

# Variable Eddington Factor Method for $S_N$ Equations with Discontinuous Galerkin Spatial Discretization and a Mixed Finite-Element Balance Equation

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## Abstract

*We present the Variable Eddington Factor (VEF) method, a non-linear Discrete Ordinates Source Iteration scheme that relaxes the consistency requirement of the transport and acceleration steps' spatial discretization. The method was applied to the 1-D, one-group neutron transport equation with Lumped Linear Discontinuous Galerkin (LLDG) transport and the constant-linear Mixed Finite Element Method (MFEM) drift diffusion acceleration. Methods for increased consistency between the transport and acceleration steps are also presented. The VEF method exhibited second-order convergence as expected from the orders of accuracy of LLDG and MFEM in isolation, accelerated source iterations as well as consistently differenced  $S_2SA$ , and survived the thick diffusion limit. In addition, the difference between the transport and acceleration steps' solution was shown to converge as the mesh was refined.*

## Keywords

Variable Eddington Factor, Source Iteration Acceleration, Lumped Linear Discontinuous Galerkin, Mixed Finite Element Method

## Running Head

Variable Eddington Factor Method

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# 1 Introduction

The Variable Eddington Factor (VEF) method, also known as Quasi-Diffusion (QD), was one of the first nonlinear methods for accelerating source iterations in  $S_N$  calculations [1]. It is comparable in effectiveness to both linear and nonlinear forms of Diffusion-Synthetic Acceleration (DSA), but it offers much more flexibility than the DSA. Stability can only be guaranteed with DSA if the diffusion equation is differenced in a manner consistent with that of the  $S_N$  equations [2]. Modern  $S_N$  codes often use advanced discretization schemes such as discontinuous Galerkin (DG) since classic discretization schemes such as step and diamond are not suitable for radiative transfer calculations in the High Energy Density Laboratory Physics (HEDLP) regime or coupled electron-photon calculations. Diffusion discretizations consistent with DG  $S_N$  discretizations cannot actually be expressed in diffusion form, but rather must be expressed in first-order or  $P_1$  form, and are much more difficult to solve than standard diffusion discretizations [3]. Considerable effort has gone into the development of “partially consistent” diffusion discretizations that yield a stable DSA algorithm with some degree of degraded effectiveness, but such discretizations are also generally difficult to develop [4, 5, 6]. A great advantage of the VEF method is that the drift-diffusion equation that accelerates the  $S_N$  source iterations can be discretized in any valid manner without concern for consistency with the  $S_N$  discretization. When the VEF drift-diffusion equation is discretized in a way that is “non-consistent,” the  $S_N$  and VEF drift-diffusion solutions for the scalar flux do not necessarily become identical when the iterative process converges. However, they do become identical in the limit as the spatial mesh is refined, and the difference between the two solutions is proportional to the spatial truncation errors associated with the  $S_N$  and drift-diffusion discretizations. In general the order accuracy of the  $S_N$  and VEF drift-diffusion solutions will be the lowest order accuracy of their respective indepen-

dent discretizations. Although the  $S_N$  solution obtained with such a “non-consistent” VEF method is not conservative, the VEF drift-diffusion solution is in fact conservative. This is particularly useful in multiphysics calculations where the low-order VEF equation can be coupled to the other physics components rather than the high-order  $S_N$  equations. Another advantage of the non-consistent approach is that even if the  $S_N$  spatial discretization scheme does not preserve the thick diffusion limit [7], that limit will generally be preserved using the VEF method.

The purpose of this paper is to investigate the application of the VEF method with the 1-D  $S_N$  equations discretized with the lumped linear-discontinuous method (LLDG) and the drift-diffusion equation discretized using the constant-linear mixed finite-element method (MFEM). To our knowledge, this combination has not been previously investigated. Our motivation for this investigation is that MFEM methods are now being used for high-order hydrodynamics calculations [8]. A radiation transport method compatible with MFEM methods is clearly desirable for developing a MFEM radiation-hydrodynamics code. Such a code would combine thermal radiation transport with hydrodynamics. However, MFEM methods are inappropriate for the standard first-order form of the transport equation. Thus the use of the VEF method with a DG  $S_N$  discretization and a MFEM drift-diffusion discretization suggests itself. Here we define a VEF method that should exhibit second-order accuracy since both the transport and drift-diffusion discretizations are second-order accurate in isolation. In addition, our VEF method should preserve the thick diffusion limit, which is essential for radiative transfer calculations in the HEDLP regime. We use the lumped rather than the standard Linear Discontinuous Galerkin discretization because lumping yields a much more robust scheme, and robustness is essential for radiative transfer calculations in the HEDLP regime. Because this is an initial study, we simplify the investigation by consid-

ering only the one-group neutron transport equation rather than the full radiative transfer equations, which include a material temperature equation as well as the radiation transport equation. The vast majority of relevant properties of a VEF method for radiative transfer can be tested with an analogous method for one-group neutron transport. Furthermore, a high-order DG-MFEM VEF method could be of interest for neutronics in addition to radiative transfer calculations. A full investigation for radiative transfer calculations will be carried out in a future study.

The remainder of this paper is organized as follows. First, we describe the VEF method analytically. Then we describe our discretized  $S_N$  equations, followed by a description of the discretized VEF drift-diffusion equation. We next give computational results. More specifically, we describe two ways to represent the  $S_N$  variable Eddington factor in the MHEM drift-diffusion equation and several ways to construct the  $S_N$  scattering source from the drift-diffusion solution for the scalar flux. Each of these options yields a different VEF method. The accuracy of these methods is then compared to that of the standard LLDG  $S_N$  solution for several test problems, and the iterative convergence rate of these methods is compared to that of the LLDG  $S_N$  equations with fully-consistent  $S_2$ SA acceleration. Finally, we give conclusions and recommendations for future work.

