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Stochastic Processes

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Abstract

Dear reader, thank you for your interest in my notes on stochastic processes. I have created these notes with care and attention to detail, attending every lecture and spending many hours outside of class to ensure their accuracy and clarity. While I do not wish to sell these notes, I would greatly appreciate any donation you may be willing to offer in exchange for their use. As a full-time student, attending these classes is a significant investment of time and resources for me, and your support would be greatly appreciated. Thank you for your consideration. If you have found these notes helpful, and would like to support my efforts in creating them, you can send a donation via PayPal to **trashajalberto@gmail.com**. Your contribution will go a long way in helping me continue to create and share high-quality educational materials. Thank you again for your support.

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Chapter 1

Introduction

1.1 Definition

The first concept that we need to present in this course is the definition of stochastic process:

Definition 1.1.1 (Stochastic process). A stochastic process is any collection of random variables

So, quite simple, no?

Let's now briefly recall what is a random variables since it will be useful to understanding what we are talking about:

A random variable is a measurable map defined as follows:

$$X : \Omega \longrightarrow S$$

where we just recall that the triplet (Ω, \mathcal{A}, P) is the well known probability space and (S, \mathcal{B}) is the measurable space.

In particular, S is a set and \mathcal{B} is a σ -field of subset of S .

It can be possible that you are not familiar with this concepts since some notion of the probability theory are needed, so if you have already problems with understanding what are we talking about pause the reading and try to comprehend the concept so far since they will be very useful for the rest of the course.

Disclaimer: in the repository
https://github.com/alberello00/stochastic_processes
you will find some notes of the course "Advanced probability"
which can be useful
to understanding some concepts.

So far we just briefly recalled the two main ingredients for defining a random variable: another ingredient that has to be mentioned when defining what is a stochastic process is the set T .

The set T can be any set, but the most natural and intuitive set that we can think of is the *Time* set. But, later on we will define better what T actually is.

To give a formal definition we can repeat that a stochastic process X is any collection of random variables namely:

$X = X_t : t \in T$
where $X_t : \Omega \longrightarrow S$ is a random variable $\forall t \in T$

Another parentheses should be opened here: in fact, if one remember what it actually is a random variable can point out that a random variable has to be a measurable map, i.e. the random variables that are in our stochastic processes has to be measurable too:

$$X_t^{-1}(B) \in \mathcal{A} \quad \forall B \in \mathcal{B}$$

where

$$X_t^{-1}(B) = \{\omega \in \Omega : X_t(\omega) \in B\}$$

1.1.1 Remarks on S and T

- S is said to be the **state space** of the process: of course, the most important case is when the state space is equal to the real line, namely

$$S = \mathbb{R} \text{ and } \mathcal{B} \text{ is the Borel } \sigma\text{-field on } \mathbb{R}$$

- T is an arbitrary set, which is usually called the **indexing set** of the **parameter space** of the process: a natural interpretation of this set can be thought as the *time* set.
- If T is a finite or countable set, X is said to be a discrete time process: for example, a sequence of random variables.
- If T is an interval of the real line \mathcal{R} , X is said to be a **continuous time process**

1.2 A process is a bivariate function

An important way to think of what is a stochastic process is that it is nothing else than a function of two variables:

$$X : \Omega \times T \longrightarrow S$$

So, we can write

$$X(\omega, t) \text{ or, equivalently, } X_t(\omega)$$

So, since a stochastic process can be thought as a function of two variables let's see what happens if we fix one of the two variables and let one be freely to change.

- Let fix $t \in T$:
Then, X_t is still a random variable, namely a measurable function on Ω .
Hence, we have

$$X_t(\omega) \quad \forall \omega \in \Omega$$

- Let fix $\omega \in \Omega$:
Then, we obtain a function of t , namely:

$$t \longrightarrow X_t(\omega)$$

This function is usually called the **path** or **trajectory** of the process.
For instance, if we choose the state process as $S = R$ and $T = [0, \infty)$: the path is a function that maps from $[0, \infty)$ into \mathcal{R}

1.2.1 Another way to see what is a stochastic process

A process can be always regarded as a random function: it is sufficient to think of X as the map $\omega \longrightarrow$ to the path associated to ω , namely:

- $\omega \longrightarrow X(\omega, *)$

1.3 Equality of the processes

One can wonder when two processes can be regarded as "equal", "similar" or, in a more general sense when two processes can be compared and see if they are in some sense indistinguishable.

In general, we can say that there are three level of "equality" of two processes:

1. $X \sim Y \Rightarrow (X_{t_1}, X_{t_2}, \dots, X_{t_n}) \sim (Y_{t_1}, Y_{t_2}, \dots, Y_{t_n})$
2. X is equal to $Y \iff P(X_t \neq Y_t) = 0 \forall t \in T$
3. X is indistinguishable from $Y \iff X_t(\omega) = Y_t(\omega) \forall t \in T$ and $\forall \omega \in A$ where $A \in \mathcal{A}$ is such that $P(A) = 1$

The natural question that can pop out in our mind is if this three type of equality are in someway related:

We can state the following implications:

$$3 \Rightarrow 2 \Rightarrow 1$$

$$1 \not\Rightarrow 2 \not\Rightarrow 3$$

1.3.1 Equivalent doesn't imply indistinguishable

Usually to say that A doesn't imply B , it is sufficient to prove that A is true while B is false.

So, let's find an example where we have two processes that satisfies the second type of equality but not the third one.

