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THE FUNCTION SPACE POINT OF VIEW IN TIME SERIES ANALYSIS

BY  
EMANUEL PARZEN

TECHNICAL REPORT NO. 47  
March 29, 1963

PREPARED UNDER CONTRACT Nonr-225(21)  
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# THE FUNCTION SPACE POINT OF VIEW IN TIME SERIES ANALYSIS\*

by

Emanuel Parzen\*\*

Stanford University

## 0. Summary

This is an expository paper which seeks to establish and show the value of the following assertion: the concept of a time series is equivalent to the idea of a probability measure on a function space.

## 1. What is a time series?

The general point of view adopted in analyzing a time series or a succession of observations  $\{X(t), t \in T\}$ , depending on the parameter  $t$  (often representing time), is the following.

A set  $T$  of values of  $t$ , called the index set of the time series, is preassigned; these are the times when observations are possible. The set  $T$  may be finite or infinite. It is possible to develop much of the theory of time series without placing any restriction on the nature of the index set  $T$ . However two important cases are when

$$T = \{0, \underline{+1}, \underline{+2}, \dots\} \quad \text{or} \quad T = \{0, 1, 2, \dots\},$$

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\*\*It is with great pleasure that I dedicate this paper to Professor Charles Loewner on his 70th birthday.

in which case the time series is said to be a discrete parameter process,  
or when

$$T = \{t: -\infty < t < \infty\} \quad \text{or} \quad T = \{t: t \geq 0\},$$

in which case the time series is said to be a continuous parameter process.

At each point  $t$  in  $T$ , a number  $X(t)$  may be observed. This number is a random variable in the sense that its value depends on chance and enjoys a probability distribution described by the one-dimensional distribution function

$$F_{X(t)}(x) = \text{Probability}[X(t) \leq x], \quad -\infty < x < \infty,$$

or the one-dimensional characteristic function

$$\varphi_{X(t)}(u) = \int_{-\infty}^{\infty} \exp[iux] dF_{X(t)}(x), \quad -\infty < u < \infty.$$

More generally, for any integer  $n$  and  $n$  points  $t_1, t_2, \dots, t_n$  in  $T$ , the  $n$  observations  $X(t_1), \dots, X(t_n)$  which can be made at these times are jointly distributed random variables whose joint probability law is specified by either (i) the joint distribution function, given for all real numbers  $x_1, \dots, x_n$  by

$$\begin{aligned} & F_{X(t_1), \dots, X(t_n)}(x_1, \dots, x_n) \\ (1.1) \quad & = \text{Probability}[X(t_1) \leq x_1, X(t_2) \leq x_2, \dots, X(t_n) \leq x_n] \end{aligned}$$

or (ii) the joint characteristic function, given for all real numbers  $u_1, \dots, u_n$  by

$$\begin{aligned}
 & \varphi_{X(t_1), \dots, X(t_n)}(u_1, \dots, u_n) \\
 &= E[\exp i(u_1 X(t_1) + \dots + u_n X(t_n))] \\
 (1.2) \quad &= \int_{-\infty}^{\infty} \dots \int_{-\infty}^{\infty} \exp i(u_1 x_1 + \dots + u_n x_n) \\
 & \quad dF_{X(t_1), \dots, X(t_n)}(x_1, \dots, x_n).
 \end{aligned}$$

The distribution function in Equation (1.1) and the characteristic function in Equation (1.2) are said to be  $n$ -dimensional, since they represent the joint probability law of  $n$  random variables.

The point of view embodied in the foregoing discussion may be summarized as follows.

A time series is a jointly distributed family of random variables  $\{X(t), t \in T\}$ , indexed by a parameter  $t$  varying in an index set  $T$ . Time series analysis is concerned with the statements that can be made about a time series, knowing only all the finite dimensional distributions of the form of (1.1).

It should be noted that the phrase "a stochastic process" is often used to describe a numerical-valued random phenomenon that arises through a process which is developing in time in a manner controlled by probabilistic laws [see Parzen (1962)]. Mathematically, a stochastic process is represented by a collection of random variables  $\{X(t), t \in T\}$ . Thus, in a

sense, the notions of a time series and of a stochastic process are equivalent; every time series is a stochastic process and vice versa.

What distinguishes the theory of time series from the theory of stochastic processes is a certain difference in emphasis.

One approach to the problem of developing mathematical models for empirical phenomena evolving in accord with probabilistic laws is to characterize such phenomena in terms of the behavior of their first and second moments. This approach has found important applications in statistical communications and control theory and in time series analysis. When a stochastic process is being studied in terms of its moments it is often called a time series. Consequently one would make the following definition.

A time series  $\{X(t), t \in T\}$  is a family of random variables (stochastic process) with finite second moments.

It is to be emphasized that it should not be assumed that a stochastic process which arises in practice necessarily has finite second moments and is therefore a time series. In particular, recent research [by Mandelbrot] has raised the question of whether certain series of economic observations involving price changes of stocks and commodities possess finite second moments.

Two important characteristics of a time series  $\{X(t), t \in T\}$  are its mean value function  $m(\cdot)$ , defined for all  $t$  in  $T$  by

$$m(t) = E[X(t)] ,$$

and its covariance kernel  $K(\cdot, \cdot)$ , defined for all  $s$  and  $t$  in  $T$



by\*

$$K(s, t) = \text{Cov}[X(s), X(t)] .$$

The importance of the mean value function and the covariance kernel derives from several facts:

(i) it is usually much easier to find the mean value function and the covariance kernel of a stochastic process than it is to find its complete probability law;

(ii) nevertheless, many important questions about a stochastic process can be answered on a basis of a knowledge only of its mean value function and covariance kernel;

(iii) for example, the continuity, differentiability and integrability properties of the covariance kernel lead to corresponding properties for the time series;

(iv) further, there exists an important class of time series, the normal stochastic processes, whose complete probability law is known once one knows its mean value function and covariance kernel.