## 2 The VEF Method

### 2.1 The Algorithm

Here, we describe the VEF method for a planar geometry, fixed-source problem:

$$\mu \frac{\partial \psi}{\partial x}(x, \mu) + \sigma_t(x) \psi(x, \mu) = \frac{\sigma_s(x)}{2} \int_{-1}^1 \psi(x, \mu') d\mu' + \frac{Q(x)}{2}, \quad (1)$$

where  $\mu = \cos \theta$  is the cosine of the angle of flight,  $\theta$ , relative to the  $x$ -axis,  $\sigma_t(x)$  and  $\sigma_s(x)$  the total and scattering macroscopic cross sections,  $Q(x)$  the isotropic fixed-source and  $\psi(x, \mu)$  the angular flux. Applying the Discrete Ordinates ( $S_N$ ) angular discretization yields the following set of  $N$  coupled, ordinary differential equations:

$$\mu_n \frac{d\psi_n}{dx}(x) + \sigma_t(x)\psi_n(x) = \frac{\sigma_s(x)}{2}\phi(x) + \frac{Q(x)}{2}, 1 \leq n \leq N, \quad (2)$$

where  $\psi_n(x) = \psi(x, \mu_n)$  is the angular flux in direction  $\mu_n$ . The  $\mu_n$  are stipulated by an  $N$ -point Gauss quadrature rule such that the scalar flux,  $\phi(x)$ , is numerically integrated as follows:

$$\phi(x) = \sum_{n=1}^N w_n \psi_n(x), \quad (3)$$

where  $w_n$  is the quadrature weight corresponding to  $\mu_n$ .

The VEF method begins by solving Eq. 2 while lagging the scattering term. This is called a source iteration, and represented as follows

$$\mu_n \frac{d\psi_n^{\ell+1/2}}{dx}(x) + \sigma_t(x)\psi_n^{\ell+1/2}(x) = \frac{\sigma_s(x)}{2}\phi^\ell(x) + \frac{Q(x)}{2}, \quad 1 \leq n \leq N, \quad (4)$$

where  $\ell$  is the iteration index. The scalar flux used in the scattering term,  $\phi^\ell$ , is assumed to be known either from the previous iteration or from the initial guess if  $\ell = 0$ . The use of a half-integral index indicates that a source iteration is the first of a two-step iteration scheme. However, if one is only doing Source Iteration (SI) without acceleration, the second step would simply be to set the final scalar flux iterate to the iterate after the source iteration:

$$\phi(x)^{\ell+1} = \phi(x)^{\ell+1/2}. \quad (5)$$

Source iteration is slow to converge in optically thick and highly scattering systems, and this is the motivation for accelerating source iteration using the VEF method. The second iterative step of the VEF method is to obtain a final “accelerated” iterate for the scalar flux by solving the VEF drift-diffusion equation using angular flux shape information from the source iteration step:

$$-\frac{d}{dx} \frac{1}{\sigma_t} \frac{d}{dx} [\langle \mu \rangle^{\ell+1/2} \phi^{\ell+1}] + \sigma_a \phi^{\ell+1} = Q, \quad (6)$$

where the Eddington factor is given by

$$\langle \mu^2 \rangle^{\ell+1/2}(x) = \frac{\int_{-1}^1 \mu^2 \psi^{\ell+1/2}(x, \mu) d\mu}{\int_{-1}^1 \psi^{\ell+1/2}(x, \mu) d\mu}. \quad (7)$$

Note that the Eddington factor depends only upon the angular shape of the angular flux, and not its magnitude. This drift diffusion equation is derived by first taking the first two angular moments of Eq. 2:

$$\frac{d}{dx} J(x) + \sigma_a(x) \phi(x) = Q(x), \quad (8a)$$

$$\frac{d}{dx} [\langle \mu^2 \rangle \phi(x)] + \sigma_t(x) J(x) = 0, \quad (8b)$$

where  $J(x)$  is the current. Then Eq. 8b is solved for  $J(x)$ , and this expression is then substituted into Eq. 8a. Performing a SI, computing the Eddington factor from the SI angular flux iterate, and then solving the drift-diffusion equation to obtain a new scalar flux iterate completes one accelerated iteration. These iterations are repeated until convergence of the scalar flux is achieved.

Acceleration occurs because the angular shape of the angular flux, and thus the Eddington factor, converges much faster than the scalar flux. In addition, solution of the drift-diffusion equation includes scattering. Thereby compensating to a large extent for lagging

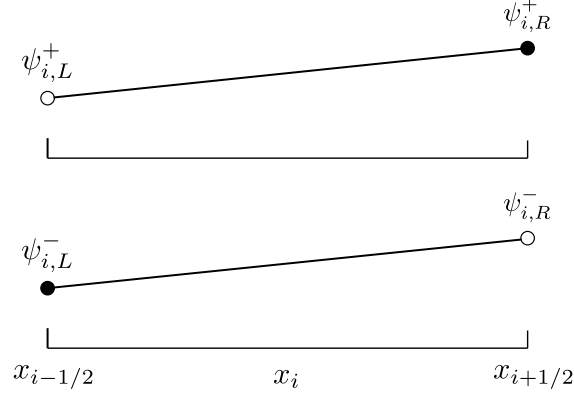


Figure 1: The distribution of unknowns in an LLDG cell. The superscript  $+$  and  $-$  indicate the angular fluxes for  $\mu_n > 0$  and  $\mu_n < 0$ , respectively.

the scattering term in the SI step.

The VEF method allows the  $S_N$  equations and drift diffusion equations to be solved with arbitrarily different spatial discretization methods. The following sections present the application of the Lumped Linear-Discontinuous Galerkin (LLDG) spatial discretization to the  $S_N$  equations and the Mixed Finite-Element Method (MFEM) to the VEF drift-diffusion equation.

