

Time Series Analysis

1. Stationary ARMA models

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- A *time series* is a sequence of data points X_t indexed a discrete set of (ordered) dates t , where $-\infty < t < \infty$.
- Each X_t can be a simple number or a complex multi-dimensional object (vector, matrix, higher dimensional array, or more general structure).
- We will be assuming that the times t are equally spaced throughout, and denote the time increment by h (e.g. second, day, month). Unless specified otherwise, we will be choosing the units of time so that $h = 1$.
- Typically, time series exhibit significant irregularities, which may have their origin either in the nature of the underlying quantity or imprecision in observation (or both).
- Examples of time series commonly encountered in finance include:
 - (i) prices,
 - (ii) returns,
 - (iii) index levels,
 - (iv) trading volumes,
 - (v) open interests,
 - (vi) macroeconomic data (inflation, new payrolls, unemployment, GDP, housing prices, ...)

- For modeling purposes, we assume that the elements of a time series are random variables on some underlying probability space.
- *Time series analysis* is a set of mathematical methodologies for analyzing observed time series, whose purpose is to extract useful characteristics of the data.
- These methodologies fall into two broad categories:
 - (i) *non-parametric*, where the stochastic law of the time series is not explicitly specified;
 - (ii) *parametric*, where the stochastic law of the time series is assumed to be given by a model with a finite (and preferably tractable) number of parameters.
- The results of time series analysis are used for various purposes such as
 - (i) data interpretation,
 - (ii) forecasting,
 - (iii) smoothing,
 - (iv) back filling, ...
- We begin with stationary time series.

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- A time series (model) is *stationary*, if for any times $t_1 < \dots < t_k$ and any τ the joint probability distribution of $(X_{t_1+\tau}, \dots, X_{t_k+\tau})$ is identical with the joint probability distribution of $(X_{t_1}, \dots, X_{t_k})$. lag
- A stationary time series model is *ergodic* if

$$\lim_{T \rightarrow \infty} \frac{1}{T} \sum_{1 \leq k \leq T} X_{t+k} = \mu, \quad (1)$$

i.e. if the time average of X_t is equal to the (ensemble) average.

- Ergodicity is a desired property of a financial time series, as we are always faced with a single realization of a process rather than an ensemble of alternative outcomes.
- The limit in (1) is usually understood in the sense of squared mean convergence.
- The notions of stationarity and ergodicity are hard to verify in practice. Luckily, there is a more practical concept.

Autocovariance and stationarity

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- A time series is covariance-stationary (a.k.a. weakly stationary), if:

- (i) $E(X_t) = \mu$ is a constant,
- (ii) For any τ , the autocovariance $\text{Cov}(X_s, X_t)$ is time translation invariant,

$$\text{Cov}(y_t, y_{t+\tau}) = E[y_t y_{t+\tau}] - \mu^2 \quad \text{Cov}(X_{s+\tau}, X_{t+\tau}) = \text{Cov}(X_s, X_t), \quad (2)$$

i.e. $\text{Cov}(X_s, X_t)$ depends only on the difference $t - s$. We will write it as Γ_{t-s} .

- For covariance stationary series, $\Gamma_{-t} = \Gamma_t$ (show it!).
- Notice that $\Gamma_0 = \text{Var}(X_t)$.
- The autocorrelation function of a time series is defined as

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$$R_{s,t} = \frac{\text{Cov}(X_s, X_t)}{\sqrt{\text{Var}(X_s)}\sqrt{\text{Var}(X_t)}}. \quad (3)$$

- For covariance-stationary time series, $R_{s,t} = R_{t-s}$, and

$$R_t = \frac{\Gamma_t}{\Gamma_0}. \quad (4)$$

Autocovariance and stationarity

- Note that μ , Γ , and R are usually unknown, and are estimated from sample data. The estimated sample mean $\hat{\mu}$, autocovariance $\hat{\Gamma}$, and autocorrelation \hat{R} are calculated as follows.
- Consider a finite sample X_0, X_1, \dots, X_T . Then

$$\begin{aligned}\hat{\mu} &= \frac{1}{T} \sum_{t=1}^T X_t, \\ \hat{\Gamma}_t &= \begin{cases} \frac{1}{T} \sum_{j=t+1}^T (X_j - \hat{\mu})(X_{j-t} - \hat{\mu}), & \text{for } t = 0, 1, \dots, T-1, \\ \hat{\Gamma}_{-t}, & \text{for } t = -1, \dots, -(T-1). \end{cases} \\ \hat{R}_t &= \frac{\hat{\Gamma}_t}{\hat{\Gamma}_0}.\end{aligned}\tag{5}$$

- Notice that this method allows us to compute up to $T - 1$ estimated sample autocorrelations.

- For practical applications, it is convenient to model a time series as a discrete-time stochastic process with a small number of parameters.
- Time series models have typically the following structure:

$$X_t = p_t + m_t + \varepsilon_t, \quad (6)$$

where the three components on the RHS have the following meaning:

- p_t is a periodic function called the *seasonality*,
 - m_t is a slowly varying process called the *trend*,
 - ε_t is a stochastic component called the *error* or *disturbance*.
- Classic linear time series models fall into three broad categories:
 - *autoregressive*,
 - *moving average*,
 - *integrated*,
- and their combinations.

White noise

- The source of randomness in the models discussed in these lectures is *white noise*. It is a process specified as follows:

$$X_t = \varepsilon_t, \quad (7)$$

where $\varepsilon_t \sim N(0, \sigma^2)$ are i.i.d. (= independent, identically distributed) normal random variables.

- Note that

$$\begin{aligned} E(\varepsilon_t) &= 0, \\ \text{Cov}(\varepsilon_s, \varepsilon_t) &= \begin{cases} \sigma^2, & \text{if } s = t, \\ 0, & \text{otherwise.} \end{cases} \end{aligned} \quad (8)$$

- The white noise process is stationary and ergodic (show it!).
- The white noise process with *linear drift*

$$X_t = at + b + \varepsilon_t, \quad a \neq 0, \quad (9)$$

is not stationary, as $E(X_t) = at + b$.

