

Non-linear (PI)DEs and affine processes

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- ① Introduction: non-linear PIDEs in Finance
- ② Numerical methods for (non-)linear PIDEs
- ③ Branching from an affine point of view
- ④ Beyond generalized Riccati equations

Credit valuation problems, utility optimization problems, super-hedging problems, American option problems lead to non-linear PIDEs.

For instance the pricing of an American Option with payoff $g : D \rightarrow \mathbb{R}$ and time horizon $T > 0$ with respect to a Markovian model X on state space D leads to a value function $v : D \times [0, T] \rightarrow \mathbb{R}$ satisfying

$$(\partial_t + \mathcal{L})v(x, t) = \mathbf{1}_{\{v(x,t) < g(x)\}} \mathbf{1}_{\{(\partial_t + \mathcal{L})v(x,t) > 0\}} \mathcal{L}g(x)$$

with boundary condition $v(x, T) = g(x)$ for $x \in D$. Here \mathcal{L} denotes the generator of the Markov process X .

In other words: $(v(X_t, t))_{0 \leq t \leq T}$ is the minimal super-martingale above $g(X_t)_{0 \leq t \leq T}$.

Linear PIDEs of the form

$$(\partial_t + \mathcal{L})u(x, t) = 0$$

with boundary condition $u(x, T) = g(x)$ allow for (Q)MC algorithms, i.e. for every $x \in D$ the value function has a stochastic representation of the form

$$u(x, t) = \mathbb{E}_{(x,t)}[g(X_T)] .$$

This expectation can be approximated by pseudo- or quasi-random samples of the Markov process X through

$$\frac{1}{N} \sum_{i=1}^N g(X_T^{(i)}).$$

This is a robust, universal and almost dimension-free method to approximate the solution u if the increments of the Markov process X can be simulated with low complexity, even though slow, i.e. with rate of convergence $1/\sqrt{N}$ which bounds the error in a probabilistic sense.

Alternatives are finite difference or finite element methods with higher convergence rates bounding the error in a deterministic way, however, their complexity depends heavily on dimension.

Non-linear PIDEs of the form

$$(\partial_t + \mathcal{L})u(x, t) + F(u(t, x)) - u(t, x) = 0$$

with boundary condition $u(x, T) = g(x)$ allow for branching Markov process representations for certain types of non-linearities F .

Generically it holds true that

$$u(t, x) = \mathbb{E}_{(x,t)} [\exp(-(T-t))g(X_T)] + \int_t^T \mathbb{E}_{(x,t)} [\exp(-(s-t))F(u(s, X_s))] ds$$

by the previous representation property. However, this is not a stochastic representation but rather a fixed point equation. Inserting the equation into itself leads towards a backwards algorithm or – under certain assumptions on F – towards a branching tree representation.

Assume that F is of the form

$$F(u) = \sum_{k=0}^M p_k u^k$$

with $p_k \geq 0$ and $\sum p_k = 1$, then the previous fixed point equation

$$\begin{aligned} u(t, x) = & \mathbb{E}_{(x,t)} [\exp(-(T-t))g(X_T)] + \\ & + \int_t^T \mathbb{E}_{(x,t)} [\exp(-(s-t)) \sum_{k=0}^M p_k (u(s, X_s))^k] ds \end{aligned}$$

leads to the short time asymptotics

$$u(t, x) = \exp(-(T - t))\mathbb{E}_{(x,t)}[g(X_T)] + \\ + (1 - \exp(-(T - t))) \sum_{k=0}^M p_k \prod_{j=1}^k \mathbb{E}_{(x,t)}[u(T, X_T^{(j)})] + o((T - t))$$

where $X^{(j)}$ denote independent copies of the Markov process X . We can now concatenate the short time asymptotics, since the expansion does not depend on u anymore.

This leads to a *branching Markov process representation*, i.e. a Markov process whose state space at time t is an integer number k of individuals in state $(x^1, \dots, x^k) \in D^k$. The particles move independently subject to the Markov process X and they die at an exponential time with parameter 1 each after giving birth to a l individuals with probability p_l (which is called branching).

