

MAP 565

Time series analysis : Lecture V

François Roueff

<http://perso.telecom-paristech.fr/~roueff/>

Telecom ParisTech – École Polytechnique

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Outline of the course

- ▷ Stochastic modeling
 - I Random processes.
 - II Spectral representation.
- ▷ Linear models
 - III Linear filtering, innovation process.
 - IV ARMA processes.
 - V Linear forecasting. ←
- ▷ Statistical inference
 - VI Overview of goals and methods.
 - VII Asymptotic statistics in a dependent context.
- ▷ Non-linear models
 - VIII Standard models for financial time series.
 - IX Complements.

← : we are here.

Outline of lectures V

- 1 Linear prediction
 - Prediction VS linear prediction
 - Linear prediction for weakly stationary processes
- 2 An illustrative example with R

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Conditional expectation

Definition : conditional expectation

Let X be a real valued random variable defined on the probability space $(\Omega, \mathcal{F}, \mathbb{P})$. Let \mathcal{G} be a sub- σ -field of \mathcal{F} .

- (a) Suppose $X \in L^2(\Omega, \mathcal{F}, \mathbb{P})$. The conditional expectation of X given \mathcal{G} is defined by

$$\mathbb{E}[X|\mathcal{G}] = \text{proj} (X | L^2(\Omega, \mathcal{G}, \mathbb{P})) .$$

- (b) It is equivalently characterized (in the a.s. sense) by

(i) $\mathbb{E}[X|\mathcal{G}] \in L^1(\Omega, \mathcal{G}, \mathbb{P})$.

(ii) For all $A \in \mathcal{G}$, we have $\mathbb{E}[X\mathbb{1}_A] = \mathbb{E}[\mathbb{E}[X|\mathcal{G}]\mathbb{1}_A]$.

If $X \in L^1(\Omega, \mathcal{F}, \mathbb{P})$, definition (b) remains valid.

Conditional expectation with respect to random variables

If $\mathcal{G} = \sigma(Z_t, t \in T)$, we denote

$$\mathbb{E}[X|Z_t, t \in T] = \mathbb{E}[X|\mathcal{G}] .$$

Basic properties

Conditional density

If (X, Y) admits a density f with respect to $\xi \otimes \xi'$, then, for all real valued g , $\mathbb{E}[g(X)|Y] = \hat{g}(Y)$ with $\hat{g}(y) = \int g(x) f(x|y) \xi(dx)$ and $f(x|y) = f(x, y) / \int f(x', y) \xi(dx')$.

Some standard properties

Let $X \in L^1(\Omega, \mathcal{F}, \mathbb{P})$.

(P-i) If X is \mathcal{G} -measurable, $\mathbb{E}[X|\mathcal{G}] = X$.

(P-ii) If X is independent of \mathcal{G} , $\mathbb{E}[X|\mathcal{G}] = \mathbb{E}[X]$

(P-iii) If Y is $\sigma(\mathcal{G})$ -meas. and $\mathbb{E}[|XY|] < \infty$, $\mathbb{E}[XY|\mathcal{G}] = Y\mathbb{E}[X|\mathcal{G}]$.

(P-iv) If $\mathcal{G} \subset \mathcal{H}$, $\mathbb{E}[\mathbb{E}[X|\mathcal{H}|\mathcal{G}] = \mathbb{E}[X|\mathcal{G}]$ (tower property).

(P-v) If $X = F(Y, Z)$ with Y \mathcal{G} -measurable and Z independent of \mathcal{G} , then $\mathbb{E}[X|\mathcal{G}] = \hat{F}(Y)$, where, for all y , $\hat{F}(y) = \mathbb{E}[F(y, Z)]$.

Prediction VS linear prediction

If $X \in L^2(\Omega, \mathcal{F}, \mathbb{P})$, then $\mathbb{E}[X|\mathcal{G}] = \text{proj}(X | L^2(\Omega, \mathcal{G}, \mathbb{P}))$, hence

$$\mathbb{E}[(X - \mathbb{E}[X|\mathcal{G}])^2] = \inf \left\{ \mathbb{E}[(X - Y)^2] : Y \in L^2(\Omega, \mathcal{G}, \mathbb{P}) \right\} .$$

We say that $\mathbb{E}[X|\mathcal{G}]$ is the **best predictor** of X given \mathcal{G} .

