } j > p, \\ \nabla(i, p) \neq 0 \quad \text{for all } i \geq q, \\ \nabla(q, j) \neq 0 \quad \text{for all } j \geq p. \end{array} \right.$$



## Corner of zeros in the table of the $\nabla(j, i)$

The minimal orders  $p$  and  $q$  are characterized by the table

$i \backslash j$	1	2	. . .	$q$	$q+1$	. . . .
1	$\rho_1$	$\rho_2$	. . .	$\rho_q$	$\rho_{q+1}$	. . . .
.						
.						
.						
$p$				$\times$	$\times$	$\times$ $\times$ $\times$ $\times$
$p+1$				$\times$	0	0 0 0 0
				$\times$	0	0 0 0 0
				$\times$	0	0 0 0 0
				$\times$	0	0 0 0 0

where  $\nabla(j, i)$  is at the intersection of row  $i$  and column  $j$ , and  $\times$  denotes a non-zero item.

## Table of the studentised statistics

In practice, only a finite number of empirical autocorrelations,  $\hat{\rho}(1), \dots, \hat{\rho}(K)$ , are available. This allows to compute the estimates,  $\hat{\nabla}(j, i)$ , of the  $\nabla(j, i)$  for  $i \geq 1$ ,  $j \geq 1$  and  $i+j \leq K+1$ . The table is thus triangular.

The orders  $p$  and  $q$  are characterized by a corner of "small values". But the determinants concern matrices of different size, and are thus not directly comparable.

We thus consider the studentised statistics defined by

$$t(i, j) = \sqrt{n} \frac{\hat{\nabla}(i, j)}{\hat{\sigma}_{\hat{\nabla}(i, j)}}.$$

## Selection of plausible values for the orders

When  $\nabla(i, j) = 0$  the statistics  $t(i, j)$  follows asymptotically a  $\mathcal{N}(0, 1)$  (provided  $EX_t^4$  exists).

One can reject the hypothesis  $\nabla(i, j) = 0$  at the level  $\alpha\%$  if  $|t(i, j)|$  is larger than the  $(1 - \alpha/2)$ -quantile,  $\Phi^{-1}(1 - \alpha/2)$ , of a  $\mathcal{N}(0, 1)$ .  
(1.96 at level 5%)

One can also **detect automatically** a corner of small values in the table of the  $t(i, j)$  if no value at this corner is larger than  $\Phi^{-1}(1 - \alpha/2)$  in absolute value.

This approach (not a formal test) allows to select a **small number of plausible values** for the orders  $p$  and  $q$ .

## Example of use of the corner methode

Simulation of size  $n = 1000$  of an ARMA(2,1):

$$X_t - 0.8X_{t-1} + 0.8X_{t-2} = \epsilon_t - 0.8\epsilon_{t-1}, \quad \epsilon_t \sim \mathcal{N}(0, 1)$$

.p.		.q.	1.	2.	3.	4.	5.	6.	7.	8.	9.	10.	11.	12.
1		17.6	-31.6	-22.6	-1.9	11.5	8.7	-0.1	-6.1	-4.2	0.5	3.5	2.1	
2		36.1	20.3	12.2	8.7	6.5	4.9	4.0	3.3	2.5	2.1	1.8		
3		-7.8	-1.6	-0.2	0.5	0.7	-0.7	0.8	-1.4	1.2	-1.1			
4		5.2	0.1	0.4	0.3	0.6	-0.1	-0.3	0.5	-0.2				
5		-3.7	0.4	-0.1	-0.5	0.4	-0.2	0.2	-0.2					
6		2.8	0.6	0.5	0.4	0.2	0.4	0.2						
7		-2.0	-0.7	0.2	0.0	-0.4	-0.3							
8		1.7	0.8	0.0	0.2	0.2								
9		-0.6	-1.2	-0.5	-0.2									
10		1.4	0.9	-0.2										
11		-0.2	-1.2											
12		1.2												