Autoregressive model $AR(1)$

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- The first class of models that we consider are the *autoregressive models* $AR(p)$. Their key characteristic is that the current observation is directly correlated with the lagged p observations.
- The simplest among them is $AR(1)$, the autoregressive model with a single lag.
- The model is specified as follows:

$$X_t = \alpha + \beta X_{t-1} + \varepsilon_t.$$

- Here, $\alpha, \beta \in \mathbb{R}$, and $\varepsilon_t \sim N(0, \sigma^2)$ is a white noise.

- A particular case of the $AR(1)$ model is the *random walk model*, namely

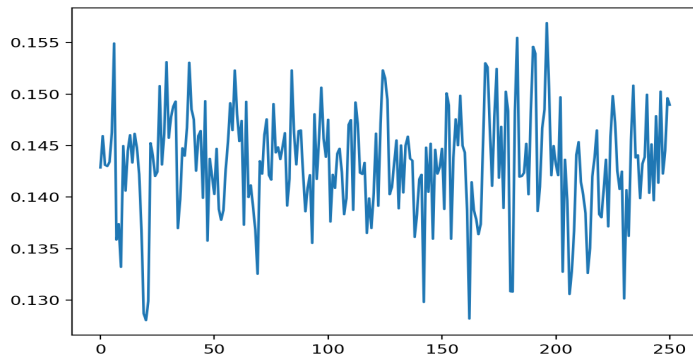
$$X_t = X_{t-1} + \varepsilon_t,$$

in which the current value of X is the previous value plus a “white noise” disturbance.

$$\begin{aligned} & \stackrel{(10)}{=} \alpha + \beta (\alpha + \beta X_{t-2} + \varepsilon_{t-1}) + \varepsilon_t \\ & \stackrel{=}{=} \frac{\alpha}{1-\beta} + \varepsilon_t + \beta \varepsilon_{t-1} + \beta^2 \varepsilon_{t-2} + \dots \end{aligned}$$

Autoregressive model $AR(1)$

- The graph below shows a simulated $AR(1)$ time series with the following choice of parameters: $\alpha = 0.1$, $\beta = 0.3$, $\sigma = 0.005$.



Autoregressive model $AR(1)$

- Here is the code snippet used to generate this graph in Python:

```
import numpy as np
import matplotlib.pyplot as plt
from statsmodels.tsa.arima_model import ARMA

alpha=0.1
beta=0.3
sigma=0.005

#Simulate AR(1)
T=250
x0=alpha/(1-beta)
x=np.zeros(T+1)
x[0]=x0
eps=np.random.normal(0.0, sigma, T)
for i in range(1, T+1):
    x[i]=alpha+beta*x[i-1]+eps[i-1]

#Take a look at the simulated time series
plt.plot(x)
plt.show()
```

Autoregressive model $AR(1)$

- Let us investigate the circumstances under which an $AR(1)$ process is covariance-stationary.
- For $\mu = E(X_t)$ to be independent of t we must have from (10):

$$\mu = \alpha + \beta\mu.$$

This equation has a solution iff $\beta \neq 1$ (except for the random walk case corresponding to $\alpha = 0, \beta = 1$). In this case,

$$\mu = \frac{\alpha}{1 - \beta}. \quad (11)$$

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- Let us now compute the autocovariance. To this end, we rewrite (10) as

$$X_t - \mu = \beta(X_{t-1} - \mu) + \varepsilon_t. \quad (12)$$

Notice that the two terms on the RHS of this equation are independent of each other.

Autoregressive model $AR(1)$

- For $\Gamma_0 = \text{Var}(X_t)$ to be independent of t , this implies that

$$\Gamma_0 = \beta^2 \Gamma_0 + \sigma^2,$$

and so

$$\Gamma_0 = \frac{\sigma^2}{1 - \beta^2}. \quad (13)$$

- Since $\Gamma_0 > 0$, this equation implies that $|\beta| < 1$.
- Multiplying (12) by $X_{t-1} - \mu$, we find that $\Gamma_1 = \beta \Gamma_0$. Iterating, we find that

$$\Gamma_k = \beta^k \Gamma_0, \quad (14)$$

with Γ_0 given by (14). The autocorrelation function is decaying exponentially fast as a function of lag between two observations.

- In conclusion, the condition for a $AR(1)$ process to be covariance-stationary is that $|\beta| < 1$.

$$\begin{aligned} \Gamma_0 &= E[(\varepsilon_t + \beta \varepsilon_{t-1} + \beta^2 \varepsilon_{t-2} + \dots) \\ &\quad (\varepsilon_{t-j} + \beta \varepsilon_{t-j-1} + \beta^2 \varepsilon_{t-j-2} + \dots)] \\ &= \sigma^2 (\phi^j + \phi^{j+2} + \phi^{j+4} + \dots) \\ &= \phi^j \frac{\sigma^2}{1 - \phi^2} = \phi^j \Gamma_0 \end{aligned}$$

Autoregressive model $AR(1)$

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- The $AR(1)$ with $|\beta| < 1$ has a natural interpretation that can be gleaned from the following “explicit” representation of X_t . Namely, iterating (10) we find that:

$$\begin{aligned} X_t &= \alpha + \beta X_{t-1} + \varepsilon_t \\ &= \alpha(1 + \beta) + \beta^2 X_{t-2} + \varepsilon_t + \beta \varepsilon_{t-1} \\ &= \dots \\ &= \alpha(1 + \beta + \dots + \beta^{L-1}) + \beta^L X_{t-L} + \varepsilon_t + \beta \varepsilon_{t-1} + \dots + \beta^{L-1} \varepsilon_{t-L+1} \\ &= \mu(1 - \beta^L) + \beta^L X_{t-L} + \sqrt{\Gamma_0(1 - \beta^{2L-1})} \xi_t \end{aligned} \tag{15}$$

where $\xi_t \sim N(0, 1)$.

