

Chapter 1

Time Series Modelling

1.1 Introduction

An observed *time series* is a set of observations $(x_t)_{t \in \mathcal{T}}$ where \mathcal{T} denotes an indexing set of finite size; each observation x_t is recorded at a specific time, related to the index t . x_t may be vector-valued if we are considering a multivariate time series, where the components may be influencing each other.

This course deals with Time Series models and their applications. We consider four areas of application:

1. The *forecasting* of future values of a time series from current and past values.
2. Computing a *transfer function* of a system, which shows the effect on the output of a system on any given series of inputs.
3. The use of *indicator* input variables in transfer function models to represent and assess the effects of unusual *intervention* events on the behaviour of a time series.
4. Examining relationships between several related time series of interest and establishing multivariate dynamic models to represent these joint relationships over time.

For now, we consider *univariate* time series; each $x_t \in \mathbb{R}$. We consider vector valued time series later.

1.2 Time Series Models

Definition 1.1 (Time Series Model). *A time series model for the observed data $\{x_t : t \in \mathcal{T}\}$ is the hypothesis that the observed data is an observation of a sequence of random variables $\{X_t : t \in \mathcal{T}\}$ and the specification of its joint probability distribution, or possibly only its expectations and covariances.*

A time series can only be observed at a finite number of times, $(x_t)_{t=1}^n$ and the n observations are a realisation of an n dimensional random vector $X = (X_1, X_2, \dots, X_n)$. These random variables may be considered to come from an infinite sequence $\{X_t, t \in \mathbb{Z}_+ \text{ or } \mathbb{Z}\}$, a *stochastic process*.

Definition 1.2 (Stochastic Process). *A stochastic process is a family of random variables $\{X_t : t \in \mathcal{T}\}$, indexed by a set \mathcal{T} , which is defined on a probability space $(\Omega, \mathcal{F}, \mathbb{P})$.*

Example 1.1 (The binary process).

A simple example of a stochastic process $\{X_t, t \in \mathbb{Z}_+\}$ is a process where the variables are i.i.d. (independent identically distributed) satisfying

$$\mathbb{P}(X_t = 1) = \mathbb{P}(X_t = -1) = \frac{1}{2}.$$

For this process, the finite dimensional marginals are well defined; for any $i_1 < \dots < i_n$,

$$\mathbb{P}(X_{i_1} = j_1, X_{i_2} = j_2, \dots, X_{i_n} = j_n) = 2^{-n}$$

for any $\{j_1, \dots, j_n\} \in \{-1, 1\}^n$. □

Definition 1.3 (IID noise). *A process $\{X_t, t \in \mathbb{Z}\}$ is said to be an IID noise with mean μ and variance σ^2 , written*

$$\{X_t\} \sim \text{IID}(\mu, \sigma^2),$$

if the random variables X_t are independent and identically distributed with $\mathbb{E}[X_t] = \mu$ and $\text{Var}(X_t) = \sigma^2$.

Usually, we are interested in $\text{IID}(0, \sigma^2)$ noise.

Notation Throughout, $\text{Var}(\cdot)$ will be used to denote variance.

The binary process is clearly an example of an $\text{IID}(0, 1)$ noise, since the variables are independent, $\mathbb{E}[X_t] = -1 \times \frac{1}{2} + 1 \times \frac{1}{2} = 0$ and $\text{Var}(X_t) = \mathbb{E}[X_t^2] - \mathbb{E}[X_t]^2 = \mathbb{E}[X_t^2] = 1$.

In many situations, the complete specification of the underlying stochastic process is not required; the methods will generally rely only on its means and covariances. Sometimes even less general assumptions are needed, but these will not be treated here.

Definition 1.4 (Mean function, Covariance function). *Let $\{X_t, t \in \mathcal{T}\}$ be a stochastic process with $\text{Var}(X_t) < \infty$ for each $t \in \mathcal{T}$. The mean function of $\{X_t\}$ is denoted by μ_X , or simply μ when there is no danger of ambiguity:*

$$\mu_X(t) := \mathbb{E}[X_t], \quad t \in \mathcal{T} \tag{1.1}$$

The covariance function of $\{X_t\}$ is denoted by C_X or C when there is no danger of ambiguity and is defined as:

$$C_X(r, s) := \text{Cov}(X_r, X_s), \quad r, s \in \mathcal{T}. \tag{1.2}$$

The symbol Cov will be used to denote covariance.