Example 1.3.1. Let V be any random variable such that $V \geq 0$ (positive-definite) and $P(V = v) = 0 \ \forall v \geq 0$. For instance, we can take the absolute value of a standard Gaussian, so $V = |Z|$ where $Z \sim N(0, 1)$. Now, let's define two random process:

- $X(t, \omega) = 0 \ \forall t \in [0, +\infty) , \ \forall \omega \in \Omega$

-

$$Y(t, \omega) = \begin{cases} 1, & \text{if } t = V(\omega). \\ 0, & \text{otherwise.} \end{cases} \quad (1.1)$$

Then, it's clear that X and Y are not indistinguishable since for $t = V(\omega)$ we get

$$Y(\omega, t) = 1 \text{ and } X(\omega, t) = 0, \text{ so } X(\omega, t) \neq Y(\omega, t)$$

However, with a close analysis, we can say that

$P(X_t \neq Y_t) = P(Y_t \neq 0)$ since $X_t = 0$ as assumption, and also $Y_t \neq 0$ when $t = V(\omega)$, therefore the $P(Y_t \neq 0) = P(V(\omega) = t)$ that we know it's equal to 0 since the probability of a point is 0 by the assumption that $P(V = v) = 0 \ \forall v \geq 0$

So, in conclusion we showed that X_t and Y_t are equivalent processes but they are not indistinguishable

1.4 Stopping times

Let take a set $T = 0, 1, 2, \dots$ and the usual sample space $(\Omega, \mathcal{A}, \mathcal{P})$; then we define a *filtration* as an increasing sequence of sub- σ fields of \mathcal{A} , namely

$$\mathcal{F}_1 \subset \mathcal{F}_\infty \subset \mathcal{F}_\epsilon \subset \dots \subset \mathcal{A}.$$

Given this ingredients we can give a formal definition of the stopping time

Definition 1.4.1. A stopping time is a map

$$T : \Omega \Rightarrow \{+\infty, 0, 1, 2, \dots\}$$

such that $\{T = n\} \in \mathcal{F}_n \ \forall n \geq 0$.

In general, a σ -field $\mathcal{G} \subset \mathcal{A}$ may be used to describe the so called state of information. By now, it's not so clear what are we referring to when we talk about the "state of information" of \mathcal{G} , so let's make some examples:

Example 1.4.1. • $\mathcal{G} = \{\emptyset, \Omega\}$ contains null information since, by definition, we always know that the \emptyset is always *False* and the Ω is always *True*.

- $\mathcal{G} = \{\emptyset, \Omega, A, A^c\}$ in this set the only information we have is whether or not A is *True*.

Now, I think that another example should clarify better what we mean by "state of the information":

Example 1.4.2. Let's take $\Omega = \{1, 2, 3, 4, 5, 6\}$ and define a σ -algebra $\mathcal{F} = \{\emptyset, \Omega, \{1, 2, 3, 4\}, \{5, 6\}\}$: now, we can see that, for example, the subset $\{1, 2\}$ is not contained, so our probabilistic model can't make conclusion in this last set, so the "information" is contained only in the set \mathcal{F} . Therefore, the only "probabilistic" conclusion we can make are only in the σ -field we constructed before; there will no exists random variable that maps the subset $\{1, 2\}$ on the interval $[0, 1]$.

T should be regarded as the first time when something we are interested actually happens: so, when we say that $T = n$ we are implicitly saying that at the time n something happened for the first time.

Remember when we defined the set where T takes values? Yes, at first can seems strange to include the $+\infty$ but now it should be more clear that saying $T = +\infty$ means that something does not happen.

Example 1.4.3. (X_n) is a sequence of real random variable and $A \in \mathcal{B}(\mathcal{R})$; $T = \inf\{n : X_n \in A\}$, so the first time n such that $X_n \in A$.

Example 1.4.4 (Casinò). Suppose you and your friends are playing at a roulette table in a casino, and you want to come up with a strategy to maximize your winnings. You decide that you will keep playing until a certain point in time, at which you will stop and walk away with your winnings. The catch is that you don't know ahead of time when this stopping time will occur - it could happen after 10 minutes of playing, or it could happen after several hours.

However, you have a secret weapon: a fortune teller who has told you that the stopping time will occur at the 10th round of the game. In other words, you will keep playing for 10 rounds, and then stop.

In this example, the stopping time is the 10th round of the game. This is a stopping time because it is determined by information that is available at or before the 10th round of the game. Specifically, the information that you have received from the fortune teller is available to you before the 10th round, and therefore the event $T = 10$ (where T is the stopping time) belongs to the sigma-algebra \mathcal{F}_{10} .

1.5 Finite dimensional distributions

Definition 1.5.1. Let X be a process indexed by $T \forall n \geq 1 \forall t_1, \dots, t_n \in T$ we have a n -dimensional random variable $(X_{t_1}, X_{t_2}, \dots, X_{t_n})$: the distributions of $(X_{t_1}, X_{t_2}, \dots, X_{t_n})$ for all $n \geq 1$ and $\forall t_1, \dots, t_n \in T$ are called finite dimensional distribution.

Example 1.5.1. Suppose that $T = \{1, 2\}$ and $X = \{X_1, X_2\}$: $X_1 \sim \text{Binomial}$ and $X_2 \sim \text{Poisson}$. Then, for obvious reasons, $(X_1, X_2) \sim N$ it's not possible and such process fail to exist.

Usually, in application we choose the finite dimensional distributions and we look for a process having such a finite dimensional distribution: the problem here arise since this process may not exist.

However there are some theorem usually called *consistency* theorems which provides conditions on the finite dimensional distributions under which the processes (with such finite dimensional distribution) exists.

Chapter 2

Conditional expectation

2.1 Conditional expectations

In order to give a proper definition of conditional expectations, three ingredients are necessary:

1. A sample space, $(\Omega, \mathcal{A}, \mathcal{P})$
2. A sub- σ -field of \mathcal{A} : $\mathcal{G} \subset \mathcal{A}$
3. A real random variable X such that $E|X| < +\infty$.