- This implies that

$$\begin{aligned} E(X_t | X_{t-L}) &= \mu(1 - \beta^L) + \beta^L X_{t-L}, \\ \text{Var}(X_t | X_{t-L}) &= \Gamma_0(1 - \beta^{2L-1}). \end{aligned} \tag{16}$$

$$y_t = \mu + \sum_{j=0}^{\infty} \psi_j \varepsilon_{t-j}$$

Autoregressive model $AR(1)$

- Since $\beta^L \rightarrow 0$ exponentially fast, for large L we have

$$X_t \approx \mu + \sqrt{\Gamma_0} \xi_t. \quad (17)$$

- In other words, the $AR(1)$ model describes a *mean reverting* time series. After a large number of observations, X_t takes the form (17), i.e. it is equal to its mean value plus a Gaussian noise.
- The rate of convergence to this limit is given by $|\beta|$: the smaller this value, the faster X_t reaches its limit behavior.
- The next question is: given a set of observations, how do we determine the values of the parameters α , β , and σ in (10)?

Maximum likelihood estimation

- *Maximum likelihood estimation* (MLE) is a commonly used method of estimating the parameters of a statistical model given a set of observations.
- It is based on the premise that the best choice of the parameter values should maximize the likelihood of making the observations given these parameters.
- Given a statistical model with parameters $\theta = (\theta_1, \dots, \theta_d)$, and a set of data $y = (y_1, \dots, y_N)$, we construct the *likelihood function* $\mathcal{L}(\theta|y)$, which links the model with the data in such a way as if the data were drawn from the assumed model.
- In practice, $\mathcal{L}(\theta|y)$ is the joint probability density function (PDF) $p(y|\theta)$ under the model, evaluated at the observed values.
- In particular, if the observations y_i are independent, then

$$\mathcal{L}(\theta|y) = \prod_{i=1}^N p(y_i|\theta), \quad (18)$$

where $p(y_i|\theta)$ denotes the PDF of a single observation.

Maximum likelihood estimation

- The value θ^* that maximizes $\mathcal{L}(\theta|y)$ serves as the best fit between the model specification and the data.
- It is usually more convenient to consider the *log likelihood function* (LLF) $-\log \mathcal{L}(\theta|y)$. Then, θ^* is the value at which the LLF attains its minimum.
- As an illustration, consider a sample $y = (y_1, \dots, y_N)$ drawn from the normal distribution $N(\mu, \sigma^2)$. Its likelihood function is given by

$$\mathcal{L}(\theta|y) = (2\pi\sigma^2)^{-N/2} \prod_{i=1}^N \exp\left(-\frac{(y_i - \mu)^2}{2\sigma^2}\right), \quad (19)$$

and the LLF is

$$-\log \mathcal{L}(\theta|y) = \frac{1}{2} N \log \sigma^2 + \frac{1}{2\sigma^2} \sum_{i=1}^N (y_i - \mu)^2 + \text{const.} \quad (20)$$

- Taking the μ and σ derivatives and setting them to 0, we readily find that the MLE estimates of μ and σ are

$$\begin{aligned}\mu^* &= \frac{1}{N} \sum_{i=1}^N y_i, \\ \sigma^* &= \frac{1}{N} \sum_{i=1}^N (y_i - \mu^*)^2.\end{aligned}\tag{21}$$

respectively.

- Note that, while μ^* is *unbiased*, the estimator σ^* is *biased* (N in the denominator above, rather than the usual $N - 1$).
- The fact that the MLE estimator of a parameter is biased is a common occurrence. One can show, however, that MLE estimators are *consistent*, i.e. in the limit $N \rightarrow \infty$ they converge to the appropriate value.
- Going forward, we will use the notation $\hat{\theta}$ rather than θ^* for the MLE estimators.

- Consider now the $AR(1)$ model and a time series of data x_0, \dots, x_T , believed to be drawn from this model. The easiest way to construct the likelihood function is to focus on the conditional PDF $p(x_1, \dots, x_T | x_0, \theta)$. This leads to the *conditional* MLE method.

- Let

$$\hat{\varepsilon}_t = x_t - \alpha - \beta x_{t-1}, \quad (22)$$

for $t = 1, \dots, T$, be the disturbances implied from the data. According to the model specification, each $\hat{\varepsilon}_t$ is independently drawn from $N(0, \sigma^2)$, and thus

$$\begin{aligned} p(x_1, \dots, x_T | x_0, \theta) &= \frac{1}{(2\pi\sigma^2)^{T/2}} \exp\left(-\frac{1}{2\sigma^2} \sum_{t=1}^T \varepsilon_t^2\right) \\ &= \frac{1}{(2\pi\sigma^2)^{T/2}} \exp\left(-\frac{1}{2\sigma^2} \sum_{t=1}^T (x_t - \alpha - \beta x_{t-1})^2\right) \end{aligned} \quad (23)$$

- Hence the LLF is given by

$$-\log \mathcal{L}(\theta | y) = \frac{1}{2} T \log \sigma^2 + \frac{1}{2\sigma^2} \sum_{t=0}^{T-1} (x_{t+1} - \alpha - \beta x_t)^2 + \text{const.} \quad (24)$$

- Minimizing this function yields:

$$\begin{pmatrix} \hat{\alpha} \\ \hat{\beta} \end{pmatrix} = \begin{pmatrix} T & \sum_{t=0}^{T-1} x_t \\ \sum_{t=0}^{T-1} x_t & \sum_{t=0}^{T-1} x_t^2 \end{pmatrix}^{-1} \begin{pmatrix} \sum_{t=0}^{T-1} x_{t+1} \\ \sum_{t=0}^{T-1} x_t x_{t+1} \end{pmatrix}, \quad (25)$$

$$\hat{\sigma}^2 = \frac{1}{T} \sum_{t=1}^T (x_t - \hat{\alpha} - \hat{\beta} x_{t-1})^2.$$