1.3 Stationarity and Strict Stationarity

A stochastic process is said to be *stationary*, if its statistical properties do not change with time. Formally, stationarity is defined in the following way.

Definition 1.5 (Stationary, Strictly Stationary, Wide sense stationary). *A time series $\{X_t, t \in \mathbb{Z}\}$ is said to be weakly stationary, or wide sense stationary, or simply stationary if*

1. $\text{Var}(X_t) < \infty$ for all $t \in \mathbb{Z}$,
2. $\mu_X(t) = \mu$ for all $t \in \mathbb{Z}$,
3. $C_X(r, r+h) = C_X(0, h)$ for all $r, h \in \mathbb{Z}$.

A process is said to be strictly stationary if any finite collection $(X_{n_1}, \dots, X_{n_k})$ has the same distribution as $(X_{n_1+t}, \dots, X_{n_k+t})$ for any $k \geq 1$ and any $(n_1, \dots, n_k, t) \in \mathbb{Z}$.

In many practical situations, only weak stationarity is considered; usually only expectation and covariance, at most, can reasonably be assessed from data. In some situations, though (for example ARCH and GARCH processes, which arise in the analysis of financial time series) it is worthwhile placing additional modelling assumptions on the data generation mechanism.

The third point in the definition of weak stationarity implies that $C_X(r, s)$ depends on r and s only through $r - s$. It is therefore convenient to define

$$\gamma_X(h) := C_X(h, 0).$$

When only one time argument appears in γ , then it denotes the autocovariance (ACVF) function of a stationary process. The value h is referred to as the *lag*.

Definition 1.6. *Let $\{X_t, t \in \mathbb{Z}\}$ be a stationary time series. The autocovariance function (ACVF) of $\{X_t\}$ is defined as*

$$\gamma_X(h) = \text{Cov}(X_{t+h}, X_t).$$

The autocorrelation function (ACF) is defined as:

$$\rho_X(h) := \frac{\gamma_X(h)}{\gamma_X(0)}.$$

A simple example of a stationary process is the so-called *white noise*.

Definition 1.7 (White noise). *A process $\{X_t, t \in \mathbb{Z}\}$ is said to be a white noise with mean μ and variance σ^2 , written*

$$\{X_t\} \sim \text{WN}(\mu, \sigma^2),$$

if $\mathbb{E}[X_t] = \mu$ for all $t \in \mathbb{Z}$ and

$$\gamma(h) = \begin{cases} \sigma^2 & \text{if } h = 0, \\ 0 & \text{if } h \neq 0. \end{cases}$$

Note that IID noise is an example of white noise, but not necessarily vice versa; the underlying distribution can be different even if the mean and covariance structures are the same; strictly stationary time series $\{X_t, t \in \mathbb{Z}\}$ with $\text{Var}(X_t) < \infty$ is stationary, but a stationary time series $\{X_t, t \in \mathbb{Z}\}$ does not need to be strictly stationary

From now on, the term ‘stationary’ will be used to denote ‘weakly’ or ‘wide sense stationary’; the term *strictly stationary* will be used for the stronger assumption.

Example 1.2 (AR(1) process).

Autoregressive (AR) processes will be considered in more detail later. A process $\{X_t, t \in \mathbb{Z}\}$ is said to be AR(1) if it stationary and satisfies:

$$X_t = \phi X_{t-1} + Z_t \quad \{Z_t\} \sim WN(0, \sigma^2).$$

For this process, the autocovariance may be computed as follows: by squaring up both sides and using $\gamma_X(0) = \text{Var}(X_t)$,

$$\gamma_X(0) = \phi^2 \gamma_X(0) + \sigma^2 \Rightarrow \gamma_X(0) = \frac{\sigma^2}{1 - \phi^2}.$$

for $h \geq 1$,

$$\gamma_X(h) = \text{Cov}(X_{t+h}, X_t) = \phi \text{Cov}(X_{t+h-1}, X_t) + \text{Cov}(Z_{t+h}, X_t) = \phi \gamma_X(h-1)$$

so that, since $\gamma_X(-h) = \gamma_X(h)$,

$$\gamma_X(h) = \frac{\sigma^2}{(1 - \phi^2)} \phi^{|h|}.$$

Its autocorrelation function (ACF) is

$$\rho_X(h) = \phi^{|h|}.$$

Note that the AR(1) process is not well defined if $|\phi| \geq 1$. □

1.4 Trends and Seasonal Components

The classical decomposition model is:

$$X_t = \mu_t + s_t + \epsilon_t,$$

where

- μ_t is a slowly changing function (the *trend*);
- s_t is a function with known period d (the ‘seasonal component’);
- ϵ_t is a stationary time series.