Normal processes: A stochastic process  $\{X(t), t \in T\}$  is said to be a normal process if for any integer  $n$  and any subset  $\{t_1, t_2, \dots, t_n\}$  of  $T$  the  $n$  random variables  $X(t_1), \dots, X(t_n)$  are jointly normally

\*In studying the general theory of time series, it is often convenient to admit complex valued random variables. The covariance kernel is then defined by

$$K(s, t) = E[X(s)\bar{X}(t)] - m(s)\bar{m}(t)$$

where  $\bar{X}(t)$  denotes the complex conjugate of  $X(t)$  and  $\bar{m}(t)$  denotes the complex conjugate of  $m(t)$ .

distributed in the sense that their joint characteristic function is given by, for any real numbers  $u_1, u_2, \dots, u_n$ ,

$$\begin{aligned} \varphi_{X(t_1), X(t_2), \dots, X(t_n)}(u_1, u_2, \dots, u_n) &= E[\exp i(u_1 X(t_1) + \dots + u_n X(t_n))] \\ &= \exp \left\{ i \sum_{j=1}^n u_j E[X(t_j)] - \frac{1}{2} \sum_{j,k=1}^n u_j u_k \text{Cov}[X(t_j), X(t_k)] \right\} \\ &= \exp \left\{ i \sum_{j=1}^n u_j m(t_j) - \frac{1}{2} \sum_{j,k=1}^n u_j u_k K(t_j, t_k) \right\}. \end{aligned}$$

Normal processes play a basic role in time series analysis for a number of reasons:

(i) Because of the central limit theorem, many random variables which arise in applications of probability theory may be considered to be approximately normally distributed; similarly, many stochastic processes can be approximated by normal processes;

(ii) Because of the mathematical tractability of normal random variables, many questions can be more simply treated for normal processes than for other kinds of time series;

(iii) Normal processes have the useful closure property that any time series [such as  $\int_0^t X(s) ds$ ,  $X'(t)$ ,  $X(t+1) - X(t)$ ] derived by means of linear operations on a normal process, is itself a normal process;

(iv) For a normal process, one obtains a knowledge of the complete probability law of the process from a knowledge of the mean value function  $m(\cdot)$  and the covariance kernel  $K(\cdot, \cdot)$ . Conversely it may be shown that if  $m(\cdot)$  and  $K(\cdot, \cdot)$  are the mean value function and covariance kernel of some time series, then there is a (unique) normal process with this mean value function and covariance kernel.

## 2. A time series as a probability measure on function space.

A time series  $\{X(t), t \in T\}$  is for many purposes best regarded as an observation on a random phenomenon each of whose possible outcomes is a real valued function with domain  $T$ . In other words, a time series  $\{X(t), t \in T\}$  is a collection of real valued functions with domain  $T$ , one of which is observed whenever a sample is taken. The observed function is therefore called a sample function, or realization, of the time series.

Given an index set  $T$ , we let  $\Omega_T$  denote the space of all real valued functions with domain  $T$ . A point  $\omega$  belonging to  $\Omega_T$  is a function on  $T$  whose value at a point  $t$  in  $T$  we denote by  $\omega(t)$ ; consequently, we may write  $\omega = \{\omega(t), t \in T\}$ .

The problem of analyzing a time series  $\{X(t), t \in T\}$  can be expressed as the problem of finding the probability function  $P[\cdot]$ , defined on suitable subsets of  $\Omega_T$  called the measurable subsets, which describes the probability distribution of possible values of the time series (in the intuitive sense that for any measurable subset  $A$  of  $\Omega_T$ ,  $P[A]$  is approximately the relative frequency of observations in a very long sequence of independent observations of the time series which are members of  $A$ .)

In this section we show how given the finite dimensional probability laws of a time series one can construct a probability measure on a suitable family of subsets of the function space  $\Omega_T$ . This probability measure will enable us to define the notion of the probability density functional of a time series which plays a central role in modern time series analysis [see Parzen (1962, 1963)].

Let  $\mathcal{A}$  be a collection of subsets of  $\Omega_T$  which contains  $\Omega_T$  as a member. A function  $P$ , with domain  $\mathcal{A}$ , is said to be a probability measure if it possesses the following properties:

Axiom 1. For every  $A \in \mathcal{A}$ ,  $P[A]$  is well defined and is a non-negative real number; in symbols,

$$P[A] \geq 0 ;$$

Axiom 2.  $P[\Omega_T] = 1$ ;

Axiom 3. For any sequence of disjoint sets  $A_1, A_2, \dots$  belonging to  $\mathcal{A}$ , whose union  $\bigcup_{n=1}^{\infty} A_n$  belongs to  $\mathcal{A}$ ,

$$P\left[\bigcup_{n=1}^{\infty} A_n\right] = \sum_{n=1}^{\infty} P[A_n] ;$$

the sets  $\{A_n\}$  are said to be disjoint (or non-overlapping) if for any two distinct indices  $j$  and  $k$  the intersection of  $A_j$  and  $A_k$  is empty,

$$A_j \cap A_k = \emptyset$$

where  $\emptyset$  denotes the empty set.

Axiom 3 is referred to as the countable additivity or sigma-additivity property of the probability function  $P$ .

The sets belonging to  $\mathcal{A}$  are called measurable sets or events. An event  $A$  is said to occur if the function representing the actual time series observed belongs to  $A$ .

In order to guarantee that the usual operations of analysis will lead to measurable sets, it is necessary to require that the family of measurable sets be a sigma-field.

A collection  $\mathcal{A}$  of subsets of  $\Omega_T$  is called a sigma-field if it has the following properties:

Axiom 1.  $\Omega_T$  belongs to  $\mathcal{A}$  (written symbolically:  $\Omega_T \in \mathcal{A}$ );

Axiom 2. If  $A$  belongs to  $\mathcal{A}$  then the complement  $A^c$  belongs to  $\mathcal{A}$  (written symbolically:  $A \in \mathcal{A}$  implies  $A^c \in \mathcal{A}$ );

Axiom 3. For any sequence  $A_1, A_2, \dots$  belonging to  $\mathcal{A}$ , the union  $\bigcup_{n=1}^{\infty} A_n$  belongs to  $\mathcal{A}$  (written symbolically:  $\{A_n\} \subset \mathcal{A}$  implies  $\bigcup_{n=1}^{\infty} A_n \in \mathcal{A}$ ).