- This can also be explicitly rewritten as

$$\hat{\beta} = \frac{\sum_{t=0}^{T-1} (x_t - \hat{x})(x_{t+1} - \hat{x}_+)}{\sum_{t=0}^{T-1} (x_t - \hat{x})^2}, \quad (26)$$

$$\hat{\alpha} = \hat{x}_+ - \hat{\beta} \hat{x},$$

where

$$\hat{x} = \sum_{t=0}^{T-1} x_t, \quad \hat{x}_+ = \sum_{t=0}^{T-1} x_{t+1}. \quad (27)$$

- The *exact* MLE method attempts to infer the likelihood of x_0 from the probability distribution. Since $x_0 \sim N(\mu, \Gamma_0)$,

$$p(x_0|\theta) = \sqrt{\frac{1-\beta^2}{2\pi\sigma^2}} \exp\left(-\frac{(x_0 - \alpha/(1-\beta))^2}{2\sigma^2/(1-\beta^2)}\right). \quad (28)$$

- On the other hand, for $t = 1, \dots, T$,

$$p(x_t|x_{t-1}, \dots, x_1, \theta) = \frac{1}{\sqrt{2\pi\sigma^2}} \exp\left(-\frac{(x_t - \alpha - \beta x_{t-1})^2}{2\sigma^2}\right). \quad (29)$$

- From the definition of conditional probability we have the following identity:

$$p(x_0, x_1, \dots, x_T|\theta) = p(x_0|\theta) \prod_{t=1}^T p(x_t|x_{t-1}, \dots, x_1, \theta). \quad (30)$$

- Therefore, the LLF is given by

$$\begin{aligned} -\log \mathcal{L}(\theta|x) = & \frac{1}{2} \log \frac{\sigma^2}{1 - \beta^2} + \frac{1}{2} T \log \sigma^2 \\ & + \frac{(x_0 - \alpha/(1 - \beta))^2}{2\sigma^2/(1 - \beta^2)} + \frac{1}{2\sigma^2} \sum_{t=1}^T (x_t - \alpha - \beta x_{t-1})^2 + \text{const.} \end{aligned} \quad (31)$$

- Unlike the conditional case, the minimum of the exact LLF cannot be calculated in closed form, and the calculation has to be done by means of a numerical search.

- Here is the Python code snippet implementing the MLE for $AR(1)$:

```
#Conditional MLE estimate
y=x[0:T]
yp=x[1:(T+1)]
m=np.sum(y)/T
mp=np.sum(yp)/T
betaCMLE=np.inner(y-m,yp-mp)/np.inner(y-m,y-m)
alphaCMLE=mp-betaCMLE*m
sigmaCMLE=np.sqrt(np.inner(yp-betaCMLE*y-alphaCMLE,
                           yp-betaCMLE*y-alphaCMLE)/T)
```

- Alternatively, one can use statsmodels functions:

```
#MLE estimate with statsmodels
model=ARMA(x,order=(1,0)).fit(method='mle')
alphaMLE=model.params[0]
betaMLE=model.params[1]
sigmaMLE=np.std(model.resid)
```


Second order autoregressive model $AR(2)$

- A second order autoregressive model $AR(2)$ model is specified as follows:

$$X_t = \alpha + \beta_1 X_{t-1} + \beta_2 X_{t-2} + \varepsilon_t, \quad (32)$$

where $\alpha, \beta_1, \beta_2 \in \mathbb{R}$, and $\varepsilon_t \sim N(0, \sigma^2)$ is a white noise.

- Under this specification, the state variable depends on its two lags (rather than one lag as in $AR(1)$).
- Let us determine the conditions under which the model is covariance-stationary.
- From the requirement that $E(X_t) = \mu$,

$$\mu = \frac{\alpha}{1 - \beta_1 - \beta_2}, \quad (33)$$

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$$X_t = \phi_1 X_{t-1} + \phi_2 X_{t-2} + \varepsilon_t$$

and so we can rewrite (32) in the following form:

$$X_t - \mu = \beta_1 (X_{t-1} - \mu) + \beta_2 (X_{t-2} - \mu) + \varepsilon_t. \quad (34)$$

$$X_{t-1} = 1 X_{t-1} + 0 X_{t-2} + 0.$$

$$\begin{pmatrix} X_t \\ X_{t-1} \end{pmatrix} = \underbrace{\begin{pmatrix} \phi_1 & \phi_2 \\ 1 & 0 \end{pmatrix}}_{\text{matrix}} \begin{pmatrix} X_{t-1} \\ X_{t-2} \end{pmatrix} + \begin{pmatrix} \varepsilon_t \\ 0 \end{pmatrix}$$

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Second order autoregressive model AR(2)

$$\Rightarrow \begin{pmatrix} X_t \\ X_{t-1} \end{pmatrix} = \bar{F}^{j+1} \begin{pmatrix} X_{t-j-1} \\ X_{t-j-2} \end{pmatrix} + \bar{F}^j \begin{pmatrix} \varepsilon_{t-j} \\ 0 \end{pmatrix} + \bar{F}^{j-1} \begin{pmatrix} \varepsilon_{t-j+1} \\ 0 \end{pmatrix} + \dots$$

- Multiplying (34) by $X_{t-j} - \mu$, for $j = 0, 1, 2$, and calculating expectations, we find that

$$\Gamma_k = \begin{cases} \beta_1 \Gamma_1 + \beta_2 \Gamma_2 + \sigma^2, & \text{if } k = 0, \\ \beta_1 \Gamma_{k-1} + \beta_2 \Gamma_{k-2}, & \text{if } k = 1, 2. \end{cases} \quad (35)$$

$$X_t = (\bar{F}^{j+1})_{1,1} X_{t-j-1} + (\bar{F}^{j+1})_{1,2} X_{t-j-2} + (\bar{F}^j)_{1,1} \varepsilon_{t-j} + \dots$$

This identity is called the *Yule-Walker equation* for the autocovariance.