The aim is to extract the deterministic components μ_t and s_t and estimate them and then check whether or not the residual component ϵ_t is a stationary time series.

1.4.1 No Seasonal Component

Assume that

$$X_t = \mu_t + \epsilon_t, \quad t = 1, \dots, n$$

where, without loss of generality, $\mathbb{E}[\epsilon_t] = 0$ (since the mean of the stationary process is systematic and is therefore considered to be part of the trend).

There are several methods for estimating μ . Three are considered here; *least squares*, *moving average* and *differencing*.

Method 1 : Least Squares estimation of μ_t The function μ_t is modelled by a function with as few parameters as necessary for accurate modelling and the parameters are estimated by the least squares technique. For example, suppose that μ_t can be modelled by a quadratic function, $\mu_t = a_0 + a_1t + a_2t^2$. The parameters $(a_k)_{k=0}^2$ are estimated by $(\hat{a}_k)_{k=0}^2$, chosen to minimise

$$\sum_{t=1}^n (x_t - a_0 - a_1t - a_2t^2)^2.$$

Method 2 : Smoothing by means of a moving average Let q be a non-negative integer and consider a smoothed version of X defined by

$$W_t := \frac{1}{2q+1} \sum_{j=-q}^q X_{t+j}, \quad q+1 \leq t \leq n-q.$$

If it turns out that μ is approximately linear over the time interval $[t-q, t+q]$ and also that q is sufficiently large so that $\frac{1}{2q+1} \sum_{j=-q}^q Y_{t+j} \simeq 0$, then

$$W_t = \frac{1}{2q+1} \sum_{j=-q}^q \mu_{t+j} + \frac{1}{2q+1} \sum_{j=-q}^q Y_{t+j} \simeq \mu_t.$$

For $t \leq q$ and $t > n-q$, W has to be defined in a different way. For example,

$$W_t = \begin{cases} \frac{1}{2t+1} \sum_{j=-t}^t X_{t+j} & t = 1, \dots, q \\ \frac{1}{2(n-t)+1} \sum_{j=-(n-t)}^{n-t} X_{t-j} & t = n-q+1, \dots, n. \end{cases}$$

Unless μ_t is a straight line and the stationary time series component Y is very small, it will not be possible to find a q satisfying both the conditions that μ is approximately linear over the interval $[t-q, t+q]$ (requiring small q) and such that $\frac{1}{2q+1} \sum_{j=-q}^q Y_s \simeq 0$ (requiring large q).

Definition 1.8 (Linear Filter). A linear filter is defined as a linear combination:

$$\hat{\mu}_t = \sum_j a_j X_{t+j},$$

where $\sum a_j = 1$ and $a_j = a_{-j}$.

A linear filter will allow a *linear trend* $\mu_t = \alpha_0 + \alpha_1 t$ to pass without distortion since

$$\sum_j a_j (\alpha_0 + \alpha_1 (t+j)) = (\alpha_0 + \alpha_1 t) \sum_j a_j + \alpha_1 \sum_j a_j j = \alpha_0 + \alpha_1 t.$$

It is possible to choose the weights $\{a_j\}$ so that a larger class of trend functions pass without distortion. For example, the *Spencer 15-point moving average*, defined as

$$\begin{cases} [a_0, a_{\pm 1}, \dots, a_{\pm 7}] = \frac{1}{320} [74, 67, 46, 21, 3, -5, -6, -3] \\ a_j = 0 \text{ for } |j| > 7 \end{cases}$$

allows a cubic trend to pass without distortion. That is, applied to $\mu_t = at^3 + bt^2 + ct + d$,

$$\hat{\mu}_t = \sum a_j X_{t+j} = \sum a_j \mu_{t+j} + \sum a_j Y_{t+j} \simeq \sum a_j \mu_{t+j} = \mu_t.$$

Conditions required for a filter to pass a trend which is polynomial of degree k without distortion are found in the exercises. Moving average methods will be considered in greater detail later.

