

Introduction to Stochastic Process

STAT 615 Course Notes

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0 Preliminaries

Definition. 0.1. A sample space Ω is the collection of individual outcomes or realizations for a random experiment.

Subsets of Ω are called events. For $A \subset \Omega$, we say “A occurs” if the outcome $\omega \in A$

Ω can be countable (finite or countably infinite) or uncountable. It need not consist of real-valued entities. Most often, its role is in the background and simply assumed.

Example. 0.1. Toss a coin infinitely many times, observing 1 (heads) or 0 (tails) for each toss. The outcomes are infinite sequences of 0's & 1's, such as $(1, 1, 0, 0, 0, 1, \dots)$. The collection Ω of all possible outcomes is uncountable in this case.

Events are subsets of Ω , often identified descriptively:

- “1st toss is H” = $\{(x_1, x_2, \dots) : x_1 = 1\}$.

- “12 H’s in the 1st 100 tosses” = $\{(x_1, x_2, \dots) : \sum_{i=1}^{100} x_i = 12\}$.
- “the long tem proportion of H’s is $\frac{1}{3}$ ” = $\{(x_1, x_2, \dots) : \frac{1}{n} \sum_{i=1}^n x_i \rightarrow \frac{1}{3}\}$.

A singleton event is a subset of just one outcome, which is not quite the same as the outcome itself.

The empty set ϕ is considered to be an event also.

Definition. 0.2. A probability measure P is a set function (i.e. P is applied to events, not to outcomes) such that

- (i) $0 \leq P(A) \leq 1$, with $P(\Omega) = 1, P(\phi) = 0$
- (ii) if A_1, A_2, \dots are disjoint then $P(\bigcup_{i=1}^{\infty} A_i) = \sum_{i=1}^{\infty} P(A_i)$ (similarly for finite unions)

when the A_i ’s are not disjoint, we have

Theorem 0.1. (*Boole’s inequality*)

$$P(\bigcup_{i=1}^n A_i) \leq \sum_{i=1}^n P(A_i)$$

.

Definition. 0.3. A collection of \mathbf{A} of subsets of Ω is called a σ -algebra (or σ -field) if

- (i) $\phi, \Omega \in \mathbf{A}$
- (ii) $A \in \mathcal{A} \implies A^c \in \mathcal{A}$
- (iii) $A_1, A_2, \dots \in \mathcal{A} \implies \bigcup_{i=1}^{\infty} A_i \in \mathcal{A}$ (Some for finite unions) and hence also $\bigcap_{i=1}^{\infty} A_i \in \mathcal{A}$

Note: the requirement is closure under countable unions and intersections. So a union of uncountably many subsets in \mathcal{A} need not itself be in \mathcal{A} . (there are important mathematical reasons for allowing this.) **Additional Note:** complement $A^c =$ “A does not occur” union of events $\bigcup_i A_i =$ “some A_i (at least one) occurs” intersection $\bigcap_i A_i =$ “every A_i occurs”

Example. 0.2. If Ω is countable (such as the integers \mathbb{Z}) then we can let \mathcal{A} be all subsets.

Example. 0.3. If $\Omega = \mathbb{R}$ (all real values), we usually use the **Borel Sets** \mathcal{B} = smallest σ -algebra that contains all singletons and all intervals. So \mathcal{B} would include countable unions of intervals, but this by no means comes close to all it has. There is no direct formula or description to characterize all arbitrary Borel sets.

On the other hand, given $B \in \mathcal{B}$ and a probability measure P , $P(B)$ can be approximated as close as we like with the probability of some finite union of intervals.

Borel sets on \mathbb{R}^d (d-dimensional vectors) are defined similarly, using rectangles.

Example. 0.4. (cond.) Let $\Omega = \{(x_1, x_2, x_3, \dots) : x_i = 0 \text{ or } x_i = 1 \text{ for each } i \geq 1\}$ be the sample space of infinite sequences of coin tosses by cylinder events (events where a finite number of the x_i 's are fixed). For example,

$$\{(x_1, x_2, \dots) : x_1 = 1, x_2 = 0\}$$

.