Now, let's define a conditional expectation:

Definition 2.1.1. A conditional expectation of X given \mathcal{G} is any real random variable $V : \Omega \Rightarrow \mathcal{R}$ such that:

1. $E(|V|) < +\infty$
2. V is \mathcal{G} -measurable
3. $E(\mathbb{1}_A \cdot X) = E(\mathbb{1}_A \cdot V) \forall A \in \mathcal{G}$

Where, on the third point we have,

$$X \cdot \mathbb{1}_A = \begin{cases} X, & \text{on } A \\ 0, & \text{on } A^c. \end{cases} \quad (2.1)$$

In particular, since $\Omega \in \mathcal{G}$, we obtain:

$$E(X) = E(X \cdot \mathbb{1}) = E(V \cdot \mathbb{1}) = E(V)$$

We need to make a second remark on the second condition stated in the definition: a random variable V is said to be \mathcal{G} -measurable if

$$V^{-1}(B) \in \mathcal{G} \forall B \in \mathcal{BR}$$

In other terms, V is a random variable not only in the "big" probability space $(\Omega, \mathcal{G}, \mathcal{P})$ but also in the "small" probability space $(\Omega, \mathcal{G}, \mathcal{P})$.

For instance, take the set $\mathcal{G} = \{\emptyset, \Omega\}$ (null information), then if X is \mathcal{G} -measurable, it must be that

$$X^{-1}(B) = \Omega \text{ or } X^{-1}(B) = \emptyset \quad \forall B \in \mathcal{B}(\mathcal{R})$$

Hence, with the set taken the only \mathcal{G} -measurable random variables are the constants. At the opposite extreme, if $\mathcal{G} = \mathcal{A}$, then every random variable is \mathcal{G} -measurable.

However, the interpretation of \mathcal{G} -measurability is under the information \mathcal{G} , so a \mathcal{G} -measurable random variable reduces to a constant: suppose we have a random variable X that is \mathcal{G} -measurable for some sub- σ -field \mathcal{G} . Intuitively, this means that X behaves like a constant, because its value can be determined from the information in \mathcal{G} : let's see with an example why this is true.

Example 2.1.1. Fix $A \in \mathcal{A}$ and define

- $\mathcal{G} = \{\emptyset, \Omega, A, A^c\}$
- $X = 3\mathbb{1}_A - 2\mathbb{1}_{A^c}$

Then, X is \mathcal{G} -measurable:

$$X^{-1}(B) = \begin{cases} \Omega, & \text{if } -2, 3 \in B \\ \emptyset, & \text{if } -2, 3 \notin B \\ A, & \text{if } 3 \in B \text{ and } -2 \notin B \\ A^c, & \text{if } 3 \notin B \text{ and } -2 \in B \end{cases} \quad (2.2)$$

But, under the information of \mathcal{G} , we know whether A is *True* or *False*, thus X becomes a constant.

Theorem 2.1.1 (Theorem). A conditional expectation V always exists, and it is almost surely unique, namely, if V_1 and V_2 are both conditional expectations, then $P(V_1 \neq V_2) = 0$ or, equivalently, $P(V_1 = V_2) = 1$

From now on, to denote a conditional expectation of X given \mathcal{G} , we adopt the notation $E(X|\mathcal{G})$, namely $V = E(X|\mathcal{G})$.

The interpretation is that the expected value is our prediction of X under the information \mathcal{G} : in the special case where $V = E(X)$, the natural interpretation is that $E(X)$ is our prediction of X without any specific information.

Now, should be more clear why $E(X|\mathcal{G})$ should be \mathcal{G} -measurable: in fact, our prediction of X under the information \mathcal{G} should be something which only depends on the information \mathcal{G} .

2.1.1 Best predictor?

In general, to make a prediction, we are not only interested in finding a good predictor, but it would be better to find the **best** predictor: don't worry, we already find it! Yes, $E(X|\mathcal{G})$ is the best prediction we can make in the following sense:

- Suppose $E(X^2) < +\infty$: then,

$$E\{(X - (E(X|\mathcal{G})))^2\} = \min E\{(X - Z)^2\}$$

We can think of Z as a predictor of X under the information \mathcal{G}

2.1.2 Properties

There are 7 useful properties of the conditional expectation:

1. Linearity: $E(aX + bY|\mathcal{G}) = aE(X|\mathcal{G}) + bE(Y|\mathcal{G})$
2. Positivity: if $X \geq 0 \rightarrow E(X|\mathcal{G}) \geq 0$
3. Constant: $E(\text{constant}|\mathcal{G}) = \text{constant}$
4. Measurability: if Z is \mathcal{G} -measurable, then $E(X \cdot Z|\mathcal{G}) = Z \cdot E(X|\mathcal{G})$
It follows that, if X is \mathcal{G} -measurable, then

$$E(X|\mathcal{G}) = E(X \cdot 1|\mathcal{G}) = X \cdot E(1|\mathcal{G}) = X \cdot 1 = X$$

5. Independence: X is said to be independent from \mathcal{G} , by definition, if

$$P(A, X \in B) = P(A) \cdot P(X \in B) \forall A \in \mathcal{G} \forall B \in \mathcal{B}(\mathcal{R})$$

If X is independent from \mathcal{G} , then $E(X|\mathcal{G}) = E(X)$

6. Chain rule: if $\mathcal{G}_\infty \subset \mathcal{G}_\epsilon \subset \mathcal{A}$, then $E(X|\mathcal{G}_\infty) = E(E(X|\mathcal{G}_\epsilon)|\mathcal{G}_\infty)$.
For instance if we know that $E(X|\mathcal{G}_\epsilon) = 0$, then
 $E(X|\mathcal{G}_\infty) = E(E(X|\mathcal{G}_\epsilon)|\mathcal{G}_\infty) = E(0|\mathcal{G}_\infty) = 0$
7. Conditional probability: if $A \in \mathcal{A}$, we can define

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

An example is the following:

Example 2.1.2. If A and $B \in \mathcal{A}$ and $P(B) > 0$,

$$P(A|B) = \frac{P(A \& B)}{P(B)}$$

Despite this condition is well-known, a nice connection between this elementary definition and the general notion of conditional probability can be done.