1.5 Linear filters

A linear process may be regarded as a *linear filter*. Let $\{X_t\}$ be a time series. A *filter* is an operation on a time series in order to obtain a new time series $\{Y_t\}$. $\{X_t\}$ is called the *input* and $\{Y_t\}$ the *output*. A linear filter \mathcal{C} is the following operation:

$$\mathcal{C}(X)_t := Y_t = \sum_{k=-\infty}^{\infty} c_{t,k} X_k. \quad (1.3)$$

We only consider the situation where $\mathbb{E}[X_t^2] < \infty$ and $\mathbb{E}[Y_t^2] < \infty$.

A linear filter is said to be *time-invariant* if $c_{t,k} = c_{t-k}$, in which case it may be written as:

$$Y_t = \sum_{j=-\infty}^{\infty} c_j X_{t-j}.$$

A time-invariant linear filter (TLF) is said to be *causal* if

$$c_j = 0 \text{ for } j < 0,$$

When the input $\{X_t\}$ of a time invariant linear filter is stationary, then the output $\{Y_t\}$ is also stationary provided $\sum_k |c_k| < +\infty$.

Definition 1.9 (Stable Linear Filter). A TLF of the form (1.3) is stable if $\sum_{k=-\infty}^{\infty} |c_k| < \infty$.

Definition 1.10 (Transfer function, Power function). Consider a stable linear filter and set

$$c(z) = \sum_{j=-\infty}^{\infty} c_j z^j.$$

The function $c(e^{-i\lambda}) := \sum_{j=-\infty}^{\infty} c_j e^{-i\lambda j}$ is known as the transfer function, while the function $|c(e^{-i\lambda})|^2$ is known as the power transfer function.

A filter may be written as $c(B)$, in the sense that

$$Y_t = c(B)X_t$$

where B as usual denotes the backward shift operator.

A linear process is a linear filter where the input is $\text{WN}(0, \sigma^2)$.

Impulse Response Function In general, for a stationary process $\{X_t : t \in \mathbb{Z}\}$, where the variables $\{X_t\}$ are functions of impulses $\{\epsilon_t : t \in \mathbb{Z}\}$, the *impulse response function* $g(s)$ is defined as:

$$g(t; s) = \frac{\partial X_{t+s}}{\partial \epsilon_t} \quad (1.4)$$

In the case of a causal linear filter $X_t = \sum_j c_j \epsilon_{t-j}$, $g(t; s) = g(s) = c_s$.

The impulse response function may be extended to vectors; if $\{\underline{X}_t : t \in \mathbb{Z}\}$ is an m -vector valued process which is a function of vector impulses $\{\underline{\epsilon}_t : t \in \mathbb{Z}\}$, then

$$g_{ij}(t, s) = \frac{\partial X_{t+s, i}}{\partial \epsilon_{t, j}}. \quad (1.5)$$

If $\{\underline{X}_t\}$ is a linear causal vector valued process satisfying $X_{t, j} = \sum_{s \geq 0} \sum_k c_{jk; s} \epsilon_{t-s, k}$ then

$$g_{ij}(t, s) = g_{ij}(s) = c_{ij; s}.$$

Method 3: Differencing to generate stationarity Let B denote the *backward shift* operator; $(BX)_t = X_{t-1}$, with powers given by $(B^j X)_t = X_{t-j}$. In other words, applying B^j to X pushes it back j time units. Strict stationarity means that $B^h X$ has the same distribution for all $h \in \mathbb{Z}_+$.

The *difference operator* ∇ is defined by

$$\nabla X_t = X_t - X_{t-1} = (1 - B)X_t,$$

where B is the *backward shift operator*. That is, $(BX)_t = X_{t-1}$. For positive integer k , ∇^k is defined by: by:

$$\nabla^k X_t = \nabla(\nabla^{k-1} X)_t.$$

For example,

$$\nabla^2 X_t = \nabla X_t - \nabla X_{t-1} = (X_t - X_{t-1}) - (X_{t-1} - X_{t-2}) = X_t - 2X_{t-1} + X_{t-2}.$$

Using the backward shift operator, this may be expressed as:

$$\nabla^2 X_t = (1 - B)^2 X_t = (1 - 2B + B^2) X_t = X_t - 2X_{t-1} + X_{t-2}.$$

