

Stochastic Modelling and Random Processes

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Chapter 1

Introduction

1.1 Motivation

Suppose we are modelling COVID. Let

- S be the number of the susceptible;
- I be the number of the infected;
- R be the number of the removed (those who have either recovered or died).

1.1.1 A Deterministic Model

A deterministic model might be

$$\begin{aligned}\dot{S} &= -\beta IS, \\ \dot{I} &= \beta IS - \gamma I, \\ \dot{R} &= \gamma I.\end{aligned}$$

But there are some problems in this model:

- S , I and R are integers, so it does not make sense to talk about \dot{S} , \dot{I} and \dot{R} .
- There is variability in when contacts are made and lead to infection.

1.1.2 A Stochastic Model

A better model might be stochastic

$$\begin{aligned}\mathbb{P}S \rightarrow S - 1 \ \& \ I \rightarrow I - 1 \text{ in } \Delta t = \beta IS \Delta t + o(\Delta t) \\ \mathbb{P}I \rightarrow I - 1 \ \& \ R \rightarrow R + 1 \text{ in } \Delta t = \gamma I \Delta t + o(\Delta t).\end{aligned}$$

The problem of this model is that contacts are usually not made uniformly in the whole population.

1.1.3 A Network Model

We can use a network model, in which nodes represent individuals and edge weights represent contact rates, to avoid uniform contacts. But this is unrealistic: the network is too big to represent 60 million people in the UK.

1.1.4 A Random Network Model

Based on the network model, we can make probability distributions on networks and derive probabilistic conclusions over the combination of stochastic dynamics and randomness of networks.

Chapter 2

Probability and Random Variables

2.1 Probability Theory

Suppose we are doing an experiment which have different random outcomes.

Definition 2.1.1 (Sample Spaces). The **sample space** of the experiment is the set of all possible outcomes, denoted as Ω .

Definition 2.1.2 (Sigma Algebra). The **σ -algebra** of subsets of Ω , denoted as \mathcal{F} , is a set of subsets of Ω which satisfies:

- $\Omega \in \mathcal{F}$;
- $A \in \mathcal{F} \implies A^c \in \mathcal{F}$;
- $\{A_i | i \in \mathcal{I}\} \subset \mathcal{F}$ with \mathcal{I} being countable $\implies \bigcup_{i \in \mathcal{I}} A_i \in \mathcal{F}$.

Remark. We say \mathcal{I} is countable if there exists a one-to-one map from \mathcal{I} into \mathbb{Z} , so “countable” includes “finite”.

Example 2.1.1. If Ω is countable, we usually take $\mathcal{F} = 2^\Omega$, which is the **power set** of Ω .

Example 2.1.2. When Ω is not countable, e.g. $[0, 1]$, if you allow **Axiom of Choice**¹, then there exist unmeasurable subsets, and we exclude them from \mathcal{F} , i.e. \mathcal{F} is the set of all Lebesgue-measurable subsets on $[0, 1]$.

Definition 2.1.3 (Events). The members of \mathcal{F} are called **events**.

Definition 2.1.4 (Probability). $\mathbb{P}[\cdot] : \mathcal{F} \mapsto \mathbb{R}$ is called a probability if

- $\mathbb{P}[A] \in [0, 1], \forall A \in \mathcal{F}$;
- $\mathbb{P}[\Omega] = 1$;
- $\mathbb{P}[\cdot]$ satisfies the **countable additivity**: $\forall \{A_i | i \in \mathcal{I}\} \subset \mathcal{F}$, where \mathcal{I} is a countable set, if A_i 's are disjoint, then

$$\mathbb{P}\left[\bigcup_{i \in \mathcal{I}} A_i\right] = \sum_{i \in \mathcal{I}} \mathbb{P}[A_i].$$

¹A Cartesian product of a collection of nonempty sets is nonempty.