## Example: automatic detection

ARMA(P,Q) MODELS FOUND WITH GIVEN SIGNIFICANCE LEVEL					
PROBA	CRIT	MODELS FOUND			
0.200000	1.28	( 2, 8)	( 3, 1)	(10, 0)	
0.100000	1.64	( 2, 1)	( 8, 0)		
0.050000	1.96	( 1,10)	( 2, 1)	( 7, 0)	
0.020000	2.33	( 0,11)	( 1, 9)	( 2, 1)	( 6, 0)
0.010000	2.58	( 0,11)	( 1, 8)	( 2, 1)	( 6, 0)
0.005000	2.81	( 0,11)	( 1, 8)	( 2, 1)	( 5, 0)
0.002000	3.09	( 0,11)	( 1, 8)	( 2, 1)	( 5, 0)
0.001000	3.29	( 0,11)	( 1, 8)	( 2, 1)	( 5, 0)
0.000100	3.72	( 0, 9)	( 1, 7)	( 2, 1)	( 5, 0)
0.000010	4.26	( 0, 8)	( 1, 6)	( 2, 1)	( 4, 0)

We find the orders  $(p,q) = (2,1)$  of the simulated model, but also other plausible values. This is not surprising: the ARMA(2,1) is well approximated by several ARMA models, e.g. AR(6), MA(11) or ARMA(1,8) (but the ARMA(2,1), which is more parcimonious in terms of parameters, can be preferred).

## Spectral density

### Definition: Fourier transform of $\gamma_X$

Let  $(X_t)$  a stationary process with **absolutely summable autocovariances**. The spectral density of  $(X_t)$  is

$$\begin{aligned}f_X(\lambda) &= \frac{1}{2\pi} \sum_{h=-\infty}^{+\infty} \gamma_X(h) e^{-i\lambda h} \\&= \frac{1}{2\pi} \sum_{h=-\infty}^{+\infty} \gamma_X(h) \cos(\lambda h) \\&= \frac{1}{2\pi} \gamma_X(0) + \frac{1}{\pi} \sum_{h=1}^{+\infty} \gamma_X(h) \cos(\lambda h), \quad \forall \lambda \in [-\pi; \pi]\end{aligned}$$

## Propositions

- $f_X$  is a real, continuous, even, nonnegative function.
- knowledge of  $f_X$  is equivalent to the knowledge of the sequence  $\{\gamma_X(h)\}$ :

$$\begin{aligned}\gamma_X(h) &= \int_{-\pi}^{\pi} f_X(\lambda) \cos(\lambda h) d\lambda \\ &= \int_{-\pi}^{\pi} f_X(\lambda) e^{-i\lambda h} d\lambda \quad \forall h \in \mathbb{Z}.\end{aligned}$$

- Intuitively, the spectral density decomposes the autocovariances: if  $f_X$  has a peak at frequency  $\lambda$  then  $\gamma_X(h)$  is large for  $h$  close to  $2\pi/\lambda$ .
- The spectral density of a WN is constant (no peak, hence no periodicity).

## Spectral density of linear transform

### Proposition

If  $(Y_t)$  has spectral density  $f_Y$  and  $\sum |a_j| < \infty$ , then  $X_t := \sum_{j=-\infty}^{\infty} a_j Y_{t-j}$  has spectral density

$$f_X(\lambda) = f_Y(\lambda) \left| \sum_j a_j e^{i\lambda j} \right|^2.$$

► Proof

Application: if  $(X_t) \sim \text{ARMA}(p, q)$  then

$$f_X(\lambda) = \frac{\psi(e^{i\lambda}) \psi(e^{-i\lambda})}{\phi(e^{i\lambda}) \phi(e^{-i\lambda})} \frac{\sigma^2}{2\pi}.$$



## Application to the canonical form

If  $(X_t) \sim \text{ARMA}$ ,

$$\Phi(B)(1 - aB)X_t = \Psi(B)(1 - bB)\epsilon_t$$

with  $\Phi(z)\Psi(z) \neq 0$  for  $|z| = 1$  and  $ab \neq 0$ , then

$$\Phi(B)(1 - \frac{1}{a}B)X_t = \Psi(B)(1 - \frac{1}{b}B)\epsilon_t^*$$

with  $\epsilon_t^*$  a WN

Indeed,

$$\epsilon_t^* = \frac{(1 - \frac{1}{a}B)(1 - bB)}{(1 - aB)(1 - \frac{1}{b}B)}\epsilon_t$$

and since  $(1 - e^{i\lambda}/a)(1 - e^{-i\lambda}/a) = |1 - ae^{i\lambda}|^2/a^2$ , we get

$$f_{\epsilon^*}(\lambda) = \frac{b^2}{a^2} \frac{\sigma^2}{2\pi}.$$

## Prediction of stationary series

**Theoretical setup:**  $(X_t)$  a stationary time series, with mean  $E(X_t) = m$  and autocovariance function  $\gamma$  (supposed to be known).