For a linear trend $\mu_t = a + bt$,

$$\nabla X_t = \nabla \mu_t + \nabla Y_t = a + bt - a - b(t-1) + \nabla Y_t = b + \nabla Y_t.$$

For the covariance,

$$\begin{aligned} \text{Cov}(\nabla Y_t, \nabla Y_s) &= \text{Cov}(Y_t, Y_s) - \text{Cov}(Y_{t-1}, Y_s) - \text{Cov}(Y_t, Y_{s-1}) + \text{Cov}(Y_{t-1}, Y_{s-1}) \\ &= \gamma_Y(t-s) - \gamma_Y(t-s-1) - \gamma_Y(t-s+1) + \gamma_Y(t-s) \\ &= 2\gamma_Y(t-s) - \gamma_Y(t-s+1) - \gamma_Y(t-s-1). \end{aligned}$$

It follows that ∇X_t is stationary with

$$\mu_{\nabla X} = b \quad \gamma_{\nabla X}(h) = 2\gamma_Y(h) - \gamma_Y(h+1) - \gamma_Y(h-1).$$

In general, if $\mu_t = \sum_{j=0}^k c_j t^j$, then

$$\nabla^k X_t = k! c_k + \nabla^k Y_t,$$

which is stationary.

1.5.1 Trend and Seasonality

Now consider the model with a seasonal component:

$$X_t = \mu_t + s_t + Y_t,$$

where $\mathbb{E}[Y_t] = 0$, $s_{t+d} = s_t$ and $\sum_{k=1}^d s_k = 0$. For simplicity in the representation, assume that n/d is an integer; in any reasonable modelling situation, n and d will be chosen so that n/d is an integer.

In models with a seasonal component, the data is often indexed by period and time-unit;

$$x_{j,k} = x_{k+d(j-1)}, \quad k = 1, \dots, d, \quad j = 1, \dots, \frac{n}{d}.$$

In this notation, $x_{j,k}$ is the observation at the k :th time-unit of the j :th period.

Three methods for dealing with seasonal components will be considered; the *small trend* method, the *moving average estimation* method and the *differencing at lag d* method.

Method S1: Small trends If the trend is considered to be constant during each period, the model may be written as:

$$X_{j,k} = \mu_j + s_k + Y_{j,k}.$$

A natural way to estimate the trend is:

$$\hat{\mu}_j = \frac{1}{d} \sum_{k=1}^d x_{j,k}$$

and a natural method for the seasonal component is:

$$\hat{s}_k = \frac{d}{n} \sum_{j=1}^{n/d} (x_{j,k} - \hat{\mu}_j).$$

Method S2: Moving average estimation For a known period d , the trend is estimated by applying a moving average to eliminate the seasonal component and to reduce the noise. For d even set $q = d/2$. The trend is estimated by:

$$\hat{\mu}_t = \frac{0.5x_{t-q} + x_{t-q+1} + \cdots + x_{t+q-1} + 0.5x_{t+q}}{d}.$$

For d odd, set $q = (d-1)/2$. The trend is estimated by:

$$\hat{\mu}_t = \frac{x_{t-q} + x_{t-q+1} + \cdots + x_{t+q-1} + x_{t+q}}{d},$$

for $q+1 \leq t \leq n-q$.

The seasonal component s_k is then estimated in the following way. Set

$$w_k = \frac{1}{\text{number of summands}} \sum_{\frac{q-k}{d} < j \leq \frac{n-q-k}{d}} (x_{k+jd} - \hat{\mu}_{k+jd}).$$

The seasonal component satisfies $\sum_{k=1}^d \hat{s}_k = 0$ and therefore the estimates are:

$$\hat{s}_k = w_k - \frac{1}{d} \sum_{i=1}^d w_i, \quad k = 1, \dots, d.$$

Method S3: Differencing at lag d Define the lag- d difference operator ∇_d by

$$\nabla_d X_t = X_t - X_{t-d} = (1 - B^d)X_t.$$

Then

$$\nabla_d X_t = \nabla_d \mu_t + \nabla_d Y_t.$$

This has no seasonal component and the methods for dealing with time series without a seasonal component may be applied.