In particular, if $\mathcal{G} = \sigma(Z_t, t \in T)$, then $\mathbb{E}[X|\mathcal{G}]$ can be **any measurable function** of $(Z_t)_{t \in T}$. Thus, in the case where $(Z_t)_{t \in T}$ is an L^2 process,

$$\mathbb{E}[(X - \mathbb{E}[X|Z_t, t \in T])^2] \leq \mathbb{E}[(X - \text{proj}(X | \overline{\text{Span}}(Z_t, t \in T)))^2] ,$$

and the equality only occurs when

$$\mathbb{E}[X|Z_t, t \in T] = \text{proj}(X | \overline{\text{Span}}(Z_t, t \in T)) \quad \text{a.s.}$$

In general, the best **linear predictor** does not achieve the same prediction error as the best **predictor** but is **much easier to determine**.

The Gaussian assumption

Theorem

Let $(X_t)_{t \in T}$ be a Gaussian process. Then, for any $t \in T$ and any countable set $I \subset T$, the conditional expectation of X_t given $(X_s)_{s \in I}$ is in $\overline{\text{Span}}(1, (X_s)_{s \in I})$, that is,

$$\mathbb{E}[X_t | X_s, s \in I] = \text{proj}(X_t | \overline{\text{Span}}(1, X_s, s \in I)) \quad \text{a.s.} \quad (1)$$

In other words,

Under the Gaussian assumption,
best predictor = best linear (or affine) predictor.

Remark

Since Gaussian processes are L^2 processes, we can rely on projections in the Hilbert space L^2 and (1) is equivalent to

$$\text{proj}(X_t | L^2(\Omega, \sigma(X_s, s \in I), \mathbb{P})) = \text{proj}(X_t | \overline{\text{Span}}(1, X_s, s \in I)) .$$

From a countable to a finite I .

The following lemma for general Hilbert spaces implies that it suffices to consider the case where I is finite.

Lemma

Let \mathcal{H} be a Hilbert space and $(E_p)_{p \geq 1}$ be a non-decreasing sequence of closed linear subspaces of \mathcal{H} . Then, for all $x \in \mathcal{H}$,

$$\lim_{p \rightarrow \infty} \text{proj}(x|E_p) = \text{proj}(x|E) \quad \text{with} \quad E = \overline{\bigcup_{p \geq 1} E_p}.$$

Let $I = \{t_1, t_2, t_2, \dots\}$. Then we use that

$$\overline{\text{Span}(1, X_s, s \in I)} = \overline{\bigcup_{p \geq 1} \text{Span}(1, X_{t_k}, k = 1, \dots, p)}$$

and

$$L^2(\Omega, \sigma(X_s, s \in I), \mathbb{P}) = \overline{\bigcup_{p \geq 1} L^2(\Omega, \sigma(X_{t_k}, k = 1, \dots, p), \mathbb{P})}.$$

Proof for a finite I .

Let $\begin{bmatrix} X & Z^T \end{bmatrix}^T$ be a Gaussian vector.

Start with the best linear predictor. Denote

$$\hat{X} = \text{proj}(X | \text{Span}(1, Z)) .$$

We may thus write $X = Y + \hat{X}$ with $Y = X - \hat{X}$, and notice that

$$\mathbb{E}[Y] = \langle Y, 1 \rangle = 0 \quad \text{and} \quad \text{Cov}(Y, Z) = \langle Y, Z \rangle = 0 .$$

On the other hand, since $\hat{X} \in \text{Span}(1, Z)$, we have that $\begin{bmatrix} Y & Z^T \end{bmatrix}^T$ is an affine function of $\begin{bmatrix} X & Z^T \end{bmatrix}^T$ and thus a Gaussian vector. We conclude that Y and Z are independent and, by (P-ii), we get

$$\mathbb{E}[X | Z] = \mathbb{E}[Y] + \mathbb{E}[\hat{X} | Z] = 0 + \hat{X} = \text{proj}(X | \text{Span}(1, Z)) .$$

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Linear prediction : general idea

Basic assumption

Let $X = (X_t)_{t \in \mathbb{Z}}$ be a weakly stationary time series.