In words, a sigma-field is a family of sets which contains the entire space  $\Omega_T$  and is closed under the operations of forming complements and countable unions. It then follows that it is closed under the operation of forming countable intersections:

Property 4. For any sequence  $A_1, A_2, \dots$  belonging to  $\mathcal{A}$ , the intersection  $\bigcap_{n=1}^{\infty} A_n \in \mathcal{A}$ .

An important example of a sigma-field is the family of all subsets of  $\Omega_T$ .

The question naturally arises: cannot all subsets of  $\Omega_T$  be made events so that it is never necessary to consider a sigma-field  $\mathcal{A}$  smaller than the family of all subsets of  $\Omega_T$ ? Unfortunately the general answer to this question is in the negative; if we desire the probability

function  $P$  to be countably additive, it is usually the case that the family  $\mathcal{A}$  of events cannot contain all subsets of  $\Omega_T$ . For example, even in the simple case that  $T$  consists of a single point so that  $\Omega_T$  is just the real line (the set of all real numbers  $\omega$  satisfying  $-\infty < \omega < \infty$ ), there is no probability function  $P$  defined on all subsets of  $\Omega_T$  that agrees with the ordinary notion of length on the subintervals of  $\Omega_T$  (see Halmos (1950), p. 70). Usually, when  $\Omega_T$  is the real line, one adopts as the family of events the family  $\mathcal{B}$  of Borel sets, where  $\mathcal{B}$  is defined as the smallest sigma-field containing as members all intervals.

The idea that, for each  $t$ ,  $X(t)$  is a random variable can now be made precise by the following definition: for each  $t$  in  $T$ ,  $X(t)$  is a function on  $\Omega_T$  whose value at a point  $\omega$  in  $\Omega_T$ , denoted by  $X(t, \omega)$ , is given by the value at  $t$  of the function  $\omega$ :

$$(2.1) \quad X(t, \omega) = \omega(t) .$$

In order to regard  $X(t)$  as a random variable, there must exist

(i) a family  $\mathcal{A}$  of subsets of  $\Omega_T$ , called the measurable sets, such that for all  $t$  in  $T$  and all real numbers  $x$

$$(2.2) \quad \{\omega: X(t, \omega) < x\} \in \mathcal{A} ,$$

and (ii) a probability measure  $P$  with domain  $\mathcal{A}$ .

A function  $X$  with domain  $\Omega_T$  is said to be  $\mathcal{A}$ -measurable if for every real number  $x$

$$\{\omega: X(\omega) < x\} \in \mathcal{A}.$$

The problem at hand is to find a sigma-field of subsets of  $\Omega_T$  such that each random variable  $X(t)$  is  $\mathcal{A}$ -measurable.

In the function space  $\Omega_T$ , there is a smallest sigma-field of events which should belong to the sigma-field of measurable sets. Let  $\mathcal{A}_T$  denote the smallest sigma-field of subsets of  $\Omega_T$  which contains all sets of the form

$$\{\omega: \omega(t) < x\}$$

where  $t$  is a point in  $T$  and  $x$  is a real number. It is clear that, for every  $t$  in  $T$ ,  $X(t)$  is  $\mathcal{A}_T$ -measurable.

Consequently, given a probability measure  $P$  on  $\mathcal{A}_T$ , one can define by (2.1) a time series  $\{X(t), t \in T\}$  consisting of random variables with domain  $\Omega_T$  and the finite dimensional probability distributions of the time series would be given by the formula

$$\begin{aligned} F_{X(t_1), \dots, X(t_n)}(x_1, \dots, x_n) &= \text{Probability}[X(t_1) \leq x_1, \dots, X(t_n) \leq x_n] \\ (2.3) \quad &= P[\{\omega \in \Omega_T: \omega(t_j) \leq x_j \quad \text{for } j = 1, 2, \dots, n\}] . \end{aligned}$$

Basic to the theory of time series analysis is the fact that the converse holds (which follows from a celebrated theorem proved by Kolmogorov (1933)).

Kolmogorov's celebrated existence theorem regarding the probability measure on function space  $\Omega_T$  induced by a stochastic process  $\{X(t), t \in T\}$ .

Let  $\{X(t), t \in T\}$  be a stochastic process with preassigned finite dimensional probability distributions. Then there exists a unique probability measure  $P_X$  on the sigma-field  $\mathcal{A}_T$  satisfying (2.3).

To prove Kolmogorov's theorem, one considers a somewhat more general problem.

Let  $T$  be an index set. Given a family of finite dimensional characteristic functions,

$$(2.4) \quad \varphi_{t_1, \dots, t_n}(u_1, \dots, u_n),$$

$n$  an integer and  $t_1, \dots, t_n$  points in  $T$ ,

what conditions need this family satisfy in order that there exist a stochastic process  $\{X(t), t \in T\}$  whose true finite dimensional characteristic functions coincide with the given set (2.4):

$$(2.5) \quad \varphi_{X(t_1), \dots, X(t_n)}(u_1, \dots, u_n) = \varphi_{t_1, \dots, t_n}(u_1, \dots, u_n).$$

In view of (2.5), it is obvious that the given set (2.4) must be mutually consistent in the sense that (i) if  $\alpha_1, \dots, \alpha_n$  is a permutation of  $1, \dots, n$ , then for any points  $t_1, \dots, t_n$  and real numbers  $u_1, \dots, u_n$ ,

$$(2.6) \quad \varphi_{t_{\alpha_1}, \dots, t_{\alpha_n}}(u_{\alpha_1}, \dots, u_{\alpha_n}) = \varphi_{t_1, \dots, t_n}(u_1, \dots, u_n),$$

since the order in which the random variables are listed is irrelevant,



and (ii) if  $m < n$ ,

$$(2.7) \quad \varphi_{t_1, \dots, t_m}(u_1, \dots, u_m) = \varphi_{t_1, \dots, t_m, t_{m+1}, \dots, t_n}(u_1, \dots, u_m, 0, \dots, 0) .$$

The content of Kolmogorov's theorem is that these consistency conditions are the only conditions that need be imposed.

In order to give a precise statement of Kolmogorov's theorem we introduce the notion of a semi-infinite interval.

