Stochastic solutions of nonlinear pde's: McKean versus superprocesses

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Stochastic solutions

- **Stochastic solution** = a stochastic process which, started from a particular point in the domain, generates after time *t* a boundary measure which, integrated over the initial condition at *t* = 0, provides the solution at *x* and time *t*.
- Example: the heat equation

$$\partial_t u(t,x) = \frac{1}{2} \frac{\partial^2}{\partial x^2} u(t,x)$$
 with $u(0,x) = f(x)$

the process is Brownian motion, $dX_t = dB_t$, and the solution

$$u(t,x) = \mathbb{E}_{x}f(X_{t}) \tag{1}$$

- The domain here is $\mathbb{R} \times [0, t)$ and the expectation value in (1) is the inner product $\langle \mu_t, f \rangle$ of the initial condition f with the measure μ_t generated by the Brownian motion at the t-boundary.
- The process should be the same for any initial condition.
- Classical results for linear pde's. Recent work in nonlinear pde's: KPP, Navier-Stokes, Poisson-Vlasov, MHD, etc.

• New exact solutions

• New numerical algorithms

Deterministic algorithms grow exponentially with the dimension d of the space, roughly N^d ($\frac{L}{N}$ the linear size of the grid). The stochastic process only grows with the dimension d.

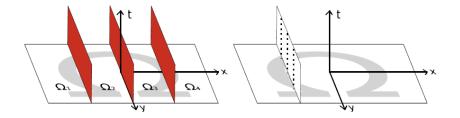
Provide localized solutions

• Sample paths started from the same point are independent. Likewise, paths starting from different points are independent from each other.

The stochastic algorithms are a natural choice for parallel and distributed computation.

- Stochastic algorithms handle equally well regular and complex boundary conditions.
- **Domain decomposition** using interpolation of localized stochastic solutions and then, in each small domain, a deterministic code. Avoids the communication time problem. Fully parallel.

Stochastic solutions and domain decomposition



Number of processors	PDD			ScaLAPACK
	T_{MC}	T_{INTERP}	T_{TOTAL}	T_{TOTAL}
64	2' 17"	<1"	3' 37"	-
128	2' 18"	<1"	2' 23"	7510' 20"
256	2' 18"	<1"	2' 18"	5223' 41"

(J. Acebrón, A. Rodríguez-Rozas, R. Spigler)

Stochastic solutions: Two construction methods

• McKean's method: a probabilistic version of the Picard series. *First* the differential equation is written as an integral equation and rearranged in a such a way that the coefficients of the successive terms in the Picard iteration obey a normalization condition *Then* the Picard iteration is interpreted as an evolution and branching process

The stochastic solution is equivalent to importance sampling of a normalized Picard series.

- **The method of superprocesses:** constructs the boundary measures of a measure-valued stochastic process and obtain the solutions of the differential equation by a scaling procedure.
- Comparison of the two methods and generalization of superprocesses (to signed measures and distribution-valued processes).

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The KPP equation: McKean's formulation

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$$\frac{\partial v}{\partial t} = \frac{1}{2} \frac{\partial^2 v}{\partial x^2} + v^2 - v$$

and initial data v(0, x) = g(x)

- $G(t, x) = \text{Green's operator for heat equation } \partial_t v(t, x) = \frac{1}{2} \frac{\partial^2}{\partial x^2} v(t, x)$ $G(t, x) = e^{\frac{1}{2}t} \frac{\partial^2}{\partial x^2}$
- KPP in integral form

$$v(t,x) = e^{-t}G(t,x)g(x) + \int_0^t e^{-(t-s)}G(t-s,x)v^2(s,x)\,ds$$
 (2)

Denoting by (ξ_t, Π_x) a Brownian motion starting from time zero and coordinate x, Eq.(2) may be rewritten

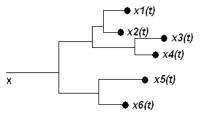
$$v(t,x) = \Pi_{x} \left\{ e^{-t}g(\xi_{t}) + \int_{0}^{t} e^{-(t-s)}v^{2}(s,\xi_{t-s}) ds \right\}$$

= $\Pi_{x} \left\{ e^{-t}g(\xi_{t}) + \int_{0}^{t} e^{-s}v_{\Box}^{2}(t-s,\xi_{s}) ds \right\}$

