Title
A high-order Eulerian Godunov method for elastic/plastic flow in solids

Permalink
https://escholarship.org/uc/item/7mc731pn

Journal
Journal of Computational Physics, 167(1)

Author
Miller, G.H.

Publication Date
2000-05-08

G.H. Miller and P. Colella

National Energy Research Scientific Computing Division

May 2000
Submitted to
Journal of Computational Physics
DISCLAIMER

This document was prepared as an account of work sponsored by the United States Government. While this document is believed to contain correct information, neither the United States Government nor any agency thereof, nor The Regents of the University of California, nor any of their employees, makes any warranty, express or implied, or assumes any legal responsibility for the accuracy, completeness, or usefulness of any information, apparatus, product, or process disclosed, or represents that its use would not infringe privately owned rights. Reference herein to any specific commercial product, process, or service by its trade name, trademark, manufacturer, or otherwise, does not necessarily constitute or imply its endorsement, recommendation, or favoring by the United States Government or any agency thereof, or The Regents of the University of California. The views and opinions of authors expressed herein do not necessarily state or reflect those of the United States Government or any agency thereof, or The Regents of the University of California.

Ernest Orlando Lawrence Berkeley National Laboratory
is an equal opportunity employer.
DISCLAIMER

This document was prepared as an account of work sponsored by the United States Government. While this document is believed to contain correct information, neither the United States Government nor any agency thereof, nor the Regents of the University of California, nor any of their employees, makes any warranty, express or implied, or assumes any legal responsibility for the accuracy, completeness, or usefulness of any information, apparatus, product, or process disclosed, or represents that its use would not infringe privately owned rights. Reference herein to any specific commercial product, process, or service by its trade name, trademark, manufacturer, or otherwise, does not necessarily constitute or imply its endorsement, recommendation, or favoring by the United States Government or any agency thereof, or the Regents of the University of California. The views and opinions of authors expressed herein do not necessarily state or reflect those of the United States Government or any agency thereof or the Regents of the University of California.

G.H. Miller and P. Colella

National Energy Research Scientific Computing Division
Ernest Orlando Lawrence Berkeley National Laboratory
University of California
Berkeley, California 94720

May 2000

This work was supported by the Director, Office of Science, Office of Advanced Scientific Computing Research, Mathematical, Information, and Computational Sciences Division, of the U.S. Department of Energy under Contract No. DE-AC03-76SF00098. Other work was supported by a subcontract from the Caltech Center for the Simulation of Dynamic Response in Materials, which in turn is supported by the Academic Strategic Alliances Program of the Accelerated Strategic Computing Initiative (ASCI/ASAP) under subcontract B341492 of the U.S. Department of Energy under Contract No. W-7405-ENG-48.
We present an explicit second-order accurate Godunov finite difference method for the solution of the equations of solid mechanics in 1, 2, and 3 spatial dimensions. The solid mechanics equations are solved in non-conservation form, with the novel application of a diffusion-like correction to enforce the gauge condition that the deformation tensor be the gradient of a vector. Physically conserved flow variables (e.g., mass, momentum, and energy) are strictly conserved; only the deformation gradient field is not. Verification examples demonstrate the accurate capturing of plastic and elastic shock waves across approximately 5 computational cells. 2D and 3D results are obtained without spatial operator splitting.

Key Words: solid mechanics, shock waves, Godunov method, elasticity, plasticity

1. INTRODUCTION

In this work, we present a higher-order Godunov method for computing in Eulerian coordinates the multidimensional dynamics of elastic-plastic solids undergoing large deformations. Our approach is based on a new formulation of the equations of solid mechanics as a first-order system of hyperbolic PDE's, a modification of that used by Trangenstein and Colella [24]. In [24], the usual conservation laws for mass, momentum and energy, plus a constitutive model, are augmented by a form of equality of mixed partial derivatives that yields conservation equations for the entries of the inverse deformation gradient. This leads to equations of the form

$$\frac{\partial U}{\partial t} + \nabla \cdot F(U) = S(U). \quad (1)$$

Work at the Lawrence Berkeley National Laboratory was sponsored by the US Department of Energy (DOE) Mathematical, Information, and Computing Sciences Division contract DE-AC03-76SF00098. Other work was supported by a subcontract from the Caltech Center for the Simulation of Dynamic Response in Materials, which in turn is supported by the Academic Strategic Alliances Program of the Accelerated Strategic Computing Initiative (ASCI/ASAP) under subcontract B341492 of DOE contract W-7405-ENG-48.
Here $S(U)$ contains source terms associated with the treatment of plasticity. These equations, by themselves, are not sufficient to specify the problem. In addition, we must impose linear constraints on the solution to guarantee that the inverse of the deformation gradient is, in fact a gradient, i.e. that the curl of the rows of the deformation gradient vanish. These constraints can be written in the following form:

$$L_C(U) = 0.$$  \hspace{1cm} (2)

