

AN INTRODUCTION TO SDES

MARIA CAMERON

CONTENTS

1. Definitions	2
2. Expected values and moments	4
3. Independence, joint distributions, covariance	6
4. Chebyshev's inequality	9
5. Types of convergence of random variables	10
6. Laws of Large Numbers and the Central Limit Theorem	13
7. Conditional probability and conditional expectation	14
8. Brownian Motion	16
8.1. Definition of Brownian Motion	16
8.2. Existence of Brownian motion	17
8.3. Construction of the Brownian motion via a refinement procedure	17
8.4. Construction of the Brownian motion by a random walk	20
8.5. Elementary properties of Brownian motion	21
9. Brownian motion and heat equation	22
10. Numerical solution of the heat equation and its solution by random walk	24
10.1. A brief introduction into the Wiener measure	27
10.2. Markov property of Brownian motion	28
11. An Introduction to SDEs.	28
11.1. Elementary properties of stochastic integral	31
11.2. Construction of the Ito integral	32
11.3. Existence and uniqueness of solutions of the Ito SDEs	32
11.4. Notation common in probability books.	33
11.5. The Ornstein-Uhlenbeck process	33
12. The Ito calculus	34
12.1. A derivation of the Ito formula	34
12.2. The geometric Brownian motion	36
12.3. Backward Kolmogorov equation	37
12.4. The expected first passage time	38
12.5. The committor equation	39
12.6. The generator of a Markov process	40
12.7. The adjoint semigroup and the forward Kolmogorov equation	41
12.8. The invariant pdf	43

13. The Langevin Dynamics	44
13.1. The full Langevin dynamics	44
13.2. The overdamped Langevin dynamics	44
14. Numerical integration of SDEs	45
14.1. The Euler-Maruyama and Milstein's methods	45
14.2. Strong and weak orders of convergence	46
References	47

1. DEFINITIONS

A probability space is a triple consisting of *the set of outcomes*, the set of subsets of the set of outcomes that we want to be able to assign probabilities to called *the σ -algebra*, and *the probability measure*, i.e. a function that assigns probabilities.

- A **sample space** Ω is the set of all possible outcomes.
- An **event** A is a subset of Ω .
- A **σ -algebra** \mathcal{B} is a subset of the set of all subsets of Ω that is closed with respect to set operations. The minimal requirements guaranteeing that the σ -algebra possesses these properties constitute the set of axioms that defines it:
 - (1) $\emptyset \in \mathcal{B}$ and $\Omega \in \mathcal{B}$;
 - (2) If $B \in \mathcal{B}$ then $B^c \in \mathcal{B}$ (B^c is the complement of B in Ω , i.e., $B^c \equiv \Omega \setminus B$).
 - (3) If $\mathcal{A} = \{A_1, \dots, A_n, \dots\}$ is a finite or countable collection in \mathcal{B} then

$$\bigcup_i A_i \in \mathcal{B}.$$

Corollary: If $\mathcal{A} = \{A_1, \dots, A_n, \dots\}$ is a finite or countable collection in \mathcal{B} then

$$\bigcap_i A_i \in \mathcal{B}.$$

Indeed,

$$\bigcap_i A_i = \left(\bigcup_i A_i^c \right)^c.$$

Example 1 Suppose you are tossing a die. For a single throw, the sample space is $\Omega = \{1, 2, 3, 4, 5, 6\}$. If you are interested in particular number on the top, the natural choice of the σ -algebra is the set of all subsets of Ω . Then $|\mathcal{B}| = 2^6 = 64$. If you are interested only in whether the outcome is odd or even, then a reasonable choice of σ -algebra is

$$\mathcal{B} = \{\emptyset, \{1, 3, 5\}, \{2, 4, 6\}, \{1, 2, 3, 4, 5, 6\}\}.$$

If you are interested only whether there is an outcome or not, you can choose the coarsest σ -algebra

$$\mathcal{B} = \{\emptyset, \{1, 2, 3, 4, 5, 6\}\}.$$

Example 2 Suppose you are doing a measurement whose outcome can be any real number. For example, you are living in a one-dimensional world, you are throwing a point object, and measuring its position with respect to a fixed point, i.e. the origin of a coordinate system in your 1D-world. The set of outcomes is \mathbb{R} . The most commonly chosen σ -algebra is the so-called *Borel σ -algebra* which is generated by all open sets in \mathbb{R} . Thanks to the properties of σ -algebra, the Borel σ -algebra can be generated by all intervals of the form $(-\infty, a]$, where $a \in \mathbb{R}$.

- A **probability measure** P is a function $P : \mathcal{B} \rightarrow [0, 1]$ such that
 - (1) $P(\Omega) = 1$;
 - (2) $0 \leq P(A) \leq 1$ for all $A \in \mathcal{B}$.
 - (3) **Countable additivity:** If $\mathcal{A} = \{A_1, \dots, A_n, \dots\}$ is a finite or countable collection in \mathcal{B} such that $A_i \cap A_j = \emptyset$ for any i, j , then

$$P\left(\bigcup_i A_i\right) = \sum_i P(A_i).$$

Corollary: $P(\emptyset) = 0$. Indeed,

$$1 = P(\Omega) = P(\Omega \cup \emptyset) = P(\Omega) + P(\emptyset) = 1 + P(\emptyset).$$

Hence, $P(\emptyset) = 0$.

- A **probability space** is the triple (Ω, \mathcal{B}, P) .
- A **random variable** η is a \mathcal{B} -measurable function $\eta : \Omega \rightarrow \mathbb{R}$.

A function is called \mathcal{B} -measurable if the preimage of any measurable subset of \mathbb{R} is in \mathcal{B} . It is proven in analysis that it suffices to check that $\{\omega \in \Omega \mid \eta(\omega) \leq x\} \in \mathcal{B}$ for any $x \in \mathbb{R}$.

- A **probability distribution function** of a random variable η is defined by

$$F_\eta(x) = P(\{\omega \in \Omega \mid \eta(\omega) \leq x\}) = P(\eta \leq x).$$

Theorem 1. *If F is a probability distribution function then*

- (1) F is nondecreasing, i.e. $x < y$ implies $F(x) \leq F(y)$.
- (2) $\lim_{x \rightarrow -\infty} F(x) = 0$, $\lim_{x \rightarrow \infty} F(x) = 1$.
- (3) $F(x)$ is continuous from the right for every $x \in \mathbb{R}$, i.e.,

$$\lim_{y \rightarrow x+0} F(y) = F(x).$$

Example 3 Suppose you are tossing a die. Consider the probability space

$$(1) \quad (\Omega = \{1, 2, 3, 4, 5, 6\}, \mathcal{B} = 2^\Omega, P(\omega) = \frac{1}{6}),$$

where 2^Ω is the set of all subsets of Ω , and $\omega \in \Omega = \{1, 2, 3, 4, 5, 6\}$. Consider the random variable $\eta(\omega) = \omega$. The probability distribution

function is given by

$$F_\eta(x) = \begin{cases} 0, & x < 1, \\ j/6, & j \leq x < j + 1, \quad j = 1, 2, 3, 4, 5 \\ 1, & x \geq 6. \end{cases}$$

- Suppose $F'_\eta(x)$ exists. Then $f_\eta(x) \equiv F'_\eta(x)$ is called the **probability density function (pdf)** of the random variable η , and

$$P(x < \eta \leq x + dx) = F_\eta(x + dx) - F_\eta(x) = f_\eta(x)dx + o(dx).$$

Example 4 The Gaussian density

$$f(x) = \frac{1}{\sqrt{2\pi\sigma^2}} e^{-\frac{(x-m)^2}{2\sigma^2}},$$

where m and σ are constants. m is the mean, while σ is the standard deviation.

Example 5 The density of an exponential random variable with parameter $a > 0$ is given by:

$$f(x) = \begin{cases} ae^{-ax}, & x \geq 0, \\ 0, & x < 0. \end{cases}$$

Example 6 The density of a uniform random variable on an interval $[a, b]$ is

$$f(x) = \frac{1}{b-a} I_{[a,b]}(x) = \begin{cases} \frac{1}{b-a}, & x \in [a, b], \\ 0, & \text{otherwise.} \end{cases}$$

Here $I_{[a,b]}(x)$ is the indicator function of the interval $[a, b]$.

- If the set of outcomes Ω is discrete (finite or countable) and the σ -algebra is the set of all subsets Ω , then the function $P(\omega)$ is often called **the probability mass function**.

2. EXPECTED VALUES AND MOMENTS

Definition 1. Let (Ω, \mathcal{B}, P) be a probability space, and η be a random variable. Then the expected value, or mean, of the random variable η is defined as

$$(2) \quad E[\eta] = \int_{\Omega} \eta(\omega) dP.$$

If Ω is a discrete set,

$$E[\eta] = \sum_i \eta(\omega_i) P(\omega_i).$$

Example 7 Suppose you are tossing a die. Consider the probability space (1) and the random variable $\eta(\omega) = \omega$, $\omega = 1, 2, 3, 4, 5, 6$. The expected value of η is

$$E[\eta] = \sum_{j=1}^6 j \frac{1}{6} = 3.5$$

Suppose that the random variable η is fixed. Then we will omit the subscript in the notation of its probability distribution function: $F_{\eta}(x) \equiv F(x)$.

The integral in Eq. (2) can be rewritten using $F(x)$:

$$E[\eta] = \int_{\mathbb{R}} x P(x < \eta \leq x + dx) = \int_{-\infty}^{\infty} x dF(x).$$

If a derivative $f(x)$ of the probability distribution function F exists, then

$$E[\eta] = \int_{-\infty}^{\infty} x f(x) dx.$$

If g is a function defined on the range of the random variable η (on $\eta(\Omega)$), then the expected value of this function is

$$E[g(\eta)] = \int_{-\infty}^{\infty} g(x) dF(x).$$

Moments: Let us take $g(x) = x^n$.

$$E[\eta^n] = \int_{-\infty}^{\infty} x^n dF(x).$$

Central moments: Let us take $g(x) = (x - E[\eta])^n$.

$$E[(\eta - E[\eta])^n] = \int_{-\infty}^{\infty} (x - E[\eta])^n dF(x).$$

Variance = 2nd central moment:

$$\text{Var}(\eta) = E[(\eta - E[\eta])^2] = \int_{-\infty}^{\infty} (x - E[\eta])^2 dF(x).$$

Example 8 Suppose you are tossing a die. Consider the probability space (1) and the random variable $\eta(\omega) = \omega$, $\omega = 1, 2, 3, 4, 5, 6$. The variance of η is

$$\text{Var}(\eta) = \frac{1}{6} \sum_{j=1}^6 (j - 3.5)^2 = \frac{35}{12} = 2.91(6).$$

The standard deviation:

$$\sigma(\eta) = \sqrt{\text{Var}(\eta)}.$$

3. INDEPENDENCE, JOINT DISTRIBUTIONS, COVARIANCE

- Two events $A, B \in \mathcal{B}$ are **independent** if

$$P(A \cap B) = P(A)P(B).$$

- Two random variables η_1 and η_2 are independent if the events

$$(3) \quad \{\omega \in \Omega \mid \eta_1(\omega) \leq x\} \quad \text{and} \quad \{\omega \in \Omega \mid \eta_2(\omega) \leq y\}$$

are independent for all $x, y \in \mathbb{R}$.

Example 9 Suppose you are tossing a die twice. Consider the probability space

$$(4) \quad (\Omega = \{1, 2, 3, 4, 5, 6\}^2, \mathcal{B} = 2^{\Omega^2}, P(\{\omega_1, \omega_2\}) = 1/36), \quad 1 \leq \omega_1, \omega_2 \leq 6.$$

Let η_1 and η_2 be random variables equal to the outcomes of the first and

TABLE 1. Two throws of a die. Values of the random variables $\xi(\omega_1, \omega_2) = \omega_1 + \omega_2$ (left) and $\beta(\omega_1, \omega_2) = \omega_1 - \omega_2$ (right).

	1	2	3	4	5	6		1	2	3	4	5	6
1	2	3	4	5	6	7	1	0	1	2	3	4	5
2	3	4	5	6	7	8	2	-1	0	1	2	3	4
3	4	5	6	7	8	9	3	-2	-1	0	1	2	3
4	5	6	7	8	9	10	4	-3	-2	-1	0	1	2
5	6	7	8	9	10	11	5	-4	-3	-2	-1	0	1
6	7	8	9	10	11	12	6	-5	-4	-3	-2	-1	0

the second throws respectively. These random variables are independent. Now consider the random variables $\eta(\omega_1, \omega_2) = \omega_1$ and $\xi(\omega_1, \omega_2) = \omega_1 + \omega_2$ (see Table 1, left). We can show that η and ξ are dependent by taking e.g., $x = 1$ and $y = 2$ in Eq. (3):

$$P(\eta \leq 1 \ \& \ \xi \leq 2) = \frac{1}{36} \neq P(\eta \leq 1)P(\xi \leq 2) = \frac{1}{6} \cdot \frac{1}{36} = \frac{1}{216}.$$

Finally, we consider the random variables $\xi(\omega_1, \omega_2) = \omega_1 + \omega_2$ and $\beta(\omega_1, \omega_2) = \omega_1 - \omega_2$ (see Table 1, right). We can show that they are dependent by taking e.g., $x = 2$ and $y = -1$ in Eq. (3):

$$P(\xi \leq 2 \ \& \ \beta \leq -1) = 0 \neq P(\xi \leq 2)P(\beta \leq -1) = \frac{1}{36} \cdot \frac{15}{36} = \frac{5}{432}.$$

- The joint distribution function of two random variables η_1 and η_2 is given by

$$F_{\eta_1 \eta_2}(x, y) = P(\{\omega \in \Omega \mid \eta_1(\omega) \leq x, \eta_2(\omega) \leq y\}) = P(\eta_1(\omega) \leq x, \eta_2(\omega) \leq y).$$

- If the second mixed derivative of $F_{\eta_1\eta_2}$ exists, it is called the **joint probability density** of η_1 and η_2 and denoted by

$$f_{\eta_1\eta_2}(x, y) := \frac{\partial F_{\eta_1\eta_2}(x, y)}{\partial x \partial y}.$$

In this case,

$$F_{\eta_1, \eta_2}(x, y) = \int_{-\infty}^x \int_{-\infty}^y f_{\eta_1\eta_2}(x, y) dx dy.$$

Exercise Show that two random variables are independent if and only if

$$F_{\eta_1\eta_2}(x, y) = F_{\eta_1}(x)F_{\eta_2}(y).$$

Furthermore, if the joint pdf $f_{\eta_1\eta_2}(x, y)$ exists, then η_1 and η_2 are independent iff

$$f_{\eta_1\eta_2}(x, y) = f_{\eta_1}(x)f_{\eta_2}(y).$$

- Given the joint pdf $f_{\eta_1\eta_2}$, one can obtain $f_{\eta_1}(x)$ by

$$f_{\eta_1}(x) = \int_{-\infty}^{\infty} f_{\eta_1\eta_2}(x, y) dy.$$

In this equation, f_{η_1} is called a **marginal** of $f_{\eta_1\eta_2}$, and the variable η_2 is **integrated out**.

- **Properties of expected value and variance** It follows from the definition, that the expected value is a linear functional:

$$(5) \quad E[a\eta_1 + b\eta_2] = aE[\eta_1] + bE[\eta_2].$$

•

$$(6) \quad \text{Var}(a\eta) = a^2\text{Var}(\eta).$$

- If η_1 and η_2 are independent, then

$$(7) \quad \text{Var}(\eta_1 + \eta_2) = \text{Var}(\eta_1) + \text{Var}(\eta_2).$$

If η_1 and η_2 are dependent, (7) is not true: take $\eta_1 = \eta_2$. In general,

$$(8) \quad \text{Var}(\eta_1 + \eta_2) = \text{Var}(\eta_1) + \text{Var}(\eta_2) + 2\text{Cov}(\eta_1, \eta_2),$$

where $\text{Cov}(\eta_1, \eta_2)$ is the covariance of η_1 and η_2 – see below. You will see below that (7) does not imply that η_1 and η_2 are independent, only that they are uncorrelated.

