Krylov Linear Solvers and Quasi Monte Carlo Methods for Transport Simulations

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Number of pages: 10 Number of tables: 4 Number of figures: 2

Abstract

 $\mathrm{QMC}\,+\,\mathrm{Krylov}$

 $\mathbf{Keywords}$ — Quasi Monte Carlo Methods, Krylov Linear Solvers

I. INTRODUCTION

II. COMPUTATIONAL RESULTS

In this section we consider an example from [1]. The formulation of the transport problem is taken from [2]. The equation for the angular flux ψ is

$$\mu \frac{\partial \psi}{\partial x}(x,\mu) + \Sigma_t(x)\psi(x,\mu) = \frac{1}{2} \left[\Sigma_s(x) \int_{-1}^1 \psi(x,\mu') \, d\mu' + q(x) \right] \text{ for } 0 \le x \le \tau$$
 (1)

The boundary conditions are

$$\psi(0,\mu) = \psi_l(\mu), \mu > 0; \psi(\tau,\mu) = \psi_r(\mu), \mu < 0.$$

II.A. Source Iteration and Linear Solvers

Source iteration is Picard iteration for the fixed point problem

$$\phi = \mathcal{S}(\phi, q, \psi_l, \psi_r)$$

To use other solvers we must convert to a linear system via

$$\mathcal{K}(\phi) = \mathcal{S}(\phi, 0, 0, 0)$$
 and $f = \mathcal{S}(0, q, \psi_l, \psi_r)$

to get

$$A\phi \equiv (I - \mathcal{K})\phi = f$$

which we can send to a linear solver.

In the computations we use the problem from [1]

$$\tau = 5, \Sigma_s(x) = \omega_0 e^{-x/s}, \Sigma_t(x) = 1, q(x) = 0, \psi_l(\mu) = 1, \psi_r(\mu) = 0,$$

and consider two cases s=1 and $s=\infty$

II.B. QMC and Krylov Linear Solvers

The linear and nonlinear solvers come from the Julia package SIAMFANLEQ.jl [3]. The documentation for these codes is in the Julia notebooks [4] and the book [5] that accompany the package.

We solve the QMC linear problem with N=1000 and Nx= 100. We use two krylov methods [6], GMRES [7] and Bi-CGSTAB [8]. Figures 1 and 2 show that the Krylov iterations take fewer than half of the number of transport sweeps that Picard iteration required.

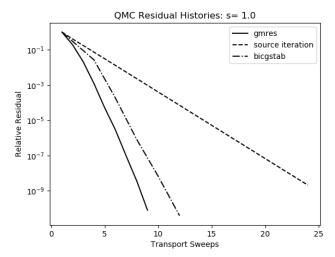


Fig. 1. s = 1

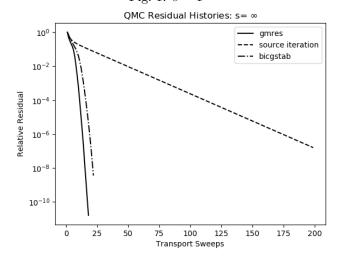


Fig. 2. $s = \infty$

II.C. Validation and calibration study

We conclude this section with a validation study. We compare the QMC results with the results from [1]. The results in [1] are exit distributions and are accurate to six figures. We have duplicated those results with an Sn computation on a fine angular and spatial mesh.

Sam, Ryan, should we use more or different values of N and Nx?

For N = 1000 and Nx = 100 we obtain the cell-average fluxes from the QMC approximation. We then use a single Sn transport sweep to recover the exit distributions from the QMC cell-average fluxes. We report the results and the corresponding results from [1] in Tables I and II.

The exit distributions, as is clear from Table I can vary by five orders of magnitude. Even so, the results from QMC agree with the benchmarks to roughly two figures.

TABLE I Exit Distributions: s = 1

	Garcia/Siewert		QMC	
$\overline{\mu}$	$\psi(0,-\mu)$	$\psi(au,\mu)$	$\psi(0,-\mu)$	$\psi(au,\mu)$
5.00e-02	5.89664 e - 01	6.07488e-06	5.71197e-01	5.85487e-06
1.00e-01	5.31120 e-01	6.92516 e- 06	5.22137e-01	6.66741 e-06
2.00e-01	4.43280 e-01	9.64232 e-06	4.41567e-01	$9.25261 \mathrm{e}\text{-}06$
3.00e-01	3.80306 e-01	1.62339e-05	3.81029 e-01	1.54416e-05
4.00e-01	3.32964 e-01	4.38580 e- 05	3.34673e-01	4.09691e-05
5.00e-01	2.96090e-01	1.69372 e-04	2.98224 e-01	1.57373e-04
6.00e-01	2.66563 e-01	5.73465e- 04	2.68871e-01	5.35989e-04
7.00e-01	2.42390 e-01	1.51282 e-03	2.44749e-01	1.42448e-03
8.00e-01	2.22235 e-01	3.24369 e-03	2.24583e-01	3.07431e-03
9.00e-01	2.05174 e-01	5.96036e-03	2.07478e-01	5.67991e-03
1.00e+00	1.90546 e-01	9.77123e-03	1.92789 e-01	9.35351 e-03