In fact, the elementary definition is actually a special case of the general definition of probability conditional to a σ -field \mathcal{G} : let us take the following σ -field $\mathcal{G} = \{\emptyset, \Omega, B, B^c\}$

It can be shown that if $f : \Omega \rightarrow \mathcal{R}$ is \mathcal{G} -measurable, then:

for some constants α and β .

Hence,

$$P(A|\mathcal{G}) = \alpha \cdot \mathbb{1}_B + \beta \cdot \mathbb{1}_{B^c},$$

but,

$$E\{P(A|\mathcal{G}) \cdot \mathbb{1}_D\} = E\{\mathbb{1}_A \cdot \mathbb{1}_D\} \quad \forall D \in \mathcal{G}.$$

So, taking $D = B$, we get

$$E\{\mathbb{1}_A \cdot \mathbb{1}_B\} = E\{\mathbb{1}_{A \cap B}\} = P(A \cap B)$$

Hence,

$$E\{P(A|\mathcal{G}) \cdot \mathbb{1}_B\} = E\alpha \cdot \mathbb{1}_B = \alpha \cdot E\{\mathbb{1}_B\} = \alpha \cdot P(B)$$

It follows that,

$$\alpha = \frac{P(A \cap B)}{P(B)} = P(A|B)$$

Similarly, one obtain

$$\beta = P(A|B^c), \text{ provided that } P(B^c) > 0$$

In short,

$$P(A|\mathcal{G}) = P(A|B) \cdot \mathbb{1}_B + P(A|B^c) \cdot \mathbb{1}_{B^c}$$

Usually, we write $E(X|Y)$ to denote $E(X|\sigma(y))$, namely $E(X|Y) = E(X|\sigma(y))$: $\sigma(y)$ is the least field which makes Y measurable.

In short, when we have a random variable Y , $\sigma(Y)$ is the smallest σ -field making Y measurable.

Let's make another (more simple) example:

Example 2.1.3. Let Y, X be *iid* with $E|X| < +\infty$. What about $E(X|X+Y)$? It can be shown that

$$E(X|X+Y) = \frac{X+Y}{2}$$

because

- $\frac{X+Y}{2}$ has the mean
- $\frac{X+Y}{2}$ is measurable with respect to $\sigma(X+Y)$
- $E(X \cdot \mathbb{1}_A) = E(\frac{X+Y}{2} \cdot \mathbb{1}_A)$

Chapter 3

Martingales

3.1 Discrete time: Martingales

Definition 3.1.1. Given a filtration $\mathfrak{F}_0 \subset \mathfrak{F}_1 \subset \dots \subset \mathcal{A}$ and the sequence of x_0, x_1, x_2, \dots real random variables: we say that (X_n) is a martingale with respect to \mathfrak{F}_n if three conditions are satisfied:

1. $E|X_n| < +\infty$
2. X_n is \mathfrak{F}_n -measurable
3. $E(X_{n+1}|\mathfrak{F}_n) = X_n$

Some remarks:

- If $E(X_{n+1}|\mathfrak{F}_n) \geq X_n$, then (X_n) is called sub-martingale.
- If $E(X_{n+1}|\mathfrak{F}_n) \leq X_n$ then (X_n) is called a super-martingale.
Obviously a martingale is both a sub-martingale and a super-martingale.
- Let $\mathfrak{F}_n^* = \sigma(X_0, X_1, \dots, X_n)$, this is the least sigma field that makes (X_0, X_1, \dots, X_n) measurable: therefore, (X_n) is still a martingale with respect to (\mathfrak{F}_n^*) .
In fact, since $\mathfrak{F}_n^* \subset \mathfrak{F}_n$, one obtains

$$E(X_{n+1}|\mathfrak{F}_n^*) = E(E(X_{n+1}|\mathfrak{F}_n)|\mathfrak{F}_n^*) = E(X_n|\mathfrak{F}_n^*) = X_n$$

In the last expression we used respectively the chain rule, the definition of conditional expectation and the notion of measurability.

A simple theorem about martingales is described above:

Theorem 3.1.1. (X_n) is a martingale if and only if $\exists C \in \mathcal{R}$ and a Z_n such that $X_n = C + \sum_{i=0}^n Z_i$.

For instance, $E|Z_i| < +\infty$, Z_i is \mathfrak{F}_i measurable and $E(Z_{i+1}|\mathfrak{F}_i) = 0$

A intuitive interpretation could be the following: imagine you are at the casinò, and Z_i represent the outcome of the i -th trial = the win at day i . X_n is the amount of money in your pocket after you play $(n + 1)$ times.
If the game is **fair**, $E(Z_{i+1}|\mathfrak{F}_i) = 0$

Let's make an example:

Example 3.1.1. If Z_i is any independent sequence of real random variables such that

$$E(Z_i) = 0 \forall i, \text{ then}$$

$X_n = c + \sum_0^n Z_i$ is a proper martingale.

In fact, letting $\mathfrak{F}_n = \sigma(X_0, X_1, \dots, X_n) = \sigma(Z_0, \dots, Z_n)$, one obtain:

$$E(Z_{i+1}|\mathfrak{F}_i) = E(Z_{i+1}|Z_0, \dots, Z_i) = \text{by independence} = E(Z_{i+1})$$

and by assumption, $E(Z_{i+1}) = 0$

3.1.1 Random walk

A random walk is a process where an object or variable takes successive random steps in some space, such as a one-dimensional number line or a two-dimensional plane. The steps can be in any direction and can be positive or negative.

