Lecture notes for Numerik IVc - Numerics for Stochastic Processes, Wintersemester 2012/2013. Instructor: Prof. Carsten Hartmann

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Course outline

1. **Probability theory**
   (a) Some basics: stochastic processes, conditional probabilities and expectations, Markov chains
   References: [MS05, Kle06]
2. **Stochastic differential equations**
   (a) Brownian motion: properties of the paths, Strong Markov Property
   References: [MS05, Øks03, Arn73]
   (b) Stochastic integrals: Itô integrals, Itô calculus, Itô isometry
   References: [MS05, Øks03, Arn73]
   (c) SDEs: existence and uniqueness of solutions, numerical discretisation, applications from physics, biolopgy and finance
   References: [Øks03, Arn73, KP92]
   (d) Misc: Kolmogorov forward and backward PDEs, infinitesimal generators, semigroup theory, stopping times, invariant distributions, Markov Chain Monte Carlo methods for PDEs and SDEs
   References: [Øks03, Arn73, KP92]
3. **Filtering theory (if time permits)**
   (a) Linear filtering: conditional expectation, best approximation, Kalman-Bucy filter for SDEs

   Reference: [Jaz07, Øks03]
4. **Approximation of stochastic processes**
   (a) Spectral theory of Markov chains: infinitesimal generator, metastability, aggregation of Markov chains
   References: [HM05, Sar11]
   (b) Markov jump processes: applications from biology, physiocs and finance, Markov decision processes, control theory
   References: [GHL09]

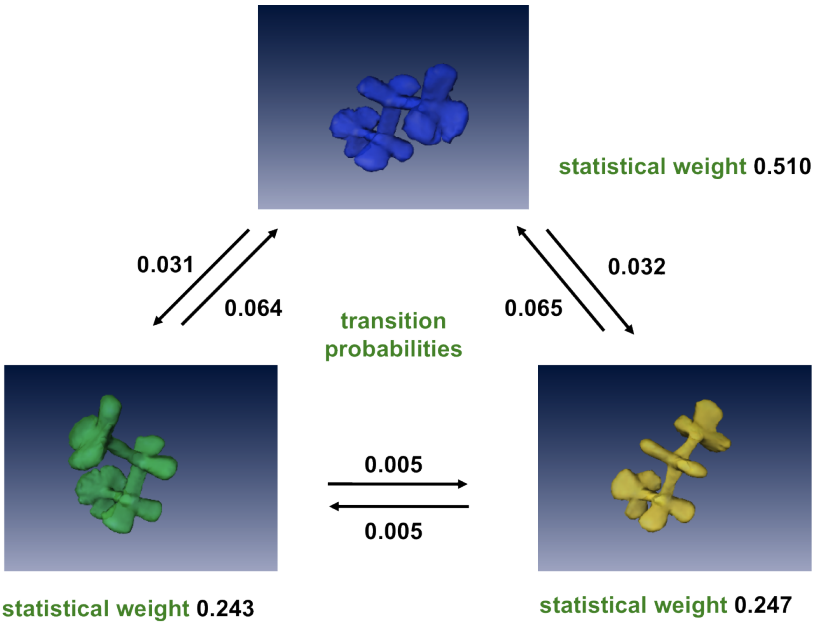

Figure 1: Simulation of a butane molecule and its approximation by a 3-state Markov chain (states in blue, green and yellow; solvent molecules not shown).

1 Day 1, 16.10.2012

1.1 Different levels of modelling

1.1.1 Time-discrete Markov chains

Time index set I is discrete, e.g. $I \subseteq \mathbb{N}$ and state space S is countable or finite, e.g. $S = \{s_1, s_2, s_3\}$ (see Figure 1). Key objects are transition probabilities. For a state space $S = \{1, \ldots, n\}$, the transition probabilities p_{ij} satisfy

$$p_{ij} = \mathbb{P}\left(X_{t+1} = j \mid X_t = i\right)$$

and yield a row-stochastic matrix $P = (p_{ij})_{i,j \in S}$.

1.1.2 Markov jump processes

These are time-continuous, discrete state-space Markov chains. Time index set $I \subseteq \mathbb{R}_+$, S discrete. For a fixed time step $h > 0$, the transition probabilities are given by (see Figure 2)

$$\mathbb{P}\left(X_{t+h} = s_j \mid X_t = s_i\right) = h\ell_{ij} + o(h)$$

where $L = (\ell_{ij})_{i,j \in S}$ and P_h are matrices satisfying $P_h = \exp(hL)$.