- Dividing (57) by Γ_0 yields the Yule-Walker equation for the autocorrelation:

$$R_k = \beta_1 R_{k-1} + \beta_2 R_{k-2}, \quad (36)$$

for $k = 1, 2$.

- This equation allows us calculate explicitly the ACF for AR(2).
- Namely, plugging in $k = 1$ and remembering that $R_{-1} = R_1$ yields $R_1 = \beta_1 + \beta_2 R_1$, or

$$R_1 = \frac{\beta_1}{1 - \beta_2}. \quad (37)$$

(Wold decomp).

eigenval of \bar{F}

$$\det \begin{pmatrix} \phi_1 - \lambda & \phi_2 \\ 1 & -\lambda \end{pmatrix} = 0 \Rightarrow \lambda^2 - \phi_1 \lambda - \phi_2 = 0 \Rightarrow 1 - \phi_1 z - \phi_2 z^2 = 0 \Rightarrow$$

$$(z = \frac{1}{\lambda})$$

(lag poly)

AR(P) is station if roots of polynomial lie outside the unit circle.

$|\lambda|$ needs to be < 1
 $\Rightarrow |z| > 1$

Second order autoregressive model $AR(2)$

- Plugging in $k = 2$ yields $R_2 = \beta_1 R_1 + \beta_2$, or

$$R_2 = \beta_2 + \frac{\beta_1^2}{1 - \beta_2}. \quad (38)$$

- Finally, substituting $k = 0$ in (34) yields

$$\Gamma_0 = (\beta_1 R_1 + \beta_2 R_2)\Gamma_0 + \sigma^2. \quad (39)$$

Solving this, we obtain

$$\Gamma_0 = \frac{(1 - \beta_2)\sigma^2}{(1 + \beta_2)((1 - \beta_2)^2 - \beta_1^2)}. \quad (40)$$

Lag operators and characteristic roots

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- We have not yet addressed the question under what condition is an $AR(2)$ time series covariance-stationary. We will now introduce the concepts that will settle this issue and will allow us to formulate criteria for stationarity for more general models,
- Let us define the *lag operator* L as a (linear) mapping:

$$LX_t = X_{t-1}. \quad (41)$$

In other words, the lag operator shifts the time index back by one unit.

- Applying the lag operator k times shifts the time index by k units:

$$L^k X_t = X_{t-k}. \quad (42)$$

We refer to L^k as the k -th power of L .

- Finally, if $\psi(z) = \psi_0 + \psi_1 z + \dots + \psi_n z^n$ is a polynomial in z , we associate with it an operator $\psi(L)$ defined by

$$\psi(L) = \psi_0 + \psi_1 L + \dots + \psi_n L^n. \quad (43)$$

$$AR(1) \quad y_t = \phi y_{t-1} + \varepsilon_t$$

$$(1 - \phi L) y_t = \varepsilon_t$$

$$y_t = \frac{1}{1 - \phi L} \varepsilon_t$$

$$= \sum_{j=0}^{\infty} (\phi L)^j \varepsilon_t$$

$$= \sum_{j=0}^{\infty} \phi^j \varepsilon_{t-j}$$

Lag operators and characteristic roots

- Notice that equation (32) can be stated as

$$\psi(L)X_t = \alpha + \varepsilon_t, \quad (44)$$

where $\psi(z) = 1 - \beta_1 z - \beta_2 z^2$.

- Solving this equation amounts to finding the inverse $\psi(L)^{-1}$ of $\psi(L)$:

$$X_t = \frac{\alpha}{\psi(1)} + \psi(L)^{-1} \varepsilon_t. \quad (45)$$

- Suppose that we can write $\psi(L)^{-1}$ as an infinite series

$$\psi(L)^{-1} = \sum_{j=0}^{\infty} \gamma_j L^j, \quad (46)$$

with

$$\sum_{j=0}^{\infty} |\gamma_j| < \infty. \quad (47)$$

Lag operators and characteristic roots

- Then

$$X_t = \frac{\alpha}{\psi(1)} + \sum_{j=0}^{\infty} \gamma_j \varepsilon_{t-j}, \quad (48)$$

with

$$E(X_t) = \frac{\alpha}{\psi(1)}, \quad (49)$$

and

$$\text{Cov}(X_t, X_{t+k}) = \sum_{j=0}^{\infty} \gamma_j \gamma_{j+k}, \text{ for } k \geq 0, \quad (50)$$

independently of t . The series is thus covariance-stationary.

- In the case of $AR(1)$, $\psi(L) = 1 - \beta L$, it is clear that the geometric series does the job:

$$(1 - \beta L)^{-1} = \sum_{j=0}^{\infty} \beta^j L^j, \quad (51)$$

- Condition (47) holds as long as $|\beta| < 1$. Another way of saying this is that the root $z_1 = 1/\beta$ of $1 - \beta z$ lies outside of the unit circle.

Lag operators and characteristic roots

- Now, if $\psi(z)$ is a polynomial with non-zero roots z_1, \dots, z_n . Then

$$\psi(L) = (-1)^n \left(\prod_{j=1}^n z_j \right) \prod_{j=1}^n (1 - z_j^{-1} L). \quad (52)$$

- If each of the roots z_j (they may be complex) lies outside of the unit circle, i.e. $|z_j^{-1}| < 1$, then we can invert $\psi(L)$ by applying (51) to each factor in the product above.
- It is not hard to verify that the convergence criterion (47), and thus the time series is stationary.
- We can summarize these arguments by stating that *a time series model given by the lag form equation (44) is covariance stationary if the roots of the polynomial $\psi(z)$ lie outside of the unit circle.*

General autoregressive model $AR(p)$

R

- The p -th order autoregressive model $AR(p)$ model is specified as follows:

$$X_t = \alpha + \beta_1 X_{t-1} + \dots + \beta_p X_{t-p} + \varepsilon_t, \quad (53)$$

where $\alpha, \beta_j \in \mathbb{R}$, and $\varepsilon_t \sim N(0, \sigma^2)$ is a white noise.