Example 10 Suppose you are tossing a die twice. Consider the probability space and random variables introduced in Example 9. Then

$$E[\xi] = E[\eta_1 + \eta_2] = E[\eta_1] + E[\eta_2] = 7.$$

$$E[\beta] = E[\eta_1 - \eta_2] = E[\eta_1] + E[-\eta_2] = 0.$$

$$\text{Var}[\xi] = \text{Var}[\eta_1 + \eta_2] = \text{Var}[\eta_1] + \text{Var}[\eta_2] = \frac{35}{6} = 5.8(3).$$

$$\text{Var}[\beta] = \text{Var}[\eta_1 - \eta_2] = \text{Var}[\eta_1] + \text{Var}[-\eta_2] = \text{Var}[\eta_1] + \text{Var}[\eta_2] = \frac{35}{6} = 5.8(3).$$

Example 11 Consider the Bernoulli random variable

$$(9) \quad \eta = \begin{cases} 1, & P(1) = p, \\ 0, & P(0) = 1 - p. \end{cases}$$

Its expected value and variance are

$$E[\eta] = 1 \cdot p + 0 \cdot (1 - p) = p,$$

$$\text{Var}(\eta) = (1 - p)^2 \cdot p + (0 - p)^2 \cdot (1 - p) = p(1 - p).$$

Now consider the sum of n independent copies of η :

$$\xi := \sum_{i=1}^n \eta_i.$$

Using Eq. (5) we calculate $E[\xi]$:

$$E[\xi] = \sum_{i=1}^n E[\eta_i] = np.$$

Since η_i , $1 \leq i \leq n$, are independent, we can calculate $\text{Var}(\xi)$ using Eq. (7):

$$\text{Var}(\xi) = \sum_{i=1}^n \text{Var}(\eta_i) = np(1 - p).$$

Finally, consider the average of n independent copies of η :

$$\zeta := \frac{1}{n} \sum_{i=1}^n \eta_i \equiv \frac{\xi}{n}.$$

Using Eqs. (5) and (6), we find

$$E[\zeta] = p,$$

$$\text{Var}(\zeta) = \text{Var}\left(\frac{\xi}{n}\right) = \frac{1}{n^2} \text{Var}(\xi) = \frac{p(1 - p)}{n}.$$

- The **covariance** of two random variables η_1 and η_2 is defined by

$$\text{Cov}(\eta_1, \eta_2) = E[(\eta_1 - E[\eta_1])(\eta_2 - E[\eta_2])].$$

Remark If η_1 and η_2 are independent, then $\text{Cov}(\eta_1, \eta_2) = 0$. If $\text{Cov}(\eta_1, \eta_2) = 0$ then η_1 and η_2 are uncorrelated. Note that uncorrelated random variables are not necessarily independent.

Example 12 Suppose you are tossing a die twice. Consider the probability space and random variables introduced in Example 9. As we have established in Example 9, ξ and β are dependent. However, they are uncorrelated. Indeed,

$$\begin{aligned} \text{Cov}(\xi, \beta) &= \sum_{1 \leq \omega_1 \leq 6, 1 \leq \omega_2 \leq 6} (\omega_1 + \omega_2 - 7)(\omega_1 - \omega_2)P(\{\omega_1, \omega_2\}) \\ &= \frac{1}{36} \left(\sum_{\omega_1 < \omega_2} (\omega_1 + \omega_2 - 7)(\omega_1 - \omega_2) + \sum_{\omega_1 > \omega_2} (\omega_1 + \omega_2 - 7)(\omega_1 - \omega_2) \right) = 0. \end{aligned}$$

Example 13 A vector-valued random variable $\eta = [\eta_1, \dots, \eta_n]$ is jointly Gaussian if

$$P(x_1 < \eta_1 \leq x_1 + dx_1, \dots, x_n < \eta_n \leq x_n + dx_n) = \frac{1}{Z} e^{-\frac{1}{2}(x-m)^\top A^{-1}(x-m)} dx + o(dx),$$

where $x = [x_1, \dots, x_n]^\top$, $m = [m_1, \dots, m_n]^\top$, $dx = dx_1 \dots dx_n$, and A is a symmetric positive definite matrix. The normalization constant Z is given by

$$Z = (2\pi)^{n/2} |A|^{1/2}, \quad \text{where } |A| = \det A.$$

In the case of jointly Gaussian random variables, the covariance matrix C whose entries are

$$C_{ij} = E[(\eta_i - E[\eta_i])(\eta_j - E[\eta_j])]$$

is equal to A . Two jointly Gaussian random variables are independent if and only if they are uncorrelated.

4. CHEBYSHEV'S INEQUALITY

Chebyshev's inequality holds for any random variable. It is a very useful theoretical tool for proving various estimates. In practice, it often gives too rough estimates which is a consequence of its universality. Chebyshev's inequality is not improvable, as we can construct a random variable for which it turns into an equality.

Theorem 2. Let η be a random variable. Suppose $g(x)$ is a nonnegative, nondecreasing function (i.e., $g(x) \geq 0$, $g(a) \leq g(b)$ whenever $a < b$). Then for any $a \in \mathbb{R}$

$$(10) \quad P(\eta \geq a) \leq \frac{E[g(\eta)]}{g(a)}.$$

Proof.

$$\begin{aligned} E[g(\eta)] &= \int_{-\infty}^{\infty} g(x) dF(x) \\ &\geq \int_a^{\infty} g(x) dF(x) \geq g(a) \int_a^{\infty} dF(x) = g(a)P(\eta \geq a). \end{aligned}$$

□

Given a random variable η we define a random variable

$$\xi := |\eta - E[\eta]|.$$

Define

$$g(x) = \begin{cases} x^2, & x \geq 0, \\ 0, & x < 0. \end{cases}$$

Plugging this into Eq. (10) we obtain

$$P(|\eta - E[\eta]| \geq a) \leq \frac{\text{Var}(\eta)}{a^2}.$$

Example 14 Suppose you are tossing a die twice. Consider the probability space and random variables introduced in Example 9. We will compare the exact probabilities with their Chebyshev estimates.

$$P(|\xi - 7| \geq 1) = P(\xi \neq 7) = 1 - \frac{6}{36} = \frac{5}{6} = 0.8(3), \quad \frac{\text{Var}(\xi)}{1} = \frac{35}{6} = 5.8(3);$$

$$P(|\xi - 7| \geq 2) = P(\xi \leq 5 \text{ or } \xi \geq 9) = \frac{20}{36} = \frac{5}{9} = 0.5(5), \quad \frac{\text{Var}(\xi)}{4} = \frac{35}{24} = 1.458(3);$$

$$P(|\xi - 7| \geq 3) = P(\xi \leq 4 \text{ or } \xi \geq 10) = \frac{12}{36} = \frac{1}{3} = 0.3(3), \quad \frac{\text{Var}(\xi)}{9} = \frac{35}{54} = 0.6(481);$$

$$P(|\xi - 7| \geq 4) = P(\xi \in \{2, 3, 11, 12\}) = \frac{6}{36} = \frac{1}{6} = 0.1(6), \quad \frac{\text{Var}(\xi)}{16} = \frac{35}{96} = 0.36458(3);$$

$$P(|\xi - 7| \geq 5) = P(\xi \in \{2, 12\}) = \frac{2}{36} = \frac{1}{18} = 0.0(5), \quad \frac{\text{Var}(\xi)}{25} = \frac{35}{150} = 0.2(3);$$

Choosing $a = k\sigma$ we get

$$P(|\eta - E[\eta]| \geq k\sigma) \leq \frac{1}{k^2}.$$

This means that for *any* random variable η defined on *any* probability space we have that the probability that η deviates from its expected value by at least k standard deviations does not exceed $1/k^2$.

The bounds given Chebyshev's inequality cannot be improved in principle, because they are exact for the random variable

$$\eta = \begin{cases} 1, & P = \frac{1}{2k^2}, \\ 0, & P = 1 - \frac{1}{k^2}, \\ -1, & P = \frac{1}{2k^2}. \end{cases}$$

It is easy to check that $E[\eta] = 0$, $\text{Var}(\eta) = \frac{1}{k^2}$. Hence

$$P(|\eta| \geq 1) = \frac{1}{k^2} = \frac{\text{Var}(\eta)}{1^2},$$

i.e. Chebyshev's inequality turns into equality.

5. TYPES OF CONVERGENCE OF RANDOM VARIABLES

Suppose we have a sequence of random variables $\{\eta_1, \eta_2, \dots\}$. In probability theory, there exist several different notions of convergence of a sequence of random variables $\{\eta_1, \eta_2, \dots\}$ to some limit random variable η .

- $\{\eta_1, \eta_2, \dots\}$ **converges in distribution** or **converges weakly**, or **converges in law** to η if

$$(11) \quad \lim_{n \rightarrow \infty} F_n(x) = F(x) \text{ for every } x \text{ where } F(x) \text{ is continuous,}$$

where F_n and F are the probability distribution functions of η_n and η respectively.

Remark Convergence of pdfs $f_n(x)$ implies convergence of $F_n(x)$. The converse is not true in general. For example, consider $F_n(x) = x - \frac{1}{2\pi n} \sin(2\pi nx)$, $x \in (0, 1)$. The corresponding pdf is $f_n(x) = 1 - \cos(2\pi nx)$, $x \in (0, 1)$. $\{F_n(x)\}$ converges to $F(x) = x$, i.e., to the uniform distribution, while $\{f_n(x)\}$ does not converge at all.

Remark In the discrete case, the convergence of probability mass functions $f(k) := P(\eta = k)$ implies the convergence of the probability distribution functions.

Example 15 Consider the sum of n independent copies of the Bernoulli random variable as in Example 11:

$$(12) \quad \xi = \sum_{i=1}^n \eta_i, \quad \text{where } \eta_i = \begin{cases} 1, & P(1) = p, \\ 0, & P(0) = 1 - p. \end{cases}$$

Its probability distribution is the binomial distribution given by

$$(13) \quad f(k; n, p) \equiv P(\xi = k) = \binom{n}{k} p^k (1-p)^{n-k},$$

where $\binom{n}{k}$ is the number of k -combinations of the set of n elements:

$$\binom{n}{k} = \frac{n!}{k!(n-k)!}.$$

Now we let $n \rightarrow \infty$ and $p \rightarrow 0$ in such a manner that the product np (i.e., the expected value of ξ) remains constant. We introduce the parameter

$$\lambda := np.$$

Consider the sequence of random variables ξ_n where ξ_n is the sum of n independent copies of Bernoulli random variable with $p = \lambda/n$, i.e.,

$$(14) \quad \xi_n = \sum_{i=1}^n \eta_i^{(n)}, \quad \text{where } \eta_i^{(n)} = \begin{cases} 1, & P(1) = \lambda/n, \\ 0, & P(0) = 1 - \lambda/n. \end{cases}$$

Plugging in $p = \lambda/n$ in the results of Example 11 we find the expected value and the variance:

$$E[\xi_n] = n \frac{\lambda}{n} = \lambda.$$

$$\text{Var}(\xi_n) = n \frac{\lambda}{n} \left(1 - \frac{\lambda}{n}\right) = \lambda \left(1 - \frac{\lambda}{n}\right).$$

We will show that the sequence ξ_n converges to the Poisson random variable with parameter λ in distribution. Consider the limit

$$\begin{aligned} \lim_{n \rightarrow \infty} f\left(k; n, \frac{\lambda}{n}\right) &= \lim_{n \rightarrow \infty} \frac{n(n-1)\dots(n-k+1)}{k!} \frac{\lambda^k}{n^k} \left(1 - \frac{\lambda}{n}\right)^{n-k} = \\ &= \frac{\lambda^k}{k!} \lim_{n \rightarrow \infty} \frac{n(n-1)\dots(n-k+1)}{n^k} \lim_{n \rightarrow \infty} \left(1 - \frac{\lambda}{n}\right)^n \lim_{n \rightarrow \infty} \left(1 - \frac{\lambda}{n}\right)^{-k} \end{aligned}$$

The first limit in the equation above is 1 as $n(n-1)\dots(n-k+1) = n^k + O(n^{k-1})$. The second limit can be calculated using the well-known fact that

$$\lim_{n \rightarrow \infty} \left(1 + \frac{1}{n}\right)^n = e.$$

Hence

$$\lim_{n \rightarrow \infty} \left(1 - \frac{\lambda}{n}\right)^n = e^{-\lambda}.$$

The third limit is 1. Therefore,

$$\lim_{n \rightarrow \infty} \frac{n!}{k!(n-k)!} \left(\frac{\lambda}{n}\right)^k \left(1 - \frac{\lambda}{n}\right)^{n-k} = \frac{\lambda^k}{k!} e^{-\lambda},$$

which is the Poisson distribution with parameter λ .

- $\{\eta_1, \eta_2, \dots\}$ **converges in probability** to η if for any $\epsilon > 0$

$$(15) \quad \lim_{n \rightarrow \infty} P(|\eta_n - \eta| \geq \epsilon) = 0$$

Remark Convergence in probability implies convergence in distribution.

Proof. We will prove this fact for the case of scalar random variables. We have $\lim_{n \rightarrow \infty} P(|\eta_n - \eta| \geq \epsilon) = 0$, we need to prove $\lim_{n \rightarrow \infty} P(\eta_n \leq x) = P(\eta \leq x)$ for every x where F_η is continuous. First we show an auxiliary fact that for any two random variables ξ and ζ , $x \in \mathbb{R}$ and $\epsilon > 0$

$$(16) \quad P(\xi \leq a) \leq P(\zeta \leq a + \epsilon) + P(|\xi - \zeta| > \epsilon).$$

Indeed,

$$\begin{aligned} P(\xi \leq a) &= P(\xi \leq a \ \& \ \zeta \leq a + \epsilon) + P(\xi \leq a \ \& \ \zeta > a + \epsilon) \\ &\leq P(\zeta \leq a + \epsilon) + P(\xi - \zeta \leq a - \zeta \ \& \ a - \zeta < -\epsilon) \\ &\leq P(\zeta \leq a + \epsilon) + P(\zeta - \xi < -\epsilon) \\ &\leq P(\zeta \leq a + \epsilon) + P(\zeta - \xi < -\epsilon) + P(\zeta - \xi > \epsilon) \\ &= P(\zeta \leq a + \epsilon) + P(|\zeta - \xi| > \epsilon). \end{aligned}$$

Applying Eq. (77) to $\xi = \eta_n$ and $\zeta = \eta$ with $a = x$ and $a = x - \epsilon$, we get

$$\begin{aligned} P(\eta_n \leq x) &\leq P(\eta \leq x + \epsilon) + P(|\eta_n - \eta| > \epsilon) \\ P(\eta \leq x - \epsilon) &\leq P(\eta_n \leq x) + P(|\eta_n - \eta| > \epsilon). \end{aligned}$$

$$P(\eta \leq x - \epsilon) - P(|\eta_n - \eta| > \epsilon) \leq P(\eta_n \leq x) \leq P(\eta \leq x + \epsilon) + P(|\eta_n - \eta| > \epsilon).$$

Taking the limit $n \rightarrow \infty$ and taking into account that $\lim_{i \rightarrow \infty} P(|\eta_n - \eta| \geq \epsilon) = 0$, we get

$$F_\eta(x - \epsilon) \leq \lim_{n \rightarrow \infty} F_{\eta_n}(x) \leq F_\eta(x + \epsilon).$$

If x is a point of continuity of F_η ,

$$\lim_{\epsilon \rightarrow 0} F_\eta(x - \epsilon) = \lim_{\epsilon \rightarrow 0} F_\eta(x + \epsilon) = F_\eta(x).$$

Therefore, taking the limit $\epsilon \rightarrow 0$ we obtain the weak convergence:

$$\lim_{n \rightarrow \infty} F_{\eta_n}(x) = F_\eta(x)$$

for any x where $F_\eta(x)$ is continuous. □

Remark The converse is, generally, not true. However, convergence in distribution to a *constant* random variable implies convergence in probability.

- $\{\eta_1, \eta_2, \dots\}$ **converges almost surely** or **almost everywhere** or **with probability 1** or **strongly** to η if

$$(17) \quad P\left(\lim_{n \rightarrow \infty} \eta_n = \eta\right) = 1.$$

Remark Convergence almost surely implies convergence in probability (by Fatou's lemma) and in distribution.

- To summarize,

$$(18) \quad \boxed{\eta_i \rightarrow \eta \text{ almost surely}} \Rightarrow \boxed{\eta_i \rightarrow \eta \text{ in probability}} \Rightarrow \boxed{\eta_i \rightarrow \eta \text{ in distribution}}$$

6. LAWS OF LARGE NUMBERS AND THE CENTRAL LIMIT THEOREM

- Let $\{\eta_1, \eta_2, \dots\}$ be a sequence of random variables with finite expected values $\{m_1 = E[\eta_1], m_2 = E[\eta_2], \dots\}$. Define

$$\xi_n = \frac{1}{n} \sum_{i=1}^n \eta_i, \quad \bar{\xi}_n = \frac{1}{n} \sum_{i=1}^n m_i.$$

Definition 2. (1) *The sequence of random variables η_n satisfies the Law of Large Numbers if $\xi_n - \bar{\xi}_n$ converges to zero in probability, i.e., for any $\epsilon > 0$*

$$\lim_{n \rightarrow \infty} P(|\xi_n - \bar{\xi}_n| > \epsilon) = 0.$$

(2) *The sequence of random variables η_n satisfies the Strong Law of Large Numbers if $\xi_n - \bar{\xi}_n$ converges to zero almost surely, i.e., for almost all $\omega \in \Omega$*

$$\lim_{n \rightarrow \infty} \xi_n - \bar{\xi}_n = 0.$$

- If the random variables η_n are independent and if $\text{Var}(\eta_i) \leq V < \infty$, then the Law of Large Numbers holds by the Chebyshev Inequality (10):

$$\begin{aligned} P(|\xi_n - \bar{\xi}_n| > \epsilon) &= P\left(\left|\sum_{i=1}^n \eta_i - \sum_{i=1}^n m_i\right| > n\epsilon\right) \\ &\leq \frac{\text{Var}(\eta_1 + \dots + \eta_n)}{\epsilon^2 n^2} \leq \frac{V}{\epsilon^2 n} \rightarrow 0 \text{ as } n \rightarrow \infty. \end{aligned}$$

•

Theorem 3. (Khinchin) *A sequence of independent identically distributed random variables $\{\eta_i\}$ with $\mathbb{E}[\eta_i] = m$ and $\mathbb{E}[|\eta_i|] < \infty$ satisfies the Law of Large Numbers.*

•

Theorem 4. (Kolmogorov) *A sequence of independent identically distributed random variables with finite expected value and variance satisfies the Strong Law of Large Numbers.*

•

Theorem 5. (The central limit theorem) *Let $\{\eta_1, \eta_2, \dots\}$ be a sequence of independent identically distributed (i.i.d.) random variables with $m = E[\eta_i]$ and $0 < \sigma^2 = \text{Var}(\eta_i) < \infty$, then*

$$(19) \quad \frac{(\sum_{i=1}^n \eta_i) - nm}{\sigma\sqrt{n}} \longrightarrow N(0, 1) \text{ in distribution,}$$

i.e., converges weakly to the standard normal distribution $N(0, 1)$ (i.e., the Gaussian distribution with mean 0 and variance 1) as $n \rightarrow \infty$.