Note: the matrix L is row sum zero, i.e. $\sum_j \ell_{ij} = 0$. The waiting times for the Markov chain in any state s_i are exponentially distributed in the sense that

$$\mathbb{P}\left(X_{t+s} = s_i,\ s \in [0, \tau) \mid X_t = s_i\right) = \exp(\ell_{ii}\tau)$$

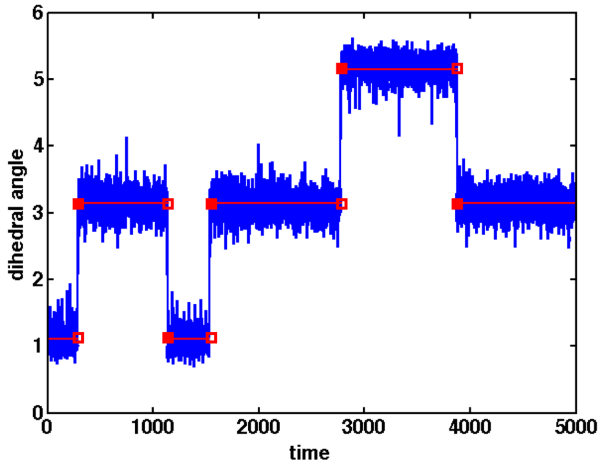

Figure 2: Simulation of butane: typical time series of the central dihedral angle (blue: metastable diffusion process, red: Markov jump process)

and the 'average waiting time' is $-\ell_{ii}$ (by definition of the exponential distribution).

Note: the spectrum of the matrix P_h is contained within the unit disk, i.e. for every eigenvalue λ of P_h, $|\lambda| \leq 1$. This property is a consequence of P_h being row-stochastic, i.e. that $\sum_j P_{h,ij} = 1$. Since $P_h = \exp(hL)$ it follows that

$$\sigma(P_h) \subset D := \left\{x \in \mathbb{R}^2 \mid |x| \leq 1\right\} \Leftrightarrow \sigma(L) \subset \mathbb{C}^- = \{y \in \mathbb{C} \mid \mathrm{Re}(y) \leq 0\}$$

Example 1.1. *Suppose one has a reversible reaction in which one has a large collection of N molecules of the same substance. The molecules can be either in state A or state B and the molecules can change between the two states. Let k^+ denote the rate of the reaction in which molecules change from state A to B and let k^- denote the rate at which molecules change from state B to A.*

For $t > 0$, consider the quantity

$$\mu_i^A(t) := \mathbb{P}\left(\textit{number of molecules in state } A \textit{ at time } t \textit{ is } i\right)$$

where $i = \{0, \ldots, N\}$. One can define quantities $\mu_i^B(t)$ in a similar way, and one can construct balance laws for these quantities, e.g.

$$\frac{d\mu_i^A(t)}{dt} = k^+\mu_{i+1}^A(t) + k^-\mu_{i-1}^A(t) - (k^+ + k^-)\mu_i^A(t).$$

The above balance law can be written in vector notation using a tridiagonal matrix L. By adding an initial condition one can obtain an initial value problem

$$\frac{d\mu^A(t)}{dt} = L^\top \mu^A(t), \quad \mu^A(0) = \mu_0.$$

The solution of the initial value problem above is

$$\mu^A(t) = \mu_0 \exp\left(tL^\top\right).$$

1.1.3 Stochastic differential equations (SDEs)

These are time-continuous, continuous state space Markov chains. SDEs may be considered to be ordinary differential equations (ODEs) with an additional noise term (cf. Figure 2). Let $b : \mathbb{R}^n \to \mathbb{R}^n$ be a smooth vector field and let $x(t)$ be a deterministic dynamical system governed by the vector field $b(\cdot)$. Then $x(t)$ evolves according to

$$\frac{dx}{dt} = b(x), \quad x(0) = x_0. \tag{1}$$

Now let $(B_t)_{t>0}$ be Brownian motion in $\mathbb{R}^d$, and let $(X_t)_{t>0}$ be a dynamical system in $\mathbb{R}^d$ which evolves according to the equation

$$\frac{dX_t}{dt} = b(X_t) + \frac{dB_t}{dt}. \tag{2}$$

The additional term $\frac{dB_t}{dt}$ represents 'noise', or random perturbations from the environment, but is not well-defined because the paths of Brownian motion are nowhere differentiable. Therefore, one sometimes writes

$$dX_t = b(X_t)dt + dB_t,$$

which is shorthand for

$$X_t = X_0 + \int_0^t b(X_t)dt + \int_0^t dB_t.$$

The most common numerical integration method for SDEs is the forward Euler method. If x is a C^1 function of time t, then

$$\left.\frac{dx}{dt}\right|_{t=s} = \lim_{h\to 0} \frac{x(s+h) - x(s)}{h}.$$

The forward Euler method for ODEs of the form (1) is given by

$$X_{t+h} = X_t + hb(X_t)$$

and for SDEs of the form (2) it is given by

$$X_{t+h} = X_t + hb(X_t) + \xi_h$$

where $0 < h \ll 1$ is the integration time step and the noise term ξ in the Euler method for SDEs is modeled by a mean-zero Gaussian random variable.