1.6 Autocovariance and Spectral Density of a stationary time series

Recall Definition 1.5 of a weakly stationary time series. It follows directly from the definition that:

$$\begin{cases} \gamma(0) \geq 0, \\ |\gamma(h)| \leq \gamma(0) & \text{for all } h \in \mathbb{Z}, \\ \gamma(h) = \gamma(-h) & \text{for all } h \in \mathbb{Z}. \end{cases} \quad (1.6)$$

An autocovariance function is clearly non-negative definite, since $\sum_{j=1}^n \sum_{k=1}^n a_j a_k \gamma(t_j - t_k)$ is the variance of $\sum_{j=1}^n a_j X_{t_j}$.

Definition 1.11. A function $\kappa : \mathbb{Z} \rightarrow \mathbb{R}$ is said to be non-negative definite, or positive semi-definite, if

$$\sum_{i,j=1}^n a_i a_j \kappa(t_i - t_j) \geq 0$$

for all n and all vectors $\underline{a} \in \mathbb{R}^n$ and $\underline{t} \in \mathbb{Z}^n$.

Theorem 1.12. The autocorrelation function of a stationary time series is a real valued, even non negative definite function defined on \mathbb{Z} . For any real valued even non negative definite function κ defined on \mathbb{Z} and for any $N \geq 1$, there exists a sequence of random variables (X_{-N}, \dots, X_N) such that $\text{Cov}(X_i, X_j) = \kappa(i - j)$.

Proof Let $\gamma(\cdot)$ be the autocovariance function of a stationary time series X_t . Then for any (t_1, \dots, t_n) and any (a_1, \dots, a_n) ,

$$0 \leq \text{Var} \left(\sum_{j=1}^n a_j X_{t_j} \right) = \sum_{j,k} a_j a_k \gamma(t_j - t_k).$$

For the other way, Let $\underline{Z} = (Z_{-N}, \dots, Z_N)$ be a vector of i.i.d. $N(0, 1)$ variables. Let K denote the $2N + 1 \times 2N + 1$ matrix with entries $K_{i,j} = \kappa(i - j)$. Then K is a non negative definite matrix. It follows that K has a decomposition $P\Lambda P^t$ where P is an orthonormal matrix and Λ is a diagonal matrix whose entries are the eigenvalues. Let $K^{1/2} = P\Lambda^{1/2}P^t$, then $(K^{1/2})^2 = K$. Let $\underline{X} = K^{1/2}\underline{Z}$, then \underline{X} is a random vector with covariance K as required. \square

1.7 Holt Winters Filtering

No trend, no seasonal component Given observations X_1, X_2, \dots, X_n from the model:

$$X_t = \mu + Z_t \quad \{Z_t\} \sim \text{WN}(0, \sigma^2)$$

where μ is considered to be approximately constant. The method of *exponential smoothing* is to compute a *smoothed* series:

$$\tilde{X}_t = \lambda X_t + (1 - \lambda)\tilde{X}_{t-1} \quad \lambda \in (0, 1) \quad (1.7)$$

where λ is the *smoothing parameter*. The forecast for time $t + h$ given the series up to time t is

$$\widehat{X}_{t+h|t} = \widetilde{X}_t.$$

The quantity \widetilde{X}_t is the estimate of μ at time t ; the assumption is that the underlying value of μ will not change between t and $t + h$.

Linear trend, no seasonal component Holt and Winters independently extended this idea (Holt (1959) and Winters (1960)) to deal with the model

$$X_t = \mu_t + Z_t \quad \{Z_t\} \sim \text{WN}(0, \sigma^2)$$

under the assumption that the trend is approximately linear. Let $m_t = \mu_t - \mu_{t-1}$. Then the equations suggested by Holt and Winters are:

$$\begin{cases} \widetilde{X}_t = \lambda_1 X_t + (1 - \lambda_1) (\widetilde{X}_{t-1} + \widetilde{m}_{t-1}) \\ \widetilde{m}_t = \lambda_2 (\widetilde{X}_t - \widetilde{X}_{t-1}) + (1 - \lambda_2) \widetilde{m}_{t-1} \end{cases} \quad (1.8)$$

where \widetilde{m}_t is the estimate of m_t at time t . The h -step ahead forecasts are then given by:

$$\widehat{X}_{t+h|t} = \widetilde{X}_t + h\widetilde{m}_t.$$

Holt Winters with linear trend and additive seasonal component Now suppose that $\{X_t\}$ is a time series with both trend and seasonal component where the seasonal component $\{s_t\}$ has period d :