## 2.2 Lumped Linear Discontinuous Galerkin $S_N$

The spatial grid and distribution of unknowns for an LLDG cell are shown in Fig. 1. We assume a computational domain of length  $x_b$  discretized into  $I$  cells. The cell centers are integral and the cell edges are half integral. The two unknowns in each cell for each discrete angle are the left and right edge discontinuous angular fluxes,  $\psi_{n,i,L}^{\ell+1/2}$  and  $\psi_{n,i,R}^{\ell+1/2}$ . The cell

edged angular fluxes are uniquely defined by upwinding:

$$\psi_{n,i-1/2}^{\ell+1/2} = \begin{cases} \psi_{n,i-1,R}^{\ell+1/2}, & \mu_n > 0 \\ \psi_{n,i,L}^{\ell+1/2}, & \mu_n < 0 \end{cases}, \quad (9a)$$

$$\psi_{n,i+1/2}^{\ell+1/2} = \begin{cases} \psi_{n,i,R}^{\ell+1/2}, & \mu_n > 0 \\ \psi_{n,i+1,L}^{\ell+1/2}, & \mu_n < 0 \end{cases}. \quad (9b)$$

The angular flux dependence within cells is linear and given in cell  $i$  by

$$\psi_n^{\ell+1/2}(x) = \psi_{n,i,L}^{\ell+1/2} B_{i,L}(x) + \psi_{n,i,R}^{\ell+1/2} B_{i,R}(x), \quad x \in (x_{i-1/2}, x_{i+1/2}), \quad (10)$$

where

$$B_{i,L}(x) = \begin{cases} \frac{x_{i+1/2}-x}{h_i}, & x \in [x_{i-1/2}, x_{i+1/2}] \\ 0, & \text{otherwise} \end{cases}, \quad (11a)$$

$$B_{i,R}(x) = \begin{cases} \frac{x-x_{i-1/2}}{h_i}, & x \in [x_{i-1/2}, x_{i+1/2}] \\ 0, & \text{otherwise} \end{cases}, \quad (11b)$$

The cell centered angular flux is the average of the left and right discontinuous edge fluxes:

$$\psi_{n,i}^{\ell+1/2} = \frac{1}{2} \left( \psi_{n,i,L}^{\ell+1/2} + \psi_{n,i,R}^{\ell+1/2} \right). \quad (12)$$

The unlumped linear-discontinuous Galerkin (LDG) discretization for Eq. 4 is obtained by substituting for  $\psi_n^{\ell+1/2}(x)$  in Eq. 4 from Eq. 10, sequentially multiplying the resultant equation by each basis function, and integrating over each cell with integration-by-parts



of the spatial derivative term. The lumped discretization equations are obtained simply by performing all volumetric integrals (after formal integration-by-parts of the spatial derivative term) using trapezoidal-rule quadrature. The LLDG discretization of Eq. 4 is given by:

$$\mu_n \left( \psi_{n,i}^{\ell+1/2} - \psi_{n,i-1/2}^{\ell+1/2} \right) + \frac{\sigma_{t,i} h_i}{2} \psi_{n,i,L}^{\ell+1/2} = \frac{\sigma_{s,i} h_i}{4} \phi_{i,L}^{\ell} + \frac{h_i}{4} Q_{i,L}, \quad (13a)$$

$$\mu_n \left( \psi_{n,i+1/2}^{\ell+1/2} - \psi_{n,i}^{\ell+1/2} \right) + \frac{\sigma_{t,i} h_i}{2} \psi_{n,i,R}^{\ell+1/2} = \frac{\sigma_{s,i} h_i}{4} \phi_{i,R}^{\ell} + \frac{h_i}{4} Q_{i,R}, \quad (13b)$$

where  $h_i$ ,  $\sigma_{t,i}$ ,  $\sigma_{s,i}$ , and  $Q_{i,L/R}$  are the cell width, total cross section, scattering cross section and discontinuous fixed source in cell  $i$ . The discontinuous scalar fluxes,  $\phi_{i,L/R}^{\ell}$ , are assumed to be known from the previous iteration or the initial guess when  $\ell = 0$ . Equations 13a, 13b, 9a, 9b, and 12 can be combined and rewritten as follows

$$\begin{bmatrix} \mu_n + \sigma_{t,i} h_i & \mu_n \\ -\mu_n & \sigma_{t,i} + \mu_n \end{bmatrix} \begin{bmatrix} \psi_{n,i,L}^{\ell+1/2} \\ \psi_{n,i,R}^{\ell+1/2} \end{bmatrix} = \begin{bmatrix} \frac{\sigma_{s,i} h_i}{2} \phi_{i,L}^{\ell} + \frac{h_i}{2} Q_{i,L} + 2\mu_n \psi_{n,i-1,R}^{\ell+1/2} \\ \frac{\sigma_{s,i} h_i}{2} \phi_{i,R}^{\ell} + \frac{h_i}{2} Q_{i,R} \end{bmatrix}, \quad (14)$$

for sweeping from left to right ( $\mu_n > 0$ ) and

$$\begin{bmatrix} -\mu_n + \sigma_{t,i} h_i & \mu_n \\ -\mu_n & -\mu_n + \sigma_{t,i} h_i \end{bmatrix} \begin{bmatrix} \psi_{n,i,L}^{\ell+1/2} \\ \psi_{n,i,R}^{\ell+1/2} \end{bmatrix} = \begin{bmatrix} \frac{\sigma_{s,i} h_i}{2} \phi_{i,L}^{\ell} + \frac{h_i}{2} Q_{i,L} \\ \frac{\sigma_{s,i} h_i}{2} \phi_{i,R}^{\ell} + \frac{h_i}{2} Q_{i,R} - 2\mu_n \psi_{n,i+1,L}^{\ell+1/2} \end{bmatrix}, \quad (15)$$

for sweeping from right to left ( $\mu_n < 0$ ), respectively. The right hand sides of Eqs. 14 and 15 are known as the scalar flux from the previous iteration, the fixed source, and the angular flux entering from the downwind cell are all known. By supplying the flux entering the left side of the first cell, the solution for ( $\mu_m > 0$ ) can be propagated from left to right by solving Eq. 14. Similarly, supplying the incident flux on the right boundary allows the solution

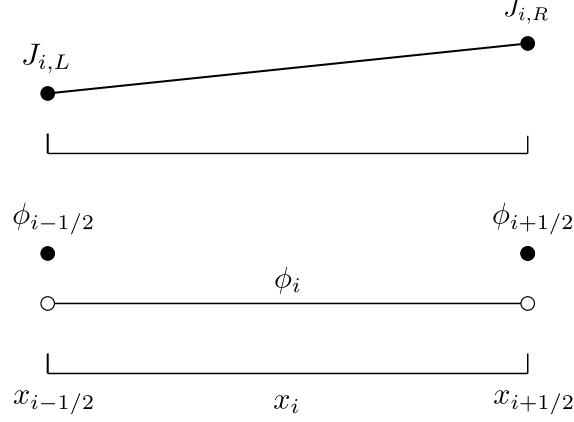


Figure 2: The distribution of unknowns in cell  $i$  for MFEM.