The KPP equation: McKean's formulation

• The stochastic solution process: a Brownian motion plus branching process with exponential holding time T, $P(T > t) = e^{-t}$. At each branching point the particle splits into two, the new particles going along independent Brownian paths. At time t > 0 one has n particles located at $x_1(t)$, $x_2(t)$, $\cdots x_n(t)$. The solution is obtained by

$$v(t,x) = \mathbb{E} \left\{ g(x_1(t)) g(x_2(t)) \cdots g(x_n(t)) \right\}$$



• An equivalent interpretation: a backwards-in-time process from time t at x. When it reaches t = 0 samples the initial condition. Generates a measure at the t = 0 boundary which is applied to $g(x) = v(0, x)_{a, c}$.

Superprocesses: Branching exit measures

- (E, \mathcal{B}) a measurable space, $M_+(E)$ the space of finite measures in Eand $(X_t, P_{0,\mu})$ a branching stochastic process with values in $M_+(E)$ and transition probability $P_{0,\mu}$ starting from time 0 and measure μ .
- \bullet The process satisfies a branching property if given $\mu=\mu_1+\mu_2$

$$P_{0,\mu} = P_{0,\mu_1} * P_{0,\mu_2}$$

that is, after the branching, (X_t^1, P_{0,μ_1}) and (X_t^2, P_{0,μ_2}) are independent and $X_t^1 + X_t^2$ has the same law as $(X_t, P_{r0,\mu})$.

• For the transition operator V_t operating on functions on E this is

$$V_t f(\mu_1 + \mu_2) = V_t f(\mu_1) + V_t f(\mu_2)$$

where $e^{-\langle V_t f, \mu \rangle} \stackrel{\circ}{=} P_{0,\mu} e^{-\langle f, X_t \rangle}$ or

$$V_{t}f\left(\mu
ight)=-\log P_{0,\mu}e^{-\left\langle f,X_{t}
ight
angle }$$

 V_t is called the *log-Laplace semigroup* associated to X_t . If the initial measure μ is δ_x one writes

$$V_t f(x) = -\log P_{0,x} e^{-\langle f_d X_t \rangle} = 0$$

Superprocesses: Branching exit measures

• In $S = [0, \infty) \times E$ consider a set $Q \subset S$ and the associated branching exit process (X_Q, P_μ) composed of a propagating Markov process in $E, \xi = (\xi_t, \Pi_{0,x})$, a set of probabilities $p_n(t, x)$ describing the branching and a parameter k defining the lifetime.

$$u(x) = V_Q f(x) = -\log P_{0,x} e^{-\langle f, X_Q \rangle}$$
(3)

 $\langle f, X_Q \rangle$ is the integral of the function f on the (space-time) boundary with the boundary exit measure generated by the process.

• This branching exit process is a (ξ, ψ) – superprocess if u(x) satisfies the equation

$$u + G_Q \psi(u) = K_Q f \tag{4}$$

where G_Q is the Green operator,

$$G_{Q}f(r,x) = \Pi_{0,x} \int_{0}^{\tau} f(s,\xi_{s}) \, ds$$

 K_Q the Poisson operator

$$K_Q f(x) = \prod_{0,x} \mathbb{1}_{\tau < \infty} f(\xi_{\hat{\tau}}) + \mathbb{P}_{\tau < \tau} + \mathbb{P}_{\tau < \tau}$$

Superprocesses: The construction

Let $\varphi(s, x; z)$ be the offspring generating function at time s and point x

$$\varphi(s,x;z) = c \sum_{0}^{\infty} p_n(s,x) z^n$$

where $\sum_{n} p_{n} = 1$ and c denotes the branching intensity.

$$P_{0,x}e^{-\langle f,X_Q\rangle} \stackrel{\circ}{=} (5)$$

$$e^{-w(0,x)} = \Pi_{0,x}\left[e^{-k\tau}e^{-f(\tau,\xi_\tau)} + \int_0^\tau dske^{-ks}\varphi\left(s,\xi_s;e^{-w(\tau-s,\xi_s)}\right)\right]$$

