

Probability and Stochastic Processes I - Lecture 10

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II.6 Stochastic Processes

Definition II.6.1 The set $\{(t, X_t) : t \in T\}$, where X_t is a random variable defined with respect to probability model (Ω, \mathcal{A}, P) for each $t \in T$, is called a *stochastic process* (sometimes a *random process*). The set T is called the *index set* of the process.

Example II.6.1 - a single random variable X_1 with $T = \{1\}$ is equivalent to the stochastic process $\{(1, X_1)\}$

- a random vector $\mathbf{X} = (X_1, \dots, X_k)'$ is equivalent to the stochastic process $\{(t, X_t) : t \in T\}$ where $T = \{1, \dots, k\}$ ■

note - T can be a very general set like the nodes of a graph and can be an infinite set

Example II.6.2 *Tossing a coin until the first head is observed.*

- suppose a coin is tossed in "independent" tosses until the first head is observed and the number of that toss is recorded
- what is Ω ?
- denote a head by 1 and a tail by 0
- it is most convenient to put

$$\begin{aligned}\Omega &= \{(\omega_1, \omega_2, \dots) : \omega_i \in \{0, 1\}\} = X_{i=1}^{\infty} \{0, 1\} \\ &= \text{the set of all sequences of 0's and 1's}\end{aligned}$$

- then define $Y : \Omega \rightarrow R$ by $Y(\omega) = i$ when $\omega_i = 1$ and $\omega_j = 0$ when $j < i$
- is Y well defined as a r.v.?
- let $p =$ probability of a head on a single toss
- if $p = 0$ then Y is not defined

- if $p > 0$, then Y is defined since

$$\begin{aligned} P(\text{"an infinite sequence of tails"}) &= \lim_{n \rightarrow \infty} P(\text{"first } n \text{ are tails"}) \\ &= \lim_{n \rightarrow \infty} (1 - p)^n = 0 \end{aligned}$$

- then putting $A_i = \{\omega : \omega_1 = 0, \dots, \omega_{i-1} = 0, \omega_i = 1\}$ and using independence

$$\begin{aligned} p_Y(i) &= P_Y(\{i\}) = P(A_i) \\ &= (1 - p)^{i-1} p P(\{\omega : \omega_{i+1} \in \{0, 1\}, \omega_{i+2} \in \{0, 1\}, \dots\}) \\ &= (1 - p)^{i-1} p \end{aligned}$$

and $Y \sim \text{geometric}(p)$

Exercise II.6.1 Prove that p_Y defines a probability distribution.

- the point is that we need an infinite dimensional Ω

- if we define $X_i(\omega) = \omega_i$, then $\{(t, X_t) : t \in \mathbb{N}\}$ is a stochastic process, called a Bernoulli(p) process ■

note - the main reason we need to consider stochastic processes is that in many applications the dependence on the index is important

Example II.6.3

- suppose $\mathbf{X} = (X_1, \dots, X_k)'$ gives various measurements of an individual's face (for facial recognition purposes)

- then the order in which the measurements is listed doesn't matter as long as we are consistent so there is no real dependence on the index

- but now suppose \mathbf{X} contains the measurements from a physical endurance test taken on an individual one month apart where i refers to month i and then the order definitely matters and we would take this into account when analyzing data

- in many contexts $t \in T \subset \mathbb{R}^1$ refers to time as for example taking $T = \{0, 1, 2, \dots\} = \mathbb{N}_0$ and X_t is the closing price of a stock on trading day t

- sometimes $T = \mathbb{Z}$ and negative times index the past

- now some stocks trade by nanoseconds so it makes sense in such a case to take, as an approximation, $T = [0, \infty)$ or $T = \mathbb{R}^1$ which is continuous time as opposed to discrete time
- stochastic processes where t is time are often referred to as *time series*
- sometimes t is multidimensional as $t = (t_1, t_2) = (\text{longitude}, \text{latitude})$ of a point on earth and $X_{(t_1, t_2)}$ measures some characteristic of that location such as its mean temperature over a particular year and such processes are referred to as *spatial processes*
- in fact to study global warming we want $t = (t_1, t_2, t_3) = (\text{longitude}, \text{latitude}, \text{time})$ for a large grid covering the earth and for years extending into the past and future ■