A proof via Fourier transform can be found in [1]. Another proof making use of characteristic functions can be found in [2].

Remark Eq. (19) can be recasted as

$$(20) \quad \frac{1}{n} \sum_{i=1}^n \eta_i \longrightarrow N\left(m, \frac{\sigma^2}{n}\right) \text{ in distribution,}$$

i.e., the average of the first n i.i.d. random variables η_i converges in distribution to the Gaussian random variable with mean $m = E[\eta_i]$ and variance σ^2/n .

7. CONDITIONAL PROBABILITY AND CONDITIONAL EXPECTATION

- The conditional probability of an event B given that the event A has happened is given by

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

Note that if A and B are independent, then $P(A \cap B) = P(A)P(B)$ and hence

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

Example 16 Suppose you are tossing a die twice. Consider the probability space (4). Let A be the event that the outcome of the first throw is even, and B be the event that the sum of the outcomes is ≥ 10 . Then (see Table 1)

$$P(B|A) = \frac{P(A \cap B)}{P(A)} = \frac{4/36}{1/2} = \frac{2}{9}.$$

Note that $P(B) = 1/6 < P(B|A)$. Hence the events A and B are dependent.

If the event A is fixed, then $P(B|A)$ defines a probability measure on (Ω, \mathcal{B}) .

- If η is a random variable on Ω , then conditional expectation of η given the event A is

$$E[\eta|A] = \int_{\Omega} \eta(\omega)P(d\omega|A) = \int_{\Omega} \eta(\omega) \frac{P(d\omega \cap A)}{P(A)} = \frac{\int_A \eta(\omega)P(d\omega)}{P(A)}.$$

Example 17 . Suppose you are tossing a die twice. Consider the probability space (4). Let A be the event that the outcome of the first throw is even, and η be the random variable whose value is the sum of outcomes, *i.e., $\eta(\{\omega_1, \omega_2\}) = \omega_1 + \omega_2$.* Then

$$E[\eta|A] = \sum_{\omega_1=1}^6 \sum_{\omega_2=1}^6 (\omega_1 + \omega_2)P(\{\omega_1, \omega_2\} | \omega_1 \in \{2, 4, 6\}).$$

Let us calculate $P(\{\omega_1, \omega_2\} \mid \omega_1 \in \{2, 4, 6\})$.

$$P(\{\omega_1, \omega_2\} \mid \omega_1 \in \{2, 4, 6\}) = \frac{P(\{\omega_1, \omega_2\} \cap (\omega_1 \in \{2, 4, 6\}))}{P(\omega_1 \in \{2, 4, 6\})}$$

$$= \begin{cases} 0, & \omega_1 \in \{1, 3, 5\}, \\ \frac{P(\{\omega_1, \omega_2\})}{P(\omega_1 \in \{2, 4, 6\})} = \frac{1/36}{1/2} = \frac{1}{18}, & \omega_1 \in \{2, 4, 6\}. \end{cases}$$

Now we continue our calculation:

$$E[\omega_1 + \omega_2 \mid \omega_1 \in \{2, 4, 6\}] = \sum_{\omega_1 \in \{2, 4, 6\}} \sum_{\omega_2=1}^6 (\omega_1 + \omega_2) \frac{1}{18} = \frac{135}{18} = 7.5.$$

Note that $E[\eta] = 7 \neq E[\eta|A] = 7.5$.

- Now we show how one can construct new random variables using conditional probability. For simplicity, we start with partitioning the set of outcomes Ω into a finite or countable number of disjoint measurable subsets:

$$\Omega = \bigcup_i A_i, \quad \text{where } A_i \in \mathcal{B}, \quad A_i \cap A_j = \emptyset.$$

Definition 3. Let η be a random variable on the probability space (Ω, \mathcal{B}, P) . Let $\mathcal{A} = \{A_i\}$ be a partition of Ω as above. Define a new random variable $E[\eta|\mathcal{A}]$ as follows:

$$(21) \quad E[\eta|\mathcal{A}] = \sum_i E[\eta|A_i] \chi(A_i),$$

where $\chi(A_i)$ is the indicator function of A_i :

$$\chi(A_i; \omega) = \begin{cases} 1, & \omega \in A_i, \\ 0, & \omega \notin A_i. \end{cases}$$

Remark Note that $E[\eta|\mathcal{A}]$ is a random variable as it is a function of the outcome ω . Indeed,

$$E[\eta|\mathcal{A}](\omega) = E[\eta|A_i] \quad \text{where } A_i \ni \omega.$$

Example 18 Suppose you are tossing a die twice. Let us partition the set of outcomes as follows:

$$\Omega = \bigcup_{i=1}^6 \{(\omega_1, \omega_2) \mid \omega_1 = i\}.$$

The corresponding partition \mathcal{A} is

$$\mathcal{A} = \{ \{(\omega_1, \omega_2) \mid \omega_1 = i\} \}_{i=1}^6.$$

Take the random variable $\xi = \omega_1 + \omega_2$ (see Table 1, left), the sum of numbers on the top. Construct a new random variable

$$\begin{aligned} E[\xi|\mathcal{A}] &= \sum_{i=1}^6 E[\xi|\omega_1 = i]\chi(\omega_1 = i) = \sum_{i=1}^6 (i + 3.5)\chi(\omega_1 = i) \\ &= 4.5\chi(\omega_1 = 1) + 5.5\chi(\omega_1 = 2) + 6.5\chi(\omega_1 = 3) \\ &\quad + 7.5\chi(\omega_1 = 4) + 8.5\chi(\omega_1 = 5) + 9.5\chi(\omega_1 = 6). \end{aligned}$$

- Now we define the conditional expectation of one random variable η given the other random variable θ . First we assume that θ assumes a finite or countable number of values $\{\theta_1, \theta_2, \dots\}$. Define the partition \mathcal{A} where

$$A_i = \{\omega \in \Omega \mid \theta = \theta_i\}.$$

Definition 4. We define a new random variable $E[\eta|\theta]$ as a the following function of the random variable θ :

$$E[\eta|\theta] := E[\eta|\mathcal{A}], \quad \text{i.e.,} \quad E[\eta|\theta] = E[\eta|A_i] \quad \text{if} \quad \theta = \theta_i.$$

Example 19 Suppose you are tossing a die twice. Let (ω_1, ω_2) be the numbers on the top. Define random variables $\xi = \omega_1 + \omega_2$ and $\theta = \omega_1$. Then it follows from our calculation from the previous example that

$$E[\xi|\theta] = 3.5 + \theta.$$

8. BROWNIAN MOTION

Various processes in nature are often modeled by stochastic differential equations of the form

$$dx = b(x, t)dt + \sigma(x, t)dw,$$

where the function $b(x, t)$ is called *the drift field*, the matrix function $\sigma(x, t)$ is called the diffusion matrix, and the factor dw is the increment of the stochastic process called *the Brownian motion*. The goal of this section is to understand what the Brownian motion is.

8.1. Definition of Brownian Motion.

Definition 5. A stochastic process (in the strict sense) is a function $v(\omega, t)$ of two arguments, where $\omega \in \Omega$, (Ω, \mathcal{B}, P) is a probability space, and $t \in \mathbb{R}$, such that

- for each ω , $v(\omega, t)$ is a function of t , and
- for each t , $v(\omega, t)$ is a random variable.

Definition 6. Brownian motion (in mathematical terminology) is a stochastic process $w(\omega, t)$, $\omega \in \Omega$, $0 \leq t < \infty$, that satisfies the following four axioms:

- (1) $w(\omega, 0) = 0$ for all ω .
- (2) For almost all ω , $w(\omega, t)$ is a continuous function of t .

- (3) For each $0 \leq s \leq t$, $w(\omega, t) - w(\omega, s)$ is a Gaussian random variable with mean 0 and variance $t - s$.
- (4) $w(\omega, t)$ has independent increments, i.e., if

$$0 \leq t_1 < t_2 < \dots < t_n$$

then

$$w(\omega, t_i) - w(\omega, t_{i-1}) \text{ for } i = 2, \dots, n \text{ are independent.}$$

Remark What is called the Brownian motion in mathematics is called the Wiener process in physics. What is called the Brownian motion in physics is called the Ornstein-Uhlenbeck process in mathematics.

Here is an equivalent definition of Brownian motion.

Definition 7. A process $w(\omega, t)$ on a probability space (Ω, \mathcal{B}, P) is called a Brownian motion if

- (1) Sample paths $w(\omega, t)$ are continuous functions of t for almost all $\omega \in \Omega$.
- (2) For any $k > 1$ and $0 \leq t_1 \leq \dots \leq t_k$, the random vector $(w(\omega, t_1), \dots, w(\omega, t_k))$ is Gaussian with mean 0 and covariance matrix

$$B(t_i, t_j) = E[w(t_i), w(t_j)] = \min\{t_i, t_j\} \equiv t_i \wedge t_j, \quad 1 \leq i, j \leq k.$$

Definition 8. A d -dimensional Brownian motion is defined as the vector process

$$w(t) = (w_1(t), \dots, w_d(t)),$$

where $w_k(t)$, $1 \leq k \leq d$ are independent Brownian motions.

8.2. Existence of Brownian motion. The question about the existence of the Brownian motion is not trivial. For example, if we upgrade axiom 2 in Definition 6 to require differentiability, such a process simply would not exist.

The original construction of Brownian motion (the Wiener process) was done by Norbert Wiener (1894 - 1964). He has shown that Fourier series

$$(22) \quad w(t) = \frac{a_0}{\sqrt{\pi}}t + \sqrt{\frac{2}{\pi}} \sum_{k=1}^{\infty} \frac{a_k}{k} \sin(kt),$$

where a_k , $k = 0, 1, 2, \dots$, are independent Gaussian random variables with mean 0 and variance 1, converges, and its sum satisfies Definition 6 for $0 \leq t \leq 1$.

In [2], the existence of Brownian motion follows from Kolmogorov's theorem about the existence of stochastic processes with covariance satisfying certain conditions.

8.3. Construction of the Brownian motion via a refinement procedure. The Brownian motion on $0 \leq t \leq 1$ can be constructed by a recursive refinement procedure [3, 2]. Consider the following collection of sets

$$\mathcal{D}_n = \left\{ \frac{k}{2^n} \mid 0 \leq k \leq 2^n \right\}$$

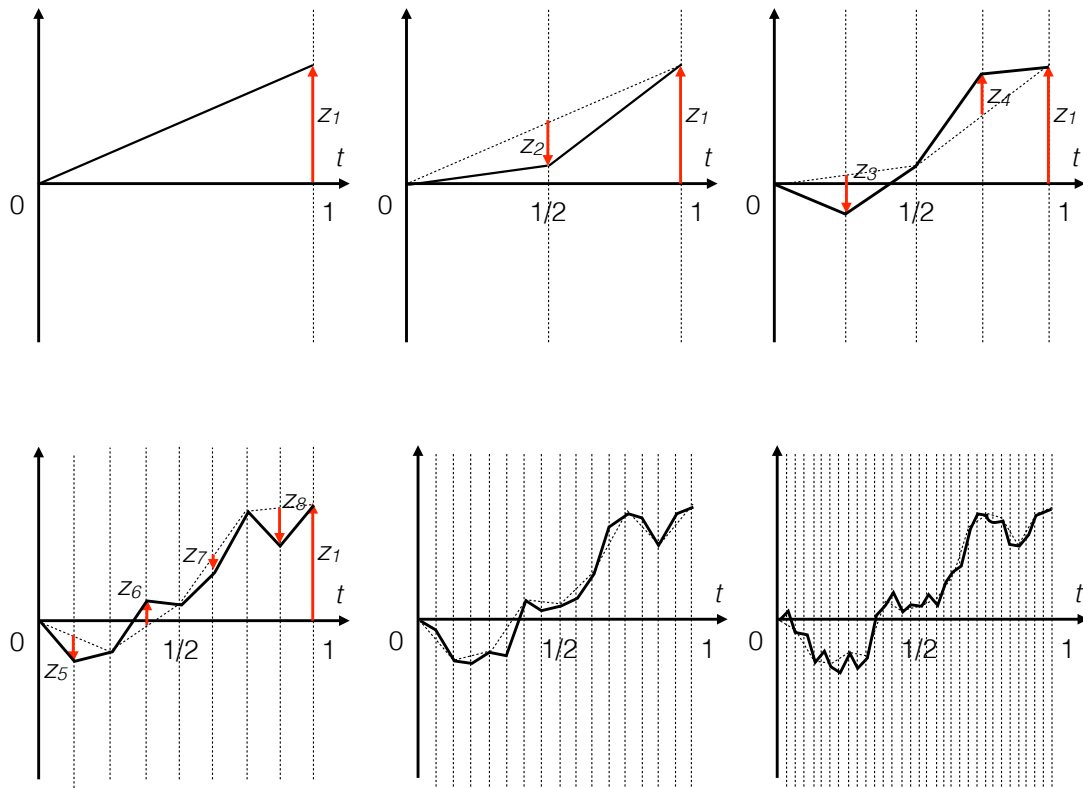


FIGURE 1. First six steps of construction a Brownian motion from Brownian random walks by a refinement procedure. Here z_1, \dots, z_8 are independent Gaussian random variables such that $z_1 \in N(0, 1)$, $z_2 \in N(0, 1/4)$, $z_3, z_4 \in N(0, \frac{1}{8})$, $z_5, z_6, z_7, z_8 \in N(0, \frac{1}{16})$.

of dyadic points. At each refinement step, we will define a Gaussian random walk satisfying the axioms of the Brownian Motion at the dyadic points (see Fig. 1).

Let $\{z_0, z_{n,j}\}$ where $n = 1, 2, 3, \dots$, $j = 1, 2, \dots, 2^{n-1}$, be a collection of independent Gaussian random variables with mean 0 and variance 1. Recall that

$$\text{if } z \in N(0, 1) \text{ then } \frac{z}{2^\alpha} \in N(0, 2^{-2\alpha}).$$

We start by constructing a Gaussian random walk on \mathcal{D}_0 by setting

$$B_0(0) = 0 \text{ and } B_0(1) = z_0.$$

Then we refine it to a Gaussian random walk B_1 on D_1 by setting

$$B_1(\mathcal{D}_0) = B_0(\mathcal{D}_0), \quad B_1(1/2) = \frac{B_0(1) + B_0(0)}{2} + \frac{z_{1,1}}{2}.$$

Note that

$$B_1(1/2) = \frac{z_0}{2} + \frac{z_{1,1}}{2}.$$

Hence

$$\text{Var}(B_1(1/2)) = \frac{1}{4} + \frac{1}{4} = \frac{1}{2}$$

as desired. Let us show that the increments of B_1 restricted to the dyadic set $\mathcal{D}_1 = \{0, 1/2, 1\}$ are independent, i.e., that $B_1(1/2) - B_1(0)$ and $B_1(1) - B_1(1/2)$ are independent. Indeed, a linear combination of independent Gaussian random variables $x_1 \sim \mathcal{N}(m_1, \sigma_1^2)$ and $x_2 \sim \mathcal{N}(m_2, \sigma_2^2)$ is Gaussian: $ax_1 + bx_2 \sim \mathcal{N}(am_1 + bm_2, a^2\sigma_1^2 + b^2\sigma_2^2)$ (check it!). Then we calculate:

$$\begin{aligned} B_1(1/2) - B_1(0) &= \frac{z_0}{2} + \frac{z_{1,1}}{2} \sim \mathcal{N}(0, 1/2), \\ B_1(1) - B_1(1/2) &= z_0 - \frac{z_0}{2} - \frac{z_{1,1}}{2} = \frac{z_0}{2} - \frac{z_{1,1}}{2} \sim \mathcal{N}(0, 1/2), \\ \mathbb{E} \left[\left(\frac{z_0}{2} + \frac{z_{1,1}}{2} \right) \left(\frac{z_0}{2} - \frac{z_{1,1}}{2} \right) \right] &= \mathbb{E} \left[\left(\frac{z_0}{2} \right)^2 - \left(\frac{z_{1,1}}{2} \right)^2 \right] = \frac{1}{2} - \frac{1}{2} = 0. \end{aligned}$$

Since uncorrelated Gaussian random variables are independent, we conclude that $B_1(1/2) - B_1(0)$ and $B_1(1) - B_1(1/2)$ are independent.