- For the covariance-stationarity, the requirement that $E(X_t) = \mu$ yields

$$\mu = \frac{\alpha}{1 - \beta_1 - \dots - \beta_p}. \quad (54)$$

- Furthermore, we require that the roots of the characteristic polynomial $\psi(z) = 1 - \alpha - \beta_1 z - \dots - \beta_p z^p$ lie outside of the unit circle.
- We can rewrite (53) in the following form:

$$X_t - \mu = \beta_1 (X_{t-1} - \mu) + \dots + \beta_p (X_{t-p} - \mu) + \varepsilon_t. \quad (55)$$

General autoregressive model $AR(p)$

- Multiplying this equation by $X_{t-j} - \mu$, for $j = 0, \dots, p$, and calculating expectations yields the Yule-Walker equation for the autocovariance:

$$\Gamma_k = \begin{cases} \beta_1 \Gamma_1 + \dots + \beta_p \Gamma_p + \sigma^2, & \text{if } k = 0, \\ \beta_1 \Gamma_{k-1} + \dots + \beta_p \Gamma_{k-p}, & \text{if } k = 1, \dots, p. \end{cases} \quad (56)$$

- Dividing (56) by Γ_0 yields the Yule-Walker equation for the autocorrelation:

$$R_k = \beta_1 R_{k-1} + \dots + \beta_p R_{k-p}, \quad (57)$$

for $k = 1, \dots, p$.

- Note that the autocorrelations satisfy essentially the same equation as the process defining X_t .
- The ACF R_k can be found as the solution to the Yule-Walker equation and are expressed in terms of the roots of the characteristic polynomial.

Choosing the number of lags in $AR(p)$

- In practice, the number of lags p is unknown, and has to be determined empirically.
- This can be done by regressing the variable on its lagged values with $p = 1, 2, \dots$, and assessing the impact of each added lag on the fit.
- It is important not to overfit the model (“torture it until it confesses”) by adding too many lags.
- Useful quantitative guides for model selection are various information criteria.
- The *Akaike information criterion* defined as follows:

$$AIC = 2k - 2 \log \mathcal{L}(\hat{\theta}|x). \quad (58)$$

Here $k = \#\theta$ is the number of model parameters, $-\log \mathcal{L}(\hat{\theta}|x)$ denotes the optimized value of the LLF.

- According to this criterion, among the candidate models the model with the lowest value of AIC is the preferred one.

Choosing the number of lags in $AR(p)$

- This is in contrast with picking the model whose optimized LLF is the lowest: this may be the result of overfitting. The AIC criterion penalizes the number of parameters, and thus discourages overfitting.
- Another popular information criteria is the *Bayesian information criterion* (a.k.a the *Schwarz criterion*), which is defined as follows:

$$\text{BIC} = \log(N)k - 2 \log \mathcal{L}(\hat{\theta}|x), \quad (59)$$

where $N = \#x$ is the number of data points.

- According to this criterion, the model with the smallest value of BIC is the preferred model.

Moving average model $MA(1)$

R

- The *moving average* model $MA(1)$ is specified as follows:

$$X_t = \mu + \varepsilon_t + \theta\varepsilon_{t-1}, \quad (60)$$

where μ and θ are constants, and ε_t is white noise.

- The key feature of the $MA(1)$ model is that its are autocorrelated with lag 1.
- The expected value of X_t is

$$E(X_t) = \mu, \quad (61)$$

as $E(\varepsilon_t) = \mu$, for all t .

- Its variance is

$$\left[\begin{aligned} E((X_t - \mu)^2) &= E((\varepsilon_t + \theta\varepsilon_{t-1})^2) \\ &= E(\varepsilon_t^2) + 2\theta E(\varepsilon_t\varepsilon_{t-1}) + \theta^2 E(\varepsilon_{t-1}^2) \\ &= (1 + \theta^2)\sigma^2. \end{aligned} \right]$$

Moving average model $MA(1)$

- For the first autocovariance, we have

$$E((X_t - \mu)(X_{t-1} - \mu)) = E((\varepsilon_t + \theta\varepsilon_{t-1})(\varepsilon_{t-1} + \theta\varepsilon_{t-2})) \\ = \theta\sigma^2.$$

- All autocovariances with lag ≥ 2 are zero (show it!).

$$E[(\varepsilon_t + \theta\varepsilon_{t-1})(\varepsilon_{t-j} + \theta\varepsilon_{t-j-1})] = 0$$

- As a result, $MA(1)$ is (unlike $AR(1)$) always covariance-stationary with

$$MA(q) : y_t = \mu + \varepsilon_t + \theta_1\varepsilon_{t-1} + \dots + \theta_q\varepsilon_{t-q}$$

$$\Gamma_t = \begin{cases} (1 + \theta^2)\sigma^2, & \text{if } t = 0, \\ \theta\sigma^2, & \text{if } |t| = 1, \\ 0, & \text{if } |t| \geq 2. \end{cases} \quad (62)$$

- As a result, the first autocorrelation $R_1 = \Gamma_1/\Gamma_0$ is given by

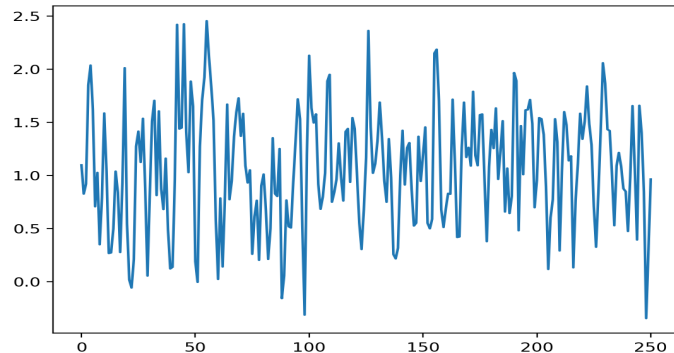
$$R_1 = \frac{\theta}{1 + \theta^2}, \quad (63)$$

with all higher order autocorrelations equal zero.