A subset  $C$  of  $\Omega_T$  is called a semi-infinite interval if it is of the form

$$(2.8) \quad C = \{\omega \in \Omega_T: \omega(t_1) \leq x_1, \dots, \omega(t_n) \leq x_n\}$$

for some integer  $n$ , points  $t_1, \dots, t_n$  belonging to  $T$ , and real numbers  $x_1, \dots, x_n$ . Note that to specify a semi-infinite rectangle one must specify an integer  $n$ ,  $n$  points  $t_1, \dots, t_n$  in the index set  $T$ , and  $n$  real numbers  $x_1, \dots, x_n$ .

Example 2A.

A semi-infinite interval. Let  $T = [0,1]$  so that  $\Omega_T$  consists of all functions on the interval 0 to 1. Let  $n = 11$ ,  $t_j = j/10$  for  $j = 0, 1, \dots, 10$ , and  $x_j = j/2$  for  $j = 0, 1, \dots, 10$ . Then

$$C = \{\omega \in \Omega(0,1): \omega(j/10) \leq j/2 \quad \text{for } j = 0, 1, \dots, 10\} .$$

Consider the following functions defined on the interval 0 to 1:

$$\begin{aligned}f_1(t) &= t + 1 \\f_2(t) &= t^2, \\f_3(t) &= \sin 5t.\end{aligned}$$

It may be shown that  $f_1(\cdot)$  does not belong to  $C$ , while  $f_2(\cdot)$  and  $f_3(\cdot)$  do belong to  $C$ .

Let  $P$  be a function defined on semi-infinite intervals in  $\Omega_T$  as follows: for  $C$  given by (2.8)

$$(2.9) \quad P[C] = F_{t_1, \dots, t_n}(x_1, \dots, x_n)$$

where  $F_{t_1, \dots, t_n}(x_1, \dots, x_n)$  is the distribution function corresponding to the characteristic function  $\varphi_{t_1, \dots, t_n}(x_1, \dots, x_n)$ .

Kolmogorov's theorem: There exists a unique probability measure  $P$  on  $\mathcal{A}_T$  whose values on semi-infinite intervals satisfies (2.9). Further, define a family of functions  $\{X(t), t \in T\}$  on  $\Omega_T$  as follows: for each  $\omega \in \Omega_T$ , the value of  $X(t)$  at  $\omega$ , denoted  $X(t, \omega)$ , is given by

$$X(t, \omega) = \omega(t), \quad \text{the value at } t \text{ of the function } \omega.$$

Then  $\{X(t), t \in T\}$  is a stochastic process, defined on the probability space  $(\Omega_T, \mathcal{A}_T, P)$  whose finite dimensional characteristic functions satisfy (2.5).

The proof of Kolmogorov's theorem requires a background in measure theory which is beyond the scope of this paper [for a proof, see Kolmogorov (1933), p. 29, or Loève (1960), p. 93].

Example 2B.

An application of Kolmogorov's theorem. Consider a time series  $\{X(t), t \in T\}$  defined by the formula

$$(2.10) \quad X(t) = \eta \varphi(t), \quad t \in T,$$

where  $\eta$  is a random variable (with finite second moment) and  $\varphi(\cdot)$  is a non-random function (for example,  $\varphi(t) = t$ ).

The statement that  $\eta$  is a random variable is not completely explicit; we also want to know the space  $\Omega$  on which  $\eta$  is defined as a function.

The random variables  $X(t)$  defined by (2.10) are functions on the same space  $\Omega$  as is  $\eta$ . Since this space may be unknown, it is sometimes useful to redefine  $X(t)$  as follows.

Given the probability distribution of  $\eta$ , the family of random variables  $\{X(t), t \in T\}$  induces a probability measure  $P$  on  $\mathcal{A}_T$  by means of the formula

$$\begin{aligned} P[\{\omega \in \Omega_T: \omega(t_1) \leq x_1, \dots, \omega(t_n) \leq x_n\}] \\ (2.11) \quad &= \text{Prob}[X(t_1) < x_1, \dots, X(t_n) < x_n] \\ &= \text{Prob}[\eta \varphi(t_1) < x_1, \dots, \eta \varphi(t_n) < x_n]. \end{aligned}$$

Next, the random variables  $\{X(t), t \in T\}$  can be redefined to be functions on  $\Omega_T$  by the formula

$$(2.12) \quad X(t, \omega) = \omega(t) .$$

The new family of random variables  $\{X(t), t \in T\}$  obtained from the definition (2.12) can be identified with the old family obtained from definition (2.10) since they have the same finite dimensional distributions.

By the procedure just described, a time series  $\{X(t), t \in T\}$  defined by an explicit formula such as (2.10) can be regarded as being equivalent to a probability measure  $P$  on the function space  $\Omega_T$ .

In particular, suppose that the random variable  $\eta$  is normal with mean  $\mu$  and variance  $\sigma^2$ . Then  $X(\cdot)$ , defined by (2.10) is a normal process with mean value function

$$(2.13) \quad m(t) = E[X(t)] = \mu \varphi(t) ,$$

and covariance kernel

$$(2.14) \quad K(s, t) = \sigma^2 \varphi(s) \varphi(t) .$$

Any normal process  $\{X(t), t \in T\}$  with the foregoing mean value function and covariance kernel induces the same probability measure  $P$  on function space  $\Omega_T$ .

Using the representation theory of time series (Parzen (1961), p. 962), it may be shown that a time series  $\{X(t), t \in T\}$  with mean value function and covariance kernel given by (2.13) and (2.14) respectively may be represented in the form of (2.10), where  $\eta$  is now a random variable

whose domain is the function space  $\Omega_T$ . Consequently when considering time series of the form of (2.10) one may assume that all random variables are functions on  $\Omega_T$ .

### 3. Probability density functionals and orthogonal measures.

The probability theory of time series is concerned with investigating the structure of a time series  $\{X(t), t \in T\}$  whose corresponding probability measure  $P$  on the function space  $\Omega_T$  is assumed known. The statistical theory of time series is concerned with a time series  $\{X(t), t \in T\}$  whose probability measure  $P$  is not known exactly but is only known to belong to a class of probability measures  $\{P_\theta, \theta \in \Phi\}$  on  $\Omega_T$ ; the class of possible probability measures  $P_\theta$  can be assumed to be indexed by a parameter  $\theta$  varying in a parameter set  $\Phi$ . Consequently an important step in developing a theory of time series analysis is to examine the relations that can exist between two probability measures  $P_1$  and  $P_2$  with a common domain  $\mathcal{A}$ .