Next, we refine B_1 to a Gaussian random walk B_2 on \mathcal{D}_2 by setting $B_2(\mathcal{D}_1) = B_1(\mathcal{D}_1)$ and

$$\begin{aligned} B_2(1/4) &= \frac{1}{2} [B_1(1/2) + B_1(0)] + \frac{z_{2,1}}{2\sqrt{2}}, \\ B_2(3/4) &= \frac{1}{2} [B_1(1) + B_1(1/2)] + \frac{z_{2,2}}{2\sqrt{2}}. \end{aligned}$$

Then,

$$\begin{aligned} \text{Var}(B_2(1/4)) &= \text{Var} \left(\frac{1}{2} [B_1(1/2) + B_1(0)] + \frac{z_{2,1}}{2\sqrt{2}} \right) = \frac{1}{4} \cdot \frac{1}{2} + \frac{1}{8} = \frac{1}{4}, \\ \text{Var}(B_2(3/4)) &= \text{Var} \left(\frac{1}{2} [B_1(1) + B_1(1/2)] + \frac{z_{2,2}}{2\sqrt{2}} \right) \\ &= \text{Var} \left(\frac{1}{2} [B_1(1) - B_1(1/2) + 2B_1(1/2)] + \frac{z_{2,2}}{2\sqrt{2}} \right) \\ &= \frac{1}{4} \left[\frac{1}{2} + 4 \cdot \frac{1}{2} \right] + \frac{1}{8} = \frac{3}{4}. \end{aligned}$$

as desired. Show that the increments of B_2 restricted to the dyadic set $\mathcal{D}_2 = \{0, 1/4, 1/2, 3/4, 1\}$ are independent. First show that $B_2(2^{-2}(k+1)) - B_2(2^{-2}k)$ and $B_2(2^{-2}(l+1)) - B_2(2^{-2}l)$

for $0 \leq k < l \leq 3$ are independent. If k is even, i.e., $k = 2p$, we have:

$$\begin{aligned} B_2\left(\frac{p}{2} + \frac{1}{4}\right) - B_2\left(\frac{p}{2}\right) &= \frac{1}{2} \left[B_1\left(\frac{p}{2}\right) + B_1\left(\frac{p+1}{2}\right) \right] + \frac{z_{2,p+1}}{2\sqrt{2}} - B_2\left(\frac{p}{2}\right) = \\ \frac{1}{2} \left[B_1\left(\frac{p+1}{2}\right) - B_1\left(\frac{p}{2}\right) \right] + \frac{z_{2,p+1}}{2\sqrt{2}} &\sim \mathcal{N}\left(0, \frac{1}{4} \cdot \frac{1}{2} + \frac{1}{8}\right) = \mathcal{N}\left(0, \frac{1}{4}\right). \end{aligned}$$

If k is odd, i.e., $k = 2p + 1$, we have:

$$\begin{aligned} B_2\left(\frac{p+1}{2}\right) - B_2\left(\frac{p}{2} + \frac{1}{4}\right) &= B_2\left(\frac{p+1}{2}\right) - \frac{1}{2} \left[B_1\left(\frac{p}{2}\right) + B_1\left(\frac{p+1}{2}\right) \right] - \frac{z_{2,p+1}}{2\sqrt{2}} = \\ \frac{1}{2} \left[B_1\left(\frac{p+1}{2}\right) - B_1\left(\frac{p}{2}\right) \right] - \frac{z_{2,p+1}}{2\sqrt{2}} &\sim \mathcal{N}\left(0, \frac{1}{4} \cdot \frac{1}{2} + \frac{1}{8}\right) = \mathcal{N}\left(0, \frac{1}{4}\right). \end{aligned}$$

Note that both $\frac{1}{2} \left[B_1\left(\frac{p+1}{2}\right) - B_1\left(\frac{p}{2}\right) \right]$ and $\frac{z_{2,p+1}}{2\sqrt{2}}$ are independent Gaussian random variables with mean 0 and variance 1/8. If $k = 2p$ and $l = 2p + 1$, using an argument similar to the one used for showing that that $B_1(1/2) - B_1(0)$ and $B(1) - B(1/2)$ are independent, we show that

$$B_2\left(\frac{p}{2} + \frac{1}{4}\right) - B_2\left(\frac{p}{2}\right) \quad \text{and} \quad B_2\left(\frac{p+1}{2}\right) - B_2\left(\frac{p}{2} + \frac{1}{4}\right)$$

are independent. If k and l are such that $\text{floor}(k/2) < \text{floor}(l/2)$, then the argument above implies that $B_2(2^{-2}(k+1)) - B_2(2^{-2}k) = x_1 + x_2$ and $B_2(2^{-2}(l+1)) - B_2(2^{-2}l) = y_1 + y_2$, where x_1, x_2, y_1, y_2 are independent Gaussian random variables with mean 0 and variance 1/8. Therefore, $B_2(2^{-2}(k+1)) - B_2(2^{-2}k)$ and $B_2(2^{-2}(l+1)) - B_2(2^{-2}l)$ for $0 \leq k < l \leq 3$ are independent. Finally, the increments over non-overlapping (no common interior points) of B_2 restricted to \mathcal{D}_2 are independent as they are sums of mutually independent increments.

Continuing inductively, we define (i) : $B_n(\mathcal{D}_{n-1}) = B_{n-1}(\mathcal{D}_{n-1})$; (ii) for $\mathcal{D}_n \setminus \mathcal{D}_{n-1}$

$$B_n\left(\frac{k}{2^{n-1}} + \frac{1}{2^n}\right) = \frac{1}{2} \left(B_{n-1}\left(\frac{k}{2^{n-1}}\right) + B_{n-1}\left(\frac{k+1}{2^{n-1}}\right) \right) + \frac{z_{n,k+1}}{2^{(n+1)/2}},$$

where $z_{n,k+1} \sim \mathcal{N}(0, 1)$. It is shown in [3] that if one continues this refinement procedure up to infinity, the resulting process satisfies the definition of Brownian motion. The first six steps on this procedure are illustrated in Fig. 1.

8.4. Construction of the Brownian motion by a random walk. This construction is left as an exercise.

Exercise Consider the mesh

$$\{t_j \mid t_j = jh, \quad h = \frac{1}{N}, \quad 0 \leq j \leq N\}.$$

Let $\{z_j\}_{j=1}^N$ be independent Gaussian random variables with mean 0 and variance 1. Consider the Gaussian random walk $B_h(t)$ defined by

$$\begin{aligned} B_h(0) &= 0, \\ B_h(t_j) &= B_h(t_{j-1}) + z_j\sqrt{h}, \quad j = 1, \dots, N, \\ B_h(t) &= \frac{1}{h} [B_h(t_{j-1})(t_j - t) + B_h(t_j)(t - t_{j-1})], \quad t_{j-1} < t < t_j, \quad j = 1, \dots, N. \end{aligned}$$

Prove that this random walk satisfies axioms (1)-(4) of Brownian motion at the points t_j , $j = 0, 1, \dots, N$.

Exercise An even simpler construction the Brownian motion can be done as follows. Consider a random walk on a mesh in the (x, t) -space, $t \geq 0$, $x \in \mathbb{R}$ in which the time-step k and the space step h are related to the time step k via $k = h^2$. Start at the origin at time 0. At any discrete moment of time $t = k, 2k, 3k, \dots$, take a step to left or to the right with probability $1/2$. Let k and h tend to zero in such a manner that the relationship $k = h^2$ is maintained. Apply the Central Limit Theorem and obtain the Brownian motion.

The last exercise will be solved in Section 9 ahead.

8.5. Elementary properties of Brownian motion.

- **The covariance function** of the Brownian motion is

$$(23) \quad E[w(t_1)w(t_2)] = \min\{t_1, t_2\} \equiv t_1 \wedge t_2.$$

Indeed, suppose $t_2 > t_1$. Then

$$\begin{aligned} E[w(t_1)w(t_2)] &= E[w(t_1)w(t_1) + w(t_1)(w(t_2) - w(t_1))] \\ &= E[w(t_1)w(t_1)] + E[w(t_1)(w(t_2) - w(t_1))] = t_1. \end{aligned}$$

- **Nowhere differentiability with probability 1** Consider the random

$$\frac{w(\omega, t + \Delta t) - w(\omega, t)}{\Delta t}.$$

It is Gaussian with mean 0 and variance $(\Delta t)^{-1}$, which tends to infinity as $\Delta t \rightarrow 0$. Hence $w(\omega, t)$ is differentiable nowhere with probability 1.

- **White noise** Despite the regular derivative of a Brownian motion does not exist, one can consider its derivative in the sense of distributions. This derivative $\eta(\omega, t)$ is called *white noise* and is defined by the property

$$\int_{t_1}^{t_2} \eta(\omega, t) dt = w(\omega, t_2) - w(\omega, t_1).$$

- **Scaling and Symmetry** If $w(t)$ is a Brownian motion then so are the processes defined by

$$\begin{aligned} x(t) &:= \frac{1}{\sqrt{c}}w(ct) \quad \text{for any positive constant } c, \\ y(t) &= -w(t). \end{aligned}$$

- **Time inversion** Let $w(t)$ be a Brownian motion. Then so is the process defined by

$$(24) \quad x(t) = \begin{cases} tw(1/t), & 0 < t < \infty, \\ 0, & t = 0. \end{cases}$$

- **Invariance under rotations and reflections (orthogonal transformations)** Let $w(t)$ be a d -dimensional Brownian motion, and T be a $d \times d$ orthogonal matrix (i.e., $T^\top = T^{-1}$). Then the process

$$x(t) = Tw(t)$$

is also a d -dimensional Brownian motion.

9. BROWNIAN MOTION AND HEAT EQUATION

In this section, we will establish a relationship between Brownian motion and the heat equation as it is done in [1]. Consider the initial-value problem for the heat equation

$$(25) \quad u_t = \frac{1}{2}u_{xx}, \quad x \in \mathbb{R}, \quad t > 0, \quad u(x, 0) = \phi(x).$$

We assume that the function $\phi(x)$ decays fast enough as $|x| \rightarrow \infty$ so that the total amount of heat is finite, i.e.,

$$\int_{-\infty}^{\infty} |\phi(x)| dx < \infty.$$

The fastest way to solve it is the one via the use of the Fourier Transform. The Fourier transform is a linear operator mapping a function $f(x)$ $x \in \mathbb{R}$, to a function $\hat{f}(k)$, $k \in \mathbb{R}$, called the Fourier Transform (FT) of f defined by

$$(26) \quad \mathcal{F}[f](k) = \hat{f}(k) := \frac{1}{\sqrt{2\pi}} \int_{-\infty}^{\infty} e^{-ikx} f(x) dx.$$

The Inverse Fourier Transform (IFT) maps $\hat{f}(k)$ back to $f(x)$:

$$(27) \quad \mathcal{F}^{-1}[\hat{f}](x) := \frac{1}{\sqrt{2\pi}} \int_{-\infty}^{\infty} e^{ikx} \hat{f}(k) dx = f(x).$$

The proof of this fact relies on Fubini's theorem and the Dominated Convergence Theorem studied in graduate analysis courses. Let us apply the Fourier Transform to $u(x, t)$ with respect to the space variable x . Then we have

$$(28) \quad u(x, t) = \frac{1}{\sqrt{2\pi}} \int_{-\infty}^{\infty} e^{ikx} \hat{u}(k, t) dx.$$

It is easy to verify by differentiating both parts of Eq. (28) with respect to x and t that

$$\begin{aligned} u_x(x, t) &= \frac{1}{\sqrt{2\pi}} \int_{-\infty}^{\infty} ik e^{ikx} \hat{u}(k, t) dx = \mathcal{F}^{-1}[ik\hat{u}] \\ u_{xx}(x, t) &= \frac{1}{\sqrt{2\pi}} \int_{-\infty}^{\infty} (ik)^2 e^{ikx} \hat{u}(k, t) dx = \mathcal{F}^{-1}[-k^2\hat{u}] \\ u_t(x, t) &= \frac{1}{\sqrt{2\pi}} \int_{-\infty}^{\infty} e^{ikx} \frac{\partial}{\partial t} \hat{u}(k, t) dx = \mathcal{F}^{-1}[\hat{u}_t] \end{aligned}$$

Plugging these identities into Eq. (25) we obtain the following initial value problem for $\hat{u}(k, t)$:

$$(29) \quad \hat{u}_t(k, t) = -\frac{k^2}{2} \hat{u}(k, t), \quad k \in \mathbb{R}, \quad t > 0, \quad \hat{u}(k, 0) = \hat{\phi}(k).$$

Therefore, we have placed a PDE with a collection of ODEs indexed by $k \in \mathbb{R}$! The solution of Eq. (25) is

$$\hat{u}(k, t) = e^{-\frac{1}{2}k^2t} \hat{\phi}(k).$$

To return to $u(x, t)$ we apply the inverse Fourier Transform:

$$\begin{aligned} u(x, t) &= \frac{1}{\sqrt{2\pi}} \int_{-\infty}^{\infty} e^{ikx} e^{-\frac{1}{2}k^2t} \hat{\phi}(k) dk \\ &= \frac{1}{\sqrt{2\pi}} \int_{-\infty}^{\infty} e^{ikx} e^{-\frac{1}{2}k^2t} dk \frac{1}{\sqrt{2\pi}} \int_{-\infty}^{\infty} e^{-iky} \phi(y) dy \\ &= \int_{-\infty}^{\infty} dy \frac{\phi(y)}{\sqrt{2\pi t}} \exp\left[-\frac{(x-y)^2}{2t}\right] \underbrace{\frac{\sqrt{t}}{\sqrt{2\pi}} \int_{-\infty}^{\infty} dk \exp\left[-\frac{1}{2}\left(k\sqrt{t} - i\frac{x-y}{\sqrt{t}}\right)^2\right]}_{=\sqrt{2\pi/t}} \end{aligned}$$

Thus, the solution of Eq. (25) is

$$(30) \quad u(x, t) = \int_{-\infty}^{\infty} dy \frac{\phi(y)}{\sqrt{2\pi t}} \exp\left[-\frac{(x-y)^2}{2t}\right].$$

To make its probabilistic interpretation more apparent, we define introduce $z = y - x$ and recast it as

$$(31) \quad u(x, t) = \int_{-\infty}^{\infty} \frac{e^{-\frac{z^2}{2t}}}{\sqrt{2\pi t}} \phi(x+z) dz.$$

The function

$$f(z) = \frac{e^{-\frac{z^2}{2t}}}{\sqrt{2\pi t}}$$

is the probability density of Gaussian random variable with mean 0 and variance t . Therefore, the solution of the heat equation at point x at time t is the expected value of the function $\phi(x + \eta(\omega))$ where $\eta \in N(0, t)$:

$$(32) \quad u(x, t) = E[\phi(x + \eta(\omega))] \quad \text{where } \eta \in N(0, t).$$

Recall that a Brownian motion $w(\omega, t)$ is a Gaussian random variable for any fixed time t . Therefore, Eq. (32) can be written as

$$(33) \quad u(x, t) = E[\phi(x + w(\omega, t))].$$

Eq. (33) suggests the following way to find the solution $u(x, t)$ at a given point (x, t) . Start Brownian motions going backward in time from the point (x, t) and run them for time t . Recall that the time-reversal of a Brownian motion is also a Brownian motion. At time t , they intersect the x -axis at the points $x + w(\omega, t)$. Find the values of ϕ at these points and average them over all Brownian motions.

10. NUMERICAL SOLUTION OF THE HEAT EQUATION AND ITS SOLUTION BY RANDOM WALK

In this section, we show that an approximate numerical solution $u(x_i, t_n)$ to the heat equation can be interpreted as the expected value of $\phi(x_i + \sum_{l=1}^n \eta_l)$ where η_l are independent Bernoulli random variables and the mesh steps in x and t are in a certain ratio [1]. This is analogous to the fact the exact analytical solution $u(x, t)$ of the heat equation (25) can be interpreted as the expected value of $\phi(x + w(\omega, t))$ as we have shown in the previous section.