The triple (Ω, \mathcal{A}, P) (i.e. sample space, σ -algebra of events probability measure) is called a probability space.

Definition. 0.4. A random variable $X = X(\omega)$ is a real-valued function applied to outcomes (i.e., $X : \Omega \rightarrow \mathbb{R}$).

An extended random variable can possibly take infinite values (i.e., $X : \Omega \rightarrow \bar{\mathbb{R}} \stackrel{def}{=} \mathbb{R} \cup \{-\infty, \infty\}$ (“ $\stackrel{def}{=}$ ” means “defined as”, the same as “=:”))

Since random variables (rvs) are functions, it is highly recommended that you take care to distinguish them from actual (possible) values. We will usually, but not always, use upper case for rvs and lower case for actual values. (Also, try to avoid using the same name, such as X or Y , for everything, and certainly make distinctions between rvs in the same problem/context.)

A random vector is a vector of rvs and a random (stochastic) process is a sequence of rvs (e.g. (x_1, x_2, \dots)) or a function that takes random values (e.g. $x(t), t \in \mathbb{R}$, where each $X(t)$ is a random variable).

Definition. 0.5. Every rv X has a cululative distribution function (cdf)

$$F_X(x) = P(X \leq x)$$

(sensible even for extended rvs)

A rv X (or its cdf) is discrete if there is a countable set of values $\{a_i\}$ such that $\sum_i P(x = a_i) = 1$, $\{P_i\} = \{P(x = a_i)\}$ is called the probability mass function (pmf).

A rv X (or its cdf) is absolutely continuous if

$$P(a < X \leq b) = F(b) - F(a) = \int_a^b f(x)dx$$

for some function $f(x)$, called the probability density function (pdf)

However, we do not wish to limit ourselves to just two types. There are, for example, rvs with neither a pmf or a pdf, and these are mixtures of types. Things become more general still for random vectors.

Furthermore, a stochastic process cannot be characterized so simply at all (More on this in the 1st chapter).

We will write $X \sim F_X$ (or $X \sim \{P_i\}$ or $X \sim f(x)$ when the context is clear) to signify that X has distribution given by F_x .

You should be familiar with all the commonly used distributions.

- ex. $Y \sim \text{Poisson}(\lambda)$ with pmf

$$P(y = k) = \frac{\lambda^k e^{-\lambda}}{k!}, k = 0, 1, 2, \dots$$

.

- ex. $T \sim \text{Gamma}(\alpha, \beta)$ with pdf

$$f(t) = \frac{\beta^\alpha t^{\alpha-1} e^{-\beta t}}{\Gamma(\alpha)}, t \geq 0$$

.

Definition. 0.6. Random variables X_1, \dots, X_m are independent if

$$P(X_1 \in A_1, X_2 \in A_2, \dots, X_m \in A_m) = \prod_{i=1}^m P(X_i \in A_i)$$

for all **Borel sets** A_1, \dots, A_m .

Equivalently,

$$P(X_1 \leq x_1, \dots, X_m \leq x_m) = \prod_{i=1}^m P(X_i \leq x_i)$$

for all real values x_1, \dots, x_m .

(The 1st definition can be applied to random elements of any kind, with appropriate A_i 's.)

A collection $\{X_t\}$ of rvs (possibly uncountably many) is independent if every finite sub-collection is. That is, if X_{t_1}, \dots, X_{t_m} are independent, for any t_1, \dots, t_m .

A sequence $\{X_n\}$ is independent and identically distributed (iid) if it is independent and all the rvs have the same distribution. We sometimes write $\{X_n\} \stackrel{iid}{\sim} F$ (or just $X_n \stackrel{iid}{\sim} F$).

Stochastic processes are generally not independent (they wouldn't be interesting if they were). But they often can be constructed or described in terms of independent rvs.

Random variables and their distributions can also be characterized in terms of expectations. As before, we do not wish to be limited to special types (and in fact it is notationally more convenient not to be).

Definition. 0.7. Here is how the expectation $E(g(x))$ is defined (if possible).

For simplicity, just let $Y = g(X)$ and we'll define $E(Y)$.

