

Numerical simulations using approximate random numbers

Oliver Sheridan-Methven

`oliver.sheridan-methven@maths.ox.ac.uk`

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Supervisors:

Prof. Michael Giles	Oxford
Dr Christopher Goodyer	Arm

Of course my code will run faster if it's vectorised.

No

Of course my code will run faster if I worked in a lower precision.

No

Well my compiler should be clever enough to decide what's best.

No

Half-precision is useless for anyone who wants accurate answers!

No

You can't design code that performs well on arbitrary vector lengths.

No



- ① What is an approximate random variable?
- ② When can we use them?
- ③ Do we still get the right answer?
- ④ What precisions can we use, and when?
- ⑤ Conclusions

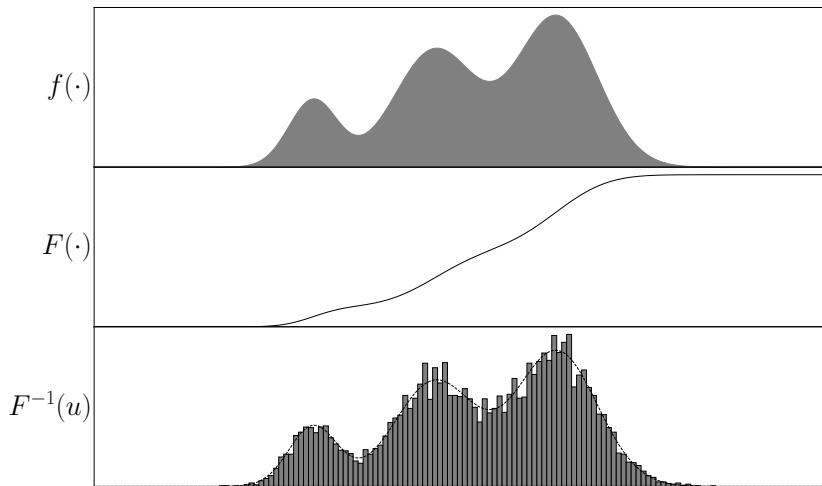
Weather Tomorrow's temperature given today's weather.

Finance Value of contract given today's prices.

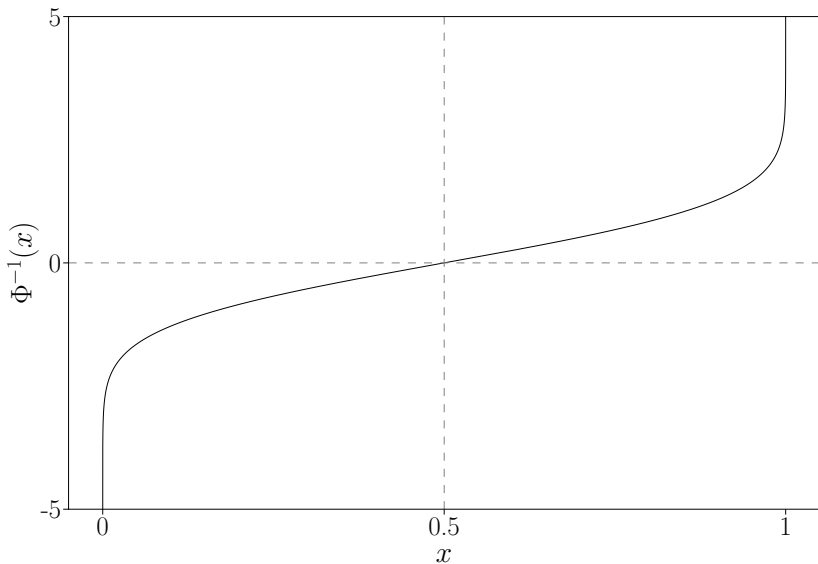
Traffic Rush hour traffic given morning congestion.