We look for the **best linear combination** of  $1, X_n, X_{n-1}, \dots, X_1$  to predict  $X_{n+h}$  for  $h \geq 1$ .

Denote by  $\hat{X}_{n+h|X_n, \dots, X_1}$  this linear prediction, which has the form:

$$\hat{X}_{n+h|X_n, \dots, X_1} = a_0 + a_1 X_n + \dots + a_n X_1$$

We want to minimize the **Mean Square Error**, MSE:

$$E(X_{n+h} - a_0 - a_1 X_n - \dots - a_n X_1)^2$$

## Examples of predictions with a fixed number of values:

- $\min_{a_0} E(X_{n+h} - a_0)^2 \implies \hat{a}_0 = E(X_t) = m.$

MSE:  $E(X_{n+h} - \hat{a}_0)^2 = \gamma(0).$

- $\min_{a_0, a_1} E(X_{n+h} - a_0 - a_1 X_n)^2?$

Cancel the derivatives with respect to  $a_1$  and  $a_2$ :

$$\begin{cases} E(X_{n+h} - a_0 - a_1 X_n) &= 0 \\ E\{(X_{n+h} - a_0 - a_1 X_n)X_n\} &= 0 \end{cases} \implies \begin{cases} \hat{a}_1 &= \rho(h), \\ \hat{a}_0 &= m(1 - \rho(h)). \end{cases}$$

Linear optimal prediction of  $X_{n+h}$  as a function of  $X_n$  and MSE:

$$\hat{X}_{n+h|X_n} = m(1 - \rho(h)) + \rho(h)X_n.$$

$$E(X_{n+h} - \hat{X}_{n+h|X_n})^2 = \gamma(0)\{1 - \rho(h)^2\}$$

## Prediction using all available values:

The solution of

$$\min_{a_0, \dots, a_n} E(X_{n+h} - a_0 - a_1 X_n - \dots - a_n X_1)^2$$

is obtained for  $\hat{a}_0$  and  $\hat{\mathbf{a}}_n = (\hat{a}_1, \dots, \hat{a}_n)'$  satisfying

$$\Gamma_n \hat{\mathbf{a}}_n = \gamma_n(h), \quad \hat{a}_0 = m \left( 1 - \sum_{i=1}^n \hat{a}_i \right)$$

where

$$\Gamma_n = [\gamma(i-j)]_{i,j=1}^n, \quad \gamma_n(h) = (\gamma(h), \gamma(h+1), \dots, \gamma(h+n-1))'.$$

$$\text{MSE} : E(X_{n+h} - \hat{X}_{n+h|X_n, \dots, X_1})^2 = \gamma(0) - \hat{\mathbf{a}}_n' \gamma_n(h).$$

## Linear prediction using the infinite past

$(X_t)$  stationary time series with mean  $E(X_t) = 0$ .

The problem is to find the **best combination** of  $X_n, X_{n-1}, \dots$ , to predict  $X_{n+h}$ .

Let  $\hat{X}_{n+h|X_n, \dots}$  this prediction, and suppose it has the forme:

$$\hat{X}_{n+h|X_n, \dots} = \sum_{j=1}^{\infty} \alpha_j X_{n+1-j}.$$

The  $\alpha_j$ 's are characterized by

$$E \left\{ \left( X_{n+h} - \sum_{j=1}^{\infty} \alpha_j X_{n+1-j} \right) X_{n+1-i} \right\} = 0, \quad i = 1, 2, \dots$$

$$\iff \gamma(h+i-1) = \sum_{j=1}^{\infty} \alpha_j \gamma(i-j), \quad i = 1, 2, \dots$$

## Using the $MA(\infty)$ representation

$$X_{n+h} = \epsilon_{n+h} + \sum_{j=1}^{\infty} c_j \epsilon_{n+h-j}, \quad \epsilon_t = \text{innovation of } X_t.$$

Project this relation over the infinite past of  $X_n$ :

$$\hat{X}_{n+h|X_n, \dots} = \sum_{j=h}^{\infty} c_j \epsilon_{n+h-j}$$

The prediction error is thus:  $X_{n+h} - \hat{X}_{n+h|X_n, \dots} = \epsilon_{n+h} + \sum_{j=1}^{h-1} c_j \epsilon_{n+h-j}$

In particular:  $X_{n+1} - \hat{X}_{n+1|X_n, \dots} = \epsilon_{n+1}$

Variance of the prediction error:

$$\text{Var}(X_{n+h} - \hat{X}_{n+h|X_n, \dots}) = \sigma^2 \left(1 + \sum_{j=1}^{h-1} c_j^2\right)$$