A simple example of a random walk is a gambler who starts with some amount of money and repeatedly flips a fair coin. If the coin lands heads, the gambler wins some amount of money, and if the coin lands tails, the gambler loses some amount of money. The gambler's wealth then follows a random walk as they make successive bets.

Definition 3.1.2. A sequence (X_n) is a random walk if:

$$X_n = X_0 + \sum_{i=1}^n Z_i, \tag{3.1}$$

where (Z_i) is a sequence of independent and identically distributed (i.i.d.) random variables, and X_0 is independent of (Z_i) .

In this notation, the sum goes from $i = 1$ to n to indicate that the first term in the sum is Z_1 , and X_0 is assumed to be a fixed constant or a random variable that is independent of (Z_i) .

A very popular example of a random walk is the following:

Example 3.1.2. • $P(X_0 \in \mathcal{Z}) = 1$

• $P(Z_i = 1) = P(Z_i = -1) = \frac{1}{2}$
when $\mathcal{Z} = \{\dots, -2, -1, 0, 1, 2, \dots\}$ is the set of integers.

3.1.2 Connection between random walk and martingale

A random walk is a martingale if and only if the following criteria are met:

- $E(|X_0|) < +\infty$
- $E(|Z_1|) < +\infty$
- $E(|Z_1|) = 0$

If X_n is a martingale and τ a stopping time, then $X_{(\tau \wedge n)}$ is still a martingale.

[Notion alert: $a \wedge b = \min(a, b)$]

In particular $X_{(\tau \wedge n)}$ is called stopped sequence:

$$X_{(\tau \wedge n)} = \begin{cases} X_n & \text{if } n \leq \tau \\ X_\tau & \text{if } n > \tau \end{cases} \quad (3.2)$$

Then the natural interpretation would be that when the event of interest occurs, the sequence became

$$X_1, X_2, \dots, X_\tau, X_\tau, \dots$$

In order to say that this is still a martingale, we have to show that

$$E(X_{\tau \wedge (n+i)} | \mathfrak{F}_n) = X_{\tau \wedge n}$$

Proof :

$$E(X_{\tau \wedge (n+i)} | \mathfrak{F}_n) = E(X_\tau \mathbb{1}_{(\tau \leq n)} + X_{n+1} \mathbb{1}_{\tau > n} | \mathfrak{F}_n)$$

$$= E(X_\tau \mathbb{1}_{\tau \leq n} | \mathfrak{F}_n) + E(X_{n+1} \mathbb{1}_{\tau > n} | \mathfrak{F}_n)$$

The first term and the second term (since this one is \mathfrak{F}_n -measurable) can be rewritten in the following way:

$$= E(\sum_{i=0}^n X_i \cdot \mathbb{1}_{\tau=i}) + \mathbb{1}_{\tau > n} \cdot E(X_{n+1} | \mathfrak{F}_n)$$

Moreover, using the definition of martingale,

$$\begin{aligned} &= \sum_{i=0}^n E(X_i \cdot \mathbb{1}_{\tau=i} | \mathfrak{F}_n) + \mathbb{1}_{\tau > n} \cdot X_n \\ &= \sum_{i=0}^n X_i \cdot \mathbb{1}_{(\tau=i)} + \mathbb{1}_{\tau > n} \cdot X_n = X_{\tau \wedge n} \end{aligned}$$

Another example follows:

Example 3.1.3. Let's take $X_n = \sum_{i=0}^n Z_i$ and $X_0 = 0$, where (Z_i) is *iid* and $E(Z_1) = 0$, $E(Z_1^2) = 1$.

Incidentally, in this case, X_n is both a martingale and a random walk.

Moreover $Y_n = X_n^2 - n$ is still a martingale. Let's prove this last point:

letting $\mathfrak{F}_n = \sigma(X_0, \dots, X_n)$ we obtain

$$E(Y_{n+1}|\mathfrak{F}_n) = E(X_{n+1}^2 - (n+1)|\mathfrak{F}_n)$$

And since $(n+1)$ is a constant

$$= E((X_n + Z_{n+1})^2|\mathfrak{F}_n) - (n+1)$$

by developing the square,

$$= E(X_n^2 + Z_{n+1}^2 + 2X_n Z_{n+1}|\mathfrak{F}_n) - (n+1)$$

applying the properties of martingales,

$$= X_n^2 + 2X_n \cdot E(Z_{n+1}|\mathfrak{F}_n) + E(Z_{n+1}^2|\mathfrak{F}_n) - (n+1)$$

$$= X_n^2 - (n+1) + 2X_n \cdot E(Z_{n+1}) + E(Z_{n+1}^2)$$

and by applying the assumptions

$$= X_n^2 - n = Y_n$$

Incidentally, it's possible to note that

$$X_n^2 = Y_n + n$$

is not a martingale.

In fact, if (M_n) is a martingale, then

$$E(M_{n+1}) = E(E(M_{n+1}|\mathfrak{F}_n)) = E(M_n)$$

Hence, for every M_n , we get

$$E(M_0) = E(M_1) = \dots$$

but, if we take the example before we can show that,

$$E(Y_n + n) = E(Y_0) + n = E(Y_{n+1} + (n+1)) \neq E(Y_n + n)$$

hence

$$E(Y_0) + (n+1) \neq E(Y_0) + n$$

For any real random variable X such that $E(|X|) < +\infty$ and any σ -field $\mathfrak{G} \subset A$, we have that

$$E(X) = E(E(X|\mathfrak{G}))$$

3.2 Gambler's ruin problem

In order to model this situation, let $X_0 = 0$ and $X_n = \sum_1^n Z_i$ where Z_i is *iid* and $P(Z_1 = 1) = P(Z_1 = -1) = \frac{1}{2}$.