for ( $\mu_n < 0$ ) to be propagated from right to left with Eq. 15. The Eddington factors are computed at cell edges as follows:

$$\langle \mu^2 \rangle_{i\pm 1/2}^{\ell+1/2} = \frac{\sum_{n=1}^N \mu_n^2 \psi_{n,i\pm 1/2}^{\ell+1/2} w_n}{\sum_{n=1}^N \psi_{n,i\pm 1/2}^{\ell+1/2} w_n}, \quad (16)$$

where the  $\psi_{n,i\pm 1/2}^{\ell+1/2}$  are defined by Eqs. 9a and 9b. The Eddington factors are computed within cell  $i$  as follows:

$$\langle \mu^2 \rangle^{\ell+1/2}(x) = \frac{\sum_{n=1}^N \mu_n^2 \psi_n^{\ell+1/2}(x) w_n}{\sum_{n=1}^N \psi_n^{\ell+1/2}(x) w_n}, \quad x \in (x_{i-1/2}, x_{i+1/2}), \quad (17)$$

where  $\psi_n^{\ell+1/2}(x)$  is defined by Eq. 10.

### 2.3 Mixed Finite-Element Method for VEF Equation

We apply the MFEM method to Eqs. 8a and 8b, and then eliminate the currents to obtain a discretization for Eq. 6. In this method, the grid is identical to that used in the LLDG  $S_N$  discretization. Separate basis functions are used for the scalar flux and current. The

unknowns in an MFEM cell are depicted in Fig. 2. The scalar flux is constant within the cell with discontinuous jumps at the cell edges and the current is a linear function defined by:

$$J_i(x) = J_{i,L}B_{i,L}(x) + J_{i,R}B_{i,R}(x), \quad (18)$$

where  $J_{i,L/R}$  are the currents at the left and right edges of the cell, and the basis functions are identical to those defined by Eqs. 11a and 11b for the LLDG  $S_N$  discretization. The constant-linear MFEM yields second order accuracy for both the scalar flux and the current.

The MFEM representation yields five unknowns per cell:  $\phi_{i-1/2}$ ,  $\phi_i$ ,  $\phi_{i+1/2}$ ,  $J_{i,L}$ , and  $J_{i,R}$ . However, each edge flux on the mesh interior is shared by two cells, so with  $I$  cells there are  $I$  cell-center scalar fluxes,  $2I$  currents, and  $2N - 1$  interior-mesh cell-edge scalar fluxes, and 2 boundary cell-edge scalar fluxes. An equation for  $\phi_i$  is found by integrating Eq. 8a over cell  $i$ :

$$J_{i,R} - J_{i,L} + \sigma_{a,i}h_i\phi_i = Q_ih_i, \quad (19)$$

where  $\sigma_{a,i}$  and  $Q_i$  are the absorption cross section and source in cell  $i$ . Equations for  $J_{i,L/R}$  are found by multiplying Eq. 8b by  $B_{i,L/R}$  and integrating over cell  $i$ :

$$-\langle\mu^2\rangle_{i-1/2}\phi_{i-1/2} + \langle\mu^2\rangle_i\phi_i + \sigma_{t,i}h_i\left(\frac{1}{3}J_{i,L} + \frac{1}{6}J_{i,R}\right) = 0, \quad (20a)$$

$$\langle\mu^2\rangle_{i+1/2}\phi_{i+1/2} - \langle\mu^2\rangle_i\phi_i + \sigma_{t,i}h_i\left(\frac{1}{6}J_{i,L} + \frac{1}{3}J_{i,R}\right) = 0, \quad (20b)$$

where the fixed source has been assumed to be isotropic. Note that  $\langle\mu^2\rangle_{i\pm 1/2}$  denotes cell edge Eddington factors, while  $\langle\mu^2\rangle_i$  denotes an average over cell  $i$  of the Eddington factors. All Eddington factors are computed using the angular fluxes from the LLDG  $S_N$  step. The edge Eddington factors are defined by Eq. 16, while the Eddington factors within each cell are

defined by Eq. 17. We stress that evaluating Eq. 17 at  $x_{i,L/R}$  does not yield  $\langle \mu^2 \rangle_{i \pm 1/2}$  because of the upwinding used to define the cell edge angular fluxes. Since the spatial dependence of the Eddington factors within each cell takes the form of a rational polynomial, the average must be computed by quadrature. We investigate three quadratures for computing these averages: midpoint quadrature,

$$\langle \mu^2 \rangle_i = \langle \mu^2 \rangle(x_i); \quad (21a)$$

trapezoidal quadrature,

$$\langle \mu^2 \rangle_i = \frac{1}{2} [\langle \mu^2 \rangle(x_{i,L}) + \langle \mu^2 \rangle(x_{i,R})]; \quad (21b)$$

and Gauss quadrature,

$$\langle \mu^2 \rangle_i = \frac{1}{2} [\langle \mu^2 \rangle(x_{i,L}^G) + \langle \mu^2 \rangle(x_{i,R}^G)]; \quad (21c)$$

where

$$x_{i,L/R}^G = \frac{h_i}{2} \mp \frac{x_{i+1/2} + x_{i-1/2}}{2\sqrt{3}}. \quad (22)$$

We believe that trapezoidal may be the most consistent with the LLDG equations because the FEM integrations within each cell associated with those equations are performed with trapezoidal quadrature.

