Efficient valuation of (non-)linear products for xVA

Felix Wolf¹ Joint work with Griselda Deelstra¹ and Lech Grzelak^{2,3}

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Expected exposure

Let $V_t = \mathbb{E}_t^{\mathbb{Q}} \Big[\sum_j \frac{B_t}{B_{T_j}} H_{T_j} \Big]$ be the discounted value of an asset (or portfolio) with payoffs H_{T_j} .

The positive exposure of V at time t is $E^+(t) = \max(0, V_t)$. At initial time t_0 we can observe the expected positive exposure at time t:

$$\mathrm{EE}(t_0, t) = \mathbb{E}_{t_0}^{\mathbb{Q}} \left[\frac{B_{t_0}}{B_t} \max(0, V_t) \right]$$

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Exposures are typically obtained by Monte Carlo simulation and the EE is obtained from the sample mean:

$$\mathrm{EE}(t_0,t) pprox rac{1}{M} \sum_{i=1}^{M} \mathrm{max}\Big(0, rac{B_{t_0}(\omega_j)}{B_{t}(\omega_j)} V_t(\omega_j)\Big).$$

Thus, we are interested in samples $V_t(\omega)$ of the random variable $V_t \mid F_{t_0}$.

Swap portfolio

Consider an interest rate (payer) swap V with price

$$V_t = \bar{N} \sum_{j=1}^m \tau_j P(t, T_j) \Big(\frac{P(t, T_{j-1}) - P(t, T_j)}{\tau_j P(t, T_j)} - K \Big).$$

In affine interest rate models, a ZCB at time t is given by

$$P(t, T) = \mathbb{E}_{t}^{\mathbb{Q}} \left[\exp\left(-\int_{t}^{T} r_{s} \, \mathrm{d}s\right) \right]$$
$$= \exp\left(A(t, T) + B(t, T)r_{t}\right),$$

where r_t is a random variable (e.g. Gaussian).

Thus, at time t_0 , V_t is a random variable in one "risk factor" r_t , following some distribution \mathcal{L} .

$$(V_t \mid \mathscr{F}_{t_0}) = (f(r_t) \mid \mathscr{F}_{t_0}) \sim \mathcal{L}$$

Approximating the portfolio

Let the portfolio Π consist of 1000 swaps $V^1,...,V^{1000}$. Simulating the price of one portfolio realisation $\Pi_t(\omega)$ requires 1000 swap evaluations:

$$\Pi_t(\omega) = V_t^1(\omega) + \cdots + V_t^{1000}(\omega) =: g(r_t(\omega)).$$

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Simplify with an approximation $\widetilde{g}_n \approx g$:

- n expensive, exact evaluations at the interpolation points: $\left(\left(r_t^1,g(r_t^1)\right),\ldots,\left(r_t^n,g(r_t^n)\right)\right)$
- (Compute the approximation)
- M cheap evaluations of the approximation $\widetilde{g}_n(r_t(\omega_j)), \ j=1,\ldots,M$.

How to interpolate between distributions?

Sample transformation

For a continuous random variable Y with cumulative distribution function F_Y , it holds $F_Y(Y) \sim \mathcal{U}[0,1]$. Proof:

$$F_{F_Y(Y)}(u) = \mathbb{P}[F_Y(Y) \le u] = \mathbb{P}[F_Y^{-1}(F_Y(Y)) \le F_Y^{-1}(u)]$$

= $\mathbb{P}[Y \le F_Y^{-1}(u)] = F_Y(F_Y^{-1}(u)) = u$
= $F_U(u)$.

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For two continuous random variables X and Y, we have $F_X(X) \sim F_Y(Y) \sim \mathcal{U}(0,1)$.

From a sample ξ of X we can obtain a sample y of Y via

$$y = F_Y^{-1}(F_X(\xi)).$$

Stochastic collocation sampling¹

X is a random variable we can easily sample from (e.g. Gaussian), Y is expensive to sample from. We can relate samples:

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This function $g := F_Y^{-1} \circ F_X$ is computationally expensive (inversion of F_Y).

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- Find interpolation points $x_1, ..., x_n$ ("collocation points") and evaluate exactly: $y_i = g(x_i), i = 1, ..., n$.
- ② Build approximation function $\widetilde{g}_n \approx g$ based on these n points.
- **3** Obtain (approximated) samples $\widetilde{y}_i = \widetilde{g}_n(\xi_i)$ of Y from (cheap) samples ξ_i of X.