Consider a rectangular mesh with step h in space and step k in time:

$$\{(x_i, t_n) \mid x_i = ih, t_n = nk\}.$$

We will denote by u_i^n the value of the exact solution u at the point (x_i, t_n) , i.e.,

$$u_i^n := u(x_i, t_n).$$

The derivatives u_t and u_{xx} can be approximated by finite differences below. Using Taylor expansions at (x_i, t_n) we get

$$(34) \quad u_t(x_i, t_n) = \frac{u_i^{n+1} - u_i^n}{k} + \frac{k}{2} u_{tt}(x_i, t^n + \alpha k), \quad \alpha \in (0, 1),$$

$$(35) \quad u_{xx}(x_i, t_n) = \frac{u_{i+1}^n - 2u_i^n + u_{i-1}^n}{h^2} + \frac{h^2}{12} u_{xxx}(x_i + \gamma h, t^n), \quad \gamma \in (-1, 1).$$

Using these finite differences, we obtain a discrete approximation to initial value problem for the heat equation:

$$(36) \quad \frac{v_i^{n+1} - v_i^n}{k} = \frac{1}{2} \frac{v_{i+1}^n - 2v_i^n + v_{i-1}^n}{h^2}, \quad v_i^0 = \phi(ih).$$

We distinguish the exact solution u of Eq. (25) and the solution v to its discretized version, Eq. (36). Starting from $n = 0$, we can solve it by the recurrence formula obtained from

Eq. (36):

$$v_i^{n+1} = v_i^n + \frac{k}{2h^2}(v_{i+1}^n - 2v_i^n + v_{i-1}^n).$$

Denote the quantity $\frac{k}{2h^2}$ by λ and rewrite the recurrence relationship as

$$(37) \quad v_i^{n+1} = (1 - 2\lambda)v_i^n + \lambda v_{i+1}^n + \lambda v_{i-1}^n.$$

Now we will find out how large is the approximation error when we replace u with v . First, we define the *local truncation error* τ_i^n , i.e., the error committed in one step starting from the exact values. In numerical PDEs, it is defined by

$$\tau_i^n := \frac{u_i^{n+1} - u_i^n}{k} - \frac{u_{i+1}^n - 2u_i^n + u_{i-1}^n}{h^2}.$$

It follows from Eqs. (34) and (35) that

$$\tau_i^n = \frac{k}{2}u_{tt}(x_i, t^n + \alpha k) - \frac{h^2}{12}u_{xxxx}(x_i + \gamma h, t^n) = O(k) + O(h^2).$$

As you can see, $\tau_i^n \rightarrow 0$ and $h, t \rightarrow 0$. This fact is called *consistency*. However, it might seem surprising at first, that the numerical solution produced by a consistent difference scheme does not necessarily converge to the exact solution at the mesh points as the mesh is refined. There is one more necessary ingredient to guarantee the convergence of the numerical solution to the exact one called *stability*. Stability means that the numerical solution is a Lipschitz function with respect to perturbations of the initial data and the difference equation. Note that the perturbations are ubiquitously present due to the local truncation error, approximations for real numbers with floating-point numbers, and the round-off error of computer arithmetics.

We will show that the numerical solution of the heat equation converges to the exact solution if $\lambda \leq 1/2$. We assume that the arithmetics is exact. We have:

$$(38) \quad u_i^{n+1} = (1 - 2\lambda)u_i^n + \lambda u_{i+1}^n + \lambda u_{i-1}^n + k\tau_i^n,$$

$$(39) \quad v_i^{n+1} = (1 - 2\lambda)v_i^n + \lambda v_{i+1}^n + \lambda v_{i-1}^n.$$

Subtract Eq. (39) from Eq. (38) and obtain the recurrence equation for the error $e_i^n := u_i^n - v_i^n$:

$$(40) \quad e_i^{n+1} = (1 - 2\lambda)e_i^n + \lambda e_{i+1}^n + \lambda e_{i-1}^n + k\tau_i^n.$$

Taking absolute values of both sides and assuming that $\lambda \leq 1/2$, i.e., $1 - 2\lambda \geq 0$, we get

$$(41) \quad |e_i^{n+1}| \leq (1 - 2\lambda)|e_i^n| + \lambda|e_{i+1}^n| + \lambda|e_{i-1}^n| + k|\tau_i^n|.$$

Define

$$E^n := \max_i |e_i^n|$$

and

$$\tau^n := \max_i |\tau_i^n|, \quad \tau := \max_{nk \leq t} |\tau^n|.$$

Then

$$E^{n+1} \leq E^n + k\tau^n \leq E^n + k\tau.$$

Therefore,

$$E^n \leq E^{n-1} + k\tau \leq E^{n-2} + 2k\tau \leq \dots \leq E^0 + nk\tau.$$

Assuming that we are starting with the exact initial condition, i.e., $E^0 = 0$. Then

$$E^n \leq nk\tau \equiv t\tau.$$

Since $\tau \rightarrow 0$ as $h, k \rightarrow 0$, we conclude that the solution to the discretized heat equation (36) converges to the solution of the heat equation (25) as we refine the mesh.

Note that we can pick any $\lambda \in (0, 1/2]$ in order to guarantee this convergence. Let us pick $\lambda = 1/2$ to make the recurrence equation Eq. (37) as simple as possible:

$$(42) \quad v_i^{n+1} = \frac{1}{2}v_{i+1}^n + \frac{1}{2}v_{i-1}^n.$$

Hence

$$\begin{aligned} v_i^n &= \frac{1}{2}v_{i+1}^{n-1} + \frac{1}{2}v_{i-1}^{n-1} \\ &= \frac{1}{4}v_{i+2}^{n-2} + \frac{1}{2}v_i^{n-2} + \frac{1}{4}v_{i-2}^{n-2} \\ &= \frac{1}{8}v_{i+3}^{n-3} + \frac{3}{8}v_{i+1}^{n-3} + \frac{3}{8}v_{i-1}^{n-3} + \frac{1}{8}v_{i-3}^{n-3} = \dots \\ &= \sum_{j=0}^n \frac{1}{2^n} \binom{n}{j} \phi((i-n+2j)h). \end{aligned}$$

Thus, the solution of the discretized heat equation Eq. (36) is

$$(43) \quad v_i^n = \sum_{j=0}^n \frac{1}{2^n} \binom{n}{j} \phi((i-n+2j)h).$$

This solution can be interpreted as follows. Consider the random walks on the mesh with step h in space and step $k = h^2$ in time starting at $(x_i = ih, t_n = nk)$ and shifting one mesh step down in time together with one mesh step left or right with probability $1/2$ each iteration. Record the values of the initial function ϕ at the points where these walks reach the x -axis (the level 0). Average these values over all walks to obtain the solution v_i^n .

The horizontal displacements of these random walks can be interpreted as sums of n independent Bernoulli random variables

$$\eta_l = \begin{cases} h, & P = 1/2 \\ -h, & P = 1/2 \end{cases}.$$

These variables have mean 0 and variance h^2 . Hence

$$E \left[\sum_{l=1}^n \eta_l \right] = 0, \quad \text{Var} \left(\sum_{l=1}^n \eta_l \right) = nh^2.$$

Let us fix t . As $k \rightarrow 0$, $n \rightarrow \infty$. Hence, by the Central Limit Theorem,

$$\sum_{l=1}^n \eta_l \rightarrow N(0, nh^2) \text{ in distribution.}$$

Recall that $k = h^2$ and hence $nh^2 = nk = t$. Set $y := 2j - n$. Then

$$P\left(\sum_{l=1}^n \eta_l = (-n + 2j)h\right) \approx \frac{e^{-y^2/2t}}{\sqrt{2\pi t}} h.$$

Therefore,

$$v_i^n = \sum_{j=0}^n \frac{1}{2^n} \binom{n}{j} \phi((i - n + 2j)h) \rightarrow \int_{-\infty}^{\infty} \frac{e^{-y^2/2t}}{\sqrt{2\pi t}} \phi(x + y) dy$$

as $n \rightarrow \infty$, $h \rightarrow 0$, $k = h^2$, $t = nk$ is fixed.

10.1. A brief introduction into the Wiener measure. Here we follow the discussion found in [1, 4]. Recall that we have derived that the solution of the heat equation Eq. (25) is given by

$$u(x, t) = E[\phi(x + w(\omega, t))],$$

where $w(\omega, t)$ is a Brownian motion. This formula says that we attach a number $\phi(x + w(\omega, t))$ to each Brownian motion and find an expectation for these numbers. Whenever we speak about an expectation, we imply some probability measure with respect to which this expectation is taken.

When we need to average some function of Brownian motion, the measure to be taken is the Wiener measure. Therefore, we need to specify the measurable space where the Wiener measure will be defined, i.e., the set of outcomes Ω and the σ -algebra on Ω .

The set of outcomes will be the set of continuous functions $y(t)$ satisfying $y(0) = 0$. The σ -algebra will contain all cylinder sets of the form

$$(44) \quad C = \{f(t) \text{ is continuous} \mid a \leq f(s) < b\},$$

where a , b , and s are parameters defining C .

The Wiener measure of each cylinder set (44) is defined so that it is equal to the probability that a Brownian motion $w(\omega, t)$ passes through the window $[a, b]$ at time s :

$$P(C) = \int_a^b \frac{e^{-x^2/2s}}{\sqrt{2\pi s}} dx.$$

The Wiener measure of the intersection of two cylinders

$$C_1 = \{f(t) \text{ is continuous} \mid a_1 \leq f(s_1) < b_1\} \text{ and } C_2 = \{f(t) \text{ is continuous} \mid a_2 \leq f(s_2) < b_2\}$$

is defined so that it is equal to the probability that a Brownian motion $w(\omega, t)$ passes through both windows:

$$a_1 \leq w(\omega, s_1) < b_1, \text{ and } a_2 \leq w(\omega, s_2) < b_2.$$

Assume $s_1 < s_2$. Taking into account that the increments of a Brownian motion are independent Gaussian random variables, we calculate

$$P(C_1 \cap C_2) = \int_{a_1}^{b_1} \frac{e^{-x^2/2s_1}}{\sqrt{2\pi s_1}} dx \int_{a_2}^{b_2} \frac{e^{-(y-x)^2/2(s_2-s_1)}}{\sqrt{2\pi(s_2-s_1)}} dy.$$

The notation for the Wiener measure is dW . Thus, the solution of the heat equation (25) can be written as

$$u(x, t) = \int \phi(x + w(\omega, t)) dW.$$

Example 20 Compute $\int F dW$ where $F(w) = \int_0^1 w^4(\omega, s) ds$.

$$\begin{aligned} \int F dW &= \int dW \int_0^1 w^4(\omega, s) ds = \int_0^1 ds \int dW w^4(\omega, s) \\ &= \int_0^1 ds \int_{-\infty}^{\infty} x^4 \frac{e^{-x^2/2s}}{\sqrt{2\pi s}} dx = \int_0^1 3s^2 ds = 1. \end{aligned}$$

10.2. Markov property of Brownian motion.

Definition 9. A stochastic process $\zeta(t)$ on $[0, T]$ is called a Markov process if for any sequences $0 \leq t_0 < \dots < t_n \leq T$ and x_0, x_1, \dots, x_n , its transition probability distribution function has the property

$$\mathbb{P}(\zeta(t_n) < x_n \mid \zeta(t_{n-1}) < x_{n-1}, \dots, \zeta(t_0) < x_0) = \mathbb{P}(\zeta(t_n) < x_n \mid \zeta(t_{n-1}) < x_{n-1}).$$

The transition probability density function, defined by

$$p(x_n, t_n \mid x_{n-1}, t_{n-1}; \dots; x_0, t_0) := \frac{\partial}{\partial x_n} \mathbb{P}(\zeta(t_n) < x_n \mid \zeta(t_{n-1}) < x_{n-1}, \dots, \zeta(t_0) < x_0)$$

then satisfies

$$(45) \quad p(x_n, t_n \mid x_{n-1}, t_{n-1}; \dots; x_0, t_0) = p(x_n, t_n \mid x_{n-1}, t_{n-1}).$$

For any three times $t > \tau > s$ and any three points x, y, z we can write the identities

$$p(y, t; z, \tau \mid x, s) = p(y, t \mid z, \tau; x, s)p(z, \tau \mid x, s) = p(y, t \mid z, \tau)p(z, \tau \mid x, s).$$

The last equality is a consequence of the Markov property. This identity implies the Chapman-Kolmogorov equation:

$$(46) \quad p(y, t \mid x, s) = \int p(y, t; z, \tau \mid x, s) dz = \int p(y, t \mid z, \tau)p(z, \tau \mid x, s) dz.$$

Theorem 6. The Brownian motion is a Markov process.

Proof. Take any sequences $0 = t_0 < \dots < t_n \leq T$ and $x_0 = 0, x_1, \dots, x_n$ and consider the joint pdf of the vector

$$w = (w(t_1), w(t_2), \dots, w(t_n)).$$

It is given by

$$(47) \quad p(x_1, t_1; x_2, t_2; \dots; x_n, t_n) = \prod_{k=1}^n \left[\frac{\exp\left\{-\frac{(x_k - x_{k-1})^2}{2(t_k - t_{k-1})}\right\}}{\sqrt{2\pi(t_k - t_{k-1})}} \right].$$

Recall that

$$p(x_n, t_n \mid x_{n-1}, t_{n-1}; \dots; x_1, t_1) = \frac{p(x_1, t_1; x_2, t_2; \dots; x_n, t_n)}{p(x_1, t_1; x_2, t_2; \dots; x_{n-1}, t_{n-1})}.$$

Using Eq. (47) we get

$$\begin{aligned} p(x_n, t_n \mid x_{n-1}, t_{n-1}; \dots; x_1, t_1) &= \frac{\prod_{k=1}^n \left[\frac{\exp\left\{-\frac{(x_k - x_{k-1})^2}{2(t_k - t_{k-1})}\right\}}{\sqrt{2\pi(t_k - t_{k-1})}} \right]}{\prod_{k=1}^{n-1} \left[\frac{\exp\left\{-\frac{(x_k - x_{k-1})^2}{2(t_k - t_{k-1})}\right\}}{\sqrt{2\pi(t_k - t_{k-1})}} \right]} = \frac{\exp\left\{-\frac{(x_n - x_{n-1})^2}{2(t_n - t_{n-1})}\right\}}{\sqrt{2\pi(t_n - t_{n-1})}} \\ &= p(x_n, t_n \mid x_{n-1}, t_{n-1}). \end{aligned}$$

Hence, the Brownian motion is a Markov process. □

11. AN INTRODUCTION TO SDES.

Here we follow the discussion found in [1, 4]. Consider a stochastic process $x(\omega, t) \equiv x(t)$ obeying the following evolution law:

$$(48) \quad dx(t) = b(x(t), t)dt + \sigma(x(t), t)dw,$$

where w is the standard Brownian motion and the functions b and σ are smooth.

This evolution law is called a *stochastic differential equation* (SDE). If $\sigma(x(t), t) = 0$, Eq. (48) becomes an ordinary differential equation (ODE)

$$(49) \quad \frac{dx}{dt} = b(x, t).$$

Suppose $x(0) = x_0$. Eq. (49) is equivalent to the following integral equation

$$(50) \quad x(t) = x_0 + \int_0^t b(x(s), s) ds.$$

A solution of an ODE is a function satisfying the ODE. If the ODE is complemented with an initial condition, then the solution to the corresponding initial-value problem is a function satisfying the given initial condition.

Now we will discuss the meaning of SDE (48). Similarly to Eq. (51) we can write

$$(51) \quad x(t) = x_0 + \int_0^t b(x(s), s) ds + \int_0^t \sigma(x(s), s) dw(s)$$

First assume that the function σ is independent of x , i.e., $\sigma(x, t) \equiv \sigma(t)$. We partition the interval $[0, t]$ into $t_0 = 0 < t_1 < t_2 < \dots < t_n = t$ and denote the fineness of the partition by Δ :

$$\Delta := \max_{1 \leq i \leq n} |t_i - t_{i-1}|.$$

Then we define

$$\int_0^t \sigma(s) dw(s) = \lim_{\substack{n \rightarrow \infty \\ \Delta \rightarrow 0}} \sum_{i=0}^{n-1} \sigma_i (w(t_{i+1}) - w(t_i))$$

where σ_i is chosen so that it approximates $\sigma(t)$ on the subinterval $[t_i, t_{i+1}]$.

The case where σ depends on x is much more difficult. We can proceed as before and write

$$\int_0^t \sigma(x(s), s) dw(s) = \lim_{\substack{n \rightarrow \infty \\ \tau \rightarrow 0}} \sum_{i=1}^n \sigma_i (w(t_{i+1}) - w(t_i)),$$

however, the value of this limit will depend on how we choose σ_i approximating $\sigma(x(t), t)$ on the interval $[t_i, t_{i+1}]$. There are two common choices. We will give their general definition. Let $f(w(t), t)$ be a smooth function depending on time and a Brownian Motion $w(t)$. In particular, $f(w(t), t)$ can be chosen to coincide with $\sigma(x(t), t)$ where $x(t)$ is a solution of SDE (48), i.e., a stochastic process depending on $w(t)$.

- **The Ito stochastic integral** is defined by the choice $f_i = f(t_i)$, i.e., f is evaluated at the left end of each subinterval:

$$\int_0^t f(w(s), s) dw = \lim_{\substack{n \rightarrow \infty \\ \Delta \rightarrow 0}} \sum_{i=0}^{n-1} f(w(t_i), t_i) (w(t_{i+1}) - w(t_i)).$$

- **The Stratonovich stochastic integral** is defined by the choice $f_i = f(t_{i+1/2})$, where $t_{i+1/2} \equiv \frac{1}{2}(t_i + t_{i+1})$. i.e., f is evaluated at the midpoint of each subinterval. The Stratonovich stochastic integral is marked by \circ :

$$\int_0^t f(w(s), s) \circ dw = \lim_{\substack{n \rightarrow \infty \\ \Delta \rightarrow 0}} \sum_{i=0}^{n-1} f(w(t_{i+1/2}), t_{i+1/2}) (w(t_{i+1}) - w(t_i)).$$

The example below demonstrates that Ito and Stratonovich stochastic integrals are different.