Definition 2.1.5 (Independence). Say $A, B \in \mathcal{F}$ are **independent** if

$$\mathbb{P}[A \cap B] = \mathbb{P}[A]\mathbb{P}[B].$$

Definition 2.1.6 (Conditional Probabilities). If $\mathbb{P}[B] > 0$, then the **conditional probability** $\mathbb{P}[A|B]$ is defined by

$$\mathbb{P}[A|B] = \frac{\mathbb{P}[A \cap B]}{\mathbb{P}[B]}, \quad \forall A \in \mathcal{F}.$$

Definition 2.1.7 (Partitions). $\{B_i | i \in \mathcal{I}\}$ is called a **partition** of the sample space Ω if:

- B_i 's are **pairwise disjoint**: $B_i \cap B_j = \emptyset, \forall i, j \in \mathcal{I}, i \neq j$;
- $B_i \neq \emptyset, \forall i \in \mathcal{I}$;
- $\{B_i | i \in \mathcal{I}\}$ **covers** Ω : $\bigcup_{i \in \mathcal{I}} B_i = \Omega$.

Theorem 2.1.1 (The Law of Total Probability). Let $\{B_i | i \in \mathcal{I}\}$ be a countable partition of Ω , with $B_i \in \mathcal{F}$ and $\mathbb{P}[B_i] > 0, \forall i \in \mathcal{I}$. Then $\forall A \in \mathcal{F}$, we have

$$\mathbb{P}[A] = \sum_{i \in \mathcal{I}} \mathbb{P}[A|B_i] \mathbb{P}[B_i].$$

Theorem 2.1.2 (Bayes' Rule). For any events A and B , if $\mathbb{P}[A] > 0$ and $\mathbb{P}[B] > 0$, then

$$\mathbb{P}[B|A] = \frac{\mathbb{P}[A|B] \mathbb{P}[B]}{\mathbb{P}[A]}.$$

Furthermore, if $\{B_i | i \in \mathcal{I}\}$ is a countable partition of Ω , with $B_i \in \mathcal{F}$ and $\mathbb{P}[B_i] > 0, \forall i \in \mathcal{I}$, then

$$\mathbb{P}[B_i|A] = \frac{\mathbb{P}[A|B_i] \mathbb{P}[B_i]}{\sum_{i \in \mathcal{I}} \mathbb{P}[A|B_i] \mathbb{P}[B_i]}.$$

Example 2.1.3. Suppose the **true positive rate** $\mathbb{P}(\text{tests positive} | \text{has COVID})$ is 0.99 and the **false positive rate** $\mathbb{P}(\text{tests positive} | \text{does not have COVID})$ is 0.01. Suppose in the population, the probability of getting contracted with COVID is 0.001, i.e. $\mathbb{P}(\text{has COVID}) = 0.001$, what is the probability that a person has COVID given his/her test is positive?

Sol.

$$\begin{aligned} & \mathbb{P}(\text{has COVID} | \text{tests positive}) \\ &= \frac{\mathbb{P}(\text{tests positive} | \text{has COVID}) \mathbb{P}(\text{has COVID})}{\mathbb{P}(\text{tests positive})} \\ &= \frac{\mathbb{P}(\text{tests positive} | \text{has COVID}) \mathbb{P}(\text{has COVID})}{\mathbb{P}(\text{tests positive} | \text{has COVID}) \mathbb{P}(\text{has COVID}) + \mathbb{P}(\text{tests positive} | \text{does not have COVID}) \mathbb{P}(\text{does not have COVID})} \\ &= \frac{0.99 \times 0.001}{0.99 \times 0.001 + 0.01 \times (1 - 0.001)} \\ &\approx 0.090. \end{aligned}$$

2.2 Random Variables

Definition 2.2.1 (Measurable Functions). Let $(\Omega, \mathcal{F}, \mathbb{P})$ and $(\mathbb{R}, \Sigma, \mathcal{L})$ be two measurable spaces, where \mathcal{L} is the Lebesgue measure. For any function $f : \Omega \mapsto \mathbb{R}$, if it satisfies $\forall A \in \Sigma, f^{-1}(A) \in \mathcal{F}$, then f is said to be **measurable**.