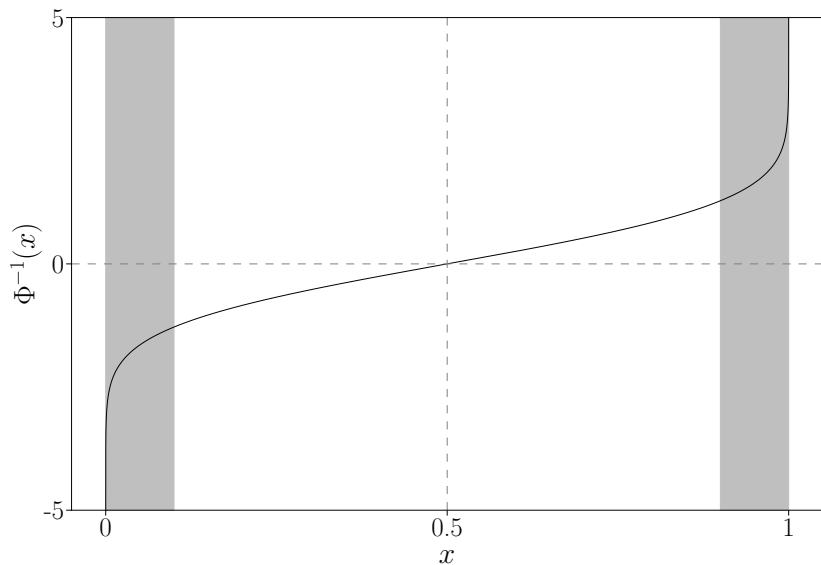
Health Risk of later secondary condition given current health.

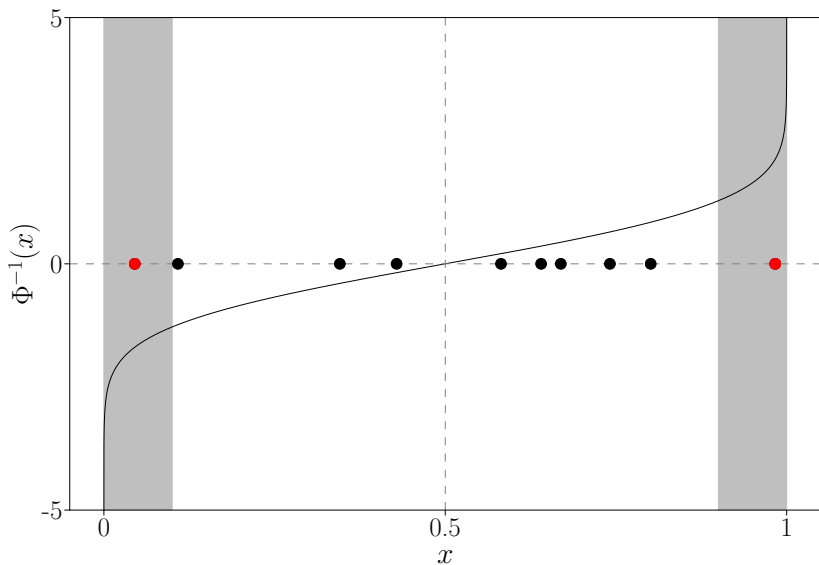
Technology Expected number of online visitors given current search trends.



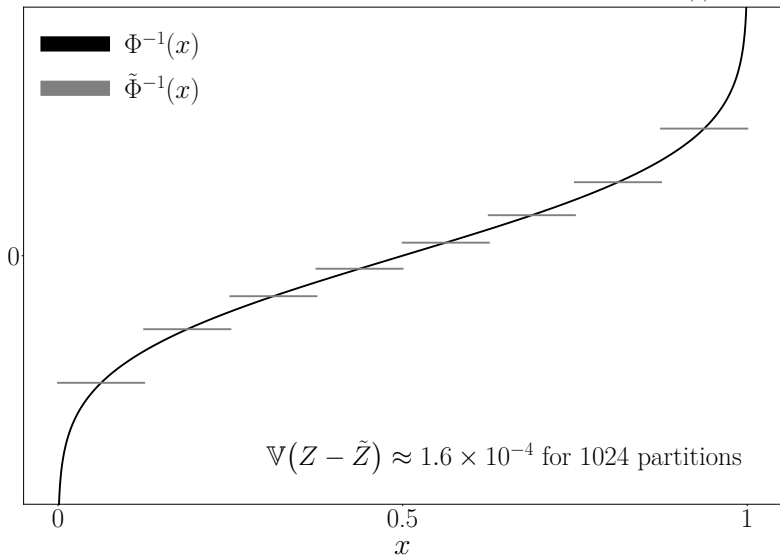
The inverse Gaussian CDF $\Phi^{-1}(\cdot)$