The number of particles in a measurable subset $A \subset D$ is an integer-measure-valued, self-exciting affine process. Let us denote the overall number of particles at time T by N_T .

Forward stochastic representation for semi-linear PIDEs

A similar consideration as before leads to the following stochastic representation formula

$$u(x, t) = \mathbb{E}_{(x,t)} \left[\prod_{j=1}^{N_T} g(X_T^{(j)}) \times \prod_{k=0}^M \left(\frac{a_k}{p_k} \right)^{\#\{\text{branchings of type } k\}} \right].$$

for equations with generic non-linearity

$$F(u) = \sum_{k=0}^M a_k u^k - u$$

and *auxiliary branching mechanism* $p_0, \dots, p_M > 0$, $\sum p_k = 1$ governing the underlying branching process.

This is a stochastic representation by a Markov process, infinite dimensional though, which can be simulated forward such as in the linear case. The result goes back to Henry-Labordere-Touzi-Wang, but roots in works of Dynkin, McKean, LeJan, Sznithman, etc.

At this point several questions arise:

- The distribution of individuals corresponds to an integer-measure-valued affine process: how is it possible that the branching measure can be replaced by a signed measure?
- Is the previous construction passing from measures to signed measures generic for affine processes on general state spaces?
- What is the relation of the auxiliary process and the functional solving the equation?

To demonstrate the essential structure let us assume for a moment $\mathcal{L} = 0$, then we are dealing with a branching process with just constant state.

We consider a Cole-Hopf transform together with time reversal $u_t = \exp(\psi_{T-t})$, then

$$\partial_t \psi_t \exp(\psi_t) = \sum_{k=0}^M p_k \exp(k\psi_t) - \exp(\psi_t)$$

with initial value $\psi_0 = \log g$. Let us denote by ν the law which takes the value $k - 1$ (one ancestor dies!) with probability p_k , for $k = 1, \dots, m$, then

$$\partial \psi_t = \mathcal{R}(\psi_t)$$

with

$$\mathcal{R}(f) := \int (\exp(f\xi) - 1) \nu(d\xi).$$

This is a generalized Riccati equation for a one-dimensional self-exciting affine (actually linear) process, namely the number of individuals at a certain point in time in the branching process picture..

- stochastically continuous, time homogenous Markov processes N with state space \mathcal{M} (usually a cone which is not necessarily pointed).
- the Fourier-Laplace transform of the marginal distribution of N is of exponential affine form

$$\mathbb{E}_{(n,0)}[\exp(\langle N_t, f \rangle)] = \exp(\phi(t, f) + \langle \psi(t, f), n \rangle)$$

where the functions ϕ and ψ satisfy so called generalized Riccati equations

$$\partial_t \phi_t = \mathcal{F}(\psi_t) \text{ and } \partial_t \psi_t = \mathcal{R}(\psi_t)$$

with $\phi(0, f) = 0$ and $\psi(0, f) = f$.

The vector fields \mathcal{F} and \mathcal{R} are of Lévy-Khintchine form. The classification of the specific form is fully understood on symmetric cones, $\mathbb{R}^m \times \mathbb{R}^n$, etc, leading to the so called admissibility conditions, *BUT* the Lévy measures are of course never signed. This can be found in the works of Cuchiero, Duffie, Filipovic, Keller-Ressel, Mayerhofer, Schachermayer, etc.

As often in mathematics it is fruitful to turn a point of view around:

- affine processes gained importance since their marginal distribution is known up to the solution of two ODEs, the generalized Riccati equations. Often the solutions are explicitly known.
- one can also apply affine processes to *represent stochastically* the solution of non-linear ODEs, which means in particular that one obtain (Q)MC algorithms for the solution of non-linear ODEs of the generalized Riccati type.

The fact that in the Henry-Labordere-Touzi-Wang representation signed Lévy measures appear must have a meaning in the world of affine processes. The result of Henry-Labordere-Touzi-Wong suggests that there are more ODE types than generalized Riccati ones, which allow for representations.