Remark. $(\Omega, \mathcal{F}, \mathbb{P})$ and $(\mathbb{R}, \Sigma, \mathcal{L})$ can be generalized:

Let (X, Σ) and (Y, T) be measurable spaces, meaning that X and Y are sets equipped with respective σ -algebras Σ and T . A function $f : X \mapsto Y$ is said to be **measurable** if for every $E \in T$ the pre-image of E under f is in Σ ; i.e. $\forall E \in T$,

$$f^{-1}(E) := \{x \in X | f(x) \in E\} \in \Sigma.$$

Definition 2.2.2 (Random Variables). A **random variable** is a measurable function $X : \Omega \mapsto \mathbb{R}$.

Definition 2.2.3 (Cumulative Distribution Functions). The **cumulative distribution function** of a random variable X is defined as

$$F(x) = \mathbb{P}[X \leq x]$$

Definition 2.2.4 (Discrete Random Variables). If $X(\Omega)$ is countable, then X is called **discrete**.

Definition 2.2.5 (Probability Mass Functions). The **probability mass function** of a discrete random variable X is defined as

$$\pi(x) = \mathbb{P}[X = x], \quad \forall x \in X(\Omega).$$

Definition 2.2.6 (Continuous Random Variables & Probability Density Functions). For a random variable X , if its cumulative distribution function satisfies

$$F(x) = \int_{-\infty}^x f(y) dy$$

for some $f \in \mathcal{L}^1(\mathbb{R})$, then X is said to be **continuous**, and f is its **probability density function**.

Remark. It is possible to have mixtures. For example, X can have a positive probability on a particular point and continuous parts on other points.

Definition 2.2.7 (Expectation). The **expectation** of a random variable X is

$$\mathbb{E}[X] := \int_{\Omega} X d\mathbb{P} = \begin{cases} \sum_{x \in X(\Omega)} x \pi(x) & X \text{ is discrete} \\ \int_{X(\Omega)} x f(x) dx & X \text{ is continuous} \end{cases}$$

Remark. The expectation may be infinite or even undefined.

Definition 2.2.8 (Variance). The **variance** of a random variable X is

$$\text{Var}[X] := \mathbb{E}[(X - \mathbb{E}[X])^2] = \mathbb{E}[X^2] - \mathbb{E}[X]^2.$$

Definition 2.2.9 (Covariance). The **covariance** of two random variables X and Y is

$$\text{Cov}[X, Y] = \mathbb{E}[(X - \mathbb{E}[X])(Y - \mathbb{E}[Y])] = \mathbb{E}[XY] - \mathbb{E}[X]\mathbb{E}[Y].$$

Definition 2.2.10 (Uncorrelated Random Variables). If $\text{Cov}[X, Y] = 0$, then X and Y are called uncorrelated.

Proposition 2.2.1. If X and Y are two independent random variables, then they are also uncorrelated. But the opposite is generally not true, except for Gaussians.

We can extend to random variables taking values in \mathbb{R}^n .

- For **cumulative distribution functions**, use the component-wise \leq instead.
- For $\text{Var}[X]$, use $\mathbb{E}[(X - \mathbb{E}[X])(X - \mathbb{E}[X])^T]$ which is a $n \times n$ matrix.
- For $\text{Cov}[X, Y]$, use $\mathbb{E}[(X - \mathbb{E}[X])(Y - \mathbb{E}[Y])^T]$.
- X and Y are independent if events $\{X \leq x\}$, $\{Y \leq y\}$ are independent, $\forall x, y$.
 - For X, Y being discrete, this is equivalent to $\pi(x, y) = \pi^X(x)\pi^Y(y)$.
 - For X, Y being continuous, this is equivalent to $f(x, y) = f^X(x)f^Y(y)$.

Theorem 2.2.1 (The Weak Law of Large Numbers). Let X_k , $k = 1, 2, \dots, X_n, \dots$ be independent and identically distributed random variables with $\mu = \mathbb{E}[X_k] < \infty$, then

$$\bar{X}_n := \frac{1}{n} \sum_{k=1}^n X_k \xrightarrow{d} \mu, \text{ as } n \rightarrow \infty,$$

where \xrightarrow{d} means **convergence in distribution**². This means the CDF of \bar{X}_n converges to the CDF of μ .

