# Krylov Linear Solvers and Quasi Monte Carlo Methods for Transport Simulations

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Number of pages: 12 Number of tables: 4 Number of figures: 3

# 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  $\psi$  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. Multigroup Equations

In general geometry the multigroup equations are

$$\mu \frac{\partial \psi_g}{\partial x}(x,\mu) + \Sigma_{t,g}(x)\psi_g(x,\mu) = \frac{1}{2} \sum_{g'=1}^G \Sigma_{s,g'\to g}(x) \int_{-1}^1 \psi_{g'}(x,\mu') d\mu' + \frac{q_g(x)}{2} \quad g = 1,\dots,G. \quad (2)$$

The boundary conditions are

$$\psi_q(0,\mu) = \psi_{l,q}(\mu), \mu > 0; \psi_q(\tau,\mu) = \psi_{r,q}(\mu), \mu < 0.$$

In matrix form, these equations are

$$\mu \frac{\partial \vec{\psi}}{\partial x}(x,\mu) + \underline{\Sigma}_t(x)\vec{\psi}(x,\mu) = \frac{1}{2}\underline{\Sigma}_s(x) \int_{-1}^1 \vec{\psi}(x,\mu') \,d\mu' + \frac{\vec{q}(x)}{2},\tag{3}$$

where

$$\vec{\psi} = (\psi_1, \psi_2, \dots, \psi_G)^{\mathrm{T}}, \qquad \vec{q} = (q_1, q_2, \dots, q_G)^{\mathrm{T}},$$
 (4)

$$\underline{\Sigma}_{t}(x) = \begin{pmatrix} \Sigma_{t,1}(x) & 0 & \dots \\ 0 & \Sigma_{t,2}(x) & 0 \dots \\ \vdots & & \ddots \\ 0 & \dots & 0 & \Sigma_{t,G}(x) \end{pmatrix},$$
 (5)

and

$$\underline{\Sigma}_{s}(x) = \begin{pmatrix} \Sigma_{s,1\to 1}(x) & \Sigma_{s,2\to 1}(x) & \dots & \Sigma_{s,G\to 1}(x) \\ \Sigma_{s,2\to 1}(x) & \Sigma_{s,2\to 1}(x) & \dots & \Sigma_{s,G\to 2}(x) \\ \vdots & \vdots & & \vdots \\ \Sigma_{s,G\to 1}(x) & \Sigma_{s,G\to 1}(x) & \dots & \Sigma_{s,G\to G}(x) \end{pmatrix},$$
(6)

## II.B. 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  $s = \infty$ 

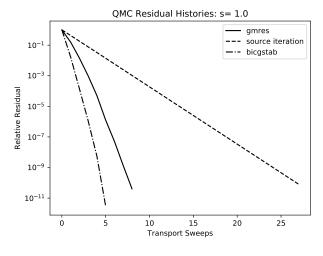


Fig. 1. s = 1

## II.C. 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=2048 and Nx= 100. We use two krylov methods [6], GMRES [7] and Bi-CGSTAB [8]. Figures 2 and 3 show that the Krylov iterations take fewer than half of the number of transport sweeps that Picard iteration required.

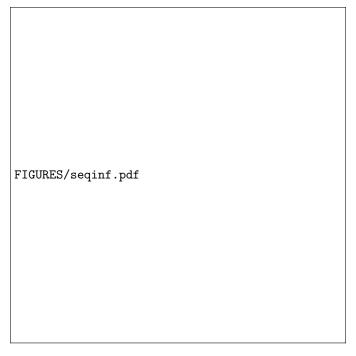


Fig. 2. s = 1

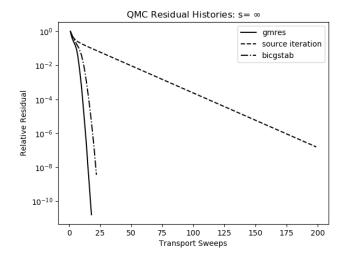


Fig. 3.  $s = \infty$ 

## II.D. 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	5.89664e-01	6.07488e-06	5.71197e-01	5.85487e-06
1.00e-01	5.31120e-01	6.92516 e- 06	5.22137e-01	$6.66741 \mathrm{e}\text{-}06$
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	1.62339 e - 05	3.81029 e-01	1.54416 e - 05
4.00e-01	3.32964 e- 01	4.38580 e- 05	3.34673 e-01	4.09691 e-05
5.00e-01	2.96090e-01	1.69372 e-04	2.98224 e-01	1.57373e-04
6.00e-01	2.66563e-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	2.05174e-01	5.96036e-03	2.07478e-01	5.67991e-03
1.00e+00	1.90546e-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		$_{ m QMC}$	
$\mu$	$\psi(0,-\mu)$	$\psi( au,\mu)$	$\psi(0,-\mu)$	$\psi( au,\mu)$
5.00e-02	8.97798e-01	1.02202 e-01	8.47454e-01	1.00663e-01
1.00e-01	8.87836e-01	1.12164e-01	8.52822 e-01	1.10325 e-01
2.00e-01	8.69581e-01	1.30419e-01	8.47710e-01	1.29064 e-01
3.00e-01	8.52299e-01	1.47701e-01	8.35879 e-01	1.46849e-01
4.00e-01	8.35503 e-01	1.64497e-01	8.22291e-01	1.64034 e-01
5.00e-01	8.18996e-01	1.81004 e-01	8.08044e-01	1.80827e-01
6.00e-01	8.02676 e-01	1.97324e-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	2.61279 e - 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	4.55590 e-02	1.73115e-02	1.63072 e-02	1.61542 e-02	1.58469 e-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	2.42755e-02	1.36521e-02	1.29509e-02	1.22769e-02
100	5.60749e-02	2.31030e-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	1.60082 e-02	4.34610 e - 03	1.41402 e-03
1600	5.99376e-02	2.15675 e-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	1.49649e-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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