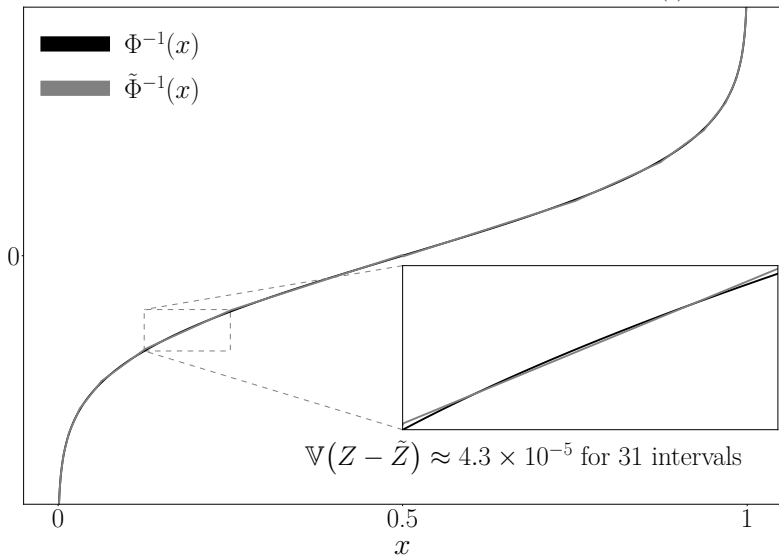


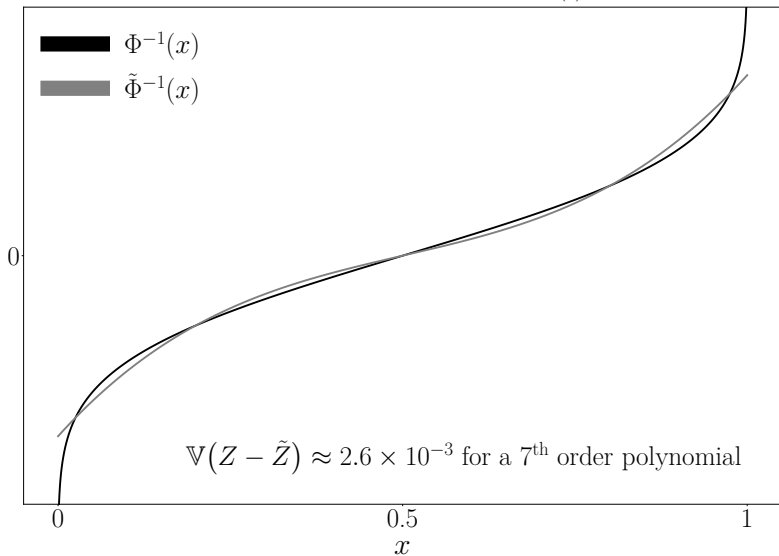


Uniform piecewise constant approximation of $\Phi^{-1}(\cdot)$



Piecewise linear dyadic approximation of $\Phi^{-1}(\cdot)$



Cubic approximation of $\Phi^{-1}(\cdot)$ 

	Average speed (clock cycles)
Intel MKL	8
Lookup table	2

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GSL (optimised)	17
Giles [1] (NVIDIA?)	12