Always cov stationary.

Moving average model $MA(1)$

- The graph below shows a simulated $MA(1)$ time series with the following choice of parameters: $\mu = 1.1$, $\beta = 0.6$, $\sigma = 0.5$.



Moving average model $MA(1)$

- Here is the code snippet used to generate this graph in Python:

```
import numpy as np
import matplotlib.pyplot as plt
from statsmodels.tsa.arima_model import ARMA

mu=1.1
theta=0.6
sigma=0.5

#Simulate MA(1)
T=250
x0=mu
x=np.zeros(T+1)
x[0]=x0
eps=np.random.normal(0.0, sigma, T+1)
for i in range(1, T+1):
    x[i]=mu+eps[i]+theta*eps[i-1]

#Take a look at the simulated time series
plt.plot(x)
plt.show()
```

MLE for $MA(1)$

- As in the case of $AR(1)$, there are two natural approaches to MLE of an $MA(1)$ model: conditional on the initial value of ε and exact.
- We begin with the *conditional* MLE method, which is somewhat easier.
- Since the value of ε_0 cannot be calculated from the observed data, we are free to set it arbitrarily; we choose $\varepsilon_0 = 0$. All the probabilities calculated below are conditional on this choice.
- We then have, for $t = 1, \dots, T$,

$$\varepsilon_t = x_t - \mu - \theta \varepsilon_{t-1}, \quad (64)$$

and so the conditional PDF of x_t is

$$p(x_t | x_{t-1}, \dots, x_1, \varepsilon_0 = 0, \theta) = \frac{1}{\sqrt{2\pi\sigma^2}} \exp\left(-\frac{\varepsilon_t^2}{2\sigma^2}\right). \quad (65)$$

- This expression is deceptively simply: in reality ε_t is a nested function of all x_s with $s \leq t$.
- The likelihood function of the sample x_1, \dots, x_T is given by the product of the probabilities above, and so

$$\mathcal{L}(\theta | x, \varepsilon_0 = 0) = \prod_{t=1}^T p(x_t | x_{t-1}, \dots, x_1, \varepsilon_0 = 0, \theta), \quad (66)$$

- The log likelihood has thus the following form:

$$-\log \mathcal{L}(\theta|x, \varepsilon_0 = 0) = \frac{1}{2} T \log \sigma^2 + \frac{1}{2\sigma^2} \sum_{t=1}^T \varepsilon_t^2 + \text{const.} \quad (67)$$

- This is a quadratic function of the x_t 's. It is cumbersome to write it down explicitly, but easy to code it in a programming language. Its minimum is easiest to find by means of a numerical search.
- In case of $|\theta| < 1$, the impact of the choice $\varepsilon_0 = 0$ phases out as we iterate through time steps. For $|\theta| > 1$ the impact of this choice accumulates, and the method cannot be used.
- For the *exact* MLE method, we notice that the joint PDF of x is given by

$$p(x|\theta) = \frac{1}{(2\pi)^{T/2} \det(\Omega)^{1/2}} \exp \left(-\frac{1}{2} (x - \mu)^\top \Omega^{-1} (x - \mu) \right), \quad (68)$$

and thus

$$-\log \mathcal{L}(\theta|x) = \frac{1}{2} \log \det(\Omega) + \frac{1}{2} (x - \mu)^\top \Omega^{-1} (x - \mu). \quad (69)$$

- Here, Ω is a band diagonal matrix:

$$\Omega = \sigma^2 \begin{pmatrix} 1 + \theta^2 & \theta & 0 & \dots & 0 \\ \theta & 1 + \theta^2 & \theta & \dots & 0 \\ 0 & \theta & 1 + \theta^2 & \dots & 0 \\ \vdots & \vdots & \vdots & \dots & \vdots \\ \vdots & \vdots & \vdots & \dots & 1 + \theta^2 \end{pmatrix} \quad (70)$$

- The numerics of minimizing (69) can be handled either by (i) a clever triangular factorization of Ω , or by the Kalman filter method (we will discuss Kalman filters later in this course).
- Unlike the conditional MLE method, the exact method does not suffer from instabilities if $|\theta| \geq 1$.

- Here is the Python code snippet implementing the MLE for $MA(1)$ using statsmodels:

```
#MLE estimate with statsmodels
model=ARMA(x,order=(0,1)).fit(method='mle')
muMLE=model.params[0]
thetaMLE=model.params[1]
sigmaMLE=np.std(model.resid)
```

General moving average model $MA(q)$

R

- A q -th order moving average model $MA(q)$ is specified as follows:

$$X_t = \mu + \varepsilon_t + \theta_1 \varepsilon_{t-1} + \dots + \theta_q \varepsilon_{t-q}, \quad (71)$$

where μ and θ_j are constants, and ε_t is white noise.

- In other words, the $MA(q)$ model fluctuates around μ with disturbances which are autocorrelated with lag q .
- The expected value of X_t is

$$E(X_t) = \mu, \quad (72)$$

while its autocovariance is

$$\Gamma_j = \begin{cases} (1 + \theta_1^2 + \dots + \theta_q^2)\sigma^2, & \text{if } j = 0, \\ (\theta_j + \theta_{j+1}\theta_1 + \dots + \theta_q\theta_{q-j})\sigma^2, & \text{if } j = 1, \dots, q, \\ 0, & \text{if } j > q. \end{cases} \quad (73)$$

ACF is great diagnostic for order of MA that you are fitting data with.

- A mixed autoregressive moving average model ARMA(p, q) is specified as follows:

$$X_t = \alpha + \beta_1 X_{t-1} + \dots + \beta_p X_{t-p} + \varepsilon_t + \theta_1 \varepsilon_{t-1} + \dots + \theta_q \varepsilon_{t-q}, \quad (74)$$

where α and β_j, θ_k are constants, and ε_t is white noise.