1. indicator rv

if

$$Y = \mathbb{1}_A = \begin{cases} 0, & \text{if } A \text{ does not occur} \\ 1, & \text{if } A \text{ does occur} \end{cases}$$

then

$$E(\mathbb{1}_A) = P(A)$$

. indicator rvs have Bernoulli distribution (which simply means discrete on $\{0, 1\}$)

2. simple rv

Y is a finite linear combination of indicator rvs

i.e. $Y = \sum_{i=1}^m c_i \mathbb{1}_{A_i}$ (hence discrete w/ finitely many values)

then

$$E(Y) = \sum_{i=1}^m c_i P(A_i)$$

. (this works even if the A_i 's are not disjoint or if the c_i 's are not distinct.)

3. nonnegative rv

$Y \geq 0$. There always exist nonnegative simple rvs $Y_n(\omega) \uparrow Y(\omega)$ for every $\omega \in \Omega$.

Then

$$E(Y) = \lim_{n \rightarrow \infty} E(Y_n) \quad (\text{which myght equal } \infty)$$

(This does not depend on the choice of sequenc $\{Y_n\}$, nor is it necessarily the best way

to compute $E(y)$.)

4. general rv Y

Define the positive part of Y , $Y_+ = \max(Y, 0)$

and the negative part, $Y_- = \max(-Y, 0)$.

(both are nonnegative)

so $Y = Y_+ - Y_-$ and $|Y| = Y_+ + Y_-$

Then if either (or both) of $E(Y_+)$ & $E(Y_-)$ are finite, define

$$E(Y) = E(Y_+) - E(Y_-)$$

. (Again, the value can be infinite - we just don't allow the case $\infty - \infty$)

The purpose of the above is just so we can define (and talk about) expectations without having to worry about the type of rv or how the expectation is to be computed. In fact, computing expectations will depend on context and often is not necessary in specific cases.

Definition. 0.8. (Lebesgue-Stieltjes notation).

If $X \sim F$, it is conventional to express

$$E(x) = \int_{-\infty}^{\infty} xF(dx)$$

, and $E(g(x)) = \int_{-\infty}^{\infty} g(x)F(dx)$, when they can be defined.

A similar usage applies to random vectors, using a joint cdf (but not to stochastic processes) .

Other terminology.

- A random variable X is finite if $|X(\omega)| < \infty$ for all $\omega \in \Omega$.
- A random variable X is bounded if there exists a constant C such that $|X(\omega)| \leq C$ for all $\omega \in \Omega$.
- Do not confuse “finite” & “bounded”.

The terms apply in similar fashion to any real valued function $g(x)$.

An event A occurs almost surely (a.s.) or with probability 1 (w.p.1) if $P(A) = 1$.

Note that “almost sure” is in the context of a specific probability measure or model, whereas the definition of A (as some event) is not.

ex. Y is almost surely finite if $P(|Y| < \infty) = 1$. However, the event $\{Y = \infty\}$ need not be empty. Indeed, we will want to discuss rvs that are almost surely finite for some models but not for others (and to identify the distinguishing condition) .

ex. X_n converges to X almost surely ($X_n \rightarrow X$ a.s.), if $P(\lim_{n \rightarrow \infty} X_n = X) = 1$,

This is stronger than convergence in probability: $P(|X_n - X| > \epsilon) \rightarrow 0$ for all positive ϵ .

The basic limit theorems

Theorem 0.2. Monotone Convergence (MCT)

For $0 \leq X_n \uparrow X : E(X_n) \rightarrow E(X)$ (even if infinite)

Likewise, if $0 \leq g_n(x) \uparrow g(x)$ for all x , then

$$\int g_n(x) dx \rightarrow \int g(x) dx$$

, and

$$\sum_{i=1}^{\infty} g_n(i) \rightarrow \sum_{i=1}^{\infty} g(i)$$

Theorem 0.3. Dominated Convergence (DCT)

For $X_n \rightarrow X$ such that $|X_n| \leq Y$ and $E(Y) < \infty$, $E(X_n) \rightarrow E(X)$.