Absolutely continuous and orthogonal probability measures. Let  $\Omega$  be a set (the sample description space) and let  $\mathcal{A}$  be a sigma-field of subsets of  $\Omega$ . Let  $P_1$  and  $P_2$  be probability measures with domain  $\mathcal{A}$ .

We say that  $P_1$  is absolutely continuous with respect to  $P_2$  if for every set  $A$  in  $\mathcal{A}$

$$(3.1) \quad P_2(A) = 0 \quad \text{implies} \quad P_1(A) = 0.$$

We write  $P_1 \ll P_2$  if  $P_1$  is absolutely continuous with respect to  $P_2$ ; the motivation for this notation is the idea that  $P_1$  is small whenever  $P_2$  is small.

We say that  $P_1$  and  $P_2$  are equivalent, denoted  $P_1 \equiv P_2$ , if each is absolutely continuous with respect to the other; in symbols,  $P_1 \equiv P_2$  if and only if  $P_1 \ll P_2$  and  $P_2 \ll P_1$ .

In order that  $P_1$  not be absolutely continuous with respect to  $P_2$  it is necessary and sufficient that there exist a set  $A$  in  $\mathcal{A}$  such that

$$(3.2) \quad P_2(A) = 0 \quad \text{and} \quad P_1(A) > 0.$$

We say that  $P_1$  and  $P_2$  are orthogonal (or perpendicular), denoted  $P_1 \perp P_2$ , if there exists a set  $A$  in  $\mathcal{A}$  such that

$$(3.3) \quad P_2(A) = 0 \quad \text{and} \quad P_1(A) = 1.$$

One can regard (3.2) as the extreme case of not being absolutely continuous.

It may be shown that if  $P_1 \ll P_2$ , then probability statements in terms of  $P_1$  can be expressed in terms of  $P_2$ ; more precisely, if  $P_1 \ll P_2$ , then there exists a function  $p$ , called the probability density function of  $P_1$  with respect to  $P_2$ , such that for any  $A$  in  $\mathcal{A}$ ,

$$(3.4) \quad P_1(A) = \int_A p \, dP_2.$$

In words, (3.4) says that to evaluate  $P_1(A)$  one integrates the function  $p$  with respect to the measure  $P_2$  over the set  $A$ . More generally,

for any  $\mathcal{A}$ -measurable function  $g$  which is integrable with respect to  $P_1$ ,

$$(3.5) \quad E_{P_1}[g] = \int_{\Omega} g \, dP_1$$

is finite, then there holds the transformation formula

$$(3.6) \quad E_{P_1}[g] = \int_{\Omega} g \, p \, dP_2 .$$

Three problems preliminary to time series analysis. To develop a general theory of time series analysis, one must begin with an understanding of the relations that can exist between two probability measures  $P_1$  and  $P_2$  on a function space. Consequently, we may speak of three problems preliminary to time series analysis:

- (i) determine whether two given probability measures  $P_1$  and  $P_2$  are orthogonal,
- (ii) determine whether one is absolutely continuous with respect to the other,
- (iii) if  $P_2 \ll P_1$ , determine the probability density functional or Radon-Nikodym derivative, denoted

$$p_{2,1} = \frac{dP_2}{dP_1} .$$

One aim of modern time series analysis is to develop ways of determining answers to these questions.



4. Signal detection and likelihood ratios.

In this section we examine the problem of detecting a signal in noise, and show how the proper formulation of this problem requires a consideration of the relations that exist between probability measures.

Many observed time series can be represented as the sum of a signal process and a noise process. More precisely, let  $T$  be a set of points, called the index set, such that at each  $t$  in  $T$  one has made (or one can make) an observation, denoted  $X(t)$ . The set of observations  $\{X(t), t \in T\}$  is a function on  $T$  which is assumed to be the sum of two other functions  $\{S(t), t \in T\}$  and  $\{N(t), t \in T\}$ :

$$X(t) = S(t) + N(t) , \quad t \in T .$$

We call  $\{S(t), t \in T\}$  the signal since it is supposed to represent the true value of the quantity being measured, while  $\{N(t), t \in T\}$  is called the noise since it represents "errors", "fluctuations", or "residuals" by which the observed function  $\{X(t), t \in T\}$  differs from the desired function  $\{S(t), t \in T\}$ .

The aim of time series analysis is to infer, from the observations, information about one or more features of the signal. The aspects of the signal in which one is interested depends on the assumptions one makes about the structure of the signal and noise processes. In this section we consider the important problem of detecting a signal in the presence of noise.

Let  $\{S(t), t \in T\}$  and  $\{N(t), t \in T\}$  be time series, called respectively the signal process and the noise process. Given an observed time series  $\{X(t), t \in T\}$ , one desires to test the hypothesis

$$H_0: X(\cdot) = N(\cdot) , \quad \text{noise alone is present} ,$$

against the alternative hypothesis

$$H_1: X(\cdot) = S(\cdot) + N(\cdot), \quad \text{signal plus noise is present,}$$

by choosing a subset  $R_1$  of the sample space  $\Omega_T$ , of possible realizations of the time series  $\{X(t), t \in T\}$ , which will be the rejection region for  $H_0$ ; that is, one says signal plus noise is present if the observed time series  $\{X(t), t \in T\}$  belongs to  $R_1$ , and one says that noise alone is present if  $\{X(t), t \in T\}$  does not belong to  $R_1$ .

In this section, we suppose that the probability distribution of  $\{X(t), t \in T\}$  under the hypotheses  $H_0$  and  $H_1$  are well defined; in this case we say that  $H_0$  and  $H_1$  are simple hypotheses. We may then introduce probability measures  $P_N$  and  $P_{S+N}$  defined on the measurable subsets  $B$  of the sample space  $\Omega_T$  by

$$\begin{aligned} P_N[B] &= \text{Prob} [\{N(t), t \in T\} \in B] \\ (4.1) \qquad &= \text{Prob} [\{X(t), t \in T\} \in B | H_0], \end{aligned}$$

$$\begin{aligned} P_{S+N}[B] &= \text{Prob} [\{S(t) + N(t), t \in T\} \in B] \\ (4.2) \qquad &= \text{Prob} [\{X(t), t \in T\} \in B | H_1]. \end{aligned}$$

Example 4A.