Equivalently,

$$\mathbb{E} \left[g(\bar{X}_n) \right] \rightarrow g(\mu), \text{ as } n \rightarrow \infty,$$

for any bounded and continuous function g . This type of convergence is called the **weak convergence**.

Or \bar{X}_n converges to μ **in probability** ($\bar{X}_n \xrightarrow{\mathbb{P}} \mu$):

$$\mathbb{P} \left[|\bar{X}_n - \mu| > \epsilon \right] \rightarrow 0, \text{ as } n \rightarrow \infty, \forall \epsilon > 0.$$

Theorem 2.2.2 (The Strong Law of Large Numbers). Let X_k , $k = 1, 2, \dots, X_n, \dots$ be independent and identically distributed random variables with $\mu = \mathbb{E}[X_k] < \infty$, then

$$\bar{X}_n \xrightarrow{a.s.} \mu,$$

where the **almost surely convergence** means

$$\mathbb{P} \left[\lim_{n \rightarrow \infty} \bar{X}_n = \mu \right] = 1.$$

Theorem 2.2.3 (Central Limit Theorem). Let X_k , $k = 1, 2, \dots, X_n, \dots$ be independent and identically distributed random variables with $\mu = \mathbb{E}[X_k] < \infty$ and $0 < \sigma^2 := \text{Var}[X_k] < \infty$, then

$$\frac{\sqrt{n}}{\sigma} (\bar{X}_n - \mu) \xrightarrow{d} \mathcal{N}(0, 1).$$

²Also called **convergence in law**.

Theorem 2.2.4 (Large Deviation Principle). *Let X_k , $k = 1, 2, \dots, X_n, \dots$ be independent and identically distributed random variables. For any interval $J \subset \mathbb{R}$,*

$$\mathbb{P} [\bar{X}_n \in J] \approx \exp \left(-n \min_{x \in J} I(x) \right),$$

meaning

$$\frac{1}{n} \log \mathbb{P} [\bar{X}_n \in J] \rightarrow - \min_{x \in J} I(x).$$

*If we know the probability distribution of X_k , an explicit expression for the rate function can be obtained. This is given by a **Legendre–Fenchel transformation**,*

$$I(x) = \sup_{\theta > 0} \theta x - \lambda(\theta),$$

*where $\lambda(\theta) = \log \mathbb{E} [e^{\theta X_k}]$ is called the **cumulant generating function (CGF)**.*

Definition 2.2.11 (Stochastic Processes). A **stochastic process** $\{X(t) | t \in T\}$ is a collection of random variables. That is, for each $t \in T$, $X(t)$ is a random variable.

- The index t is often interpreted as time and, as a result, we refer to $X(t)$ as the **state** of the process at time t .
- The set T is called the **index set** of the process.
 - When T is a countable set, the process is said to be a **discrete-time** process.
 - If T is an interval of the real line, the stochastic process is said to be a **continuous-time** process.
- The **state space** of a stochastic process is defined as the set of all possible values that the random variables $X(t)$ can assume.

Chapter 3

Discrete-Time Markov Chain

3.1 Discrete-Time Markov Chains

Definition 3.1.1 (Discrete-Time Stochastic Processes). A **discrete-time stochastic process** with state space S is a sequence $\{Y_n | n \in \mathbb{N}\}$ of random variables taking values in S .

Definition 3.1.2 (Discrete-Time Markov Chains). Let $\{X_n | n \in \mathbb{N}\}$ be a discrete-time stochastic process with a discrete state space S . The process is called a **Markov chain**, if for all $A \subset S$, $n \in \mathbb{N}$ and $s_0, \dots, s_n \in S$,

$$\mathbb{P}[X_{n+1} \in A | X_n = s_n, \dots, X_0 = s_0] = \mathbb{P}[X_{n+1} \in A | X_n = s_n].$$

Proposition 3.1.1. *For any Markov chain $\{X_n | n \in \mathbb{N}\}$, conditional on the present, the past and the future are independent, i.e. $\forall n \in \mathbb{N}_+, \forall s_n \in S, X_{n+1} | X_n = s$ and $X_{n-1} | X_n = s$ are independent.*