Moreover, we suppose that player 1 has $a > 0$ euros and player 2 has $b > 0$ euros. There are three possibilities

1. The game does not end, namely

$$-a < X_n < b \quad \forall n$$

It can be shown that $P(-a < X_n < b \forall n) = 0$. (For now we can take this fact as given)

2. The winner is player 1, namely

$$X_n = b \text{ happens before } X_n = -a$$

3. The winner is player 2, namely

$$X_n = -a \text{ happens before } X_n = b$$

Let $\tau = \text{time when the game ends} = \inf n : X_n = -a \text{ or } X_n = b$. Now, let's evaluate

$$E(X_\tau) = -a \cdot P(X_\tau = -a) + b \cdot P(X_\tau = b)$$

by denoting the last term with α , we get

$$= -a \cdot (1 - \alpha) + b \cdot \alpha$$

So, it's easy to show that

$$E(X_\tau) + a = \alpha(a + b)$$

and

$$\alpha = \frac{E(X_\tau) + a}{a + b}$$

where α is the probability that player 1 is the winner. In the end, we note that X_n is a martingale since $E(Z_1) = 0$, so that $X_{\tau \wedge n}$ is still a martingale.

Therefore, since $P(\tau < \infty) = 1$ and $n \rightarrow +\infty$ at a certain point

$$\tau = \min \tau, n$$

and so we get that

$$E(X_\tau) = E(\lim_{n \rightarrow \infty} X_{\tau \wedge n})$$

and by some mathematical fact, which we take as given, we have

$$= \lim_{n \rightarrow \infty} E(X_{\tau \wedge n})$$

and since $(X_{\tau \wedge n})$ is a martingale and $E(X_{\tau \wedge n}) = E(X_{\tau \wedge 0}) \forall n$ we conclude

$$= E(X_{\tau \wedge 0}) = E(X_0) = E(0) = 0$$

So, in conclusion $\alpha = \frac{a}{a+b}$.

3.3 Doob-Mayer decomposition

Let $f : \mathbb{R} \rightarrow \mathbb{R}$ be a function such that

$$E(|f(X_n)|) < +\infty \quad \forall n$$

Then $f(X_n)$ is a sub-martingale provided

- X_n is a martingale and f is convex or,
- X_n is a sub-martingale and f is convex and increasing

Here, what we are trying to say is that, take X_n and you are interested to know if $f(X_n)$ is a martingale or a sub-martingale.

For instance, if X_n is a martingale, then, for example, X_n^2 or $|X_n|$ are sub-martingales.

Definition 3.3.1 (Doob-Meyer decomposition). (X_n) is a sub-martingale if and only if it's possible to write the sub-martingale as follows:

- $X_n = M_n + A_n$

where M_n is a martingale and A_n is an increasing sequence, namely

$$0 = A_0 \leq A_1 \leq A_2 \leq \dots$$

and A_n is \mathfrak{F}_{n-1} -measurable ("predictability").

Also, $E(A_n) < +\infty$.

Furthermore, the sequences A_n and M_n are almost surely unique.

Example 3.3.1. Suppose $X_n = \sum_1^n Z_i$, $X_0 = 0$ where (Z_i) is iid with $E(Z_i) = 0$ and $E(Z_i^2) = 1$. X_n is a martingale but X_n^2 is not a martingale.

However X_n^2 is a sub-martingale. Moreover, since $Y_n = X_n^2 - n$ is martingale, we can say that $M_n = Y_n = X_n^2 - n$ and $A_n = n$

To prove this theorem we have to show both the implication (from left to right and to right to left):

- Let first explain the "easy implication":

$$\begin{aligned} E(X_{n+1} | \mathfrak{F}_n) &= E(M_{n+1} + A_{n+1} | \mathfrak{F}_n) \\ &= E(M_{n+1} | \mathfrak{F}_n) + E(A_{n+1} | \mathfrak{F}_n) \end{aligned}$$

since the first term is a martingale and the second is \mathfrak{F}_n -measurable

$$= M_n + A_{n+1}$$

and since it is an increasing sequence

$$\geq M_n + A_n = X_n$$

- Now we need to show the other way round, the "harder implication":

Let $A_0 = 0$ and $A_n = \sum_{j=0}^{n-1} \{E(X_{j+1}|\mathfrak{F}_j) - X_j\} \forall n > 0$.

Now that we defined A_n , we have to show that A_n is increasing and predictable:

A_n is predictable because it involves only X_0, X_1, \dots, X_{n-1} and $E(X_1|\mathfrak{F}_0), \dots, E(X_n|\mathfrak{F}_{n-1})$ and all such random variables are measurable with respect to \mathfrak{F}_{n-1} . Hence, A_n is \mathfrak{F}_{n-1} -measurable.

Moreover, A_n is increasing as

$$\begin{aligned} A_{n+1} &= \sum_{j=0}^n \{E(X_{j+1}|\mathfrak{F}_j) - X_j\} \\ &= A_n + \{E(X_{n+1}|\mathfrak{F}_n) - X_n\} \geq A_n \end{aligned}$$

Remind me to explain better this steps please.

Next, let's define $M_n = X_n - A_n$:

it remains only to see that M_n is a martingale.