A specified signal in normal noise. One important case in which  $H_0$  and  $H_1$  are simple hypotheses is when the following assumptions hold.

Assumption on the index set: The index set  $T$  is a finite set  $\{t_1, t_2, \dots, t_n\}$ .

Assumption on the noise process: The noise process  $\{N(t), t \in T\}$  is assumed to possess finite second moments, to have zero means:

$$(4.3) \quad E[N(t)] = 0,$$

and known covariance kernel  $K$ :

$$(4.4) \quad E[N(s) N(t)] = K(s, t).$$

Further,  $\{N(t), t \in T\}$  is a normal process; that is, the  $n$  random variables  $N(t_1), \dots, N(t_n)$  are jointly normally distributed so that their joint characteristic function is given by

$$(4.5) \quad \begin{aligned} \varphi_{N(t_1), \dots, N(t_n)}(u_1, \dots, u_n) &= E \left[ \exp \left\{ i \sum_{j=1}^n u_j X(t_j) \right\} \right] \\ &= \exp \left[ -\frac{1}{2} \sum_{j,k=1}^n u_j K(t_j, t_k) u_k \right]. \end{aligned}$$

Assumption on the signal process: The signal  $\{S(t), t \in T\}$  is a known non-random function. The signal plus noise process  $\{X(t) = S(t) + N(t), t \in T\}$  is then a normal process with mean value function

$$(4.6) \quad E[X(t)] = S(t)$$

and covariance kernel

$$(4.7) \quad \text{Cov}[X(s), X(t)] = K(s, t) .$$

Under the assumptions of this example,  $P_N$  is the probability measure on function space induced by a normal process with mean value function identically zero and covariance kernel  $K$ , while  $P_{S+N}$  is the probability measure induced by a normal process with mean value function equal to the signal function  $S(t)$  and covariance kernel  $K$ .

Perfect detectability and the singular detection problem: We say that the hypotheses  $H_0$  and  $H_1$  are perfectly detectable, or that the problem of detecting the signal process  $S(\cdot)$  in the presence of the noise process  $N(\cdot)$  is singular, if there exists a set  $A$  in the sample space  $\Omega_T$  such that

$$(4.8) \quad P_N[A] = 0 , \quad P_{S+N}[A] = 1 .$$

By choosing  $A$  as the rejection region  $R_1$  for  $H_0$  one has probability zero of incorrectly identifying noise as signal plus noise or signal plus noise as noise. Note that the probability measures  $P_N$  and  $P_{S+N}$  are orthogonal if (4.8) holds. By definition, then, the hypotheses  $H_0$  and  $H_1$  are perfectly detectable if and only if  $P_N$  and  $P_{S+N}$  are orthogonal.

The regular detection problem: The problem of detecting the signal process in the presence of the noise process  $N(\cdot)$  is called regular if

$P_N$  and  $P_{S+N}$  are not orthogonal. In this case, new principles must be introduced in order to optimally choose the rejection region  $R_1$ . First, one distinguishes two types of errors that can occur:

(i) a false alarm (or error of type I) occurs when one says that a signal plus noise is present when in fact noise alone is present;

(ii) a detection failure (or error of type II) occurs when one says that noise alone is present when in fact signal plus noise is present.

A rejection region  $R$  is then characterized by two numbers  $\alpha$  and  $\beta$ , defined by

$$\begin{aligned} \alpha &= \text{Prob} [\text{false alarm}] \\ (4.9) \quad &= \text{Prob} [(X(t), t \in T) \in R | H_0] , \end{aligned}$$

$$\begin{aligned} \beta &= \text{Prob} [\text{detection failure}] \\ (4.10) \quad &= 1 - \text{Prob} [(X(t), t \in T) \in R | H_1] . \end{aligned}$$

In certain cases it may be possible to assign a numerical measure to the seriousness of a false alarm and of a detection failure; one denotes these costs by\*  $L_1$  and  $L_2$  respectively. Further it may be possible to determine the fraction  $\pi_s$  of experimental situations in which signal plus noise is present; one calls  $\pi_s$  the prior probability that signal

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\*In the general theory of hypothesis testing, a false alarm is called a type I error, and  $L_1$  is the cost of a type I error, while a detection failure is called a type II error and  $L_2$  is the cost of a type II error.

is present. To each critical region one can assign a risk  $\rho$ , called the Bayes risk and defined as the expected cost of an incorrect decision:

$$(4.11) \quad \rho = \alpha(1 - \pi_S) L_1 + \beta\pi_S L_2 .$$

The Bayes rejection region  $R$  (or optimum rejection region according to the Bayes criterion) is defined as the region which minimizes  $\rho$ , the expected cost of an incorrect decision.\*\*

It may be difficult to use a Bayes rejection region for one or both of the following reasons:

- (i) because of difficulty in assigning the losses  $L_1$  and  $L_2$ ,
- (ii) because of difficulty in assigning the prior probability  $\pi_S$ .

In these circumstances one may use the Neyman-Pearson rejection region (or optimum rejection region according to the Neyman-Pearson criterion) which is defined as the rejection region  $R$  minimizing  $\beta$ , the detection failure probability, subject to the restriction that  $\alpha$ , the false alarm probability, is less than or equal to some desired level  $\alpha_0$ .

We next show how one may determine the Bayes rejection region and the Neyman-Pearson rejection region by introducing probability density functionals.