Eliminating  $J_{i,R}$  from Eq. 20a and  $J_{i,L}$  from Eq. 20b yields:

$$J_{i,L} = \frac{-2}{\sigma_{t,i} h_i} \left\{ 2 [\langle \mu^2 \rangle_i \phi_i - \langle \mu^2 \rangle_{i-1/2} \phi_{i-1/2}] - [\langle \mu^2 \rangle_{i+1/2} \phi_{i+1/2} - \langle \mu^2 \rangle_i \phi_i] \right\}, \quad (23a)$$

$$J_{i,R} = \frac{-2}{\sigma_{t,i} h_i} \left\{ 2 [\langle \mu^2 \rangle_{i+1/2} \phi_{i+1/2} - \langle \mu^2 \rangle_i \phi_i] - [\langle \mu^2 \rangle_i \phi_i - \langle \mu^2 \rangle_{i-1/2} \phi_{i-1/2}] \right\}. \quad (23b)$$

An equation for  $\phi_{i+1/2}$  on the mesh interior is found by enforcing continuity of current at

the cell edges:

$$J_{i,R} = J_{i+1,L} . \quad (24)$$

Using the definitions of  $J_{i,L}$  and  $J_{i,R}$  from Eqs. 23a and 23b in the balance equation (Eq. 19) and continuity equation (Eq. 24) yields equations for all cell-center fluxes and interior-mesh cell-edge fluxes. The resulting balance and continuity equations are:

$$-\frac{6}{\sigma_{t,i}h_i}\langle\mu^2\rangle_{i-1/2}\phi_{i-1/2} + \left(\frac{12}{\sigma_{t,i}h_i}\langle\mu^2\rangle_i + \sigma_{a,i}h_i\right)\phi_i - \frac{6}{\sigma_{t,i}h_i}\langle\mu^2\rangle_{i+1/2}\phi_{i+1/2} = Q_i h_i , \quad (25a)$$

$$\begin{aligned} -\frac{2}{\sigma_{t,i}h_i}\langle\mu^2\rangle_{i-1/2}\phi_{i-1/2} + \frac{6}{\sigma_{t,i}h_i}\langle\mu^2\rangle_i\phi_i - 4\left(\frac{1}{\sigma_{t,i}h_i} + \frac{1}{\sigma_{t,i+1}h_{i+1}}\right)\langle\mu^2\rangle_{i+1/2}\phi_{i+1/2} \\ + \frac{6}{\sigma_{t,i+1}h_{i+1}}\langle\mu^2\rangle_{i+1}\phi_{i+1} - \frac{2}{\sigma_{t,i+1}h_{i+1}}\langle\mu^2\rangle_{i+3/2}\phi_{i+3/2} = 0 . \end{aligned} \quad (25b)$$

The equations for the outer boundary fluxes,  $\phi_{1/2}$  and  $\phi_{I+1/2}$ , involve boundary conditions together with continuity conditions. For instance, the equation for  $\phi_{1/2}$  is

$$J_{1,L} = J_{1/2} , \quad (26)$$

where  $J_{1,L}$  is defined in Eq. 23a, and  $J_{1/2}$  is the left boundary current defined by a boundary condition. For a reflective condition,

$$J_{1/2} = 0 . \quad (27)$$

For a source condition,

$$J_{1/2} = 2 \sum_{\mu_n > 0} \mu_n \psi_{n,1/2} w_n - B_{1/2} \phi_{1/2} , \quad (28)$$

where

$$B_{1/2} = \frac{\sum_{n=1}^N |\mu_n| \psi_{n,1/2} w_n}{\sum_{n=1}^N \psi_{n,1/2} w_n} \quad (29)$$

is the boundary Eddington factor [9]. The equation for  $\phi_{I+1/2}$  is

$$J_{I,R} = J_{I+1/2}. \quad (30)$$

where  $J_{I,R}$  is defined in Eq. 23b, and  $J_{I+1/2}$  is the right boundary current. For a reflective condition,

$$J_{I+1/2} = 0. \quad (31)$$

For a source condition,

$$J_{I+1/2} = B_{I+1/2} \phi_{I+1/2} - 2 \sum_{\mu_n < 0} |\mu_n| \psi_{n,I+1/2} w_n, \quad (32)$$

where

$$B_{I+1/2} = \frac{\sum_{n=1}^N |\mu_n| \psi_{n,I+1/2} w_n}{\sum_{n=1}^N \psi_{n,I+1/2} w_n}. \quad (33)$$

These transport-consistent Marshak-like source boundary conditions are derived starting with the identity

$$J_{1/2} = j^+ - j^-, \quad (34)$$

where  $j^\pm$  denotes the positive half-range currents associated with  $\mu > 0$  and  $\mu < 0$ , respectively. For the left boundary condition, we simply perform the following algebraic manipulations:

$$J_{1/2} = j^+ - j^- = 2j^+ - (j^+ + j^-) = 2j^+ - \frac{j^+ + j^-}{\phi} \phi = 2j^+ - B_{1/2} \phi. \quad (35)$$

For the right boundary condition, we similarly obtain

$$J_{I+1/2} = j^+ - j^- = (j^+ + j^-) - 2j^- = \frac{j^+ + j^-}{\phi} \phi - 2j^- = B_{I+1/2} \phi - 2j^- . \quad (36)$$

Note that these source boundary conditions become equivalent to the standard Marshak boundary conditions if the  $S_N$  angular flux is isotropic. The system of  $2I + 1$  equations for the cell-center and cell-edge fluxes can be assembled into a matrix of both cell centered and cell edge scalar fluxes and solved with a banded matrix solver of bandwidth five.

## 2.4 Increased Consistency Between LLDG and MFEM

The MFEM representation for the scalar flux is constant within a cell, but the LLDG representation for the scalar flux is linear. This suggests that improved accuracy of the  $S_N$  solution could be achieved by somehow constructing a linear scalar flux dependence from the MFEM solution. One simple method for doing this is to use the MFEM cell-edge scalar fluxes to compute a slope. This works quite well, for neutronics. However, it will be inadequate in a radiative transfer calculation because slopes must also be generated for the material temperatures, and an MFEM approximation for the temperatures will not include edge temperatures. We have chosen to use a more generally applicable approach based upon standard data reconstruction techniques that require only cell-centered values to compute slopes [10]. We also limit such slopes to avoid non-physical scalar fluxes. For example, the reconstructed left and right scalar fluxes in a cell  $i$  are given by

$$\phi_{i,L/R} = \phi_i \mp \frac{1}{4} \xi_i (\Delta \phi_{i+1/2} + \Delta \phi_{i-1/2}) , \quad (37)$$

where  $\xi$  is a van Leer-type slope limiter [10]:

$$\xi_i = \begin{cases} 0, & \text{if } r_i \leq 0, \\ \max \left\{ \frac{r_i}{1+r_i}, \frac{1}{1+r_i} \right\}, & \text{if } r_i > 0, \end{cases} \quad (38)$$

$$r_i = \frac{\Delta\phi_{i-1/2}}{\Delta\phi_{i+1/2}}, \quad (39)$$

and

$$\Delta\phi_{i+1/2} = \phi_{i+1} - \phi_i. \quad (40)$$

On the boundaries, we use

$$\phi_{1,L/R} = \phi_1 \mp \frac{1}{2}\Delta\phi_{3/2}, \quad (41a)$$

$$\phi_{I,L/R} = \phi_I \mp \frac{1}{2}\Delta\phi_{I-1/2}. \quad (41b)$$

We also set any negative left or right flux values in the boundary cells to zero by appropriately rotating the slopes. I'm not sure if this was how the slopes were really calculated, but we need to fully explain what was done, and we should do something to prevent negativities from arising.