The measure-valued process starts from δ_x at time 0, τ is the first exit time from Q and $f(\tau, \xi_{\tau})$ the value of a function in the boundary ∂Q . Using $\int_0^{\tau} k e^{-ks} ds = 1 - e^{-k\tau}$ and the Markov property $\Pi_{0,x} \mathbb{1}_{s < \tau} \Pi_{s,\xi_s} = \Pi_{0,x} \mathbb{1}_{s < \tau}$, Eq.(5) for $e^{-w(0,x)}$ is converted into

$$e^{-w(0,x)} = \Pi_{0,x} \left[e^{-f(\tau,\xi_{\tau})} + k \int_0^{\tau} ds \left[\varphi\left(s,\xi_s; e^{-w(\tau-s,\xi_s)}\right) - e^{-w(\tau-s,\xi_s)} \right] \right]$$

Superprocesses: The limiting procedure

Replace w(0, x) by $\beta w_{\beta}(0, x)$ and f by βf . β may be interpreted as the mass of the particles and when the measure-valued process $X_Q \rightarrow \beta X_Q$ then $P_{\mu} \rightarrow P_{\frac{\mu}{\beta}}$.

$$e^{-\beta w(0,x)} = \Pi_{0,x} \left[e^{-\beta f(\tau,\xi_{\tau})} + k_{\beta} \int_{0}^{\tau} ds \left[\varphi_{\beta} \left(s, \xi_{s}; e^{-\beta w(\tau-s,\xi_{s})} \right) - e^{-\beta w(\tau-s,\xi_{s})} \right] \right]$$

Defining

$$u_{\beta} = \left(1 - e^{-\beta w_{\beta}}\right) / \beta \quad ; \quad f_{\beta} = \left(1 - e^{-\beta f}\right) / \beta$$
$$\psi_{\beta} \left(r, x; u_{\beta}\right) = \frac{k_{\beta}}{\beta} \left(\varphi \left(r, x; 1 - \beta u_{\beta}\right) - 1 + \beta u_{\beta}\right)$$

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One obtains

$$u_{\beta}(0,x) + \Pi_{0,x} \int_{0}^{\tau} ds \psi_{\beta}\left(s,\xi_{s};u_{\beta}\right) = \Pi_{0,x} f_{\beta}\left(\tau,\xi_{\tau}\right)$$

that is

$$u_{\beta} + G_{Q}\psi_{\beta}\left(u_{\beta}\right) = K_{Q}f_{\beta} \tag{6}$$

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When $\beta \to 0$, $f \to f_{\beta}$ and if ψ_{β} goes to a well defined limit ψ then u_{β} tends to a limit u, solution of (4) associated to a superprocess. Also in the $\beta \to 0$ limit

$$u_{eta}
ightarrow w_{eta} = -\log P_{0,x} e^{-\langle f, X_Q
angle}$$

If to reproduce with (6) the equation we want it must be $\beta \to 0$ and $k_{\beta} \to \infty$, the superprocess would correspond to a cloud of particles for which both the mass and the lifetime tend to zero.

The KPP equation as a superprocess

The KPP equation

$$v(t,x) = e^{-t}G(t,x)g(x) + \int_0^t e^{-s}G(t-s,x)v^2(s,x)\,ds$$

is identical to

$$P_{0,x}e^{-\langle f,X_Q\rangle} \stackrel{\circ}{=} e^{-w(0,x)} =$$

$$= \Pi_{0,x} \left[e^{-k\tau}e^{-f(\tau,\xi_{\tau})} + \int_0^{\tau} dske^{-ks}\varphi\left(s,\xi_s;e^{-w(\tau-s,\xi_s)}\right) \right]$$
with $k = 1$, $e^{-w(0,x)} = v(\tau,x)$, $e^{-f(\tau,\xi_{\tau})} = g(\xi_{\tau})$,
 $\varphi\left(s,\xi_s;e^{-w(\tau-s,\xi_s)}\right) = v^2(\tau-s,\xi_s)$.

• The McKean probabilistic construction corresponds to an intermediate step in the superprocess construction. Summing over the exit measure, the solution is

$$v(t,x) = e^{-\langle f, X_Q \rangle} = e^{-\sum_i f\left(\xi_{\tau_i}\right)} = e^{\sum_i \log g\left(\xi_{\tau_i}\right)} = \prod_i g\left(\xi_{\tau_i}\right)$$

essentially the same as before.