Recall the **Wold decomposition**

$$X_t = \text{mean} + \underbrace{\sum_{k \geq 0} \psi_k \epsilon_{t-k}}_{\text{purely non-det. process}} + \text{deterministic process},$$

where $(\epsilon_t)_{t \in \mathbb{Z}}$ is the innovation (white noise) defined by

$$\begin{aligned} \epsilon_t &= X_t - \text{proj}(X_t | \mathcal{H}_{t-1}^X) \\ &= X_t - \lim_{p \rightarrow \infty} \text{proj}(X_t | \mathcal{H}_{t-1,p}^X). \end{aligned}$$

From now on, we only consider the **centered purely non-deterministic part**, so we assume that X is **centered and purely non-deterministic**.

Linear prediction : general idea (cont.)

There are two different ways to consider the problem of linear prediction :

- ▶ **Model-based** linear prediction : a parametric model has been determined (either because the system producing the data is well known and understood or because a model has been **statistically inferred** from historical data). In this case, use the **best linear predictor** of the corresponding model. **Examples of models** :
 - ▶ ARMA(p, q) models.
 - ▶ Dynamic linear models.
 - ▶ Extension to non-linear models (and non-linear prediction).
- ▶ **Direct calculation** of linear prediction coefficients using the autocovariance function γ .

AR(p) model

Consider an AR(p) model

$$X_t = \sum_{k=1}^p \phi_k X_{t-k} + \epsilon_t .$$

Many successful applications:

- ▶ Statistical inference : estimate the parameters of the model and use them for time series analysis.
- ▶ Forecasting : use the AR(p) best linear predictor.
- ▶ Coding : use the AR(p) representation to code, transmit and reconstruct a signal (as in speech coding in the standards of GSM).

Yule Walker equations

Definition : linear prediction coefficients of order p

Let $p \geq 1$. The forward linear prediction coefficients of order p , denoted by $\phi_p^+ = [\phi_{1,p}^+ \ \dots \ \phi_{p,p}^+]$, are defined by

$$\text{proj} \left(X_t \mid \mathcal{H}_{t-1,p}^X \right) = \sum_{k=1}^p \phi_{k,p}^+ X_{t-k} ,$$

which is equivalent to

$$\Gamma_p^+ \phi_p^+ = \gamma_p^+ , \quad (2)$$

$$\text{where } \gamma_p^+ = \text{Cov} \left(X_t, \begin{bmatrix} X_{t-1} \\ \vdots \\ X_{t-p} \end{bmatrix} \right)^T \quad \text{and} \quad \Gamma_p^+ = \text{Cov} \left(\begin{bmatrix} X_{t-1} \\ \vdots \\ X_{t-p} \end{bmatrix} \right)^T$$

Linear prediction coefficients (cont)

Note that we have

$$\gamma_p^+ = \begin{bmatrix} \gamma(1) \\ \vdots \\ \gamma(p) \end{bmatrix} \text{ and } \Gamma_p^+ = \begin{bmatrix} \gamma(0) & \gamma(-1) & \cdots & \gamma(-p+1) \\ \gamma(1) & \gamma(0) & \gamma(-1) & \vdots \\ \vdots & \ddots & \ddots & \ddots \\ \vdots & & & \gamma(-1) \\ \gamma(p-1) & \gamma(p-2) & \cdots & \gamma(1) & \gamma(0) \end{bmatrix}.$$

Moreover the variance of the error is given by

$$\sigma^2(p) := \text{Var} \left(X_t - \text{proj} \left(X_t | \mathcal{H}_{t-1,p}^X \right) \right) = \gamma(0) - (\phi_p^+)^T \overline{\gamma_p^+}. \quad (3)$$

Eq. (2) and (3) are called the **Yule-Walker** equations.

Innovation algorithm

Let $\epsilon_{1,0}^+ = X_1$ and, for all $t \geq 2$,

$$\epsilon_{t,t-1}^+ = X_t - \text{proj}(X_t | \text{Span}(X_s, s = 1, \dots, t-1)) .$$

Then $(\epsilon_{t,t-1}^+)_{t \geq 1}$ is an orthogonal sequence such that

$$\|\epsilon_{t,t-1}^+\|^2 = \sigma_{t-1}^2$$

(which decreases with t) and

$$\text{Span}(X_s, s = 1, \dots, t) = \text{Span}(\epsilon_{s,s-1}^+, s = 1, \dots, t) .$$

Innovation algorithm (cont.)