### 3 Computational Results

We need to describe the problem here. Unless otherwise noted, all calculations were performed using midpoint quadrature to compute the cell average Eddington factors. Figure 3a shows the iterative convergence as a function of unaccelerated iteration number for the scalar flux and the Eddington factor. The Eddington factor's large drop in relative norm we need to describe the norm between the first and second iterations supports the claim that the



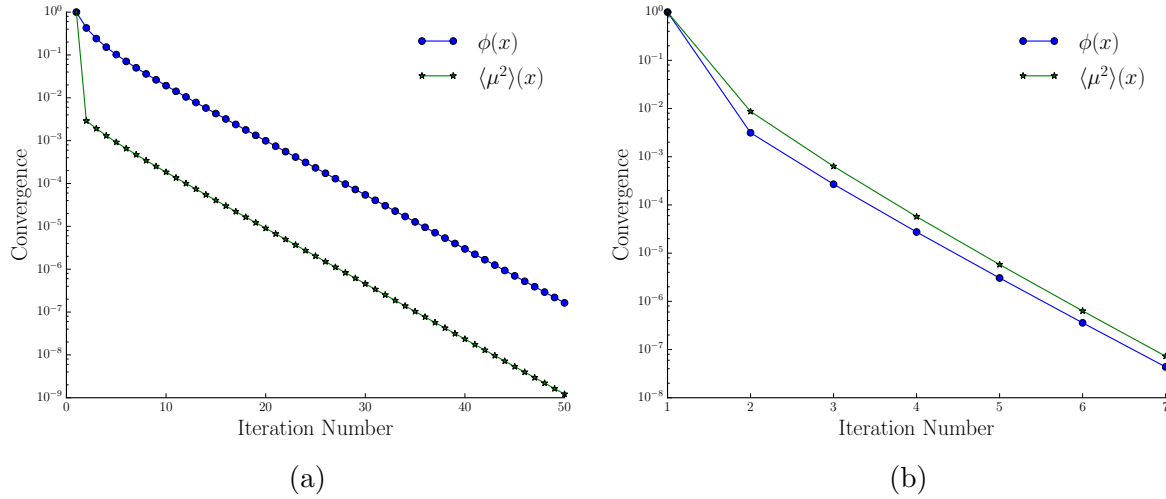


Figure 3: The convergence rate for  $\phi(x)$  and  $\langle \mu^2 \rangle(x)$  for (a) unaccelerated and (b) VEF accelerated SI.

angular shape of the angular flux, and thus the Eddington factor, converges rapidly. When compared to Fig. 3b, a plot of the iterative convergence for the VEF method, it is clear that the VEF method transfers the fast rate of convergence of the Eddington factor to the scalar flux.

To compare SI, VEF, and consistently differenced S<sub>2</sub>SA, a homogeneous test problem with a reflecting left boundary, a vacuum right boundary, and a total thickness of 10 cm, was used. This system was discretized into 50 spatial cells. The total macroscopic cross section,  $\sigma_t$ , was set to  $1 \text{ cm}^{-1}$  leading to an optical thickness per cell of 0.2 mfp and a problem thickness of 10 mfp. The convergence tolerance was set to  $10^{-6}$ . [We need to explain the convergence tolerance in full detail.](#) Figure 4 shows the number of iterations required for convergence for SI, VEF, and S<sub>2</sub>SA for varying ratios of  $\sigma_s$  to  $\sigma_t$ . The ratio of unaccelerated to VEF accelerated iterations ranged from 1.6 to 7. This suggests that acceleration is occurring and that the VEF method is not just doing twice the amount of work per iteration. In

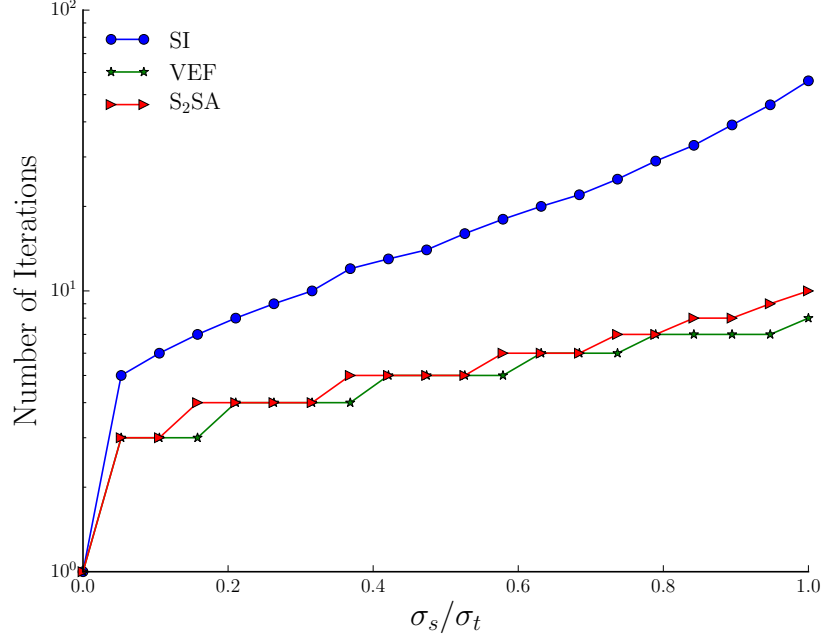


Figure 4: A comparison of the number of iterations required for Source Iteration, VEF acceleration, and S<sub>2</sub>SA to converge for varying ratios of  $\sigma_s$  to  $\sigma_t$ .

addition, the VEF method performed similarly to S<sub>2</sub>SA. I am skeptical of the numbers for the unaccelerated case. You need to include an estimate of the spectral radius to properly converge the error when the spectral radius is near one. You need a tighter tolerance -  $\text{tol} = \text{tol} * (1 - \rho)$ , where  $\rho$  is the spectral radius estimate. Otherwise you will not actually converge the error to  $\text{tol}$  but to  $\text{tol}/(1 - \rho)$ .