In fact,

$$\begin{aligned} E(M_{n+1}|\mathfrak{F}_n) &= E(X_{n+1} - A_{n+1}|\mathfrak{F}_n) \\ &= E(X_{n+1}|\mathfrak{F}_n) - E(A_{n+1}|\mathfrak{F}_n) \\ &= E(X_{n+1}|\mathfrak{F}_n) - A_{n+1} \\ &= E(X_{n+1}|\mathfrak{F}_n) - \{A_n + E(X_{n+1}|\mathfrak{F}_n) - X_n\} \\ &= X_n - A_n \end{aligned}$$

Now, let's prove that A_n and M_n are actually almost surely unique:
to prove this fact suppose

$$M_n + A_n = X_n = M'_n + A'_n$$

Then, $M_n - M'_n$ is a martingale and $A'_n - A_n$ is \mathfrak{F}_{n-1} -measurable: thus,

$$\begin{aligned} M_{n-1} - M'_{n-1} &= E(M_n - M'_n|\mathfrak{F}_{n-1}) = E(A'_n - A_n|\mathfrak{F}_{n-1}) \\ &= A'_n - A_n \end{aligned}$$

Hence,

$$A'_n - A_n = M_{n-1} - M'_{n-1}$$

Thus, for $n < 1$,

$$A'_1 - A_1 = M_0 - M'_0 = X_0 - X_0 = 0$$

Hence $A'_1 = A_1$, and by induction $A'_n = A_n \forall n$: since

$$M_n + A_n = X_n = M'_n + A'_n$$

we obtain that $M_n = M'_n \forall n$

3.3.1 Convergence theorem of martingales

Theorem 3.3.1. Let X_n be a submartingale. If $\sup_n E[|X_n|] < \infty$, then:

1. X_n converges almost surely (a.s.) to some limit X : $X_n \xrightarrow{a.s.} X$
2. If X_n is uniformly integrable, then X_n converges a.s. and in L^1 to X : $X_n \xrightarrow{a.s., L^1} X$

• Remark

– X_n is said to be uniformly integrable if:

$$\lim_{c \rightarrow +\infty} \sup_n E\{|X_n| \cdot \mathbb{1}_{|X_n| > c}\} = 0$$

If the $\sup_n E\{|X_n|^{1+\epsilon}\}$ is finite for some $\epsilon > 0 \Rightarrow X_n$ is uniformly integrable. The contrary is not true.

It is also possible to show that if X_n is uniformly integrable $\Rightarrow \sup_n E|X_n| < +\infty$

Let's show an example where the $\sup_n E|X_n| < +\infty$ but X_n is not uniformly integrable.

Let X_n be such that

$$P(X_n = 0) = \frac{n-1}{n} \text{ and } P(X_n = n) = \frac{1}{n}$$

Then, $E|X_n| = n \cdot P(X_n = n) + 0 \cdot P(X_n = 0) = n \cdot \frac{1}{n} = 1$.

Hence, $\sup_n E|X_n| = \sup_n 1 = 1 < +\infty$.

Fix $c > 0$, then $\sup_n E(|X_n| \cdot \mathbb{1}_{|X_n| > c})$

$$= \sup_n [E\{\mathbb{1}_{X_n=0} \cdot \mathbb{1}_{|X_n| > c} \cdot (|X_n|)\} + E\{\mathbb{1}_{X_n=n} \cdot \mathbb{1}_{|X_n| > c} \cdot (|X_n|)\}] \quad (3.3)$$

$$= \sup_n E\{\mathbb{1}_{(X_n=0)} \cdot \mathbb{1}_{(X_n=c)} \cdot n\}$$

$$= \sup_n n \cdot E\{\mathbb{1}_{(X_n=n)}\}$$

$$= \sup_n n \cdot \frac{1}{n} = 1$$

then

$$\lim_{c \rightarrow +\infty} \sup_n E(|X_n| \cdot \mathbb{1}_{|X_n| > c}) = 1 \neq 0$$

Has to be completed

Chapter 4

Markov Processes - Markov Chains

4.1 Definition

Definition 4.1.1. A sequence X_0, X_1, X_2, \dots of real i.i.d. random variables to be a markov chain with respect to a filtration $\mathfrak{F}_0 \subset \mathfrak{F}_1 \subset \dots \subset \mathfrak{A}$ if two conditions are satisfied:

1. X_n is \mathfrak{F}_n -measurable $\forall n$
2. $P(X_{n+1} \in A | \mathfrak{F}_n) = P(X_{n+1} \in A | X_n) \forall n \geq 0 \forall A \in \mathfrak{B}(\mathfrak{R})$

Roughly speaking, we are saying that $P(\text{future} | \text{Past and present}) = P(\text{future} | \text{present})$

4.2 Homogeneous Markov Chain

Definition 4.2.1. Let $\alpha(x, A) = P(X_{n+1} \in A | X_n = x) \forall x \in \mathbb{R}, A \in \mathfrak{B}(\mathbb{R})$ and $n \geq 0$

\Downarrow

then the sequence (X_n) is homogeneous $\leftrightarrow a_n = a_0 \forall n \geq 0$.

If (X_n) is an homogeneous markov chain, the probability distribution of the sequence is determined by:

- The probability distribution of X_0 , usually called initial distribution;
- The conditional distribution α_0 , usually called transition kernel.

If we write $\alpha = \alpha_0$ and we call Π the probability distribution of X_0 , we have, by applying the definition, that the

$$\begin{aligned} P(X_0 \in A_0, X_1 \in A_1) &= E_x \{ P(X_1 \in A_1 | X_0 = x) \cdot \mathbb{1}_{(x)} \}_{A_0} \\ &= E_x \alpha(x, A_1) \cdot \mathbb{1}_{x \in A_0} \\ &= \sum_x \alpha(x, A_1) \cdot \mathbb{1}_{x \in A_0} \cdot \Pi(x) \text{ if } X_0 \text{ is discrete.} \end{aligned}$$

Instead, if X_0 is absolutely continuous,

$P(X_0 \in A_0, X_1 \in A_1) = \int \alpha(x_j, A_1) \cdot \mathbb{1}_{x \in A_0} \cdot f(x)$, where f is the density of X_0

Suppose for instance, that X_n is discrete for every n : then,

$$\begin{aligned} P(X_0 = x_0, X_1 = x_1, X_2 = x_2) &= P(X_2 = x_2 | X_0 = x_0, X_1 = x_1) \cdot P(X_0 = x_0, X_1 = x_1) \\ &= P(X_2 = x_2 | X_0 = x_0, X_1 = x_1) \cdot P(X_1 = x_1 | X_0 = x_0) \cdot P(X_0 = x_0) \\ &= P(X_2 = x_2 | X_1 = x_1) \cdot P(X_1 = x_1 | X_0 = x_0) \cdot P(X_0 = x_0) \\ &= \alpha(x_1, \{x_2\}) \cdot \alpha(x_0, \{x_1\}) \cdot \Pi(x_0) \end{aligned}$$

4.2.1 Stationary sequence

In general, a sequence (X_n) is said to be stationary if

$$(X_1, X_2, X_3, \dots) \sim (X_0, X_1, X_2, X_3, \dots).$$

If (X_n) is stationary, then $X_n \sim X_0 \forall n \geq 0$, but not conversely.