A general result on non-linear ODEs

Consider a state space $\mathcal{M} \subset \mathbb{R}^d$ and four vectors of Lévy measures $\nu_+^{re}, \nu_-^{re}, \nu_+^{im}, \nu_-^{im}$ corresponding to the characteristic vector fields $\mathcal{R}_{\pm}^{r/i}$. Only ν_+^{re} is a generic Lévy measure of finite variation, all the others are assumed to be of finite activity. We assume the constant part \mathcal{F} to vanish here since it is not important for the argument to come.

Assume furthermore that the sum over all measures

$$\nu = \nu_+^{re} + \nu_-^{re} + \nu_+^{im} + \nu_-^{im}$$

satisfies the admissibility conditions and describes a self-exciting pure jump affine (actually linear) process N taking values in \mathcal{M} . Then one can construct a second affine process \tilde{N} , actually a pure jump linear process, with state space $\mathcal{M} \times \mathbb{Z}^{2d}$ and corresponding Lévy measures $\tilde{\nu}$ again being decomposable in four measures, too.

$$\tilde{\nu} = \tilde{\nu}_+^{re} + \tilde{\nu}_-^{re} + \tilde{\nu}_+^{im} + \tilde{\nu}_-^{im}$$

Fix $i = 1, \dots, d$: coordinate i of the measure $\tilde{\nu}_+^{re}$ corresponds to the push forward along $\mathcal{M} \ni m \mapsto (m, 0, 0) \in \mathcal{M} \times \mathbb{Z}^{2d}$ of coordinate i of ν_+^{re} ; coordinate i of the measure $\tilde{\nu}_-^{re}$ corresponds to the push forward along $\mathcal{M} \ni m \mapsto (m, e_i, 0) \in \mathcal{M} \times \mathbb{Z}^{2d}$ of coordinate i of ν_-^{re} ; coordinate i of the measure $\tilde{\nu}_+^{im}$ corresponds to the push forward along $\mathcal{M} \ni m \mapsto (m, 0, e_i) \in \mathcal{M} \times \mathbb{Z}^{2d}$ of coordinate i of ν_+^{im} ; whereas coordinate i of the measure $\tilde{\nu}_-^{im}$ corresponds to the push forward along $\mathcal{M} \ni m \mapsto (m, e_i, e_i) \in \mathcal{M} \times \mathbb{Z}^{2d}$ of coordinate i of ν_-^{im} . All other jump measures necessary to fully specify the affine process \tilde{N} vanish.

The non-trivial components of the ψ function of \tilde{N} started at $(f, i\pi, \dots, i\pi, i\pi/2, \dots, i\pi/2)$ actually solve

$$\partial\psi_t = \mathcal{R}_+^{re}(\psi_t) - \mathcal{R}_i^{re}(\psi_t) + i\mathcal{R}_+^{im}(\psi_t) - i\mathcal{R}_-^{im}(\psi_t) \quad (1)$$

$$= \int (\exp(\langle \psi_t, \xi \rangle) - 1) \eta(d\xi) = \mathcal{R}(\psi_t), \quad (2)$$

where $\eta = \nu_+^{re} - \nu_-^{re} + i\nu_+^{im} - i\nu_-^{im}$ is a complex measure.

In other words loosely speaking we have stochastic representations for non-linear ODEs with vector fields being Fourier-Laplace transforms of finite complex-valued measures on a certain state space.

Notice that we can also add an additive noise W to the equation

$$d\psi_t = \mathcal{R}(\psi_t) dt + dW_t,$$

which finally leads to the stochastic representation

$$\exp(\psi^i(t, f)) = \tag{3}$$

$$= \mathbb{E}_{(e_i, 0)} \left[\exp(\langle \tilde{N}_t, (f, i\pi, i\pi/2) \rangle) \exp\left(\int_0^t \langle N_s, dW_s \rangle\right) | \sigma(W)_t \right]. \tag{4}$$