Denote the prediction coefficients in the innovation basis $(\epsilon_{k,k-1}^+)_{k=1,\dots,p}$ by $\theta_p = (\theta_{k,p})_{k=1,\dots,p}$, that is,

$$\text{proj}(X_{p+1} | \text{Span}(X_s, s = 1, \dots, p)) = \sum_{k=1}^p \theta_{k,p} \epsilon_{k,k-1}^+.$$

Then one gets back to the prediction coefficients in the observation basis recursively by identifying

$$\begin{aligned} \sum_{k=1}^p \phi_{k,p}^+ X_{p+1-k} &= \sum_{k=1}^p \theta_{k,p} \left(X_k - \sum_{j=1}^{k-1} \phi_{j,k-1}^+ X_{k-j} \right) \\ &= \theta_{p,p} X_p + \sum_{k=2}^p \left(\theta_{p+1-k,p} - \sum_{j=1}^{k-1} \theta_{p+1-j,p} \phi_{k-j,p-j}^+ \right) X_{p+1-k} \end{aligned}$$

Algorithm 1: Innovation algorithm

Data: $\gamma(k, j)$, $1 \leq j \leq k \leq K + 1$, X_1, \dots, X_{K+1}

Result: $\epsilon_{1,0}^+, \dots, \epsilon_{K+1,K}^+$, θ_p and σ_p^2 for $p = 1, \dots, K$.

Initialization: set $\sigma_0^2 = \gamma(1, 1)$ and $\epsilon_{1,0}^+ = X_1$.

for $p = 1, \dots, K$ **do**

for $m = 1, \dots, p$ **do**

 Set $\theta_{m,p} = \sigma_{m-1}^{-2} \left(\gamma(p+1, m) - \sum_{j=1}^{m-1} \overline{\theta_{j,m-1}} \theta_{j,p} \sigma_j^2 \right)$

end

 Set

$$\sigma_p^2 = \gamma(p+1, p+1) - \sum_{m=1}^p |\theta_{m,p}|^2 \sigma_{m-1}^2$$

$$\epsilon_{p+1,p}^+ = X_{p+1} - \sum_{m=1}^p \theta_{m,p} \epsilon_{m,m-1}^+.$$

end

Innovation algorithm : numerical complexity

- ▷ $O(p^3)$ operations are needed to compute θ_p and ϕ_p^+ .
- ▷ If X is known to be an $MA(q)$ process, then for all $p \geq q$, we have

$$\theta_p = \begin{bmatrix} 0 & \dots & 0 & \theta_{p-q+1,p} & \dots & \theta_{p,p} \end{bmatrix}^T$$

Hence, in this special case, the innovation algorithm can be performed in $O(p)$ operations.

- ▷ The innovation algorithm is also valid for a non-stationary L^2 sequence $(X_t)_{t \geq 1}$.
- ▷ Exploiting the stationarity to compute ϕ_p^+ , Levinson's Algorithm can be performed in $O(p^2)$ operations.

Partial auto-correlation function

Recall that the sequence $\kappa := (\phi_{p,p}^+)_{p \geq 1}$ is called the **partial autocorrelation** function of X .

$$\text{Then we have } \kappa(1) = \frac{\gamma(1)}{\gamma(0)} = \rho(1) = \frac{\langle X_t, X_{t-1} \rangle}{\|X_t\| \|X_{t-1}\|}.$$

If $p \geq 2$, this formula can be extended as follows.

Denote the **forward** and **backward** linear prediction **errors** by

$$\epsilon_{t,p}^+ = X_t - \text{proj}(X_t | \mathcal{H}_{t-1,p}^X) \quad \text{and} \quad \epsilon_{t,p}^- = X_t - \text{proj}(X_t | \mathcal{H}_{t+p,p}^X)$$

Then we have

$$\kappa(p) = \frac{\langle \epsilon_{t,p-1}^+, \epsilon_{t-p,p-1}^- \rangle}{\|\epsilon_{t,p-1}^+\| \|\epsilon_{t-p,p-1}^-\|} = \frac{\text{Cov}(\epsilon_{t,p-1}^+, \epsilon_{t-p,p-1}^-)}{\sigma_{p-1}^2}.$$
$$\begin{array}{ccccccc} X_{t-p} & & X_{t-p+1} & \dots & X_{t-1} & & X_t \\ & & \underbrace{\hspace{10em}} & & & & \\ & & \mathcal{H}_{t-1,p-1}^X & & & & \end{array}$$

Algorithm 2: Levinson-Durbin algorithm.