Example 21 Calculate $\int_a^b w dw$ (the Ito stochastic integral) and $\int_a^b w \circ dw$ (the Stratonovich stochastic integral). Let the partition be uniform, i.e.,

$$\Delta t = \frac{b-a}{n}.$$

We start with the Ito stochastic integral.

$$\begin{aligned} \int_a^b w dw &= \lim_{\Delta \rightarrow 0} \sum_i w(t_i) [w(t_{i+1}) - w(t_i)] \\ &= \frac{1}{2} \lim_{\Delta \rightarrow 0} \sum_i [w^2(t_{i+1}) - w^2(t_i) - (w(t_{i+1}) - w(t_i))^2] \end{aligned}$$

Note that

$$\begin{aligned} \sum_i (w^2(t_{i+1}) - w^2(t_i)) &= w^2(t_n) - w^2(t_{n-1}) + w^2(t_{n-1}) - \dots + w^2(t_1) - w^2(t_0) \\ &= w^2(b) - w^2(a). \end{aligned}$$

We also compute

$$E \left[\sum_i (w(t_{i+1}) - w(t_i))^2 \right] = \sum_i (t_{i+1} - t_i) = b - a.$$

$$\begin{aligned} \text{Var} \left(\sum_i (w(t_{i+1}) - w(t_i))^2 \right) &= \sum_i [E[(w(t_{i+1}) - w(t_i))^4] - (E[(w(t_{i+1}) - w(t_i))^2])^2] \\ &\leq \sum_i E [(w(t_{i+1}) - w(t_i))^4] = \frac{n}{\sqrt{2\pi\Delta t}} \int_{-\infty}^{\infty} x^4 e^{-x^2/2\Delta t} dx \\ &= 3n(\Delta t)^2 = \frac{3n(b-a)^2}{n^2} = \frac{3(b-a)^2}{n} \rightarrow 0 \text{ as } n \rightarrow \infty. \end{aligned}$$

Here we have used the fact that the fourth central moment of $\mathcal{N}(\mu, \sigma^2)$ is $3\sigma^4$. Hence

$$\int_a^b w dw = \frac{w^2(b) - w^2(a)}{2} - \frac{b-a}{2}.$$

The expected value of this integral is zero. Indeed, $w(t_i) [w(t_{i+1}) - w(t_i)]$ is a product of two independent Gaussian random variables with mean 0, hence $E [w(t_i) [w(t_{i+1}) - w(t_i)]] = 0$ for all i . Therefore,

$$E \left[\int_a^b w dw \right] = E \left[\lim_{\Delta \rightarrow 0} \sum_i w(t_i) [w(t_{i+1}) - w(t_i)] \right] = 0.$$

Now we will calculate the Stratonovich stochastic integral:

$$\begin{aligned}
 \int_a^b w \circ dw &= \lim_{\Delta \rightarrow 0} \sum_i w(t_{i+1/2})(w(t_{i+1}) - w(t_i)) \\
 &= \lim_{\Delta \rightarrow 0} \sum_i [w(t_{i+1/2}) - w(t_i) + w(t_i)][w(t_{i+1}) - w(t_{i+1/2}) + w(t_{i+1/2}) - w(t_i)] \\
 &= \lim_{\Delta \rightarrow 0} \sum_i (w(t_{i+1/2}) - w(t_i))^2 \\
 &+ \lim_{\Delta \rightarrow 0} \sum_i (w(t_{i+1/2}) - w(t_i))(w(t_{i+1}) - w(t_{i+1/2})) \\
 &+ \lim_{\Delta \rightarrow 0} \sum_i w(t_i) [w(t_{i+1}) - w(t_i)].
 \end{aligned}$$

The first limit can be evaluated by finding its mean $(b - a)/2$ and showing that its variance tends to zero as $n \rightarrow \infty$. The second limit is zero as it is the sum of products of independent Gaussian random variables with mean zero and vanishing variance as $n \rightarrow \infty$. The third limit is the Ito stochastic integral that we have just evaluated. Hence,

$$\int_a^b w \circ dw = \frac{b - a}{2} + \frac{w^2(b) - w^2(a)}{2} - \frac{b - a}{2} = \frac{w^2(b) - w^2(a)}{2}.$$

As you see,

$$\int_a^b w \circ dw \neq \int_a^b w dw.$$

11.1. Elementary properties of stochastic integral. Here we follow the discussions in [4, 5]. We will use a shorter notation denoting $f(w(t), t)$ by $f(t, \omega)$ where ω is the stochastic argument of the Brownian Motion.

We will consider stochastic process $f(t, \omega)$ on $0 \leq t \leq T$ satisfying the following conditions:

Condition (1): $f(t, \omega)$ is independent of the increments of the Brownian motion $w(t, \omega)$ in the future, i.e., $f(t, \omega)$ is independent of $w(t + s, \omega) - w(t, \omega)$ for all $s > 0$. Such processes are called *adapted to the Brownian filtration* \mathcal{F}_t .

Condition (2):

$$\int_0^T E[f^2(s, \omega)] ds < \infty.$$

Now we list some useful elementary properties. The first two properties are similar to those of the Riemann integral. The other ones are specific for the Ito integral. Let $f(t, \omega)$ and $g(t, \omega)$ be any functions satisfying conditions (1) and (2) above.

(1) Linearity:

$$\int_0^t (af(s, \omega) + bg(s, \omega))dw(s, \omega) = a \int_0^t f(s, \omega)dw(s, \omega) + b \int_0^t g(s, \omega)dw(s, \omega).$$

(2) Additivity. Let $0 < T_1 < T$. Then

$$\int_0^T f(s, \omega)dw(s, \omega) = \int_0^{T_1} f(s, \omega)dw(s, \omega) + \int_{T_1}^T f(s, \omega)dw(s, \omega).$$

(3) If f is a deterministic function, i.e., $f(s, \omega) \equiv f(s)$, then

$$\int_0^t f(s)dw(s, \omega) \sim N\left(0, \int_0^t f^2(s)ds\right).$$

(4) For any $0 \leq \tau \leq t \leq T$,

$$\begin{aligned} E \left[\int_0^t f(s, \omega) dw(s, \omega) \right] &= 0; \\ E \left[\int_0^t f(s, \omega) dw(s, \omega) \mid \int_0^\tau f(s, \omega) dw(s, \omega) = x \right] &= x; \\ E \left[\left(\int_0^T f(s, \omega) dw(s, \omega) \right)^2 \right] &= \int_0^T E[f^2(s, \omega)] ds. \end{aligned}$$

(5)

$$E \left[\int_0^T f(s, \omega) dw(s, \omega) \int_0^T g(s, \omega) dw(s, \omega) \right] = \int_0^T E[f(s, \omega)g(s, \omega)] ds.$$

11.2. Construction of the Ito integral. First note that for any constant random function $f(\omega)$

$$\int_a^b f(\omega) dw(s) = f(\omega)(w(b) - w(a)).$$

Therefore, it is easy to construct the Ito integral for any simple random function $h(t, \omega)$ that assumes a finite number of values. The integrals for simple functions can be extended to integrals for any functions satisfying conditions 1 and 2.

Theorem 7. *For every function $f(t, \omega)$ satisfying Conditions 1 and 2, there is a sequence of step functions $f_n(t, \omega)$ satisfying Conditions 1 and 2 such that*

$$(52) \quad \lim_{n \rightarrow \infty} \int_0^T |f(t, \omega) - f_n(t, \omega)|^2 dt = 0$$

for almost all $\omega \in \Omega$, and the limit

$$(53) \quad I(t, \omega) := \lim_{n \rightarrow \infty} \int_0^t f_n(s, \omega) dw(s, \omega)$$

is uniform in T for almost all $\omega \in \Omega$ and is independent of the sequence $f_n(t, \omega)$ satisfying conditions 1 and 2.

11.3. Existence and uniqueness of solutions of the Ito SDEs. Consider the SDE

$$(54) \quad dx(t) = b(x(t), t)dt + \sigma(x(t), t)dw, \quad x(0) = x_0 \in \mathbb{R}^d, \quad t \in [0, T],$$

where w is the standard Brownian motion. We assume that the functions b and σ satisfy the following conditions. There exists a constant C such that

$$(55) \quad \|b(x, t)\| + \|\sigma(x, t)\|_F \leq C(1 + \|x\|), \quad \text{for all } x \in \mathbb{R}^d, \quad t \in [0, T],$$

where $\|\cdot\|_F$ denotes the Frobenius matrix norm:

$$\|A\|_F := \sqrt{\text{tr}(A^\top A)},$$

and

$$(56) \quad \|b(x, t) - b(y, t)\| + \|\sigma(x, t) - \sigma(y, t)\|_F \leq C\|x - y\|, \quad \text{for all } x, y \in \mathbb{R}^d, \quad t \in [0, T].$$

The first condition says that b and σ do not grow faster than linearly in x , and the second condition is an analogue of the Lipschitz condition. In this case, Eq. (54) with $E[\|x_0\|^2] < \infty$ has a unique solution such that

$$E \left[\int_0^t \|x(s)\|^2 ds \right] < \infty \quad \text{for all } t \in [0, T].$$

From now on, we will autonomize SDEs to save some writing. If b and/or σ explicitly depend on t , we introduce a new independent variable s , declare that t is a new dependent variable, and add the equation $dt = ds$. This is called “autonomization”. Therefore, the assumption that the SDE is autonomous does not lead to the loss of generality.

11.4. Notation common in probability books. It is common in the probability community to denote stochastic processes by capital letters with subscripts specifying their time arguments:

$$(57) \quad dX_t = b(X_t)dt + \sigma(X_t)dw, \quad X_0 = x, \quad t \in [0, T], \quad x \in \mathbb{R}^d.$$

Terminology:

- X_t satisfying Eq. (57) is called a *diffusion process*;
- $b(x)$ is called a *drift*;
- the matrix $\Sigma(x) = \sigma(x)\sigma^\top(x)$ is called a *diffusion matrix*.

Exercise Show that b and Σ satisfy:

$$(58) \quad \lim_{t \rightarrow s} E \left[\frac{X_t - X_s}{t - s} \mid X_s = x \right] = b(x, s)$$

$$(59) \quad \lim_{t \rightarrow s} E \left[\frac{[X_t - X_s][X_t - X_s]^\top}{t - s} \mid X_s = x \right] = \Sigma(x, s).$$

11.5. The Ornstein-Uhlenbeck process. The Ornstein-Uhlenbeck process models the velocity of a heavy particle pushed around by light particles. The variable X_t is the velocity of the particle. For simplicity, we consider it in 1D:

$$(60) \quad dX_t = -\gamma X_t dt + \sqrt{2D}dw, \quad X_0 = x \in R$$

where x is a fixed number, γ is the friction coefficient, and D is the diffusion coefficient.

The exact solution of Eq. (60) can be written in the closed form involving a stochastic integral. We proceed as we do when we solve a linear non-homogeneous first order ODE. Switch the term $-\gamma X_t$ to the left-hand side and multiply the equation by the integrating factor $\exp(\gamma t)$. Then we get

$$d(e^{\gamma t} X_t) = \sqrt{2D}e^{\gamma t} dw.$$

Suppose $X_0 = x$. Integrating from 0 to t we obtain:

$$e^{\gamma t} X_t - x = \sqrt{2D} \int_0^t e^{\gamma s} dw_s.$$

Hence the solution of Eq. (60) is

$$X_t = xe^{-\gamma t} + \sqrt{2D} \int_0^t e^{-\gamma(t-s)} dw_s.$$

The solution X_t is a Gaussian random variable with mean $xe^{-\gamma t}$ and variance $\frac{D}{\gamma}(1 - e^{-2\gamma t})$. The variance is found as follows. We partition the interval $[0, t]$ into n equal subintervals

of length $h = t/n$ and let $n \rightarrow \infty$:

$$\begin{aligned} \text{Var} \left(\sqrt{2D} \int_0^t e^{-\gamma(t-s)} dw_s \right) &= 2De^{-2\gamma t} E \left[\left(\int_0^t e^{\gamma s} dw_s \right)^2 \right] \\ &= 2De^{-2\gamma t} \lim_{n \rightarrow \infty} E \left[\sum_{j=0}^{n-1} e^{\gamma 2jh} [w((j+1)h) - w(jh)]^2 \right] \\ &= 2De^{-2\gamma t} \int_0^t e^{2\gamma s} ds = 2De^{-2\gamma t} \frac{1}{2\gamma} (e^{2\gamma t} - 1) \\ &= \frac{D}{\gamma} (1 - e^{-2\gamma t}). \end{aligned}$$

As $t \rightarrow \infty$, the velocity X_t of the particle becomes a Gaussian random variable with mean 0 and variance $\frac{D}{\gamma}$.

Now let us calculate the covariance function of the Ornstein-Uhlenbeck process. Suppose $t \geq s$.

$$\begin{aligned} \text{Cov}(X_s, X_t) &= E \left[\left(\sqrt{2D} \int_0^s e^{-\gamma(s-\tau)} dw_\tau \right) \left(\sqrt{2D} \int_0^t e^{-\gamma(t-\tau)} dw_\tau \right) \right] \\ &= 2De^{-\gamma(t+s)} E \left[\left(\int_0^s e^{\gamma\tau} dw_\tau \right) \left(\int_0^s e^{\gamma\tau} dw_\tau + \int_s^t e^{\gamma\tau} dw_\tau \right) \right] \\ &= \frac{D}{\gamma} e^{-\gamma(t+s)} (e^{2\gamma s} - 1) \\ &= \left[-\frac{D}{\gamma} \right] e^{-\gamma(t+s)} + \frac{D}{\gamma} e^{-\gamma(t-s)}. \end{aligned}$$

Note that as t and s tend to infinity, the first term decays to zero. Then the covariance function depends only on the difference $t - s$ and is given by

$$R(t - s) = \frac{D}{\gamma} e^{-\gamma|t-s|}.$$

The Ornstein-Uhlenbeck process (60) with the initial condition $x \sim N(0, \frac{D}{\gamma})$ can serve as a model for a colored noise.

12. THE ITO CALCULUS

SDEs can be solved analytically only in special cases. Even if so, we often need not the formula for the solution but the expected value of some function defined on the random trajectories. This function can be the first passage time to a given region of the phase space or the probability to reach first one given region rather than the other given region. To answer such kind of questions, it is handy to be able to calculate the time evolution of functions defined on trajectories. The Ito formula provides us with a tool to do it.

12.1. A derivation of the Ito formula. The most important result in the Ito calculus is the Ito formula.

12.1.1. *1D case.* For simplicity, we will start with the 1D case. Let $x(t)$ be trajectory obeying the ODE

$$\frac{dx}{dt} = b(x).$$

Then any function $f(x(t), t)$ evolves in time according to

$$(61) \quad \frac{df}{dt} = \frac{\partial f}{\partial t} + \frac{\partial f}{\partial x} \frac{dx}{dt} = \frac{\partial f}{\partial t} + b(x) \frac{\partial f}{\partial x}.$$

Now let $x(t, \omega)$ be a trajectory of the SDE

$$dX_t = b(X_t)dt + \sigma(X_t)dw.$$

One could naively write

$$df = \frac{\partial f}{\partial t} dt + \frac{\partial f}{\partial x} dX_t = \frac{\partial f}{\partial t} dt + \frac{\partial f}{\partial x} (b(X_t)dt + \sigma(X_t)dw)$$

but this would be **WRONG**. This is because $dw = O(\sqrt{dt})$ and $(dw)^2 = O(dt)$. The correct differential of f is given by Eq. (62) below.

Let us derive it. We want to find the differential, i.e., the part of the increment of the order of dt or larger for a function $f(X_t, t)$, where $dX_t = b(X_t)dt + \sigma(X_t)dw$. We will write a formal Taylor expansion of $f(X_t, t)$ and keep all terms of the order of dt or larger. To save some space, the arguments (X_t, t) in all derivatives will be omitted.

$$f(X_t + dX_t, t + dt) = f(X_t, t) + \frac{\partial f}{\partial t} dt + \frac{\partial f}{\partial x} dX_t + \frac{1}{2} \frac{\partial^2 f}{\partial t^2} (dt)^2 + \frac{\partial^2 f}{\partial x \partial t} dX_t dt + \frac{1}{2} \frac{\partial^2 f}{\partial x^2} (dX_t)^2 + \dots$$

The term dw is $O(\sqrt{dt})$. The term $dX_t dt$ is $O((dt)^{3/2})$. The term $(dw)^2$ contained in $(dX_t)^2$ is $O(dt)$. Hence we need to keep only the term $\frac{1}{2} \frac{\partial^2 f}{\partial x^2} (dX_t)^2$ out of the second order terms. Therefore,

$$df(X_t, t) = \frac{\partial f}{\partial t} dt + \frac{\partial f}{\partial x} dX_t + \frac{1}{2} \frac{\partial^2 f}{\partial x^2} (dX_t)^2.$$

Writing dX_t explicitly we get:

$$\begin{aligned} (dX_t)^2 &= (b(X_t)dt + \sigma(X_t)dw)(b(X_t)dt + \sigma(X_t)dw) \\ &= b^2(X_t)(dt)^2 + 2b(X_t)\sigma(X_t)dt dw + \sigma^2(X_t)(dw)^2 = \sigma^2(X_t)dt + o(dt). \end{aligned}$$

Hence,

$$(62) \quad df(X_t, t) = \left[\frac{\partial f}{\partial t} + b(X_t) \frac{\partial f}{\partial x} + \frac{1}{2} \sigma^2(X_t) \frac{\partial^2 f}{\partial x^2} \right] dt + \sigma(X_t) \frac{\partial f}{\partial x} dw.$$

Note that Eq. (62) can be rewritten as

$$(63) \quad df(X_t, t) = \left[\frac{\partial f}{\partial t} + Lf \right] dt + \sigma(x) \frac{\partial f}{\partial x} dw, \quad \text{where } L = b(X_t) \frac{\partial}{\partial x} + \frac{1}{2} \sigma^2(X_t) \frac{\partial^2}{\partial x^2}.$$

L is called the infinitesimal generator of the process.