Accuracy and precision

HIGH PRECISION
HIGH ACCURACY



LOW PRECISION
HIGH ACCURACY



HIGH PRECISION
LOW ACCURACY



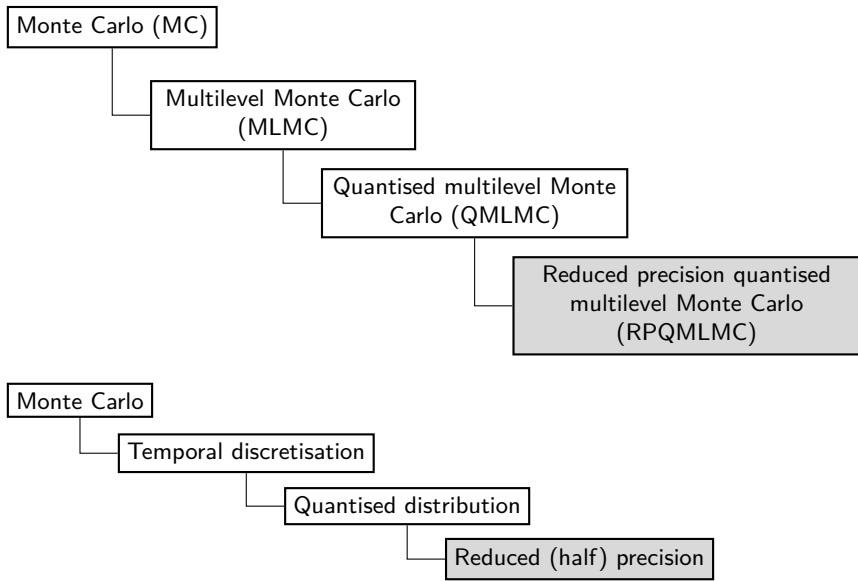
NO PRECISION
NO ACCURACY
NO SPECTATORS



J. HARRIS

$$\mathbb{E}(P) \approx \mathbb{E}(\hat{P}_{\text{Accurate}}) = \mathbb{E}(\hat{P}_{\text{Crude}}) + \mathbb{E}(\hat{P}_{\text{Accurate}} - \hat{P}_{\text{Crude}})$$

Nested multilevel Monte Carlo I



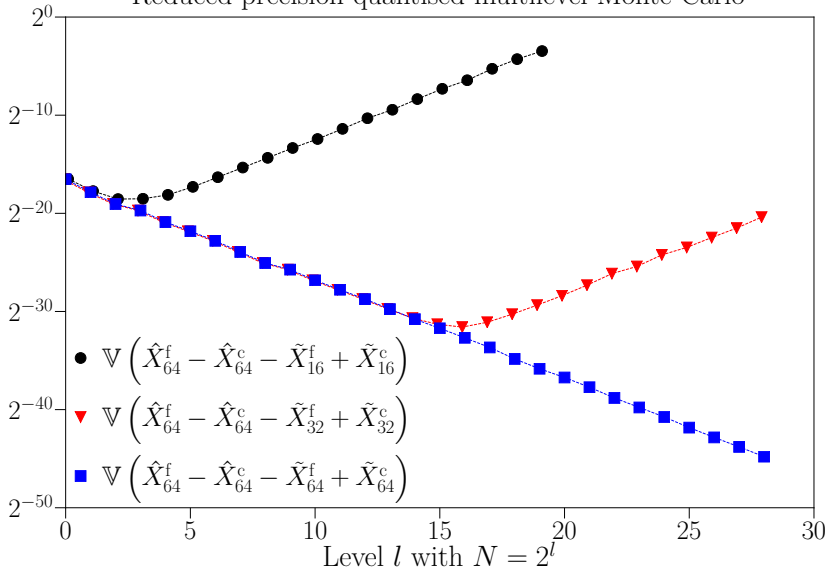
Correction = Discretisation \times Quantisation

$$\text{Correction} = \text{Discretisation} \times \text{Quantisation} \\ + \text{Roundoff}$$

Correction = Discretisation \times Quantisation

$$+ \frac{\underbrace{\text{Roundoff}}_1}{\text{Discretisation}}$$

Reduced precision quantised multilevel Monte Carlo



Running our QMLMC estimator with a single quantisation level (1024 bins) for finely discretised paths gives the following average time per path:

Relative accuracy $\varepsilon = 10^{-3}$ Times per path 10^{-4} s	Memory intensive	Work intensive
Original MLMC	24.8	17.0
Quantised MLMC	13.2	3.99

Level	paths
Original	1 920 000
Quantised	1 980 000
Correction	14 000

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↓ ×2

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$\downarrow \times 4$

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Quantised	1 980 000
Correction	14 000

Errors from using a cheap proxy distribution can be **quantified** and controlled by the introduction of a nested multilevel Monte Carlo framework.

There is a degree of freedom in the construction of this proxy. Put the results in the low level **cache**, or use a very cheap (piece-wise) **polynomial**.

The resultant approximations **converge**.

The **approximate schemes scale** as we move to wider vectors (SVE) and lower precisions (FP16), benefiting from greater SIMD parallelisation and faster FP16 calculations.

Half-precision can be used in the coarsest calculations (which consume most of the computer time). (BFloat16 may require Kahan summation...).

- [1] Mike Giles. Approximating the erfinv function. In *GPU Computing Gems, Jade Edition*, volume 2, pages 109–116. Elsevier, 2011.