Let us assume that there exists a measure  $Q$  on the measurable subsets of the sample space  $\Omega_T$ , and functions  $p_N$  and  $p_{S+N}$  with domain  $\Omega_T$ , with the property that, for every measurable subset  $B$  of  $\Omega_T$ ,

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\*\*Note that a rejection region which minimizes the average probability of error  $\pi_e$ ,  $\pi_e = \alpha(1 - \pi_S) + \beta\pi_S$ , is the same as the Bayes rejection region with unit costs,  $L_1 = L_2 = 1$ .

$$(4.12) \quad P_N[B] = \int_B p_N dQ$$

$$(4.13) \quad P_{S+N}[B] = \int_B p_{S+N} dQ .$$

In order to emphasize that its argument is a function  $(X(t), t \in T)$ , we call  $p_N$  a functional, and sometimes denote it by  $p_N(X(t), t \in T)$ . It is called the probability density functional of  $P_N$  with respect to  $Q$ . The function  $p_N$  may be written symbolically as a derivative,

$$(4.14) \quad p_N = \frac{dP_N}{dQ} ,$$

and is then called the Radon-Nikodym derivative of  $P_N$  with respect to  $Q$ . Similarly,  $p_{S+N}$  is called the probability density functional, or Radon-Nikodym derivative, of  $P_{S+N}$  with respect to  $Q$ .

In order for the probability density functionals  $p_N$  and  $p_{S+N}$  to exist, there must exist a measure  $Q$  with respect to which both  $P_N$  and  $P_{S+N}$  are absolutely continuous; recall that  $P_N$  is absolutely continuous with respect to  $Q$  if

$$(4.15) \quad Q(A) = 0 \quad \text{implies} \quad P_N(0) \quad \text{for all } A \subset \Omega .$$

Such a measure is given by

$$(4.16) \quad Q = P_N + P_{S+N} .$$



Consequently there is no loss of generality in assuming that there exist functions  $p_N$  and  $p_{S+N}$  satisfying (4.12) and (4.13) respectively.

In terms of probability density functionals, the false alarm probability  $\alpha$  and detection failure probability  $\beta$  of a rejection region are given by

$$(4.17) \quad \alpha = \int_R p_N dQ$$

$$(4.18) \quad \beta = 1 - \int_R p_{S+N} dQ .$$

Consequently, the Bayes risk  $\rho$  of a rejection region  $R$  is given by

$$(4.19) \quad \rho = \int_R \{ (1 - \pi_S) L_1 p_S - \pi_S L_2 p_{S+N} \} dQ + \pi_S L_2 .$$

To minimize  $\rho$ , one should choose  $R$  as the set of observations  $\{X(t), t \in T\}$  for which the integrand in (4.19) is negative, so that the Bayes rejection region  $R$  is given by

$$(4.20) \quad R = \left\{ \left( X(t), t \in T \right) : \frac{p_{S+N}}{p_N} > \frac{(1 - \pi_S) L_1}{\pi_S L_2} \right\} .$$

The ratio

$$(4.21) \quad \frac{p_{S+N}}{p_N}$$

of probability density functionals is called in classical statistical literature the likelihood ratio since  $p_N(X(t), t \in T)$  is defined to be

the likelihood that noise alone is present given that the observed time series was  $\{X(t), t \in T\}$ , and  $p_{S+N}(X(t), t \in T)$  is the likelihood that signal plus noise is present given that the observed time series was  $\{X(t), t \in T\}$ .

The likelihood ratio (4.21) has a probabilistic meaning which shows that the measure  $Q$  used in defining the likelihood ratio plays no role. Let

$$(4.22) \quad A = \left\{ \left( X(t), t \in T \right) : p_{S+N} > 0 \quad \text{and} \quad p_N = 0 \right\}.$$

If  $p_{S+N}[A] = 0$ , then  $p_{S+N}$  is absolutely continuous with respect to  $p_N$ , and the Radon-Nikodym derivative of  $p_{S+N}$  with respect to  $p_N$  is given by the likelihood ratio:

$$(4.23) \quad \frac{dp_{S+N}}{dp_N} = \frac{p_{S+N}}{p_N}.$$

The Radon-Nikodym derivative of  $p_{S+N}$  with respect to  $p_N$  is called the probability density functional of signal plus noise with respect to noise; where no ambiguity can arise it is denoted by  $p$ . To summarize, the probability density functional  $p$  of signal plus noise with respect to noise is a function on the sample space  $\Omega_T$  satisfying

$$(4.24) \quad p_{S+N}[B] = \int_B p \, dp_N, \quad B \subset \Omega,$$

and exists if and only if  $p_{S+N}$  is absolutely continuous with respect to  $p_N$ , which holds if and only if  $p_{S+N}[A] = 0$ , where  $A$  is defined

by (4.22); it then follows that  $p$  is given by the likelihood ratio

$$(4.25) \quad p = \frac{P_{S+N}}{P_N} .$$

If  $P_{S+N}[A] > 0$ , instead of (4.24), one has for every measurable subset  $B$  of  $\Omega$ ,

$$(4.26) \quad P_{S+N}[B] = \int_B p \, dP_N + P_{S+N}[AB] ,$$

where  $p$  is still given by (4.25); (4.26) is an example of the Lebesgue decomposition theorem which states that to the probability measures  $P_N$  and  $P_{S+N}$  there exists a function  $p$  on  $\Omega$  and set  $A$  such that  $P_N[A] = 0$  and (4.26) holds. We have made this assertion concrete by showing how one may find  $p$  and  $A$ .

Optimum detectors: A random variable  $U$  (that is, a function on the sample space  $\Omega$ ) which has the property that the rejection region which is optimum according to a certain criterion may be expressed in terms of the values of  $U$  is called an optimum detector according to that criterion.

From (4.20) it follows that the likelihood ratio is an optimum Bayes detector. Indeed the optimum rejection region according to the Bayes criterion consists of all observations  $\{X(t), t \in T\}$  for which the likelihood ratio is above a certain threshold value  $\Lambda_0$  given by

$$(4.27) \quad \Lambda_0 = \frac{(1 - \pi_s) L_1}{\pi_s L_2} .$$

Similarly, it may be shown that the likelihood ratio is an optimum Neyman-Pearson detector. Indeed the optimum Neyman-Pearson rejection region according to the Neyman-Pearson criterion consists of all observations  $\{X(t), t \in T\}$  for which the likelihood ratio is above a certain threshold  $\Lambda_0$ , determined by the condition that the false alarm probability be equal to  $\alpha_0$ :

$$(4.28) \quad P_N \left[ \frac{p_{S+N}}{p_N} > \Lambda_0 \right] = \alpha_0 .$$

To prove this assertion, one uses the fact that the Neyman-Pearson rejection region  $R$  is that region which subject to the condition

$$(4.29) \quad \int_R p_N dQ \leq \alpha$$

maximizes

$$(4.30) \quad \int_R p_{S+N} dQ .$$

Intuitively, one sees that the optimum rejection region  $R$  should contain those sample points which have the highest value of the likelihood ratio (4.21) since for a given contribution to the integral in (4.29) these points make a maximum contribution to the integral in (4.30). A formal proof of this assertion is easily given, using the fundamental lemma of Neyman and Pearson.