12.1.2. *Multidimensional case.* Now we turn to the multidimensional case is stated below. We define the generator by

$$(64) \quad L = b(X_t) \cdot \nabla + \frac{1}{2} \Sigma(X_t) : \nabla \nabla,$$

where the symbol “:” means

$$A : B := \sum_{i=1}^n \sum_{j=1}^n A_{ij} B_{ij} \equiv \text{tr} \left(A^\top B \right)$$

and

$$(\nabla \nabla)_{ij} = \frac{\partial^2}{\partial x_i \partial x_j}.$$

In the coordinate form, L can be rewritten as

$$(65) \quad L = \sum_{j=1}^d b_j(X_t) \frac{\partial}{\partial x_j} + \frac{1}{2} \sum_{i,j=1}^d \Sigma_{ij}(X_t) \frac{\partial^2}{\partial x_i \partial x_j}.$$

Theorem 8. (Ito’s formula) Assume that b and σ satisfy Eqs. (55) and (56) and $E[\|X_0\|^2] < \infty$. Let $f(x, t)$ be twice continuously differentiable in x in \mathbb{R}^d and continuously differentiable in t on $[0, T]$. Then the process $f(X_t, t)$ satisfies:

$$(66) \quad \boxed{f(X_t, t) = f(X_0, 0) + \int_0^t \frac{\partial f(X_s, s)}{\partial s} ds + \int_0^t Lf(X_s, s) ds + \int_0^t \nabla f(X_s, s)^\top \sigma(X_s) dw_s.}$$

Eq. (66) is equivalent to

$$(67) \quad \frac{df}{dt} = \frac{\partial f}{\partial t} + \nabla f \cdot \frac{dX_t}{dt} + \frac{1}{2} \frac{dX_t}{dt} \cdot \nabla \nabla f \frac{dX_t}{dt},$$

or

$$(68) \quad df(X_t, t) = \frac{\partial f}{\partial t} dt + \sum_{i=1}^d \frac{\partial f}{\partial x_i} dX_i + \frac{1}{2} \sum_{i,j=1}^d \frac{\partial^2 f}{\partial x_i \partial x_j} dX_i dX_j,$$

Ito’s formula is proven e.g. in [4].

12.2. **The geometric Brownian motion.** The geometric Brownian motion is a stochastic process satisfying the following SDE

$$(69) \quad dX_t = \mu X_t dt + \sigma X_t dw,$$

where w is the standard Brownian motion, μ (the percentage drift) and σ (the percentage volatility) are constants. The SDE (69) is used in mathematical finance to model the stock prices in the Black-Scholes model.

Eq. (69) has an analytic solution that can be found as follows. Introduce the new dependent variable $Y_t = \log X_t$. Using Ito's formula (62) and taking into account that $\frac{dY}{dX} = \frac{1}{X}$ and $\frac{d^2Y}{dX^2} = -\frac{1}{X^2}$ we write the differential of Y_t :

$$dY_t = \left(\mu X_t \frac{dY}{dX} + \frac{1}{2} \sigma^2 X_t^2 \frac{d^2Y}{dX^2} \right) dt + \sigma X_t \frac{dY}{dX} dw = \left(\mu - \frac{\sigma^2}{2} \right) dt + \sigma dw.$$

The right-hand side of the SDE for Y_t is independent of Y_t and hence Y_t is found just by integration of the right-hand side:

$$Y_t = Y_0 + \left(\mu - \frac{\sigma^2}{2} \right) t + \sigma w_t.$$

Returning to $X_t = \exp(Y_t)$ we get the exact solution of Eq. (69):

$$(70) \quad X_t = X_0 e^{\left(\mu - \frac{\sigma^2}{2} \right) t + \sigma w_t}.$$

We say that a random variable X is *lognormal* if its logarithm is Gaussian, i.e. if $\log X \sim N(m, s^2)$.

Exercise Check that if $Y \sim N(m, s^2)$ then

$$E[e^Y] = e^{m + \frac{s^2}{2}}, \quad \text{Var}(e^Y) = e^{2m + s^2} (e^{s^2} - 1),$$

and the pdf of $X = e^Y$ is given by

$$f_{e^Y}(x) = \frac{1}{x \sqrt{2\pi s^2}} e^{-\frac{(\log x - m)^2}{2s^2}}.$$

Note that

$$Y_t = \log X_t \sim N \left(Y_0 + \left(\mu - \frac{\sigma^2}{2} \right) t, \sigma^2 t \right),$$

hence the geometric Brownian motion X_t has a log-normal distribution. The mean and the variance of X_t are

$$E[X_t] = X_0 e^{\mu t}, \quad \text{Var}(X_t) = X_0^2 e^{2\mu t} (e^{\sigma^2 t} - 1).$$

12.3. Backward Kolmogorov equation. Imagine that we are interested in some quantity f that depends on X_t evolving according to the SDE

$$dX_t = b(X_t)dt + \sigma(X_t)dw.$$

Suppose that we want to find the expected value of f at a future time T given that at the present time t , $X_t = x$. For example, you can think of f being an option price that depends on the stock price X . We denote the expected value of f at time T conditioned on $X_t = x$ by $u(x, t)$. Let us find the time evolution of

$$u(x, t) := E[f(X_T) | X_t = x] = \int_{\mathbb{R}^d} f(y) p(y, T | x, t) dy,$$

In words, imagine that we start a stochastic process X_s at time t at the point x . We stop it at a fixed time T . We want to find the expected value of $f(X_T)$. This expected value

$u(x, t)$ depends on the initial time t and the initial point x . Obviously, for the terminal time $t = T$ we have: $u(x, T) = f(x)$. Using Ito's formula (66) we calculate

$$u(X_T, T) - u(X_t, t) = \int_t^T \frac{\partial}{\partial s} u(X_s, s) ds + \int_t^T Lu(X_s, s) ds + \int_t^T \nabla u(X_s, s)^\top \sigma(X_s) dw_s.$$

Now we will take expected values of both parts of this equation conditioned on $X_t = x$. Note that

$$E[u(X_T, T) \mid X_t = x] = E[f(X_T) \mid X_t = x] = u(x, t)$$

and

$$E[u(X_t, t) \mid X_t = x] = u(x, t).$$

Hence the conditional expectation of the left-hand side is 0. Also note that by property (4) in Section 11.1,

$$E \left[\int_t^T \nabla u(X_s, s)^\top \sigma(X_s) dw_s \right] = 0.$$

Therefore, for all x, t and T we have

$$\int_t^T \left[\frac{\partial}{\partial s} u(x, s) + Lu(x, s) \right] ds = 0.$$

Hence for all $t \leq T$, $u(x, t)$ satisfies the PDE with the final condition:

$$\frac{\partial}{\partial t} u(x, t) + Lu(x, t) = 0 \quad u(x, T) = f(x).$$

Re-defining u as

$$u(x, t) = E[f(X_t) \mid X_0 = x], \quad \text{i.e., } u(x, t) = u_{\text{old}}(x, T - t),$$

we obtain that for $0 \leq t \leq T$

$$(71) \quad \boxed{\frac{\partial}{\partial t} u(x, t) = Lu(x, t), \quad u(x, 0) = f(x).}$$

Eq. (71) is called the *Backward Kolmogorov Equation*. It describes the time evolution of expected values. Note that in the re-definition of u we used the fact that X_t evolves according to an autonomous SDE which is invariant with respect to a time shift.

12.4. The expected first passage time. Let $A \subset \mathbb{R}^d$ be some region. The first passage time to A is defined as

$$\tau_A = \inf\{t \geq 0 \mid X_t \in A\}.$$

Let $u(x, t)$ be the expected first passage time to A for the process X_t starting at x , i.e.,

$$u(x, t) = E[\tau_A \mid X_t = x].$$

We calculate:

$$u(X_{\tau_A}, \tau_A) - u(X_t, t) = \int_t^{\tau_A} \frac{\partial}{\partial s} u(X_s, s) ds + \int_t^{\tau_A} Lu(X_s, s) ds + \int_t^{\tau_A} \nabla u(X_s, s)^\top \sigma(X_s) dw_s.$$

Next, we take the expected values of the left- and right-hand side of this equation conditioned on $X_t = x$. Taking into account that

$$E[u(X_{\tau_A}, \tau_A) \mid X_t = x] = E[\tau_A \mid X_t = x] = u(x, t) \quad \text{and} \quad E[u(X_t, t) \mid X_t = x] = u(x, t)$$

we get for all t and x

$$\int_t^{\tau_A} \frac{\partial}{\partial s} u(x, s) ds + \int_t^{\tau_A} Lu(x, s) ds = 0.$$

We also note that

$$\frac{\partial}{\partial t} u(x, t) = 1.$$

Finally, if $x \in A$ we have $\tau_A = 0$. Hence the mean first passage time $u(x)$ satisfies the following boundary value problem

$$(72) \quad \boxed{Lu = -1, \quad x \notin A, \quad u(\partial A) = 0.}$$

Example 22 Let $dX_t = \sqrt{2\beta^{-1}}dw$, $X_0 = x$, i.e., a scaled one-dimensional Brownian motion starting at the point x . Let us find the expected exit time from the interval $[-1, 1]$. As we have shown, the expected exit time $u(x)$ satisfies Eq. (72), which in our case becomes

$$\beta^{-1}u'' = -1, \quad -1 \leq x \leq 1, \quad u(1) = u(-1) = 0.$$

Solving this equation we obtain

$$u(x) = \frac{\beta}{2}(1 - x^2).$$

12.5. The committor equation. Let $A \subset \mathbb{R}^d$ and $B \subset \mathbb{R}^d$ be some regions. The committor function $q(x)$ is defined as the probability that the process starting at the point x first reaches B rather than A [6, 7]. Let us derive a boundary-value problem for the committor. It is clear that $q(\partial A) = 0$ and $q(\partial B) = 1$. For $x \in (A \cup B)^c$ let us define the first passage time to $A \cup B$, i.e.,

$$\tau_{AB} = \inf\{t \geq 0 \mid X_t \in A \cup B\}.$$

We calculate:

$$q(X_{\tau_{AB}}) - q(X_0) = \int_0^{\tau_{AB}} Lq(X_s) ds + \int_0^{\tau_{AB}} \nabla q(X_s)^\top \sigma(X_s) dw_s.$$

Take the expected values of the left- and right-hand side of this equation conditioned on $X_0 = x$. We get that for all x

$$q(x) - q(x) = \int_0^{\tau_{AB}} Lq(x) ds = 0$$

Therefore, the solution of the boundary-value problem

$$Lq = 0, \quad x \in (A \cup B)^c, \quad q(\partial A) = 0, \quad q(\partial B) = 1$$

is the committor function.

Example 23 Let $dX_t = -V'(x)dt + \sqrt{2\beta^{-1}}dw$, $X_0 = x$, i.e., a particle moving according to the overdamped Langevin dynamics in the potential force field $V(x)$. In the 1D case, the committor equation can be solved exactly. We have:

$$-V'(x)q'(x) + \beta^{-1}q''(x) = 0, \quad a \leq x \leq b, \quad q(a) = 0, \quad q(b) = 1.$$

Multiply this equation by $\beta e^{-\beta V(x)}$. Then its left-hand side becomes a complete differential:

$$\left(e^{-\beta V(x)} q'(x) \right)' = 0.$$

Integrating this equation and taking the boundary conditions into account we get

$$q(x) = \frac{\int_a^x e^{\beta V(y)} dy}{\int_a^b e^{\beta V(y)} dy}.$$

12.6. The generator of a Markov process. In Section 12.3, we fixed a function f and we considered the expectation of f at time t as a function of the initial point x and time t :

$$u(x, t) := E[f(X_t) \mid X_0 = x].$$

Now we fix the time t and consider the same expectation as a map applied to the set of continuous and bounded functions $f(x)$, $x \in \mathbb{R}^d$. Therefore, we define the family of operators indexed by t :

$$(P_t f)(x) := E[f(X_t) \mid X_0 = x] = \int_{\mathbb{R}^d} f(y) p(y, t \mid x, 0) dy.$$

We will call the operator P_t the *transfer operator*. It is analogous to the stochastic matrix P in the discrete-time Markov chains.

The operator P_t possesses the semigroup properties:

$$P_0 = I, \quad P_{t+s} = P_t \circ P_s \quad \text{for all } t, s \geq 0.$$

Indeed,

$$(P_0 f)(x) := E[f(X_0) \mid X_0 = x] = f(x),$$

Hence P_0 is the identity. Recall the Chapman-Kolmogorov equation (Eq. (46)) expressing the Markov property. Using it, we write:

$$\begin{aligned} (P_{t+s} f)(x) &= \int_{\mathbb{R}^d} f(y) p(y, t+s \mid x, 0) dy = \int_{\mathbb{R}^d} f(y) dy \int_{\mathbb{R}^d} p(y, t+s \mid z, t) p(z, t \mid x, 0) dz \\ &= \int_{\mathbb{R}^d} dz p(z, t \mid x, 0) \int_{\mathbb{R}^d} p(y, t+s \mid z, t) f(y) dy = (P_t P_s f)(x). \end{aligned}$$

Due to this, the operator P_t is often referred to as the *Markov semigroup*.

Now consider the limit as $t \downarrow 0$:

$$(73) \quad (L f)(x) := \lim_{t \rightarrow 0^+} \frac{(P_t f)(x) - f(x)}{t}.$$

Assume that this limit exists. This limit is called the *infinitesimal generator* of the transfer operator P_t or the *generator* of the Markov process X_t .

Recall that $(P_t f)(x) = u(x, t)$. Eq. (73) implies that

$$\frac{\partial u}{\partial t} = \lim_{s \rightarrow 0} \frac{P_{t+s} f - P_t f}{s} = L P_t f = L u, \quad u(x, 0) = f(x).$$

This is the Backward Kolmogorov equation that we have obtained in Section 12.3. Matching it with Eq. (71) we see that the operator L must be given by Eq. (64).

12.7. The adjoint semigroup and the forward Kolmogorov equation. In Section 12.6 we considered the expected value of $f(X_t)$ conditioned on $X_0 = x$. Now we assume that X_0 does not start at x with probability 1 but the starting point is distributed according to a pdf $\mu_0(x)$. Then the expected value of $f(X_t)$ is

$$E[f(X_t)] = \int_{\mathbb{R}^d} P_t f(x) \mu_0(x) dx.$$

Writing $P_t f$ explicitly and switching the order of integration we obtain

$$\begin{aligned} E[f(X_t)] &= \int_{\mathbb{R}^d} P_t f(x) \mu_0(x) dx = \int_{\mathbb{R}^d} \mu_0(x) dx \int_{\mathbb{R}^d} dy f(y) p(y, t | x, 0) \\ &= \int_{\mathbb{R}^d} dy f(y) \int_{\mathbb{R}^d} \mu_0(x) p(y, t | x, 0) dx \\ (74) \quad &=: \int_{\mathbb{R}^d} dy f(y) P_t^* \mu_0. \end{aligned}$$

In the original order of integration, we froze the pdf $\mu_0(x)$ while evolved $f(X_t)$ in time. After switching the order of integration, we froze f and evolved the pdf μ in time. Finally, we have defined the evolution operator for the pdf:

$$(75) \quad \mu_t(x) := (P_t^* \mu_0)(x) := \int_{\mathbb{R}^d} \mu_0(x) p(y, t | x, 0) dx.$$

The operator P_t^* is adjoint to the transfer operator P_t . Indeed, consider the inner product

$$(f, g) := \int_{\mathbb{R}^d} f(x) g(x) dx.$$

Then Eq. (74) shows that

$$E[f(X_t)] = (P_t f, \mu_0) = (f, P_t^* \mu_0).$$

The infinitesimal generator for the adjoint semigroup P_t^* is defined by

$$(76) \quad (L^* \mu_0)(x) := \lim_{t \rightarrow 0} \frac{(P_t^* \mu_0)(x) - \mu_0(x)}{t}.$$

It is easy to check that L and L^* are adjoint, i.e., for all admissible f and g ,

$$(77) \quad (L f, g) = (f, L^* g).$$

Eqs. (75) and (76) show that the time evolution of the probability density function is given by

$$(78) \quad \frac{\partial \mu(x, t)}{\partial t} = L^* \mu(x, t), \quad \mu(x, 0) = \mu_0(x).$$

Eq. (78) is called the *forward Kolmogorov equation* or the *Fokker-Planck equation*.