In Tables III and IV we look at the relative errors in the QMC exit distributions as compared to a highly accurate SN result. We compensate for the widely varying scales by tabulating, for each value of N and Nx

$$R = \max(R^0, R^\tau)$$

where

$$R^0 = \max_{\mu} \frac{|\psi^{SN}(0, -\mu) - \psi^{QMC}(0, -\mu)|}{\psi^{SN}(0, -\mu)}$$

and

$$R^{\tau} = \max_{\mu} \frac{|\psi^{SN}(\tau, \mu) - \psi^{QMC}(\tau, \mu)|}{\psi^{SN}(\tau, \mu)}.$$

	Garcia/Siewert		QMC		
$\overline{\mu}$	$\psi(0,-\mu)$	$\psi(au,\mu)$	$\psi(0,-\mu)$	$\psi(\tau,\mu)$	
5.00e-02	8.97798e-01	1.02202e-01	8.47454e-01	1.00663e-01	
1.00e-01	8.87836e-01	1.12164 e-01	8.52822 e-01	1.10325 e-01	
2.00e-01	8.69581 e-01	1.30419e-01	8.47710e-01	1.29064 e-01	
3.00e-01	8.52299 e-01	1.47701e-01	8.35879 e- 01	1.46849 e-01	
4.00e-01	8.35503 e-01	1.64497e-01	8.22291e-01	1.64034 e-01	
5.00e-01	8.18996 e-01	1.81004 e-01	8.08044 e-01	1.80827e-01	
6.00 e-01	8.02676 e-01	1.97324 e-01	7.93459e-01	1.97336e-01	
7.00e-01	7.86493e-01	2.13507e-01	7.78672e-01	2.13625 e-01	
8.00e-01	7.70429e-01	2.29571e-01	7.63768e-01	2.29725 e-01	
9.00e-01	7.54496e-01	2.45504 e-01	7.48818e-01	2.45642 e-01	
1.00e+00	7.38721e-01	2.61279e-01	7.33889e-01	2.61361e-01	

Ryan, for large Nx I see convergence as N increases. Is it clearly 1/N? Am I missing something? Am I tabulating the wrong thing?

$$\label{eq:table_interpolation} \begin{split} & \text{TABLE III} \\ & \text{Exit Distributions Errors: } s = 1.0 \end{split}$$

$Nx \setminus N$	1000	2000	4000	8000	16000
50	1.41162e-01	1.36428e-01	1.34747e-01	1.35736e-01	1.35577e-01
100	7.08438e-02	6.60744 e - 02	6.52017e-02	6.51605 e-02	6.49914 e-02
200	4.17171e-02	3.30480 e-02	3.23088e-02	3.21432 e-02	3.17467e-02
400	4.55590e-02	1.73115e-02	1.63072 e-02	1.61542 e-02	1.58469e-02
800	4.83754e-02	1.93087e-02	1.29178e-02	8.30117e-03	7.96562e-03
1600	5.07584e-02	2.03691e-02	1.44681e-02	4.52388e-03	4.11350e-03
3200	5.09694e-02	2.13418e-02	1.48086e-02	2.88667e-03	2.18194e-03

$Nx \setminus N$	1000	2000	4000	8000	16000
50	5.95648e-02	2.42755 e-02	1.36521e-02	1.29509e-02	1.22769e-02
100	5.60749e-02	2.31030 e-02	1.31680 e-02	6.45949 e - 03	6.59550 e-03
200	5.62864e-02	2.32524 e-02	1.42149e-02	4.77319e-03	3.55246 e-03
400	5.30954e-02	2.17854e-02	1.48225 e - 02	4.73260 e-03	2.05558e-03
800	7.66264e-02	1.88155e-02	1.60082e-02	4.34610 e-03	1.41402 e-03
1600	5.99376e-02	2.15675e-02	1.56784 e - 02	4.21636e-03	1.29138e-03
3200	5.74319e-02	1.89482 e-02	2.00195 e-02	3.26688e-03	1.49649 e - 03

III. CONCLUSION

ACKNOWLEDGMENTS

The research of CTK was supported by Department of Energy grant DE-NA003967, and National Science Foundation Grants DMS-1745654, and DMS-1906446.

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