Data: $\gamma(k)$, $k = 0, \dots, K$

Result: $\{\phi_{m,p}^+\}_{1 \leq m \leq p, 1 \leq p \leq K}$, $\kappa(1), \dots, \kappa(K)$

Initialization: set $\kappa(1) = \phi_{1,1}^+ = \gamma(1)/\gamma(0)$ and $\sigma_1^2 = \gamma(0)(1 - \kappa(1)^2)$.

for $p = 1, 2, \dots, K - 1$ **do**

Set

$$\kappa(p+1) = \sigma_p^{-2} \left(\gamma(p+1) - \sum_{k=1}^p \phi_{k,p}^+ \gamma(p+1-k) \right)$$

$$\sigma_{p+1}^2 = \sigma_p^2 (1 - \kappa(p+1)^2)$$

$$\phi_{p+1,p+1}^+ = \kappa(p+1)$$

for $m \in \{1, \dots, p\}$ **do**

Set

$$\phi_{m,p+1}^+ = \phi_{m,p}^+ - \kappa(p+1) \overline{\phi_{p+1-m,p}^+}.$$

end

end

1 Linear prediction

2 An illustrative example with R


```
#####
#           Linear Predictive Coding           #
#####
quantize <- function(x,corder=3){
# code input x with quantiz. levels given by corder
if (length(corder)>1) # corder are quant. levels
{
  ql <- corder
} else # set quant. levels from normal quantiles
{
  ql <- qnorm(seq(from=0,to=1,by=1/corder),
              mean = mean(x), sd = sqrt(3*var(x)))
}
xc <- NULL
for (t in 1:(length(x))){
  xc= c(xc,ql[which.min(abs(ql-x[t]))])}
return(list(xc=xc,ql=ql))
}

lpcoding <- function(x,corder=3,rorder=10){
# estimation of ar parameters
ac <- acf(x,type=c('covariance'),plot= FALSE)
arc <- acf2AR(ac$acf[1:(rorder+1)])
arc <- arc[nrow(arc),]
# residuals computation
res <- NULL
xinitz <- c(rep(0,rorder),x)
for (t in ((rorder+1):length(xinitz))){
  res <- c(res,xinitz[t]-
           t(as.vector(xinitz[(t-1):(t-rorder)]))
           %*% as.vector(arc))
}
# coded residuals
resc <- quantize(res, corder=corder)
# reconstructed time series from coded residuals
return(list(xc=filter(resc$xc,arc,method='recursive'),
           ar=arc))
}
```

```

}
lpcodingblocks <- function(x,bt=0.02,freq,
                           corder=3,rorder=10){
  bl <- floor(bt*freq) # block length
  rorder <- min(c(rorder,floor(bl/5)))
  xc <- NULL
  for (k in 1:floor(length(x)/bl)){
    xc <- c(xc,lpcoding(x[((k-1)*bl+1):(k*bl)],
                        corder=corder,rorder=rorder)$xc)
  }
  return(xc)
}
# get the original speech audio sample
require(audio)
X <- load.wave('/home/roueff/data/dataset/audio/speech/3meninaboat.wav')
fr <- X$rate
subsamp <- 2**3
extract <- ts(X[seq(from=1,to=length(X),by=subsamp)]/max(abs(X)),
              frequency=fr/subsamp)
ts.plot(extract)

# AR Coding and direct coding
extractc <- ts(lpcodingblocks(extract,
                              freq=frequency(extract),
                              corder=5,rorder=20),
              start=start(extract),
              frequency=frequency(extract))

extractbadc <- ts(quantize(extract,corder=5)$xc,
                 start=start(extract),
                 frequency=frequency(extract))

# plot coded signals
lines(extractc,col=2)
lines(extractbadc,col=3)
# plot within a time window

```

```

for (term in c('', 'c', 'badc')){
  eval(parse(text=paste('we', term,
    ' <- window(extract',
    term, ', start=3.5, end=4)',
    sep=''))))
}
ts.plot(we)
lines(wec, col=2)
lines(webadc, col=3)
# save wav files
require(audio)
save.wave(audioSample(extract,
  rate=frequency(extract)),
  '/tmp/3.wav')
save.wave(audioSample(extractc,
  rate=frequency(extract)),
  '/tmp/3c.wav')
save.wave(audioSample(extractbadc,
  rate=frequency(extract)),
  '/tmp/3badc.wav')

```