The previous prediction formula is essentially of theoretical interest because the  $\epsilon$  are not observable, but it allows for simple updating formulas:

$$\hat{X}_{n+h|X_n, \dots} - \hat{X}_{n+h|X_{n-1}, \dots} = c_h \epsilon_n = c_h [X_n - \hat{X}_{n|X_{n-1}, \dots}]$$

## Using the $AR(\infty)$ representation

Since

$$X_{n+h} = \sum_{i=1}^{\infty} b_i X_{n+h-i} + \epsilon_{n+h}$$

we get

$$\hat{X}_{n+h|X_n, \dots} = \sum_{i=1}^{\infty} b_i \hat{X}_{n+h-i|X_n, \dots}$$

with  $\hat{X}_{n+h-i|X_n, \dots} = X_{n+h-i}$  if  $h \leq i$ .

Neglecting the values anterior to  $t = 1$ , we deduce

$$\hat{X}_{n+h|X_n, \dots} \approx \sum_{i=1}^{n+h-1} b_i \hat{X}_{n+h-i|X_n, \dots}.$$



## Using the ARMA form

Suppose  $n > p$  and  $n > q$

$$X_{n+h} = \sum_{i=1}^p \phi_i X_{n+h-i} + \epsilon_{n+h} - \sum_{j=1}^q \psi_j \epsilon_{n+h-j}$$

$$\rightarrow \hat{X}_{n+h|X_n, \dots} = \sum_{i=1}^p \phi_i \hat{X}_{n+h-i|X_n, \dots} - \sum_{j=1}^q \psi_j \hat{\epsilon}_{n+h-j|X_n, \dots}$$

with

$$\begin{aligned} \hat{\epsilon}_{n+h-j|X_n, \dots} &= 0, \quad \text{if } h > j \\ &= \epsilon_{n+h-j}, \quad \text{otherwise} \end{aligned}$$

and  $\hat{X}_{n+h-i|X_n, \dots} = X_{n+h-i} \quad \text{if } h \leq i.$

## Joint use of these formulas:

We aim at computing predictions at horizon  $H$ .

At time  $n$ : we have to compute  $\hat{X}_{n+1|X_n,\dots}, \dots, \hat{X}_{n+H|X_n,\dots}$ .

At time  $n+1$ : we have to  
update the previous predictions:

$$\hat{X}_{n+2|X_{n+1},\dots}, \dots, \hat{X}_{n+H|X_{n+1},\dots}$$

compute a new prediction:  $\hat{X}_{n+H+1|X_{n+1},\dots}$ .

## Several steps

- One can use the  $AR(\infty)$  representation to compute  $\hat{X}_{n+1|X_n, \dots}, \dots, \hat{X}_{n+H|X_n, \dots}$   
 (we will no longer use the observations  $X_1, \dots, X_n$ )
- At time  $n+1$  we observe  $X_{n+1}$ : one can use the updating formula deduced from the  $MA(\infty)$  representation.  
 $\Rightarrow \hat{X}_{n+2|X_{n+1}, \dots}, \dots, \hat{X}_{n+H|X_{n+1}, \dots}$
- Remains to compute  $\hat{X}_{n+H+1|X_{n+1}, \dots}$ : one can use the formula deduced ARMA representation, provided that:
  - $H > q$  ( $\rightarrow \hat{\epsilon}_{n+H+1-j|X_{n+1}, \dots} = 0, \quad \forall j \leq q$ )
  - $H > p$  (we do not need to keep the observations)
- The procedure can be iterated when  $X_{n+2}$  is observed.

## Example: ARMA (1,1)

$$X_t - 0.4X_{t-1} = \epsilon_t - 0.5\epsilon_{t-1}$$

horizon:  $H = 3$

8 Observations (obtained by simulation):

0.480   -0.458   0.427   -0.159   -0.006   0.516   -0.499   0.566

AR( $\infty$ ) representation:

$$X_{n+h} = -0.1 \sum_{i=1}^{\infty} 0.5^{i-1} X_{n+h-i} + \epsilon_{n+h}$$

$$\rightarrow \hat{X}_{9|8,7,\dots} = 0.541, \quad \hat{X}_{10|8,\dots} = 0.592 \quad \hat{X}_{11|8,\dots} = 0.668$$

Then we observe  $X_9 = -0.309$ .