- The equation above has the following lag operator representation:

$$\psi(L)X_t = \alpha + \varphi(L)\varepsilon_t, \quad (75)$$

where

$$\begin{aligned} \psi(z) &= 1 - \beta_1 z - \dots - \beta_p z^p, \\ \varphi(z) &= 1 + \theta_1 z + \dots + \theta_q z^q. \end{aligned} \quad (76)$$

- The process (45) is covariance stationary if the roots of ψ lie outside of the unit circle.

- In this case, we can write the model in the form

$$X_t = \mu + \gamma(L)\varepsilon_t, \quad (77)$$

where $\mu = \alpha/\psi(1)$, and $\gamma(L) = \psi(L)^{-1}\varphi(L)$. Explicitly, $\gamma(L)$ is an infinite series:

$$\gamma(L) = \sum_{j=0}^{\infty} \gamma_j L^j, \quad (78)$$

with

$$\sum_{j=0}^{\infty} |\gamma_j|^2 < \infty. \quad (79)$$

- This form of the model specification is called the *moving average form*.
- The parameters ARMA models are estimated by means of the MLE method. The complexity of computation required to minimize the LLF increases with the number of parameters.
- Information criteria, such as AIC or BIC, remain useful quantitative guides for model selection.

Forecasting time series with $ARMA(p, q)$

- An important function of time series analysis is making predictions about future values of the observed data, i.e. *forecasting*.
- Data based forecasting problem can be formulated as follows: given the observations $X_{1:t} = X_1, \dots, X_t$, what is the best forecast $X_{t+1|1:t}^*$ of X_{t+1} ?
- In mathematical terms, the problem requires minimizing a suitable loss function. We choose to minimize the *mean squared error* (MSE) given by

$$E((X_{t+1} - X_{t+1|1:t}^*)^2). \quad (80)$$

- We claim that $X_{t+1|1:t}^*$ is, indeed, given by the conditional expected value:

$$X_{t+1|1:t}^* = E_t(X_{t+1}). \quad (81)$$

Here E_t denotes expectation, conditional on the information up to time t ,

$$E_t(\cdot) = E(\cdot | X_{1:t}). \quad (82)$$

Forecasting time series with $ARMA(p, q)$

- Indeed, if Z is any random variable measurable with respect to the information set generated by $X_{1:t}$, then

$$\begin{aligned} E((X_{t+1} - Z)^2) &= E((X_{t+1} - E_t(X_{t+1}) + E_t(X_{t+1}) - Z)^2) \\ &= E((X_{t+1} - E_t(X_{t+1}))^2) + E((E_t(X_{t+1}) - Z)^2) \\ &\quad + 2E((X_{t+1} - E_t(X_{t+1}))(E_t(X_{t+1}) - Z)). \end{aligned}$$

- We argue that the cross term above is zero. Indeed

$$\begin{aligned} E_t((X_{t+1} - E_t(X_{t+1}))(E_t(X_{t+1}) - Z)) &= E_t(X_{t+1} - E_t(X_{t+1}))(E_t(X_{t+1}) - Z) \\ &= (E_t(X_{t+1}) - E_t(X_{t+1}))(E_t(X_{t+1}) - Z) \\ &= 0. \end{aligned}$$

Since $E(\cdot) = E(E_t(\cdot)|X_t)$, the claim follows.

Forecasting time series with $ARMA(p, q)$

- As a result

$$E((X_{t+1} - Z)^2) = E((X_{t+1} - E_t(X_{t+1}))^2) + E((E_t(X_{t+1}) - Z)^2),$$

which has its minimum at $Z = E_t(X_{t+1})$. This proves (81).

- The argument above is, in fact, quite general, and it easily extends to general k -period forecasts $X_{t+k|1:t}^*$. Minimizing the corresponding MSE yields:

$$X_{t+k|1:t}^* = E_t(X_{t+k}). \quad (83)$$

- Later we will generalize this method to time series models with more complex structure.

Forecasting time series with $ARMA(p, q)$

- As an example, a single period forecast in an $AR(1)$ model is

$$\begin{aligned}X_{t+1|1:t}^* &= E_t(X_{t+1}) \\&= E_t(\alpha + \beta X_t + \varepsilon_{t+1}) \\&= \alpha + \beta X_t.\end{aligned}\tag{84}$$

- The forecast error is ε_{t+1} , and so the variance of the forecast error is σ^2 .
- Likewise, a single period forecast in an $AR(p)$ model is

$$X_{t+1|1:t}^* = \alpha + \beta_1 X_t + \dots + \beta_p X_{t-p+1}.\tag{85}$$

with forecast error is ε_{t+1} , and the variance of the forecast error is σ^2 .

Forecasting time series with $ARMA(p, q)$

- A two-period forecast in an $AR(1)$ model is given by

$$\begin{aligned}X_{t+2|1:t}^* &= E_t(X_{t+2}) \\&= E_t(\alpha + \beta X_{t+1} + \varepsilon_{t+2}) \\&= (1 + \beta)\alpha + \beta^2 X_t.\end{aligned}\tag{86}$$

- The error of the two period forecast is $\varepsilon_{t+2} + \beta\varepsilon_{t+1}$; its variance is $(1 + \beta^2)\sigma^2$.
- A one period forecast in an $MA(1)$ model is

$$\begin{aligned}X_{t+1|1:t}^* &= E_t(X_{t+1}) \\&= E_t(\mu + \varepsilon_{t+1} + \theta\varepsilon_t) \\&= \mu + \theta\varepsilon_t.\end{aligned}\tag{87}$$

- The forecast error is ε_{t+1} , and its variance is σ^2 .
- These calculations can be generalized to produce a general formula for a multi-period forecast in an $ARMA(p, q)$ model. This result is known as the *Wiener-Kolmogorov prediction formula* and its discussion can be found in [1].

References



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