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The KPP equation as a superprocess

Let u(t, x) = 1 - v(t, x), which satisfies the equations

$$\frac{\partial u}{\partial t} = \frac{1}{2} \frac{\partial^2 u}{\partial x^2} - u^2 + u$$

 $u(t,x) + \Pi_{x} \int_{0}^{t} \left(u^{2} \left(t - s, \xi_{s} \right) - u \left(t - s, \xi_{s} \right) \right) ds = \Pi_{x} \left(1 - g \left(\xi_{t} \right) \right)$

that is, for KPP, $\psi(\mathbf{0}, x; u) = u^2 - u$

$$\begin{split} \psi_{\beta}\left(0,x;u_{\beta}\right) &= \frac{k_{\beta}}{\beta}\left(\varphi\left(0,x;1-\beta u_{\beta}\right)-1+\beta u_{\beta}\right) \\ &= \frac{k_{\beta}}{\beta}\left(c\sum p_{n}\left(1-\beta u_{\beta}\right)^{n}-1+\beta u_{\beta}\right) \\ &= \frac{k_{\beta}c}{\beta}\left(\beta^{2}u_{\beta}^{2}-\beta u_{\beta}\right)=u^{2}-u \end{split}$$

with $p_n = \delta_{n,2}$. Therefore $c = \beta = 1$ and $k_\beta = 1$. That is, for KPP the superprocess is not a scaling limit. It coincides with the McKean process. In this case, because $\beta = 1$ instead of $\beta \to 0$, the solution is $(1 = e^{-w})_{n \to \infty}$.

Superprocesses and nonlinear heat equation

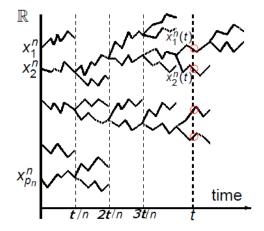
With other limiting choices for β , stochastic solutions are constructed for other equations, in particular for equations without the natural Poisson clock which is present in the KPP equation. For example for

$$\frac{\partial u}{\partial t} = \frac{1}{2} \frac{\partial^2 u}{\partial x^2} - u^2$$
$$\psi(0, x; u) = u^2$$

$$\psi_{\beta}\left(0,x;u_{\beta}\right) = \frac{k_{\beta}}{\beta} \left(\beta u_{\beta} - 1 + \sum_{n=0}^{2} p_{0} + p_{1}\left(1 - \beta u_{\beta}\right) + p_{2}\left(1 - \beta u_{\beta}\right)^{2}\right)$$
$$= u_{\beta}^{2}$$

leads to $p_1 = 0$; $p_0 = p_2 = \frac{1}{2}$; $k_\beta = \frac{2}{\beta}$ In this case, with $\beta \to 0$, the solution is given by (3) and the superprocess is a scaling limit $(n \to \infty$ in the figure) where both mass and lifetime of the particles tend to zero and at each bifurcation one has equal probability of either dying without offspring or having two children, k = k = 0

Superprocesses and nonlinear heat equation



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Superprocesses for more general interactions

The construction may be generalized for interactions u^{α} with $1 < \alpha \leq 2$. With $z = 1 - \beta u_{\beta}$ one has

$$\varphi(0, x; z) = \sum_{n} p_{n} z^{n} = z + u_{\beta}^{\alpha} = z + \frac{\beta}{k_{\beta}} \frac{(1-z)^{\alpha}}{\beta^{\alpha}} = z + \frac{1}{k_{\beta}\beta^{\alpha-1}}$$
$$\times \left(1 - \alpha z + \frac{\alpha (\alpha-1)}{2} z^{2} - \frac{\alpha (\alpha-1) (\alpha-2)}{3!} z^{3} + \cdots\right)$$

Choosing $k_{\beta} = \frac{\alpha}{\beta^{\alpha-1}}$ the terms in *z* cancel and for $1 < \alpha \le 2$ the coefficients of all the remaining *z* powers are positive and may be interpreted as branching probabilities. It would not be so for $\alpha > 2$.

$$p_0 = \frac{1}{\alpha}; \quad p_1 = 0; \quad \cdots \quad p_n = \frac{(-1)^n}{\alpha} \begin{pmatrix} \alpha \\ n \end{pmatrix} \quad n \ge 2$$

with $\sum_{n} p_{n} = 1$. With this choice of branching probabilities, $k_{\beta} = \frac{\alpha}{\beta^{\alpha-1}}$ and $\beta \to 0$ one obtains a superprocess which provides a solution to the equation $\frac{\partial u}{\partial t} = \frac{1}{2} \frac{\partial^{2} u}{\partial x^{2}} - u^{\alpha}$ for $1 < \alpha \leq 2$.

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