- think of a realized value of a stochastic process as a function $X(\omega) : T \rightarrow R^1$, called the *sample function*, with value $X_t(\omega)$ at index t
- so a stochastic processes is in effect a probability measure on functions from T into R^1
- a stochastic process is a generalization of random variable where we start with (Ω, \mathcal{A}, P) and get P_X as follows
- $X : \Omega \rightarrow R^T =$ set of functions with domain T mapping into R^1
 $= \{g : g : T \rightarrow R^1\}$
- $\mathcal{B}^T =$ smallest σ -algebra on R^T containing all sets of the form

$$\{g : g(t_1) \in (a_1, b_1], \dots, g(t_n) \in (a_n, b_n]\}$$

for any $\{t_1, \dots, t_n\} \subset T$ and intervals $(a_1, b_1], \dots, (a_n, b_n]$

- then require $X^{-1}B \in \mathcal{A}$ for any $B \in \mathcal{B}^T$ which implies

$$P_X(B) = P(X^{-1}B)$$

Example II.6.4

- suppose $T = [0, \infty)$, $\omega \sim N(0, 1)$ and define $X_t(\omega) = \omega t$ so X gives a ray from the origin with $N(0, 1)$ distributed slope
- suppose $T = [0, 1]$, $\omega \sim \text{Uniform}(0, 10)$ and define $X_t(\omega) = \cos(\omega t)$ so X gives a cosinusoid with random frequency

- the definition of a stochastic process immediately implies that the finite dimensional distributions are consistent in the sense that the distribution of $(X_{s_1}, \dots, X_{s_m})$ can be obtained from that of $(X_{t_1}, \dots, X_{t_n})$ by marginalizing whenever $\{s_1, \dots, s_m\} \subset \{t_1, \dots, t_n\}$
- as a general way of defining stochastic processes we have the following

Proposition II.6.1 (*Kolmogorov's Consistency Theorem*)

Suppose $T \subset \mathbb{R}$ and a probability model $(R^n, B^n, P_{(t_1, \dots, t_n)})$ is given for each $\{t_1, \dots, t_n\} \subset T$. If the probability models are consistent, then there exists probability model (Ω, \mathcal{A}, P) and random variables X_t such that $\{(t, X_t) : t \in T\}$ is a stochastic process with $P_{(X_{t_1}, \dots, X_{t_n})} = P_{(t_1, \dots, t_n)}$.

Example II.6.5 *Bernoulli(p) process*

- let $T = \{1, 2, \dots\}$ and $P_{(t_1, \dots, t_n)}$ be the discrete probability measure concentrated on $\{0, 1\}^n$ given by the probability function

$$P_{(t_1, \dots, t_n)}(x_1, \dots, x_n) = \begin{cases} \prod_{i=1}^n p^{x_i} (1-p)^{1-x_i} & \text{if } (x_1, \dots, x_n) \in \{0, 1\}^n \\ 0 & \text{otherwise} \end{cases}$$

- these distributions are clearly consistent and so KCT indicates that this provides a valid definition of a s.p. $\{(t, X_t) : t \in T\}$ ■

Definition II.6.2 A s.p. $\{(t, X_t) : t \in T\}$ is a *Gaussian process* whenever

$$(X_{t_1}, \dots, X_{t_n}) \sim N_n(\mu(t_1, \dots, t_n), \Sigma(t_1, \dots, t_n))$$

where $\mu(t_1, \dots, t_n) \in R^n$, $\Sigma(t_1, \dots, t_n) \in R^{n \times n}$ p.d. for every $\{t_1, \dots, t_n\} \subset T$. ■

- does a Gaussian process exist?

Example II.6.6 *Gaussian white noise process*

- specify $\sigma^2 : T \rightarrow (0, \infty)$ and for $\{t_1, \dots, t_n\} \subset T$

$$\begin{aligned}\mu(t_1, \dots, t_n) &= (0, \dots, 0)' \\ \Sigma(t_1, \dots, t_n) &= \text{diag}(\sigma^2(t_1), \dots, \sigma^2(t_n))\end{aligned}$$

- then putting $(X_{t_1}, \dots, X_{t_n}) \sim N_n(\mathbf{0}, \text{diag}(\sigma^2(t_1), \dots, \sigma^2(t_n)))$ defines a valid s.p. (**Exercise II.6.2**) ■