For stochastic dynamical systems which evolve according to SDEs as in (2), one can consider the probability that a system at some point $x \in \mathbb{R}^d$ will be in a set $A \subset \mathbb{R}^d$ after a short time $h > 0$:

$$\mathbb{P}\left(X_{t+h} \in A \,|X_t = x\right).$$

The associated transition probability density functions of these stochastic dynamical systems are Gaussian because the noise term in (2) is Gaussian.

What has been the generator matrix L in case of a Markov jump process is an infinite-dimensional operator acting on a suitable Banach space. Specifically,

$$Lf(x_0) = \lim_{t\to 0} \frac{\mathbb{E}_{x_0}[f(X_t)] - f(x_0)}{t},$$

provided that the limit exists. Here $f\colon \mathbb{R}^n \to \mathbb{R}$ is any measurable function and $\mathbb{E}_{x_0}[\cdot]$ denotes the expectation over all random paths of X_t satisfying $X_0 = x_0$. L is a second-order differential operator if f is twice differentiable.

2 Day 2, 23.10.2012

Preliminaries from probability theory

Let $(\Omega, \mathcal{E}, \mathbb{P})$ be a probability space, where Ω is a set and $\mathcal{E} \subseteq 2^{\Omega}$ is a σ-field or σ-algebra on Ω, and $\mathbb{P}$ is a probability measure (i.e., $\mathbb{P}$ is a nonnegative, countably additive measure on $(\Omega, \mathcal{E})$ with the property $\mathbb{P}(\Omega) = 1$).

2.1 Conditioning

Let $A \in \mathcal{E}$ be a set of nonzero measure, i.e. $\mathbb{P}(A) > 0$ and define $\mathcal{E}_A$ to be the set of all subsets of A which are elements of $\mathcal{E}$, i.e.

$$\mathcal{E}_A := \{E \subset A \mid E \in \mathcal{E}\}.$$

Definition 2.1 (Conditional probability, part I)**.** *For an event A and an event $E \in \mathcal{E}_A$, the* conditional probability *of E given A is*

$$\mathbb{P}(E|A) := \frac{\mathbb{P}(E \cap A)}{\mathbb{P}(A)}.$$

Remark 2.2. *Think of $\mathbb{P}_A := \mathbb{P}(\cdot\,|A)$ as a probability measure on the measurable space $(A, \mathcal{E}_A)$.*

Given a set $B \in \mathcal{E}$, the *characteristic* or *indicator* function $\chi_B : \Omega \to \{0, 1\}$ satisfies

$$\chi_B(x) = \begin{cases} 1 & x \in B \\ 0 & x \notin B. \end{cases}$$

Definition 2.3 (Conditional expectation, part I)**.** *Let $X : \Omega \to \mathbb{R}$ be a random variable with finite expectation with respect to $\mathbb{P}$. The* conditional expectation *of X given an event A is*

$$\mathbb{E}(X|A) = \frac{\mathbb{E}[X\chi_A]}{\mathbb{P}(A)}.$$

Remark 2.4. *We have*

$$\mathbb{E}(X|A) = \frac{1}{\mathbb{P}(A)} \int_A X d\mathbb{P} = \int X d\mathbb{P}_A.$$

Remark 2.5. *Observe that $\mathbb{P}(E|A) = \mathbb{E}[\chi_E|A]$.*

Up to this point we have only considered the case where A satisfies $\mathbb{P}(A) > 0$. We now consider the general case.

Definition 2.6 (Conditional expectation, part II)**.** *Let $X : \Omega \to \mathbb{R}$ be an integrable random variable with respect to $\mathbb{P}$ and let $\mathcal{F} \subset \mathcal{E}$ be any sub-sigma algebra of $\mathcal{E}$. The* conditional expectation of X given $\mathcal{F}$ *is a random variable $Y := \mathbb{E}[X|\mathcal{F}]$ with the following properties:*

- *Y is measurable with respect to $\mathcal{F}$: $\forall B \in \mathcal{B}(\mathbb{R})$, $Y^{-1}(B) \in \mathcal{F}$.*
- *We have*

$$\int_F X d\mathbb{P} = \int_F Y d\mathbb{P} \quad \forall F \in \mathcal{F}.$$

Remark 2.7. *The second condition in the last definition amounts to the projection property as can be seen by noting that*

$$\mathbb{E}\left[X\chi_F\right] = \int_F X d\mathbb{P} = \int_F Y d\mathbb{P} = \mathbb{E}\left[Y\chi_F\right] = \mathbb{E}\left[\mathbb{E}\left[X|\mathcal{F}\right]\chi_F\right].$$

By the Radon-Nikodym theorem [MS05], the conditional expectation exists and is unique up to $\mathbb{P}$-null sets.