$$X_t = \mu_t + s_t + Z_t \quad \{Z_t\} \sim \text{WN}(0, \sigma^2)$$

The Holt-Winters algorithm accommodates the seasonal component in the following way: let $Y_t = X_t - s_t$, then \widetilde{Y}_t is an approximation of μ_t and

$$\begin{cases} \widetilde{Y}_t = \lambda_1 (X_t - \widetilde{s}_{t-d}) + (1 - \lambda_1) (\widetilde{Y}_{t-1} + \widetilde{m}_{t-1}) \\ \widetilde{m}_t = \lambda_2 (\widetilde{Y}_t - \widetilde{Y}_{t-1}) + (1 - \lambda_2) \widetilde{m}_{t-1} \\ \widetilde{s}_t = \lambda_3 (X_t - \widetilde{Y}_t) + (1 - \lambda_3) \widetilde{s}_{t-d} \end{cases}$$

The initial conditions are:

$$\begin{cases} \widetilde{Y}_{d+1} = X_{d+1} \\ \widetilde{m}_{d+1} = \frac{1}{d} (X_{d+1} - X_1) \\ \widetilde{s}_i = X_i - (X_1 + \widetilde{m}_{d+1}(i-1)) \quad i = 1, \dots, d+1 \end{cases}$$

The predictors are:

$$\widehat{X}_{t+h|t} = \widetilde{Y}_t + h\widetilde{m}_t + \widetilde{s}_{t+h} \quad h = 1, 2, \dots$$

The parameters $\lambda_1, \lambda_2, \lambda_3 \in (0, 1)$ may be chosen by minimising the sum of squares of the one-step prediction error on data that has already been observed:

$$\sum_{i=d+2}^n \left(X_i - \widehat{X}_{i|i-1} \right)^2$$

Holt Winters Seasonal Multiplicative The *multiplicative* model is simply: $X_t = \exp\{\mu_t + s_t + \epsilon_t\}$ so that

$$\log X_t = \mu_t + s_t + \epsilon_t \quad \{\epsilon_t\} \sim WN(0, \sigma^2).$$

If P denotes the period, then the forecasting model for time $t + \tau$ with forecasting origin t is:

$$\widehat{x}_{t,t+\tau} = (\widehat{a}_t + \tau \widehat{b}_t) \widehat{S}_{t+\tau-P}$$

where:

$$\begin{cases} \widehat{a}_t = \lambda_1 \left(\frac{x_t}{\widehat{S}_{t-P}} \right) + (1 - \lambda_1) (\widehat{a}_{t-1} + \widehat{b}_{t-1}) \\ \widehat{b}_t = \lambda_2 (\widehat{a}_t - \widehat{a}_{t-1}) + (1 - \lambda_2) \widehat{b}_{t-1} \\ \widehat{S}_t = \lambda_3 \left(\frac{x_t}{\widehat{a}_t} \right) + (1 - \lambda_3) \widehat{S}_{t-P}. \end{cases}$$

The justification of this is left as an exercise.

1.8 Time Series in R

1.8.1 Extracting the Trend, Seasonal Component and Noise in R

The `stl` command may be used to decompose a time series into trend, seasonal component and noise. The computation of ‘trend’ is based on moving average. For illustration, consider the carbon dioxide data from Mauna Loa in the file `atmospheric-carbon-dioxide-recor.csv` .

```
> www =
"https://www.mimuw.edu.pl/~noble/courses/TimeSeries/data/atmospheric-
carbon-dioxide-recor.csv"
> carbon = read.csv(www)
```

Delete observation 611 which is ‘na’:

```
> carbon = carbon[-611,]
```

(this deletes the last row, which is ‘na’).

```
> y = carbon$MaunaLoaCO2
> MaunLoaCo2 = ts(data = y, frequency = 12)
```

(this gets it into an appropriate format - each row represents a year)

```
> output.stl = stl(MaunLoaCo2, s.window = "periodic")
> plot(output.stl)
```

This gives a plot of the original data, the seasonal component, the trend and the ‘remainder’.

```
> a <- output.stl$time.series
> acf(a)
```

The time.series part of the stl output gives a decomposition into trend, the seasonal and the noise. The acf gives the autocorrelation for each of these; the trend, seasonal and noise, while the off-diagonals show the cross autocorrelations.