Instead, if X_n is an homogeneous markov chain, then

$$(X_n) \text{ is stationary} \Leftrightarrow X_n \sim X_0 \forall n \geq 0$$

Example 4.2.1. Any random walk is an homogeneous markov chain: in fact, let $X_n = X_0 + \sum_{i=1}^n Z_i$ where

- X_0 is independent of (Z_i)
- (Z_i) is iid

Let $\mathfrak{F}_n = \sigma(X_0, X_1, x_2, \dots, X_n)$, Then,

$$\begin{aligned} P(X_{n+1} \in A | X_0 = x_0, \dots, X_n = x_n) &= P(X_n + Z_{n+1} \in A | X_0 = x_0, \dots, X_n = x_n) \\ &= P(x_n + Z_{n+1} \in A | X_0 = x_0, \dots, X_n = x_n) \end{aligned}$$

and since independence

$$= P(x_n + Z_{n+1} \in A)$$

and since they are identically distributed

$$= P(X_n + Z_1 \in A)$$

So, similarly

$$P(X_{n+1} \in A | X_n = x_n) = P(x_n + Z_1 \in A)$$

In this case

$$\alpha(x, A) = P(X_1 \in A | X_0 = x) = P(x + Z_1 \in A)$$

In particular, let us consider the symmetric random walk on \mathbb{Z} :

$$X_0 \in \mathbb{Z} \text{ and } P(Z_1 = 1) = P(Z_1 = -1) = \frac{1}{2}$$

In this case $\alpha(i, \{j\}) = P(X_1 = j | X_0 = i)$, that can be written as

$$= P(Z_1 + i = j) = P(Z_1 = j - i)$$

$$= \begin{cases} \frac{1}{2} & \text{if } j - i = 1 \text{ or } j - i = -1 \\ 0 & \text{otherwise} \end{cases} \quad (4.1)$$

If you have reach this point, well done! Don't give up, we are almost at the half of our journey!!

4.2.2 Recurrent chains

So far we considered a general homogeneous markov chain: from now on, we deal with an important special case, namely we assume X_n is discrete $\forall n \geq 0$.

In this case, we have

$$P(X_n \in S, \forall n \geq 0) = 1$$

where S is a finite or countable set, called the **state space** of the chain and $x \in S$ is called **state**.

A state x can be **recurrent** or **transient**:

- is said to be **recurrent** if $P(\exists n > 0 \text{ with } X_n = x | X_0 = x) = 1$
- is said to be **transient** if $P(\exists n > 0 \text{ with } X_n = x | X_0 = x) < 1$

Let's recall the definition of a stopping time:

$$\tau_x = \inf\{n > 0 : X_n = x\} \begin{cases} \text{first time } n > 0 \text{ such that } X_n = x \\ +\infty \text{ if } X_n \neq x \forall n > 0 \end{cases} \quad (4.2)$$

Now, we can show that if x is recurrent, then

- x is null recurrent if

$$E(\tau_x | X_0 = x) = +\infty$$

- x is positive recurrent if

$$E(\tau_x | X_0 = x) < +\infty$$

4.2.3 Irreducible chains

An homogeneous markov chain is **irreducible** if $\forall x, y \in S \exists m \geq 1 \text{ and } n \geq 1$ such that

- $P(X_n = y | X_0 = x) > 0$
- $P(X_n = x | X_0 = y) > 0$

revise the notation above. Now, let α be the transition kernel:

$$\alpha(x, A) = P(X_1 \in A | X_0 = x) \quad \forall x \in S, \forall A \in \mathcal{B}(\mathbb{R})$$

In this case, since S is countable we will write

$$\alpha(x, y) \text{ instead of } \alpha(x, \{y\}) \quad \forall x, y \in S$$

Hence, $\alpha(x, y) = P(X_1 = y | X_0 = x)$ In this notation, let us suppose $S = \mathbb{Z}$ and $\alpha(n, n+1) = 1$ In this case, $\alpha(n, n+1)$ represents the probability of transitioning from state n to state $n+1$ in one step, given that the current state is n . Since $\alpha(n, n+1) = 1$, it means that the chain can only move to state $n+1$ from state n and cannot stay in state n or move to any other state in one step.

Next, suppose

- $S = \{a, b\}$
- $\alpha(a, b) = \alpha(b, a)$

if the chain is irreducible, all states are of the same type. If S is finite we can conclude that \exists a positive recurrent state but \nexists any null recurrent state.

Theorem 4.2.1. Let $x \in S$. Then,

$$x \text{ is recurrent} \Leftrightarrow \sum_{n=1}^{+\infty} P(X_n = x | X_0 = x) = +\infty \Leftrightarrow P(X_n = x \text{ for infinitely many } n | X_0 = x) = 1$$

The first equivalence follows from the definition of recurrent state, which is that starting from x , the chain returns to x with probability 1. This means that the probability of returning to x in n steps, given that the chain started at x is $P(X_n = x | X_0 = x)$. So if x is recurrent, then the probability of ever returning to x is infinite, which implies that the sum of these probabilities over all n is infinite as well.