Updating formula:

$$\hat{X}_{9+h|9,\dots} - \hat{X}_{9+h|8,\dots} = c_h[X_9 - \hat{X}_{9|8,7,\dots}],$$

with  $c_1 = -0.1$ ,  $c_2 = -0.04$

$$\rightarrow \hat{X}_{10|9,\dots} = 0.677, \quad \hat{X}_{11|9,\dots} = 0.702$$

Using the ARMA(1,1) representation :

$\hat{X}_{12|9,\dots} = 0.4\hat{X}_{11|9,\dots} = 0.281$  and then repeat ...

◀ Return to the theoretical predictions of an ARMA

## Transformation linéaire d'une série stationnaire

### Propriété

Soit  $(X_t)$  une série **stationnaire** (au second ordre),  $(a_i)_{i \in \mathbb{Z}}$  une suite de réels telle que  $\sum_{i=-\infty}^{\infty} |a_i| < \infty$ . Alors

$$a(B)X_t = \sum_{i=-\infty}^{\infty} a_i B^i X_t = \sum_{i=-\infty}^{\infty} a_i X_{t-i}$$

converge en moyenne quadratique and presque sûrement, [► Proof](#) et définit une série stationnaire.

**Remarque:** En particulier, une combinaison linéaire **finie** des variables  $X_{t-i}$  définit toujours une série stationnaire (donc la stationnarité au second ordre des MA est automatique). [◀ Return](#)

## Autocovariance d'une transformation linéaire

### Propriété

Soit  $(X_t)$  une série stationnaire,  $(a_i)_{i \in \mathbb{Z}}$  une suite de réels telle que  $\sum_{i=-\infty}^{\infty} |a_i| < \infty$ . Alors

$$Y_t = a(B)X_t = \sum_{i=-\infty}^{\infty} a_i B^i X_t = \sum_{i=-\infty}^{\infty} a_i X_{t-i}$$

est une série stationnaire de fonction d'autocovariance

$$\gamma_Y(h) = \sum_{i=-\infty}^{\infty} \sum_{j=-\infty}^{\infty} a_i a_j \gamma_X(h+i-j).$$

► Preuve

## Manipulation des séries en $B$

Les séries en  $B$  appliquées à des processus stationnaires se manipulent **comme des séries entières**. En particulier, if

$\sum_{i=-\infty}^{\infty} |a_i| < \infty$ ,  $\sum_{i=-\infty}^{\infty} |b_i| < \infty$ ,  $a(B) = \sum_{i=-\infty}^{\infty} a_i B^i$  and  $b(B) = \sum_{i=-\infty}^{\infty} b_i B^i$  alors for all processus stationnaire  $(X_t)$

$$\begin{aligned} a(B)b(B)X_t &= \sum_{i=-\infty}^{\infty} a_i \sum_{j=-\infty}^{\infty} b_j B^{i+j} X_t = b(B)a(B)X_t \\ &= \sum_{k=-\infty}^{\infty} \left( \sum_{j=-\infty}^{\infty} a_{k-j} b_j \right) B^k X_t := c(B)X_t. \end{aligned}$$

De plus  $a(B) = b(B) \iff$

$$a_i = b_i, \forall i \iff a(z) = b(z), \forall 1 - \delta < |z| < 1 + \delta$$

for un  $\delta > 0$ .



## Preuve de la convergence de la série

- Par Fubini,

$$E \sum_{i=-\infty}^{\infty} |a_i X_{t-i}| < \infty,$$

donc la série est finie avec probabilité 1

- Pour  $0 < p < q$ ,

$$\left\| \sum_{p \leq |i| \leq q} a_i X_{t-i} \right\|_2 \leq \sum_{p \leq |i| \leq q} |a_i| \|X_{t-i}\|_2 \rightarrow 0$$

quand  $p, q \rightarrow \infty$  donc la série converge en moyenne quadratique par le critère de Cauchy.

## Preuve de la stationnarité

En justifiant les interversions de  $E$  et  $\sum$  par Lebesgue et Fubini, on obtient

$$EY_t = a(B)EX_t = a(1)EX_1 = \sum_{i=-\infty}^{\infty} a_i EX_1$$

puis

$$E(Y_t - EY_1)(Y_{t-h} - EY_1) = \sum_{i=-\infty}^{\infty} \sum_{j=-\infty}^{\infty} a_i a_j E(X_{t-i} - EX_1)(X_{t-j-h} - EX_1).$$

◀ Retour