Definition 2.8 (Conditional probability, part II)**.** *Define the conditional probability of an event $E \in \mathcal{E}$ given A by $\mathbb{P}(E|A) := \mathbb{E}\left[\chi_E|A\right]$*

Exercise 2.9. *Let X, $Y : \Omega \to \mathbb{R}$ and scalars $a, b \in \mathbb{R}$. Prove the following properties of the conditional expectation:*

- *(Linearity):*
$$\mathbb{E}\left[aX + bY|A\right] = a\mathbb{E}\left[X|A\right] + b\mathbb{E}\left[Y|A\right].$$
- *(Law of total expectation):*
$$\mathbb{E}\left[X\right] = \mathbb{E}\left[X|A\right] + \mathbb{P}(A) + \mathbb{E}\left[X|A^c\right]\mathbb{P}(A^c)$$
- *(Law of total probability):*
$$\mathbb{P}(B) = \mathbb{P}(B|A)\mathbb{P}(A) + \mathbb{P}(B|A^c)\mathbb{P}(A^c).$$

Example 2.10. *The following is a collection of standard examples.*

- *Gaussian random variables: Let X_1, X_2 be jointly Gaussian with distribution $N(\mu, \Sigma)$, where*
$$\mu = \begin{pmatrix}\mathbb{E}[X_1]\\ \mathbb{E}[X_2]\end{pmatrix}, \quad \Sigma = \begin{pmatrix}a & b\\ b & c\end{pmatrix}$$
such that Σ is positive definite. The density of the distribution is
$$\rho(x) = \frac{1}{\sqrt{det(2\pi\Sigma)}}\exp\left[-\frac{1}{2}\left(x-\mu\right)^\top \Sigma\left(x-\mu\right)\right]$$
(Ex.: Compute the distribution of X_1 given that $X_2 = a$ for some $a \in \mathbb{R}$.)
- *(Conditioning as coarse-graining): Let $Z = \{Z_i\}_{i=1}^M$ be a partition of Ω, i.e. $\Omega = \cup_{i=1}^M Z_i$ with $Z_i \cap Z_j = \emptyset$ and define*
$$Y(\omega) = \sum_{i=1}^M \mathbb{E}\left[X\ |Z_i\right]\chi_{Z_i}(\omega).$$
Then $Y = \mathbb{E}\left[X|Z\right]$ is a conditional expectation (cf. Figure 3)
- *(Exponential waiting times): exponential waiting times are random variables $T : \Omega \to [0, \infty)$ with the* memoryless property*:*
$$\mathbb{P}\left(T > s + t\ |T > s\right) = \mathbb{P}\left(T > t\right).$$
This property is equivalent to the statement that T has an exponential distribution, i.e. that $\mathbb{P}(T > t) = \exp(-\lambda t)$ for a parameter value $\lambda > 0$.

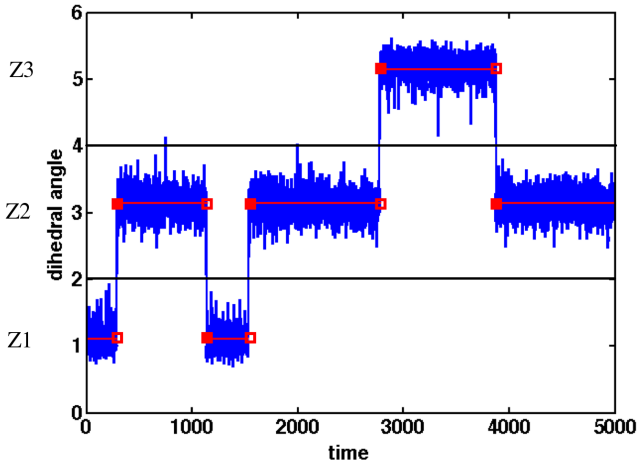

Figure 3: Simulation of butane, coarse-grained into three states Z_1, Z_2, Z_3.

2.2 Stochastic processes

Definition 2.11 (Stochastic process)**.** *A* stochastic process $X = \{X_t\}_{t\in I}$ *is a collection of random variables on a probability space* $(\Omega, \mathcal{E}, \mathbb{P})$ *indexed by a parameter* $t \in I \subseteq [0, \infty)$*. We call* X

- discrete *in time if* $I \subseteq \mathbb{N}_0$
- continuous *in time if* $I = [0, T]$ *for any* $T < \infty$*.*

How does one define probabilities for X? We provide a basic argument to illustrate the possible difficulties in defining the probability of a stochastic process in an unambiguous way. By definition of a stochastic process, $X_t = X_t(\omega)$ is measurable for every fixed $t \in I$, but if one has an event of the form

$$E = \{\omega \in \Omega \mid X_t(\omega) \in [a, b] \ \forall t \in I\}$$

how does one define the probability of this event? If t is discrete, the σ-additivity of $\mathbb{P}$ saves us, together with the measurability of X_t for every t. If, however, the process is time-continuous, X_t is defined only almost surely (a.s.) and we are free to change X_t on a set A_t with $\mathbb{P}(A_t) = 0$. By this method we can change X_t on $A = \cup_{t\in I} A_t$, The problem now is that $\mathbb{P}(A)$ need not be equal to zero even though $\mathbb{P}(A_t) = 0 \ \forall t \in I$. Furthermore, $\mathbb{P}(E)$ may not be uniquely defined. So what can we do? The solution to the question of how to define probabilities for stochastic processes is to use finite-dimensional distributions or marginals.