(again, similar statements for integrals & sums)

Theorem 0.4. Fatou's Lemma $X_n \geq 0, X_n \rightarrow X$

$$\liminf_{n \rightarrow \infty} E(X_n) \geq E(X)$$

(& similarly for sums and integrals)

(Note:

$$\liminf_{n \rightarrow \infty} a_n = \lim_{n \rightarrow \infty} (\inf_{k \geq n} a_k) \quad \& \quad \limsup_{n \rightarrow \infty} a_n = \lim_{n \rightarrow \infty} (\sup_{k \geq n} a_k)$$

.)

Theorem 0.5. Strong Law of Large Numbers (SLLN) For iid X_n ,

$$\frac{1}{n} \sum_{i=1}^n X_i \rightarrow E(X) \quad (\text{if defined})$$

almost surely, and

$$E\left(\left|\frac{1}{n} \sum_{i=1}^n X_i - E(X)\right|\right) \rightarrow 0 \quad (\text{if } E(|x|) < \infty)$$

.

Equalities & inequalities

Theorem 0.6.

- If $X \geq 0$ then $E(X) = \int_0^\infty P(X > x) dx$, even if infinite (& valid even if $P(X = \infty) > 0$).
- If $X \geq 0$ and is integer-valued then

$$E(X) = \sum_{n=0}^{\infty} P(X > n)$$

Theorem 0.7.

(i) (Markov) If $g(x) \geq 0$ & nondecreasing then

$$P(X \geq x) \leq \frac{E(g(X))}{g(x)}$$

e.g.

$$P(|X| > x) \leq \frac{E(|x|)}{x} \quad \& \quad P(|X| > x) \leq \frac{E(x^2)}{x^2}$$

.

(ii) $|E(x)| \leq E(|x|)$

(iii) (Lyapunov) $E(|X|^r) \leq (E(x^2)E(y^2))^{\frac{1}{2}}, \quad \text{if } 0 < r < s.$

(iv) (Cauchy-Schwaiz) $|E(xy)| \leq (E(x^2)(y^2))^{\frac{1}{2}}$

(v) (Jenson) If $g(x)$ is convex then $E(g(x)) \geq g(E(x))$ with equality only if $g(x) = a + bx$ for some constants a, b .

Note: “convex” means $g(ax + (1 - a)y) \leq ag(x) + (1 - a)g(y), 0 < a < 1$

Equivalently, $g(b) - g(a) = \int_a^b g'(u)du$ all a, b and $g'(u)$ is nondecreasing (the derivative need not exist at all values of u for this to hold)

Theorem 0.8. (Kolmogorov maximal inequality)

If X_n are independent, each $E(X_n^2) < \infty$ & $S_n = X_1 + \cdots + X_n$

then

$$P(\max_{k \leq n} |S_k - E(S_k)| > x) \leq \frac{\text{Var}(S_n)}{x^2}$$

The next result takes many guises, but you should set the idea from examples shown.

Theorem 0.9. (Fubini-tonelli)

Suppose either that $g(x, y) \geq 0$ for all x, y or the quantities below are finite when $g(x, y)$ is replaced with $|g(x, y)|$.

(i)

$$\iint g(x, y) dx dy = \iint g(x, y) dy dx$$

.

(ii)

$$\int \sum_{n=0}^{\infty} g(n, y) dy = \sum_{n=0}^{\infty} \int g(n, y) dy$$

.

(iii)

$$\int E(g(X, y)) dy = E\left(\int g(x, y) dy\right)$$

.

(iv)

$$\sum_{n=0}^{\infty} E(g(X, n)) = E\left(\sum_{n=0}^{\infty} g(X, n)\right)$$

.

(v)

$$\sum_{n=0}^{\infty} \sum_{k=0}^{\infty} g(n, k) = \sum_{k=0}^{\infty} \sum_{n=0}^{\infty} g(n, k)$$

.

(vi)

$$\iint g(x, y) F(dx) G(dy) = \iint g(x, y) G(dy) F(dx)$$

, where the latter are lebesgue-stieltjes integrals. (When $g(x, y) \geq 0$ all x, y the double integrals/sums/etc above can equal ∞ , on both sides.)

We rely heavily on conditional probabilities and conditional expectation. Like expectation, however, we want to discuss & use them without being restricted to joint discrete distributions or joint continuous distributions. We provide here, without elaboration, a suitably general definition.