Here $L_C$ is a system of linear differential operators with constant coefficients. The constraint equation is an initial-value constraint: if (1) is satisfied, and $L_C(U)$ is identically zero at some initial time, then $L_C(U)$ vanishes identically for all later times. The constraint (2) plays an essential role in the analysis of the characteristic structure of the system (1). In order to get the physically correct eigenvectors and eigenvalues from the quasilinear form of the equations, one must use the constraint to replace some of the spatial derivatives by others. In general, solutions to (1), without imposing (2), give rise to unphysical wave propagation properties, even for linearized waves as was observed in [24].

A difficulty arises when one attempts to compute solutions to (1,2) using a conservative finite difference method. To the extent that a modified equation analysis is valid, we expect the behavior of the numerical solution to behave very similarly to the solution to the following system of PDE's:

$$\frac{\partial U^{\text{Mod}}}{\partial t} + \nabla \cdot F(U^{\text{Mod}}) = S(U^{\text{Mod}}) + \tau_U(U^{\text{Mod}})$$

$$L_C(U^{\text{Mod}}) = \tau_C(U^{\text{Mod}}).$$  \hspace{1cm} (3)

Here $\tau_U$ and $\tau_C$ are truncation error terms, which are nonzero. In general, these terms, and in particular $\tau_C$, cannot be eliminated. The practice of enforcing a discretized form of the constraint (2) at the end of each time step using a Hodge projection would guarantee that a discretized form of (2) is satisfied identically. However, that will change the form of $\tau_C$, but not set it to zero. The observation that the truncation error terms are a small perturbation to the equations is not sufficient to guarantee that $U^{\text{Mod}}$ is close to $U$. There is much less known about the well-posedness of systems of equations that are combinations of evolution equations and constraints than there is about pure evolution equations, and unexpected pathologies are known to occur [16].

The approach we want to take on this problem starts with an analysis due to Godunov [8, 9]. Numerical methods based on this approach have been recently investigated for the MHD equations in [19], the case for which Godunov first applied this analysis. Godunov modifies (1) in the following way:

$$\frac{\partial U}{\partial t} + \nabla \cdot F(U) = S(U) + \xi L_C(U).$$  \hspace{1cm} (4)

Here $\xi = \xi(U)$ can be chosen so that the system has the physically correct linearized eigenstructure, independent of whether $L_C$ vanishes. In addition, $L_C$ satisfies a transport equation such that if $L_C(U)$ is identically zero at some time, then it remains so for all later times.

The numerical method we present here is based on the form of the equations given by (4). Thus we are discretizing a well-posed initial value problem without constraints, independent of the whether or not the constraint (2) is satisfied. This gives us a high degree
of confidence that a stable and consistent method can be developed. Of course, the extent to which we compute a solution to the original physical problem (1,2) depends strongly on whether the constraint is satisfied, but now that is purely an accuracy issue, without any impact on the stability of the method. In fact, we will investigate the use of various methods of limiting discrete measures of $L_c$, similar to filtering methods developed for incompressible flow [20, 11].