Definition 2.12. *(Finite dimensional distributions): Fix* $d \in \mathbb{N}$*,* $t_1, \ldots, t_d \in I$*. The* finite-dimensional distributions *of the stochastic process* X *for* $(t_1, \ldots, t_d)$ *are defined as*

$$\mu_{t_1,\ldots,t_d}(B) := \mathbb{P}_{(X_{t_k})_{k=1,\ldots,d}}(B) = \mathbb{P}\left(\{\omega \in \Omega \mid (X_{t_1}(\omega), \ldots, X_{t_d}(\omega)) \in B\}\right)$$

for $B \in \mathcal{B}(\mathbb{R}^d)$*.*

Here and in the following we use the shorthand notation $\mathbb{P}_Y := \mathbb{P} \circ Y^{-1}$ to denote the *push forward* of $\mathbb{P}$ by the random variable Y.

Theorem 2.13. *(Kolmogorov Extension Theorem): Fix $d \in \mathbb{N}$, $t_1, \ldots, t_d \in I$, and let $\mu_{t_1,\ldots,t_d}$ be a* consistent *family of finite-dimensional distributions, i.e.*

- *for any permutation π of $(1, \ldots, d)$,*

$$\mu_{t_1,\ldots,t_d}(B_1 \times \ldots B_d) = \mu_{(t_{\pi(1)},\ldots,t_{\pi(d)}}(B_{\pi(1)} \times \ldots \times B_{\pi(d)})$$

- *For $t_1, \ldots, t_{d+1} \in I$, we have that*

$$\mu_{t_1,\ldots,t_{d+1}}(B_1 \times \ldots B_d \times \mathbb{R}) = \mu_{t_1,\ldots,t_d}(B_1 \times \ldots \times B_d).$$

Then there exists a stochastic process $X = (X_t)_{t\in I}$ with $\mu_{t_1,\ldots t_d}$ as its finite-dimensional distribution.

Remark 2.14. *The Kolmogorov Extension Theorem does not guarantee uniqueness, not even $\mathbb{P}$-a.s. uniqueness, and, as we will see later on, such a kind of uniqueness would not be a desirable property of a stochastic process.*

Definition 2.15. *(Filtration generated by a stochastic process X): Let $\mathcal{F} = \{\mathcal{F}_t\}_{t\in I}$ with $\mathcal{F}_s \subset \mathcal{F}_t$ for $s < t$ be a filtration generated by $\mathcal{F}_t = \sigma(\{X_s \mid s \leq t\})$ is called the filtration generated by X.*

2.3 Markov processes

Definition 2.16. *A stochastic process X is a* Markov process *if*

$$\mathbb{P}(X_{t+s} \in A \mid \mathcal{F}_s) = \mathbb{P}(X_{t+s} \in A \mid X_s) \tag{3}$$

where

$$\mathbb{P}(\cdot|X_s) := \mathbb{P}(\cdot|\sigma(X_s)),$$
$$\mathbb{P}(E|\sigma(X_s)) := \mathbb{E}[\chi_E \mid \sigma(X_s)]$$

for some event E.

Remark 2.17. *If I is discrete, then X is a Markov process if*

$$\mathbb{P}(X_{n+1} \in A \mid X_0 = x_0, \ldots, X_n = x_n) = \mathbb{P}(X_{n+1} \in A \mid X_n = x_n)$$

Example 2.18. *Consider a Markov Chain $(X_t)_{t\in\mathbb{N}_0}$ on a continuous state space $S \subset \mathbb{R}$ and let $\mathcal{S}$ be a σ-algebra on S. Let the evolution of $(X_t)_{t\in\mathbb{N}_0}$ be described by the* transition kernel *$p(\cdot,\cdot) : S \times \mathcal{S} \to [0,1]$ which gives the single-step transition probabilities:*

$$\begin{aligned} p(x, A) &:= \mathbb{P}(X_{t+1} \in A \mid X_t = x) \\ &= \int_A q(x, y)dy. \end{aligned}$$

In the above, $A \in \mathcal{B}(S)$ and $q = \frac{d\mathbb{P}}{d\lambda}$ is the density of the transition kernel with respect to Lebesgue measure. The transition kernel has the property that

$\forall x \in S$, $p(x, \cdot)$ is a probability measure on $\mathcal{S}$, while for every $A \in \mathcal{S}$, $p(\cdot, A)$ is a measurable function on S.