Proof.

$$\begin{aligned} & \mathbb{P}[X_{n+1} = s_{n+1}, X_{n-1} = s_{n-1} | X_n = s_n] \\ &= \frac{\mathbb{P}[X_{n-1} = s_{n-1}, X_n = s_n, X_{n+1} = s_{n+1}]}{\mathbb{P}[X_n = s_n]} \\ &= \mathbb{P}[X_{n-1} = s_{n-1}] \cdot \mathbb{P}[X_n = s_n | X_{n-1} = s_{n-1}] \cdot \mathbb{P}[X_{n+1} = s_{n+1} | X_n = s_n, X_{n-1} = s_{n-1}] \cdot \frac{1}{\mathbb{P}[X_n = s_n]} \\ &= \mathbb{P}[X_{n-1} = s_{n-1}] \cdot \mathbb{P}[X_n = s_n | X_{n-1} = s_{n-1}] \cdot \mathbb{P}[X_{n+1} = s_{n+1} | X_n = s_n] \cdot \frac{1}{\mathbb{P}[X_n = s_n]} \\ &= \mathbb{P}[X_{n-1} = s_{n-1}] \cdot \frac{\mathbb{P}[X_{n-1} = s_{n-1} | X_n = s_n] \cdot \mathbb{P}[X_n = s_n]}{\mathbb{P}[X_{n-1} = s_{n-1}]} \cdot \mathbb{P}[X_{n+1} = s_{n+1} | X_n = s_n] \cdot \frac{1}{\mathbb{P}[X_n = s_n]} \\ &= \mathbb{P}[X_{n-1} = s_{n-1} | X_n = s_n] \cdot \mathbb{P}[X_{n+1} = s_{n+1} | X_n = s_n] \end{aligned}$$

□

3.1.1 Homogeneity

Definition 3.1.3 (Homogeneity). A Markov chain $\{X_n | n \in \mathbb{N}\}$ is **homogeneous** if for all $A \subset S$, $n \in \mathbb{N}$ and $s \in S$,

$$\mathbb{P}[X_{n+1} \in A | X_n = s] = \mathbb{P}[X_1 \in A | X_0 = s].$$

Example 3.1.1 (Random Walk with Boundaries). Let $\{X_n | n \in \mathbb{N}\}$ be a **simple random walk** on $S = \{1, \dots, L\}$ with $p(x, y) = p\delta_{y, x+1} + q\delta_{y, x-1}$. The boundary conditions are

- **periodic** if $p(L, 1) = p, p(1, L) = q$,
- **absorbing** if $p(L, L) = 1, p(1, 1) = 1$,
- **closed** if $p(L, L) = p, p(1, 1) = q$,
- **reflecting** if $p(L, L - 1) = 1, p(1, 2) = 1$.

3.1.2 Transition Matrices and Transition Functions

Definition 3.1.4 (One-Step Transition Matrices). For a homogeneous discrete-time Markov chain $\{X_n | n \in \mathbb{N}\}$ taking values in $\{s_1, s_2, s_3, \dots, s_n, \dots\}$, its **one-step transition matrix** P is defined as

$$P_{i,j} = \mathbb{P} [X_{n+1} = s_j | X_n = s_i] .$$

Remark. The sum of each row of a one-step transition matrix is 1, i.e.

$$P |1\rangle = |1\rangle .$$

Proposition 3.1.2. Let $\pi_0(\cdot)$ be the probability mass function of X_0 , then

$$\mathbb{P} [X_0 = s_0, X_1 = s_1, \dots, X_n = s_n] = \pi_0(s_0) P_{s_0, s_1} \cdots P_{s_{n-1}, s_n} .$$