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Why collocation?

"Classic interpolation" framework:

$$\Pi_t(\omega) = g(r_t(\omega)) \approx \widetilde{g}_n(r_t(\omega))$$

Collocation framework:

$$y = (F_Y^{-1} \circ F_X)(\xi) = g(\xi) \approx \widetilde{g}_n(\xi).$$

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Example Lagrange polynomial over interpolation points x_1, \ldots, x_n :

$$\widetilde{g}_n(\xi) = \sum_{i=1}^n g(x_i)\ell_i(\xi),$$

where

$$\ell_i(\xi) = \prod_{j=1, i \neq i}^n \frac{\xi - x_j}{x_i - x_j}$$

Connection to Gaussian Quadrature

Let p_i be an orthogonal, polynomial basis in $L^2(X)$, i.e.

$$\int_{a}^{b} p_{i}(x)p_{j}(x)f_{X}(x)dx = \delta_{ij}\mathbb{E}[p_{i}(X)^{2}].$$

We want to find weights w_i and collocation points x_i , i = 1, ..., n, so that

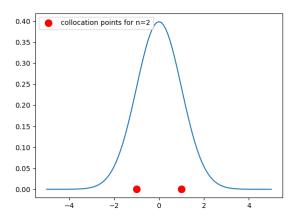
$$\int_{a}^{b} g(x)f_{X}(x)dx \approx \sum_{i=1}^{n} g(x_{i})w_{i}.$$

Find weights and points by consideration of enough exact integrals:

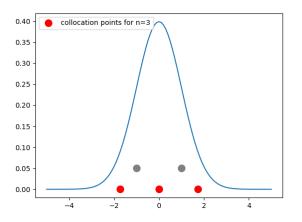
$$\int_{a}^{b} p_{0}(x)dx = w_{1}p_{0}(x_{1}) + \dots + w_{n}p_{0}(x_{n})$$
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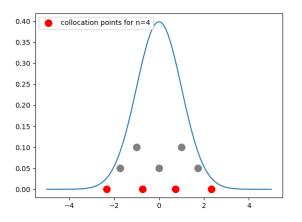
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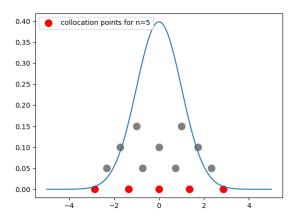
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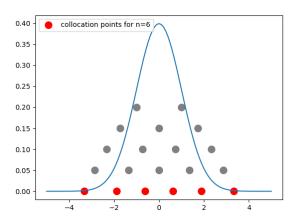
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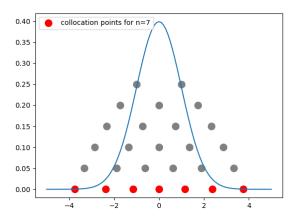
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Choice of interpolation function

The approximation \widetilde{g}_n ...

- must be cheap to evaluate:
 "n exact + M cheap evaluations ≪ M exact valuations"
- must offer high accuracy
- may preserve properties of g (e.g. monotonicity)
- may be differentiable

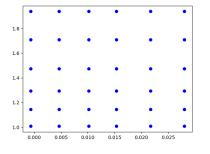
There are many options:

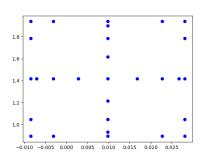
Lagrange polynomials, Chebyshev polynomials, Hermite polynomials, ...

Higher dimensions²

The number of interpolation points should not grow too fast.

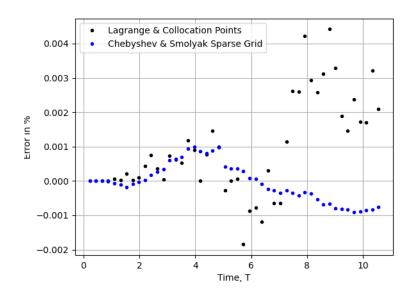
Cartesian grid of (optimal) collocation points vs. Smolyak sparse grid





 $^{^2}$ L.A. Grzelak. Sparse Grid Method for Highly Efficient Computation of Exposures for xVA. arXiv:2104.14319, 2021.

Hybrid portfolio of (many) stock contracts and swaps.



Many directions to investigate

- Error bounds for different interpolation methods (in higher dimensions)
- Interplay between interpolation points and interpolation methods
- Effects on portfolios of non-linear derivatives

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Thank you for listening!