For a concrete example, consider the Euler-Maruyama discretization of an SDE for a fixed time step Δt,

$$X_{n+1} = X_n + \sqrt{\Delta t}\xi_{n+1}, \quad X_0 = 0,$$

where $(\xi_i)_{i\in\mathbb{N}}$ are independent, identically distributed (i.i.d) Gaussian $\mathcal{N}(0,1)$ random variables. The process $(X_i)_{i\in\mathbb{N}}$ is a Markov Chain on $\mathbb{R}$. The transition kernel $p(x, A)$ has the Gaussian transition density

$$q(x,y) = \frac{1}{\sqrt{2\pi\Delta t}} \exp\left[-\frac{1}{2}\frac{|y-x|^2}{\Delta t}\right].$$

Thus, if $X_n = x$, then the probability that $X_{n+1} \in A \subset \mathbb{R}$ is given by

$$\mathbb{P}\left(X_{n+1} \in A | X_n = x\right) = \int_A q(x,y)dy.$$

3 Day 3, 30.10.2012

Recapitulation:

- A stochastic process $X = (X_t)_{t\in I}$ is a collection of random variables $X_t : \Omega \to \mathbb{R}$ indexed by $t \in I$ (e.g. $I = [0,\infty)$) on some probability space $(\Omega, \mathcal{E}, \mathbb{P})$.
- A filtration $\mathcal{F} := (\mathcal{F}_t)_{t\in I}$ is a collection of increasing sigma-algebras satisfying $\mathcal{F}_t \subset \mathcal{F}_s$ for $t < s$. A stochastic process X is said to be *adapted to* $\mathcal{F}$ if $(X_s)_{s\le t}$ is $\mathcal{F}_t$-measurable. For example, if we define $\mathcal{F}_t := \sigma(X_s \ : \ s \le t)$, then X is adapted to $\mathcal{F}$.
- The probability distribution of a random variable X is given in terms of its finite dimensional distributions.

Example 3.1 (Continued from last week)**.** *Let $I = \mathbb{N}_0$ and consider a sequence $(X_n)_{n\in\mathbb{N}_0}$ of random variables $X_n = X_n^{\Delta t}$ governed by the relation*

$$X_{n+1}^{\Delta t} = X_n^{\Delta t} + \sqrt{\Delta t}\xi_{n+1}, \quad X_0^{\Delta t} = 0 a.s. \tag{4}$$

where $\Delta t > 0$, and $(\xi_k)_{k\in\mathbb{N}_0}$ are i.i.d random variables with $\mathbb{E}[\xi_k] = 0$ and $\mathbb{E}\left[\xi_k^2\right] = 1$ (not necessarily Gaussian). To obtain a continuous-time stochastic process, the values of the stochastic process on non-integer time values may be obtained by linear interpolation (cf. Figure 4 below). We want to consider the limiting behaviour of the stochastic process in the limit as Δt goes to zero. Set $\Delta t = t/N$ for a fixed terminal time $t < \infty$ and let $N \to \infty$ ($\Delta t \to 0$). Then, by the central limit theorem,

$$X_N^{\Delta t} = \sqrt{\frac{t}{N}} \sum_{k=1}^{N} \xi_k \rightharpoonup \sqrt{t}Z \tag{5}$$

where $Z \sim \mathcal{N}(0,1)$, and "$\rightharpoonup$" means "convergence in distribution", i.e., weak convergence of the induced probability measure; equivalently, the limiting random variable is distributed according to $\mathcal{N}(0,t)$. In other words the limiting

distribution of the random variable $X_N^{\Delta t}$ *for fixed* $t = N\Delta t$ *is the same as the distribution of a centered Gaussian random variable with variance* t*. As this is true for any* $t > 0$*, we can think of the limiting process as a continuous-time Markov process* $B = (B_t)_{t>0}$ *with Gaussian transition probabilities,*

$$\begin{aligned}\mathbb{P}\left(B_{t+s} \in A \,|B_s = x\right) &= \int_A q_{s,t}(x,y)dy \\ &= \frac{1}{\sqrt{2\pi|t-s|}} \int_A \exp\left(-\frac{|y-x|^2}{2|t-s|}\right) dy.\end{aligned}$$