The Method of Manufactured Solutions (MMS) was used to compare the accuracy of the VEF method as the cell width was decreased. We need to explain how you constructed the MMS solution. I can't tell anything about the problem. In general, you are supposed to provide enough information for the reader to reproduce your calculations. The L2 norm of the difference between the numerical and MMS solutions was compared at five logarithmically

Reconstruction Method	Eddington Representation	Order	$C$	$R^2$
None	Average	1.997	0.682	$9.9999 \times 10^{-1}$
None	Rational Polynomial	1.998	0.687	1.0000
Center	Average	2.007	0.726	$9.9992 \times 10^{-1}$
Center	Rational Polynomial	2.009	0.732	$9.9991 \times 10^{-1}$

Table 1: The order of accuracy, error, and  $R^2$  values for the permutations of the two Eddington representation methods and two slope reconstruction methods.

spaced cell widths between 0.5 mm and 0.01 mm. A line of best fit of the form

$$E = Ch^n \quad (42)$$

was used to find the order of accuracy,  $n$ , and the constant of proportionality,  $C$ , of the numerical error,  $E$ . These values are provided in Table 1 for the permutations of average and rational polynomial Eddington representation and slope reconstruction and no reconstruction. All of the permutations are second order accurate and have similar overall accuracy. This suggests that slope reconstruction and Eddington representation do not affect the *order* of numerical accuracy.

The convergence between unaccelerated SI and the VEF method was compared as a function of cell width for a simple homogeneous slab and for Reed’s problem. In both cases, the left boundary was reflecting and the right boundary was vacuum. The homogeneous slab had a scattering ratio of 0.75. The cross sections and source for Reed’s problem are provided in Table 2. The L2 norm of the difference between the SI solution and VEF solution is plotted for the four permutations of no reconstruction, van Leer slope limited reconstruction, average Eddington representation, and rational polynomial Eddington representation in Figures 5a and 5b for the homogeneous slab problem and Reed’s problem.

In the homogeneous problem, VEF with van Leer limited slope reconstruction was five

	Region 1	Region 2	Region 3	Region 4	Region 5
$q$	10	0	0	0	1
$\Sigma_t$	10	0.001	1	5	1
$\Sigma_a$	10	0	0.1	0	0.1
Domain	$0 \leq x < 2$	$2 \leq x < 4$	$4 \leq x < 6$	$6 \leq x < 7$	$7 \leq x \leq 8$

Table 2: The cross sections and source used for Reed's problem.

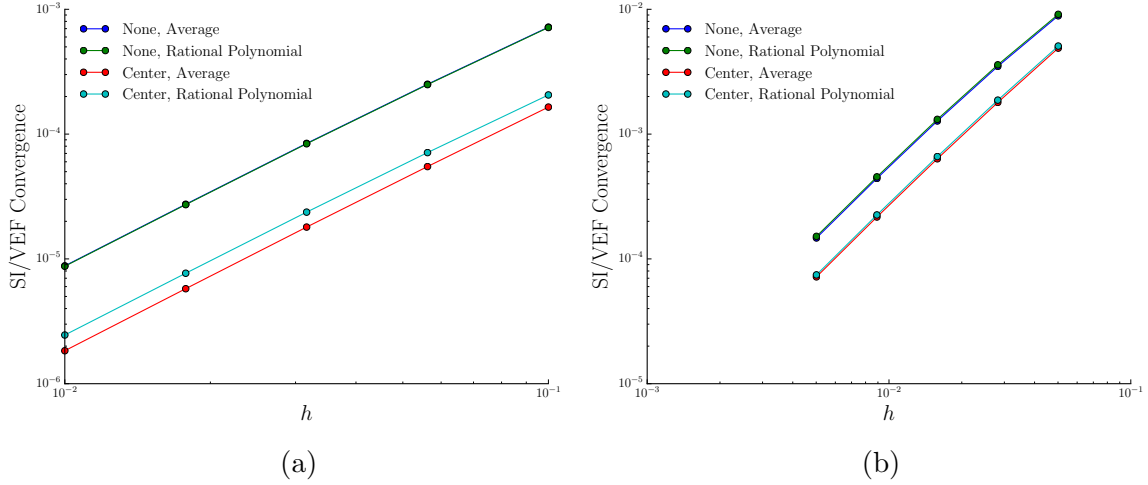


Figure 5: The L2 norm of the difference between SI and the four permutations of the VEF method as the cell spacing is decreased for (a) the homogeneous slab problem and (b) Reed's problem.

times more convergent [what are we talking about here - iterative convergence or error in the final solution?](#) than VEF without reconstruction. Use of Gauss quadrature for computing the Eddington factors decreased the van Leer reconstruction convergence by 30%. [Wow, I don't get this.](#) We are computing a more accurate average and getting a less accurate solution. [We should add the trapezoidal quadrature results to this](#) In Reed's problem, reconstruction was twice as accurate as no reconstruction. The calculations using trapezoidal and Gauss quadratures to compute the average Eddington factors were XXX% and 4% (more/less) accurate than those using midpoint quadrature. [We need to add trapezoidal results.](#) Note

that Average and Rational Polynomial are somewhat confusing terms because the MFEM formalism always calls for an average value of the Eddington factor in the gradient term. It is just a question of how you compute the average.