Definition. 0.9. Let Y be a r.v. such that $E(Y)$ exists, and X be a random element (r.v., r. vector, stochastic process). The conditional expectation of Y given X , denoted $E(y|X)$ is

a function of the random variable X , say $h(x)$, (Note: this also means $E(Y|X)$ is itself a r.v.) such that

$$E(g(x)Y) = E(g(x)E(Y|X)) \quad (*)$$

for all (measurable) bounded real-valued functions $g(x)$. (“measurable” just means $\{x : g(x) \leq y\}$ is a Borel set, which is a technical restriction not usually of much concern to us.)

Conditional Probability $P(y \in A|X) = E(\mathbb{1}_A(y)|X)$.

The definition does not tell us how to compute $E(y|X)$; it merely says that we can define it. Note that $E(y|X)$ is a random variable (but it must be a function of X) . Often we may have a way to define or compute $h(x) = E(y|X = x)$ suitably and then set $E(y|X) = h(X)$.

Properties of conditional expectation

Theorem 0.10. *Given X , and r.v.'s Y, Y_1, Y_2, \dots (w/ means)*

- (i) $y_1 \leq y_2 \implies E(y_1|X) \leq E(y_2|X)$.
- (ii) $E(aY_1 + bY_2|X) = aE(Y_1|X) + bE(y_2|X)$.
- (iii) *MCT, DCT & Fatou's lemma all hold, cond. on X .*
- (iv) *The formula (*) in **Theorem.0.6.** holds whenever the expectations exist (not just for bounded $g(x)$).*
- (v) *(Pull-Out) $E(g(x)y|X) = g(X)E(y|x)$, if it exists.*
- (vi) *(In)equalities for expectation (Jensen, etc.) all hold, cond. on X .*
- (vii) *If Y is independent of X then $E(Y|X) = E(Y)$.*
- (viii) *$y = h(x)$ then $E(y|x) = y$*

(ix) (Tower) $E(y|x) = E(E(y|x, z)|x) = E(E(y|x)|x, z)$.

Basically, we can work with conditional expectations in an intuitive fashion based on familiarity with definitions for the case (x, y) is jointly discrete or jointly absolutely cont.

But now we can do so without those restrictions, including the case X is a stochastic process (or a portion of one) consisting of infinitely many rv's.

Take care to use the tower property, (ix) above, correctly. It is not usually true that

$$E(E(y|x)|z) = E(E(y|z)|x)$$

One other caution: $E(y|x)$ is not originly defined since $P(bx = 0) = 1$ implies $E(y|x) + b(x)$ also satisfies the definition. We don't usually have a problem w/ this if there are only countably many r.v.s.

1 Introduction

- definitions
- examples (simple random walk, simple branching proc.)
- probability generating functions
- some analytic results

Definition. 1.1. A stochastic process is a collection of random variables, $\{X_t\}_{t \in \mathbb{T}}$, defined on the same probability space. The state space \mathbb{X} is a set of possible values for X_t and the index set \mathbb{T} is the set on which X_t evolves.

(a single instance of values $\{X_t\}$ is called a sample path.)

For us, \mathbb{X} will usually be the integers \mathbb{Z} or the real line \mathbb{R} (or some other subset of \mathbb{R} but it can be \mathbb{R}^m or something more exotic.

\mathbb{T} will usually be the nonnegative integers \mathbb{Z}_+ or the nonnegative real numbers $\mathbb{R}_+ = [0, \infty)$, but it also can be more general.

Example. 1.1. $X_t = \#$ of people infected on day t of an epidemic

$X_{s,t}$ = temperature at point s on the globe and time t

$X(A) = \#$ of fireant colonies in a region A (here \mathbb{T} is a collection of sets, including all open & closed sets)

Stochastic processes have many applications:

e.g. signal process, random networks, queues, climate, earthquakes, insect populations, stock market, economics, epidemics, and so on.

Note: I generally use upper case for random variables and lower case for actual values.

Just as we do for a random variable & a random vector, we want to discuss the distribution of a stochastic process. But, with infinitely many rv's, we cannot simply express it as an ordinary function (such as a cdf). Instead we have