The stochastic process B *is* homogeneous *or* time-homogeneous *because the transition probability density* $q_{s,t}(\cdot,\cdot)$ *does not depend on the actual values of* t *and* s*, but only on their difference, i.e.,*

$$q_{s,t}(\cdot,\cdot) = \tilde{q}_{|s-t|}(\cdot,\cdot) \tag{6}$$

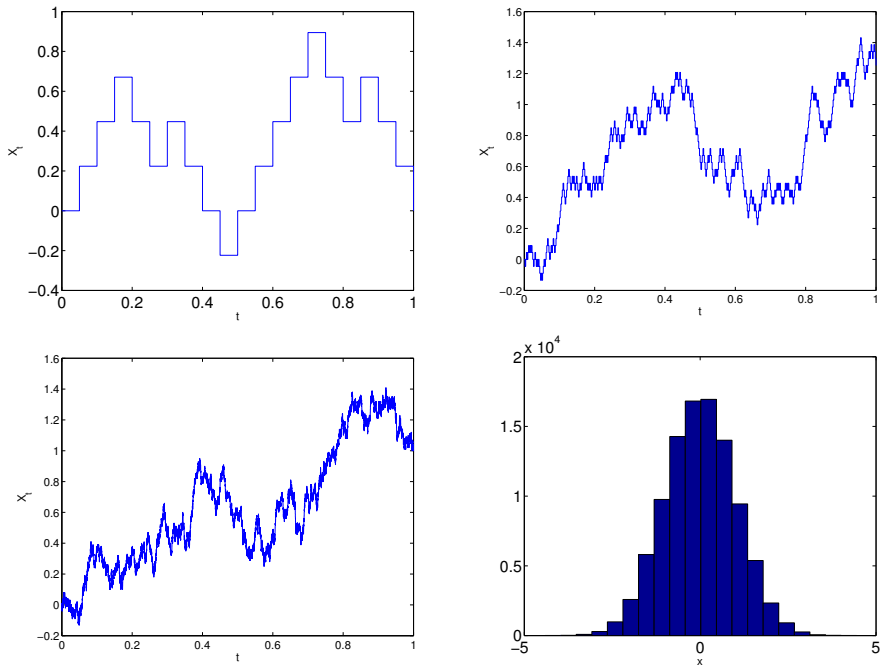

Figure 4: Sample paths of $(X_n^{\Delta t})_n$ for $\Delta t = 0.05,\ 0.002,\ 0.001$ over the unit time interval $[0,1]$, with piecewise constant interpolation. The lower right plot shows the histogram (i.e., the unnormalized empirical distribution) of $(X_{1000}^{\Delta t})$ at time $t = 1$, averaged over 10 000 independent realizations.

Remark 3.2. *The choice of exponent* $1/2$ *in* $\sqrt{\Delta t} = (\Delta t)^{1/2}$ *in (5) is unique. For* $(\Delta t)^\alpha$ *with* $\alpha \in (0, \frac{1}{2})$*, the limit of* $X_n^{\Delta t}$ *"explodes" in the sense that the variance of the process blows up, i.e.,* $\mathbb{E}[(X_N^{\Delta t})^2] \to \infty$ *as* $N \to \infty$*. On the other hand, for* $(\Delta t)^\alpha$ *with* $\alpha > 1/2$*,* $X_N^{\Delta t} \to 0$ *in probability as* $N \to \infty$*.*

3.1 Brownian motion

Brownian motion is named after the British botanist, Robert Brown (1773-1858), who first observed the random motion of pollen particles suspended in water. Einstein called the Brownian process "Zitterbewegung" in his 1905 paper, *Über die von der molekularkinetischen Theorie der Wärme gefordete Bewegung von in ruhenden Flüssigkeiten suspendierten Teilchen.* The Brownian motion is a continuous-time stochastic process which is nowhere differentiable. It is also a *martingale* in the sense that on average, the particle stays in the same location at which it was first observed. In other words, the best estimate of where the particle will be after a time $t > 0$ is its initial location.

Definition 3.3. *(Brownian motion) The stochastic process $B = (B_t)_{t>0}$ with $B_t \in \mathbb{R}$ is called the 1-dimensional* Brownian motion *or the 1-dimensional* Wiener process *if it has the following properties:*

(i) $B_0 = 0$ $\mathbb{P}$*-a.s.*

(ii) B *has* independent increments*, i.e., for all $s < t$, $(B_t - B_s)$ is a random variable which is independent of B_r for $0 \leq r \leq s$.*

(iii) B *has* stationary, Gaussian increments*, i.e., for $t > s$ we have*[1]

$$B_t - B_s \overset{D}{=} B_{t-s} \tag{7a}$$

$$\overset{D}{=} \mathcal{N}(0, t-s). \tag{7b}$$

(iv) Trajectories of Brownian motion are continuous functions of time.