Example 4B.

Detection of a specified signal in normal noise. Consider the signal and noise processes described in example 4A. Assume that the covariance matrix

$$K = \begin{bmatrix} K(t_1, t_1) & \dots & K(t_1, t_n) \\ \vdots & & \vdots \\ K(t_n, t_1) & \dots & K(t_n, t_n) \end{bmatrix}$$

is non-singular with inverse denoted

$$K^{-1} = \begin{bmatrix} K^{-1}(t_1, t_1) & \dots & K^{-1}(t_1, t_n) \\ \vdots & & \vdots \\ K^{-1}(t_n, t_1) & \dots & K^{-1}(t_n, t_n) \end{bmatrix}.$$

In words,  $K^{-1}(t_i, t_j)$  does not denote the reciprocal of  $K(t_i, t_j)$ , but rather denotes the  $(i, j)$ -th element of the inverse matrix  $K^{-1}$ . The signal and noise process and the noise process then both possess probability density functionals  $p_{S+N}$  and  $p_N$  with respect to Lebesgue measure  $Q$ :

$$p_N(X(t), t \in T) = \left\{ (2\pi)^n |K| \right\}^{-\frac{1}{2}} \exp \left[ -\frac{1}{2} \sum_{i,j=1}^n X(t_i) K^{-1}(t_i, t_j) X(t_j) \right],$$

$$p_{S+N}(X(t), t \in T) = \left\{ (2\pi)^n |K| \right\}$$

$$\exp \left[ -\frac{1}{2} \sum_{i,j=1}^n \{X(t_i) - S(t_i)\} K^{-1}(t_i, t_j) \{X(t_j) - S(t_j)\} \right]$$

where  $|K|$  is the determinant of the matrix  $K$ .

The likelihood ratio is given by

$$\frac{P_{S+N}}{P_N} = \exp \left[ \sum_{i,j=1}^n X(t_i) K^{-1}(t_i, t_j) S(t_j) - \frac{1}{2} \sum_{i,j=1}^n S(t_i) K^{-1}(t_i, t_j) S(t_j) \right].$$

In order to write this expression more compactly, let us introduce the notation (which will play an important role in the sequel)

$$(f, g)_K = \sum_{i,j=1}^n f(t_i) K^{-1}(t_i, t_j) g(t_j),$$

defined for any two functions  $f$  and  $g$  on  $T$ . Then

$$\frac{P_{S+N}}{P_N} = \exp \left[ (X, S)_K - \frac{1}{2} (S, S)_K \right].$$

Since the likelihood ratio is a monotone increasing function of  $(X, S)_K$ , it follows that  $(X, S)_K$  is an optimum (Bayes or Neyman-Pearson) detector. Indeed, the rejection region for testing  $H_0$  against  $H_1$  can be expressed as the set of observations  $\{X(t), t \in T\}$  for which  $(X, S)_K$  is above a certain threshold  $\Lambda_1$ , say. A detector of the form

$$(X, S)_K = \sum_{i=1}^n X(t_i) \left\{ \sum_{j=1}^n K^{-1}(t_i, t_j) S(t_j) \right\}$$

is said to be a "correlation detector" or a "matched filter" since it is obtained by "correlating" or "matching" the specified signal shape  $S(t)$  with the observed time series  $X(t)$ .

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### Exercises

1. In the notation defined in Example 2A, show that  $f_1(\cdot)$  does not belong to  $C$ , while  $f_2(\cdot)$  and  $f_3(\cdot)$  do belong to  $C$ .
2. Let  $H_0$  and  $H_1$  be simple hypotheses about a time series  $(X(t), 0 \leq t \leq 1)$ . Under both hypotheses,  $(X(t), 0 \leq t \leq 1)$  is a normal process with covariance kernel

$$K(s, t) = \varphi(s) \varphi(t)$$

and mean value function respectively given by

$$H_0: E[X(t)] \equiv 0 ,$$

$$H_1: E[X(t)] = S(t) .$$

Assume that  $S(\cdot)$  and  $\varphi(\cdot)$  are orthogonal,

$$\int_0^1 S(t) \varphi(t) dt = 0 .$$

Show that  $H_0$  and  $H_1$  are perfectly detectable.

Hint:  $U = \int_0^1 S(t) X(t) dt$  is a perfect detector, since

$$P_1[U = 0] = 1, \quad P_2[U = \int_0^1 S^2(t) dt \neq 0] = 1.$$



3. Let  $H_0$  and  $H_1$  be simple hypotheses about a time series  $\{X(t), 0 \leq t \leq 1\}$ . Under both hypotheses,  $\{X(t), 0 \leq t \leq T\}$  is of the form

$$X(t) = \eta \varphi(t)$$

where  $\varphi(t)$  is a non-random function and  $\eta$  is a normal random variable,

$$\text{Under } H_0: E[\eta] = 0, \quad \text{Var}[\eta] = 1$$

$$\text{Under } H_1: E[\eta] = m, \quad \text{Var}[\eta] = 1.$$

Show that an optimum detector for testing  $H_0$  against  $H_1$  is given by

$$\exp \left[ -\frac{1}{2} \left\{ (\eta - m)^2 - \eta^2 \right\} \right] = \exp \left\{ m\eta - \frac{1}{2} m^2 \right\}$$

and therefore is given by  $\eta$ . A possible formula for  $\eta$  is

$$\eta = \frac{X(t_1)}{\varphi(t_1)}$$

for any point  $t_1$  in  $0 \leq t_1 \leq 1$  such that  $\varphi(t_1) \neq 0$ .

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