Lastly, slope reconstruction and Eddington factor averaging were tested in the diffusion limit. The cross sections and source were scaled as follows [11]:

$$\sigma_t(x) \rightarrow \sigma_t(x)/\epsilon, \quad (43a)$$

$$\sigma_s(x) \rightarrow \epsilon\sigma_s(x), \quad (43b)$$

$$Q(x) \rightarrow \epsilon Q(x). \quad (43c)$$

As  $\epsilon \rightarrow 0$ , the system becomes diffusive. The number of iterations for convergence within a tolerance of  $10^{-8}$  as  $\epsilon \rightarrow 0$  is plotted in Fig. 6. The error between the VEF solution and the exact diffusion solution is provided in Fig. 7. This supports the claim that the VEF method is robust as all four permutations properly preserved the diffusion limit.

## 4 Conclusions and Future Work

We have presented the VEF method for one-group neutron transport in slab geometry and the pairing of Lumped Linear Discontinuous Galerkin for the  $S_N$  transport step and the constant-linear Mixed Finite-Element Method for the drift diffusion acceleration step. We have numerically demonstrated that the LLDG/MFEM VEF method accelerates Source Iteration by transferring the rapid convergence of the angular shape of the angular flux to the scalar flux. In addition, the VEF method performed similarly to consistently differenced  $S_2SA$ .

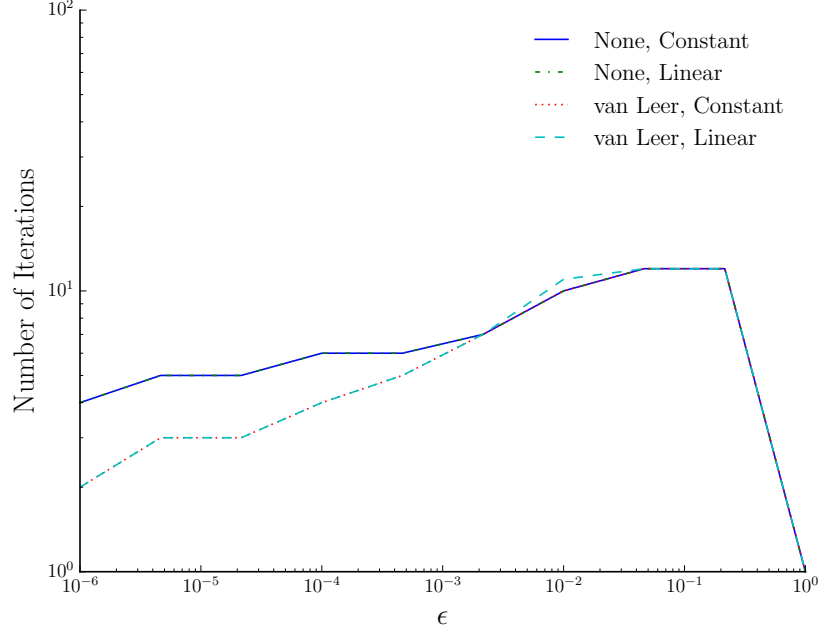


Figure 6: The number of iterations required for convergence for the permutations of slope reconstruction and angular flux representation in the diffusion limit.

Methods for increased consistency between LLDG and MFEM were also presented. This included a more consistent method for computing the Eddington factor on the cell centers and a cell-centered slope reconstruction. It was shown that both the VEF method without and without added consistency measures was second-order accurate as expected from the orders of accuracy of LLDG and MFEM in isolation and that all of the VEF methods were robust in the diffusion limit. In addition, while this nonlinear scheme produces two solutions, one from  $S_N$  and one from drift diffusion, the solutions were shown to converge as the mesh was refined for both homogeneous and inhomogeneous systems. Furthermore, the drift-diffusion solution for scalar flux and current is conservative and thus is the preferred solution to compute quantities of interest.

Slope reconstruction significantly increased the convergence between the  $S_N$  and drift

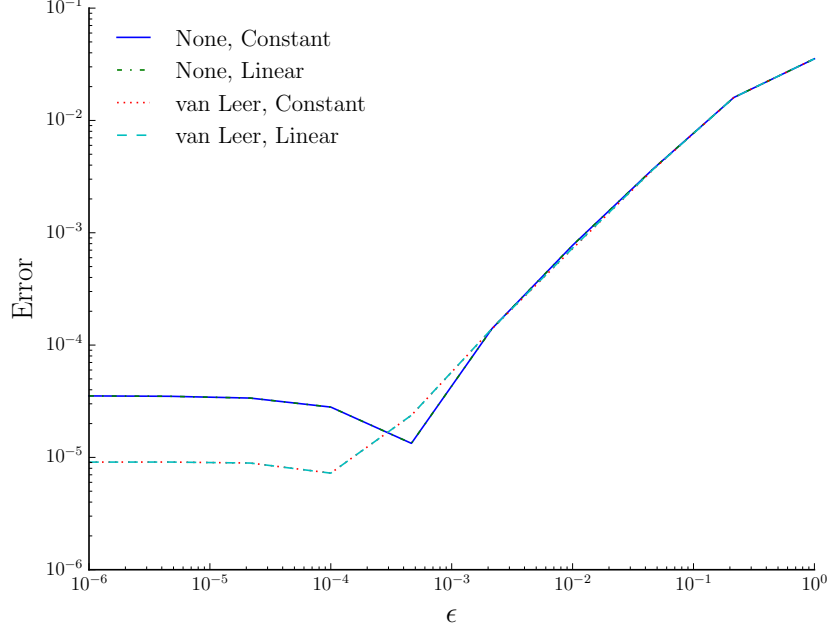


Figure 7: The error between the VEF methods and the exact diffusion solution as  $\epsilon \rightarrow 0$ .

diffusion solutions. While consistent representation of the Eddington factor did not significantly impact numerical accuracy or solution convergence, it is expected to be important in radiative transfer calculations due to the non-linearities in temperature and the spatial dependence of the cross sections. [We need to take a look at this. I want to make sure I understood what you did. I don't understand why we see less accuracy with the rational polynomial representation.](#)

The VEF method is especially suited for multiphysics calculations as VEF drift diffusion is conservative and much less expensive than a transport sweep. In addition, the multiphysics and transport discretizations can be arbitrarily different.

Future work includes extending the VEF method presented in this paper to the radiative transfer equations, verifying the VEF method in higher dimensions, and generalizing the

method to higher order finite-elements.

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