Definition 3.4. *(Gaussian process) A 1-dimensional process $G = (G_t)_{t>0}$ is called a* Gaussian process *if for any collection $(t_1, \ldots, t_m) \subset I$ for arbitrary $m \in \mathbb{N}_0$, the random variable $(G_{t_1}, \ldots, G_{t_m})$ has a Gaussian distribution, i.e. it has a density*

$$f(g) = \frac{1}{\sqrt{det(2\pi\Sigma)}} \exp\left[-\frac{1}{2}(g-\mu)^\top \Sigma^{-1}(g-\mu)\right] \tag{8}$$

where $g = (g_1, \ldots, g_m)$, $\mu \in \mathbb{R}^m$ is a constant vector of means and $\Sigma = \Sigma^\top \in \mathbb{R}^{m\times m}$ is a symmetric positive semi-definite matrix.

Remark 3.5. *The Brownian motion process is a Gaussian process with the vector of means $\mu = 0$ and covariance matrix*

$$\Sigma = \begin{pmatrix} t_1 & 0 & \ldots & 0 \\ 0 & t_2 - t_1 & \ddots & \vdots \\ \vdots & \ddots & \ddots & 0 \\ 0 & \ldots & 0 & t_m - t_{m-1} \end{pmatrix} \tag{9}$$

The covariance matrix is diagonal due to the independence of the increments of Brownian motion.

[1] The notation "$X \overset{D}{=} Y$" means "X has the same distribution as Y".

Remark 3.6. *Some further remarks are in order.*

(a) Conditions (i)-(iii) define a consistent family of finite-dimensional distributions. Hence, the existence of the process B is guaranteed by the Kolmogorov Extension Theorem.

(b) Conditions (i)-(iii) imply that $\mathbb{E}[B_t] = 0$ and $\mathbb{E}[B_t B_s] = \min(t,s)$ $\forall s,t \in \mathbb{R}$. The proof is left as an exercise.

(c) The discrete process $(X_n^{\Delta t})_{n\in\mathbb{N}_0}$ converges in distribution to a Brownian motion $(B_t)_{t\geq 0}$ if the time discrete is linearly interpolated between two successive points. In other words, if we consider the continuous-time stochastic processes $(X_t^{\Delta t})_{t>0}$ (which is obtained by linear interpolation between the $X_N^{\Delta t}$) and B as random variables on the space of continuous trajectories ($C(\mathbb{R}_+)$ and $\mathcal{B}(C(\mathbb{R}_+))$), then the process $(X_t^{\Delta t})_{t>0}$ converges in distribution to B.

(d) We have that

$$\begin{aligned}\mathbb{E}\left[(B_t - B_s)^2\right] &= \mathbb{E}\left[(B_{t-s})^2\right] \text{ by (7a) in Definition 3.3}\\ &= |t-s| \text{ by (7b) in Definition 3.3.}\end{aligned}$$

(e) Brownian motion enjoys the following scaling invariance, also known as self-similarity of Brownian motion*: for every $t > 0$ and $\alpha > 0$,*

$$B_t \overset{D}{=} \alpha^{-1/2} B_{\alpha t}.$$

An alternative construction of Brownian motion

Observe that we have constructed Brownian motion by starting with the scaled random walk process and using the Kolmogorov Extension Theorem. Now we present an alternative method for constructing Brownian motion that is useful for numerics, called the Karhunen-Loève expansion of Brownian motion. We will consider this expansion for Brownian motion on the unit time interval $[0,1]$.

Let $\{\eta_k\}_{k\in\mathbb{N}}$ be a collection of independent, identically distributed (i.i.d) Gaussian random variables distributed according to $\mathcal{N}(0,1)$, and let $\{\phi_k(t)\}_{k\in\mathbb{N}}$ be an orthonormal basis of

$$L^2([0,1]) = \left\{ u : [0,1] \to \mathbb{R} \,:\, \int_0^1 |u(t)|^2 dt < \infty \right\}. \tag{10}$$

By construction, the basis functions satisfy

$$\langle \phi_i, \phi_j \rangle = \int_0^1 \phi_i(t)\phi_j(t)dt = \delta_{ij}\,,$$

and we can represent any function $\forall f \in L^2([0,1])$ by

$$f(t) = \sum_{k\in\mathbb{N}} \alpha_k \phi_k(t)$$

for $\alpha_k = \langle f, \phi_k \rangle$. We have the following result.

Theorem 3.7. *(Karhunen-Loève): The process* $(W_t)_{0\leq t\leq 1}$ *defined by*

$$W_t = \sum_{k\in\mathbb{N}} \eta_k \int_0^t \phi_k(s)ds \tag{11}$$

is a Brownian motion.

Proof. We give only a sketch of the proof. For details, see the Appendix in [MS05], or [KS91]). The key components of the proof are to show the following:

(i) The infinte sum which defines the Karhunen-Loève expansion is absolutely convergent, uniformly on $[0, 1]$.

(ii) It holds that $\mathbb{E}[W_t] = 0$ and $\mathbb{E}[W_t W_s] = \min(s, t)$.

□

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