Eq. (64) allows us to find the adjoint generator L^* explicitly. Consider the process governed by the SDE

$$dX_t = b(X_t)dt + \sigma(X_t)dw, \quad X(0) = x, \quad t \geq 0, \quad x \in \mathbb{R}^d.$$

For it, the generator is the differential operator given by

$$L = b(x) \cdot \nabla + \frac{1}{2} \Sigma(x) : \nabla \nabla.$$

To find the adjoint generator L^* , we consider the identity Eq. (77):

$$\begin{aligned} (Lf, g) &= \int_{\mathbb{R}^d} (b \cdot \nabla f + \frac{1}{2} \Sigma : \nabla \nabla f) g dx \\ &= \int_{\mathbb{R}^d} f \left(-\nabla \cdot (gb) + \frac{1}{2} \nabla \nabla : (\Sigma g) \right) dx = (f, L^*g). \end{aligned}$$

Here we have integrated by parts the first term once and the second term twice. Hence,

$$(79) \quad L^*g = -\nabla \cdot (gb) + \frac{1}{2} \nabla \nabla : (\Sigma g)$$

Example 24 We will elaborate the procedure of obtaining L^* in 2D. The extension to higher dimensions is straightforward. Consider a 2D stochastic process of the form

$$\begin{aligned} dX_t &= b_1(X_t, Y_t)dt + \sigma_{11}(X_t, Y_t)dw_1 + \sigma_{12}(X_t, Y_t)dw_2 \\ dY_t &= b_2(X_t, Y_t)dt + \sigma_{21}(X_t, Y_t)dw_1 + \sigma_{22}(X_t, Y_t)dw_2 \end{aligned}$$

In the vector notations it looks as

$$d \begin{pmatrix} X_t \\ Y_t \end{pmatrix} = \begin{pmatrix} b_1 \\ b_2 \end{pmatrix} dt + \begin{pmatrix} \sigma_{11} & \sigma_{12} \\ \sigma_{21} & \sigma_{22} \end{pmatrix} \begin{pmatrix} dw_1 \\ dw_2 \end{pmatrix}.$$

The matrix $\Sigma = \sigma \sigma^\top$ is

$$\Sigma \equiv \begin{pmatrix} \Sigma_{11} & \Sigma_{12} \\ \Sigma_{12} & \Sigma_{22} \end{pmatrix} \equiv \begin{pmatrix} \sigma_{11}^2 + \sigma_{12}^2 & \sigma_{11}\sigma_{21} + \sigma_{12}\sigma_{22} \\ \sigma_{11}\sigma_{21} + \sigma_{12}\sigma_{22} & \sigma_{21}^2 + \sigma_{22}^2 \end{pmatrix}.$$

The generator L applied to a function f is:

$$Lf = b_1 \partial_x f + b_2 \partial_y f + \frac{1}{2} [\Sigma_{11} \partial_{xx} f + 2\Sigma_{12} \partial_{xy} f + \Sigma_{22} \partial_{yy} f].$$

The adjoint generator L^* is found from the identity $(Lf, g) = (f, L^*g)$:

$$\begin{aligned} (Lf, g) &= \int_{\mathbb{R}^2} \left[gb_1 \partial_x f + gb_2 \partial_y f + \frac{1}{2} g [\Sigma_{11} \partial_{xx} f + 2\Sigma_{12} \partial_{xy} f + \Sigma_{22} \partial_{yy} f] \right] dx dy \\ &= \int_{\mathbb{R}^2} f \left[-\partial_x (b_1 g) - \partial_y (b_2 g) + \frac{1}{2} [\partial_{xx} (\Sigma_{11} g) + 2\partial_{xy} (\Sigma_{12} g) + \partial_{yy} (\Sigma_{22} g)] \right] dx dy \\ &= (f, L^*g). \end{aligned}$$

Hence

$$L^*g = -\partial_x(b_1g) - \partial_y(b_2g) + \frac{1}{2} [\partial_{xx}(\Sigma_{11}g) + 2\partial_{xy}(\Sigma_{12}g) + \partial_{yy}(\Sigma_{22}g)].$$

12.8. The invariant pdf. Recall irreducible continuous-time Markov chains. The invariant measure is the solution of $\pi L = 0$ or, equivalently, $L^\top \pi^\top = 0$. π and π^\top are row and column vectors respectively. If the invariant measure is normalizable, we normalize it so that $\sum_i \pi_i = 1$ and call it the invariant distribution. For irreducible Markov chains with a finite number of states, any probability distribution converges over time to the unique invariant distribution π , i.e., for any initial distribution p_0 , the solution of

$$\frac{dp}{dt} = pL, \quad p(0) = p_0$$

converges to π . Such Markov chains are called *ergodic*. The Markov Chain Monte Carlo methods employ this property. Also recall that irreducibility of a continuous-time Markov chain with a finite number of states implies that the eigenvalue 0 of L has multiplicity one.

Suppose that the equation $L^*f = 0$ for a Markov process X_t has a unique positive solution μ up to a multiplicative constant, and this solution is normalizable so that $\int \mu(x)dx = 1$, then μ is the unique invariant pdf. In this case, for any initial pdf μ_0 , the pdf μ_t converges to μ as $t \rightarrow \infty$:

$$\lim_{t \rightarrow \infty} P_t^* \mu_0 = \mu.$$

Such Markov processes are also called *ergodic*.

The unique invariant pdf $\mu(x)$ satisfies the stationary forward Kolmogorov equation (stationary Fokker-Planck equation):

$$(80) \quad L^* \mu = 0, \quad \int_{\mathbb{R}^d} \mu(x) dx = 1.$$

Exercise (1) Show that the generator of the 1D Ornstein-Uhlenbeck process (60) is given by

$$L = -\gamma x \frac{d}{dx} + D \frac{d^2}{dx^2}.$$

(2) Integrating by parts, derive the expression for the adjoint generator

$$L^*g = \frac{d}{dx}(\gamma xg) + D \frac{d^2g}{dx^2}.$$

(3) Solve the equation

$$L^* \mu = 0, \quad \int_{-\infty}^{\infty} \mu dx = 1$$

and find that the invariant pdf for the 1D Ornstein-Uhlenbeck process is

$$(81) \quad \mu(x) = \sqrt{\frac{\gamma}{2\pi D}} e^{-\frac{\gamma x^2}{2D}}.$$

13. THE LANGEVIN DYNAMICS

13.1. The full Langevin dynamics. The Langevin equation models the dynamics of heavy particles in the potential force field pushed around by light particles:

$$(82) \quad \begin{aligned} dq &= \frac{p}{m} dt \\ dp &= (-\nabla V(q) - \gamma p) dt + \sqrt{2\gamma m \beta^{-1}} dw, \end{aligned}$$

where $(q, p) \in \mathbb{R}^{2d}$ are the positions and momenta of the heavy particles, γ is the friction coefficient, m is the mass of the heavy particles, and $-\nabla V(q)$ is the potential force acting on the heavy particles. Eq. (82) can be written in the form (57) by introducing

$$X_t = \begin{bmatrix} q \\ p \end{bmatrix}, \quad b(X_t) = \begin{bmatrix} p/m \\ -\nabla V(q) - \gamma p \end{bmatrix}, \quad \sigma = \sqrt{2\gamma m \beta^{-1}} \begin{bmatrix} 0 & 0 \\ 0 & I \end{bmatrix},$$

where I is the $d \times d$ identity matrix.

Exercise (1) Show that the infinitesimal generator for Eq. (82) is given by

$$L = \frac{p}{m} \cdot \nabla_q - \nabla_q V \cdot \nabla_p + \gamma (-p \nabla_p + m \beta^{-1} \Delta_p).$$

(2) Derive the expression for the adjoint generator

$$L^* g = -\frac{p}{m} \cdot \nabla_q g + \nabla_q V \cdot \nabla_p g + \gamma (\nabla_p \cdot (p g) + m \beta^{-1} \Delta_p g).$$

(3) Solve the stationary Fokker-Planck equation and show that the invariant pdf is given by

$$\mu(q, p) = \frac{1}{Z} e^{-\beta H(q, p)}, \quad \text{where } H(q, p) = \frac{|p|^2}{2m} + V(q).$$

13.2. The overdamped Langevin dynamics. Suppose the friction coefficient γ in Eq. (82) is large and/or the mass m is small, i.e., $m\gamma^{-1}$ is small. We divide the equation for p by γ :

$$\gamma^{-1} dp = \gamma^{-1} m dv = (-\gamma^{-1} \nabla V(q) - p) dt + \gamma^{-1} \sqrt{2m\gamma\beta^{-1}} dw.$$

We use that assumption that $m\gamma^{-1}$ is small and set the left-hand side of the SDE above to 0. Then we replace p with $m \frac{dq}{dt}$ in the right-hand side and multiply both sides by γ :

$$m\gamma \frac{dq}{dt} = -\nabla V(q) + \sqrt{2m\gamma\beta^{-1}} \frac{dw}{dt}$$

or

$$m\gamma dq = -\nabla V(q) dt + m\gamma \sqrt{\frac{2\beta^{-1}}{m\gamma}} dw(t)$$

Now we want to cancel out $m\gamma$. To do so, we rescale the time by introducing $\tau = (m\gamma)^{-1}t$. Then $dt = m\gamma d\tau$. Recall that if $w(t)$ is a Brownian motion, then for any $c > 0$, $c^{-1/2}w(ct)$ is also a brownian motion, i.e., $c^{1/2}w(t) = w(ct)$. Hence

$$dw(t) = dw(m\gamma\tau) = (m\gamma)^{1/2} dw(\tau).$$

Therefore, choosing the new time τ and canceling $m\gamma$ we get

$$dq = -\nabla V(q)d\tau + \sqrt{2\beta^{-1}}dw(\tau).$$

For the overdamped Langevin dynamics, the generator L is given by

$$L = -\nabla V \cdot \nabla + \beta^{-1}\Delta.$$

The adjoint generator L^* is

$$L^*g = -\nabla \cdot (g\nabla V) + \beta^{-1}\Delta g.$$

The invariant pdf is

$$\mu = \frac{1}{Z}e^{-\beta V(x)}.$$

14. NUMERICAL INTEGRATION OF SDEs

Analytical solutions to SDEs can be obtained only in special cases. Therefore, it is important to develop methods for numerical integration of SDEs. Here we will present the Euler-Maruyama method (1955), a higher order Milstein's method (1974), and MALA (1996) (Metropolis Adjusted Langevin Method), and discuss some basic concepts of numerical analysis of SDE solvers. An introductory text on numerical integrators for SDEs that includes Matlab codes is the paper by D. Higham [8].

14.1. The Euler-Maruyama and Milstein's methods. Consider the initial-value problem

$$(83) \quad dX_t = b(X_t)dt + \sigma(X_t)dw, \quad 0 \leq t \leq T, \quad X_0 = x.$$

We discretize the time interval $[0, T]$ into n equal subintervals of length $h = \frac{T}{n}$ and generate a vector $w = [w_1, \dots, w_n]$ of n independent Gaussian random variables with mean 0 and variance h to simulate the Brownian motion. In Matlab, it is convenient to do it using the command `randn`. In C, I use the Box-Muller algorithm. We will denote the exact and numerical solutions to Eq. (83) at the mesh points by X_j and Y_j respectively, $j = 0, 1, \dots, n$. Then one can compute the numerical solution using the recurrence:

$$(84) \quad Y_{j+1} = Y_j + b(Y_j)h + \sigma(Y_j)w_{j+1}, \quad Y_0 = x.$$

This method is called the *Euler-Maruyama* method. It was proposed by G. Maruyama in 1955 [?].

A more accurate is Milstein's method (1974) [?]:

$$(85) \quad Y_{j+1} = Y_j + b(Y_j)h + \sigma(Y_j)w_{j+1} + \frac{1}{2}\sigma(Y_j)\sigma'(Y_j)[w_{j+1}^2 - h], \quad Y_0 = x.$$

As one can see, Milstein's method coincides with the Euler-Maruyama if σ is constant. Milstein's method is derived using the Ito formula [?]. The exact expression for X_{j+1} is

$$(86) \quad X_{j+1} = X_j + \int_{t_j}^{t_{j+1}} b(X_s)ds + \int_{t_j}^{t_{j+1}} \sigma(X_s)dw_s.$$

By the Ito formula (62) we have the following expressions for $b(X_s)$ and $\sigma(X_s)$ for $s \in [t_j, t_{j+1}]$:

$$\begin{aligned} b(X_s) &= b(X_j) + \int_{t_j}^s \left[b(X_u) \frac{db(X_u)}{dX} + \frac{1}{2} \sigma^2(X_u) \frac{db^2(X_u)}{dX^2} \right] du + \int_{t_j}^s \sigma(X_u) \frac{db(X_u)}{dX} dw_u \\ &= b(X_j) + \int_{t_j}^s Lb(X_u) du + \int_{t_j}^s L^1 b(X_u) dw_u, \\ \sigma(X_s) &= \sigma(X_j) + \int_{t_j}^s L\sigma(X_u) du + \int_{t_j}^s L^1 \sigma(X_u) dw_u, \\ &\text{where } L^1 := \sigma(X_u) \frac{d}{dX}. \end{aligned}$$

We plug these expressions into Eq. (86) and get:

$$\begin{aligned} X_{j+1} &= X_j + \int_{t_j}^{t_{j+1}} \left[b(X_j) + \int_{t_j}^s Lb(X_u) du + \int_{t_j}^s L^1 b(X_u) dw_u \right] ds \\ &\quad + \int_{t_j}^{t_{j+1}} \left[\sigma(X_j) + \int_{t_j}^s L\sigma(X_u) du + \int_{t_j}^s L^1 \sigma(X_u) dw_u \right] dw_s \\ &= X_j + b(X_j)h + \sigma(X_j)(w_{j+1} - w_j) + L^1 \sigma(X_j) \int_{t_j}^{t_{j+1}} w_s dw_s + O(h^{3/2}) \end{aligned}$$

Recall that

$$\int_a^b w dw = \frac{1}{2} ((w_b - w_a)^2 - (b - a)).$$

Using the fact that

$$L^1 \sigma(X) = \sigma(X) \sigma'(X),$$

and neglecting terms of order $h^{3/2}$ or smaller we obtain Milstein's method.

14.2. Strong and weak orders of convergence.

Definition 10. We say that a method for solving the initial value problem for SDEs on the time interval $[0, T]$ has a strong order of convergence γ if there is a constant C such that

$$(87) \quad E \left[\sup_{0 \leq t_j \leq T} |X_j - Y_j| \right] \leq Ch^\gamma$$

if $h = T/n$ is sufficiently small.

In other words, a method has a strong order of convergence γ , if the mean of errors over realizations of Brownian motions decays as $O(h^\gamma)$.

Definition 11. We say that a method for solving the initial value problem for SDEs on the time interval $[0, T]$ has a weak order of convergence γ if there is a constant C such that

$$(88) \quad \sup_{0 \leq t_j \leq T} |E[X_j] - E[Y_j]| \leq Ch^\gamma$$

if $h = T/n$ is sufficiently small.

In other words, a method has a weak order of convergence γ , if the error of the mean over realizations of Brownian motions decays as $O(h^\gamma)$.

One can show that if the functions b and σ in SDE (83) are nice enough (in particular, globally Lipschitz) then the Euler-Maruyama method has a strong order 1/2 and a weak order 1, while Milstein's method has both weak and strong orders 1. One can check these orders by applying the given method to an SDE with an analytical solution, e.g., the geometric Brownian motion, plotting weak and strong errors versus h in the log-log scale, and estimating the slopes of the lines (see Fig. 4 in [8]) using least squares fits.

The proof of the strong orders of convergence of the Euler-Maruyama and Milstein's methods is not short and simple unlike basic methods for ODEs. Exact formulations and exact proofs of the strong convergence results for Euler-Maruyama and Milstein's methods can be found e.g. in [?].

REFERENCES

- [1] A. Chorin and O. Hald, *Stochastic Tools in Mathematics and Science*, 3rd edition, Springer 2013
- [2] L. Koralov and Ya. Sinai, *Theory of probability and stochastic processes*, 2nd edition, Springer, 2007
- [3] Peter Moertens and Yuval Peres, *Brownian Motion*, electronic book, 2008
- [4] Zeev Schuss, *Theory and Applications of Stochastic Processes, An analytical approach*, Springer, 2010
- [5] Grigorios Pavliotis, *Stochastic processes and Applications, Diffusion Processes, the Fokker-Planck, and Langevin Equations*, Springer, 2014
- [6] W. E, E. Vanden-Eijnden, Towards a theory of transition paths, *Journal of statistical physics* 123 (3) (2006) 503.
- [7] W. E, E. Vanden-Eijnden, Transition-path theory and path-finding algorithms for the study of rare events, *Annu. Rev. Phys. Chem.* 61 (2010) 391–420.
- [8] Desmond J. Higham, *An Algorithmic Introduction to Numerical Simulation of Stochastic Differential Equations*, *SIAM Review*, **43**, 3, (2001) 525-546
- [9] Wikipedia