The dotted blue lines indicate ‘error’ bars. The plot of interest is the residual (or ‘remainder’). The acf indicates clear correlations between the residuals; they are not $WN(0, \sigma^2)$. The plot is in Figure 1.1

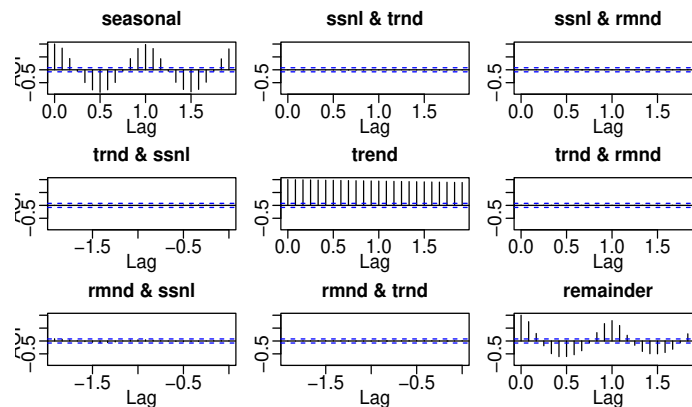


Figure 1.1: Mauna Loa: estimated acf for decomposition

To get the sample standard deviation of each column in the time series, try:

```
> apply(a,2,sd)
seasonal      trend  remainder
2.0402413 21.0085895  0.2735003
```

This indicates that the remainder is small compared with the trend and seasonal components.

1.8.2 Holt Winters Filtering: Implementation in R

The ‘Air Passengers’ data set is included in the data sets that come with R. Implementation of Holt-Winters can be carried out as follows: Type

```
> data(AirPassengers)
> AP <- AirPassengers
> str(AP)
```

```
Time-Series [1:144] from 1949 to 1961: 112 118 132 129 121 135 148 148 136 119
...
```

The data set is now loaded. Type

```
> ?HoltWinters
```

to obtain the syntax for the command. Note that values for λ_1, λ_2 and λ_3 may be given; if the user does not give the values, then they are computed by minimising the sum of squares of the one-step prediction errors as outlined above. To make a *multiplicative seasonal* Holt Winters, try:

```
> AP.hw <- HoltWinters(AP,seasonal="mult")
> plot(AP.hw)
> legend("topleft",c("observed","fitted"),lty=1,col=1:2)
```

This gives a plot showing both the original data and the one-step predictors. The plot is in Figure 1.2.

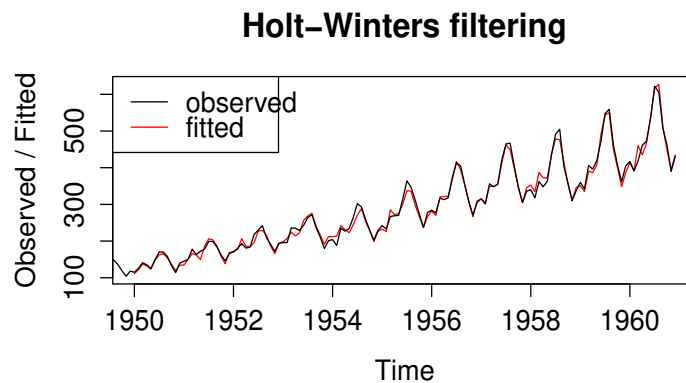


Figure 1.2: Air Passenger data with Holt Winters filtering

Prediction is made quite simply using the ‘predict’ command, which makes the arithmetical computations from the Holt Winters object. The following shows the predictions for the next four years.

```
> AP.predict <- predict(AP.hw,n.ahead=4*12)
> ts.plot(AP,AP.predict,lty=1:2)
```

The plot is found in Figure 1.3.

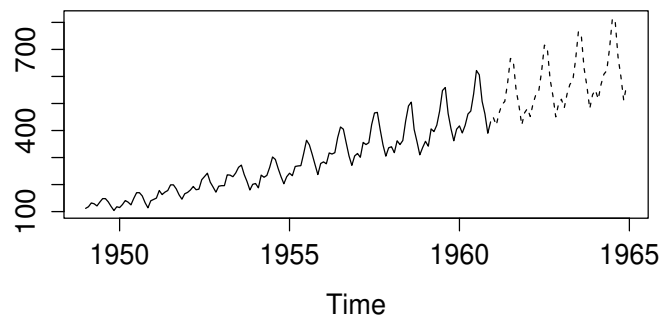


Figure 1.3: Air Passenger data: Holt Winters prediction