If we use a row vector $\langle \pi_0 |$ to represent the probability distribution of X_0 , such that $\langle \pi_0 |_i = \mathbb{P} [X_0 = s_i]$, then the probability distribution of X_n can be represented as

$$\langle \pi_n | = \langle \pi_0 | P^n .$$

Definition 3.1.5 (Transition Functions). The transition matrix of $\{X_n | n \in \mathbb{N}\}$ can be written into the **transition function** $p_n(x, y)$ instead:

$$p_n(x, y) := \mathbb{P} [X_n = y | X_0 = x] .$$

3.1.3 Chapman-Kolmogorov Equations

Theorem 3.1.1 (Chapman-Kolmogorov Equations). For a homogeneous discrete-time Markov chain $\{X_n | n \in \mathbb{N}\}$, its transition function fulfills the **Chapman-Kolmogorov equations**

$$p_{k+n}(x, y) = \sum_{z \in S} p_k(x, z) p_n(z, y) \quad \text{for all } k, n \geq 0, x, y \in S .$$

Remark. In matrix form, the Chapman-Kolmogorov equations read

$$P_{n+k} = P_n P_k \quad \text{and in particular} \quad P_{n+1} = P_n P_1 .$$

Corollary 3.1.1. Let P_n be the n -step transition matrix of a homogeneous discrete-time Markov chain $\{X_n | n \in \mathbb{N}\}$, then

$$P_n = P^n \quad \& \quad P_0 = I .$$

3.1.4 Stationary Distributions

Definition 3.1.6 (Stationarity). Let $\{X_n | n \in \mathbb{N}\}$ be a homogeneous discrete-time Markov chain with state space S . The distribution $\pi(x)$, $x \in S$, is called **stationary** if for all $y \in S$

$$\sum_{x \in S} \pi(x) p(x, y) = \pi(y),$$

or

$$\langle \pi | P = \langle \pi |.$$

Remark. If π is a stationary distribution, then it is a left eigenvector with eigenvalue 1.

Remark. To solve the stationary distributions, we can solve

$$\begin{cases} \langle \pi | P &= \langle \pi | \\ \langle \pi | \mathbf{1} &= 1 \end{cases}$$

Theorem 3.1.2. *Every homogeneous finite discrete-time Markov chain has a stationary distribution.*

Proof. Let

$$\Delta = \left\{ \langle \pi | \mid \pi_i \geq 0, \langle \pi | \mathbf{1} = 1 \right\}$$

Then $P_{i,j} \geq 0$ and $P | \mathbf{1} = | \mathbf{1}$, so $\pi \in \Delta \implies \langle \pi | P \in \Delta$. Notice that Δ is compact and convex, and P is continuous (linear), so by the [Brouwer's Fixed-Point Theorem](#), P has a fixed point $\langle \pi^* | \in \Delta$, such that $\langle \pi^* | P = \langle \pi^* |$. \square

Remark. There can be more than one stationary distributions. For example, if a Markov chain has two parts with no transitions between them, then let $\langle \pi_1 |$ and $\langle \pi_2 |$ be stationary probabilities of them, and any convex combination of $\langle \pi_1 |$ and $\langle \pi_2 |$ is a stationary distribution.

Definition 3.1.7 (Cycles). A **cycle** is a closed path in S along the graph of allowed transitions by P , and its length is greater than 0.

Definition 3.1.8 (Transience & Recurrence). Say $i \in S$ is **transient** if there does not exist any cycle through i . Otherwise, i is **recurrent**.

Definition 3.1.9 (Communication). Say $i, j \in S$ communicate with each other if there exist a cycle through i and j , denoted as $i \leftrightarrow j$.

Proposition 3.1.3. *Communication is an equivalent relation on the set of all recurrent states:*

- $i \leftrightarrow i, \forall i \in S$;
- $i \leftrightarrow j \iff j \leftrightarrow i, \forall i, j \in S$;
- $i \leftrightarrow j \leftrightarrow k \implies i \leftrightarrow k, \forall i, j, k \in S$.

Definition 3.1.10 (Classes, Communicating Components). A **class** (also called a **communicating component**) is a set of all communicating states in the state space.

Chapter 4

Continuous-Time Markov Chain

Chapter 5

Continuous State Space Markov Processes

Chapter 6

Stochastic Particle Systems

Chapter 7

Networks

Chapter 8

Random Networks