2. GOVERNING EQUATIONS

The mechanical behavior of solids is described by observable variables (e.g., density $\rho$, momentum $\rho v$, internal energy $E$, and the deformation $F$ with respect to a chosen reference state), and also unobservable internal parameters which describe the response of the material to deviatoric stress. One constitutive representation of this behavior is through the multiplicative decomposition of the total deformation into elastic and inelastic components,

$$F = F^eF^p.$$  \hfill (5)

Here $F$ is the Lagrangian coordinate deformation which relates the spatial coordinate frame $x = x(a,t)$ to the material coordinate frame $a$:

$$F_{\alpha\beta} = \frac{\partial x_\alpha}{\partial a_\beta}. \hfill (6)$$

We refer to $F^p$ as the plastic deformation tensor, although the numerical scheme we will present applies to more general inelastic deformations. According to (5), $F^p$ is a fictitious state of total deformation in which there is no elastic deformation: given an initial total deformation $F$, and a purely elastic relaxation path $F^e \rightarrow I$, the total observable deformation will evolve to $F^p$, $F \rightarrow F^p$. The state $F^p$ is a function of the deformation history of the material. We represent this history through a single scalar parameter $\kappa$, a work hardening measure, and constitutive flow rules

$$\dot{F}^p = h(\rho, g, F^p, E, \kappa) \hfill (7)$$

$$\kappa = K(\rho, g, F^p, E, \kappa) \hfill (8)$$

which depend on the state variables but not their gradients.

The equations of solid mechanics are then given by

$$\frac{\partial}{\partial t} \begin{pmatrix} \rho \\ \rho v \\ \rho E \\ g e_x \\ g e_y \\ g e_z \\ \rho F^e e_x \\ \rho F^e e_y \\ \rho F^e e_z \\ \rho \kappa \end{pmatrix} + \frac{\partial}{\partial x_\alpha} \begin{pmatrix} \rho v_\alpha \\ \rho v_\alpha - \sigma_\alpha \\ \rho E v_\alpha - v_\beta \sigma_\beta \\ g v_\alpha x_\alpha \\ g v_\alpha y_\alpha \\ g v_\alpha z_\alpha \\ \rho F^e v_\alpha x_\alpha \\ \rho F^e v_\alpha y_\alpha \\ \rho F^e v_\alpha z_\alpha \\ \rho \kappa v_\alpha \end{pmatrix} = \begin{pmatrix} 0 \\ \rho f \\ \rho(\Phi + v \cdot f) \\ (v \times (\nabla \times g^T))^T e_x \\ (v \times (\nabla \times g^T))^T e_y \\ (v \times (\nabla \times g^T))^T e_z \\ \rho e_x \\ \rho e_y \\ \rho e_z \\ \rho K \end{pmatrix} \hfill (9)$$

where $e_x, e_y, e_z$ are the Cartesian unit vectors, and $E$ is the sum of internal energy and kinetic energy, $(E = E + \frac{1}{2}v \cdot v)$. For generality, we include a heat source term $\Phi$, and a
body force vector \( f \). The system of equations (9) is abbreviated

\[
\frac{\partial U}{\partial t} + \frac{\partial F_\alpha(U)}{\partial x_\alpha} = S(U),
\]

where \( U \) is the vector of quasi-conservation-form variables \((\rho, \rho v, \rho E, \text{ etc ...})\), \( F_\alpha(U) \) is the flux in direction \( e_\alpha \), and where \( S(U) \) is the vector of source terms. Here we follow the treatment in [24] and use the inverse deformation gradient \( g = \mathcal{F}^{-1} \) as dependent variables. However, we introduce additional nonconservative terms in the evolution equations of \( g \).

We will show below that the addition of these terms leads to a well-behaved hyperbolic structure for the equations, independent of whether the curl of \( g^T \) vanishes. However, we note here that \( g = \nabla \times g^T \) satisfies the following evolution equation:

\[
\frac{\partial g}{\partial t} + \nabla \cdot (vg - g_v) = 0.
\]

In particular, if \( g \) vanishes identically at time \( t = 0 \), it vanishes at all later times.

To solve these equations we adopt a predictor-corrector strategy. For each time step, we first solve the conservative flux differencing left-hand side of (10) using fluxes derived (by solution to Riemann problems) from edge- and time-centered variables that include time-centered contributions from the source terms. The solution obtained by flux differencing is then modified by addition of the source terms, evaluated using time-centered and cell-centered variables, and acting over the full time step \( \Delta t \).

The solution to the flux differencing equations is based upon the standard high-order Godunov strategy. This strategy begins with a characteristic analysis of the equations, which makes use of the linearized 1D equations in direction \( e_\alpha \):

\[
\begin{pmatrix}
\frac{\rho}{F_{p_{ex}}} \\
\frac{\mu e_x}{F_{p_{ex}}} \\
\frac{\mu e_y}{F_{p_{ex}}} \\
\frac{\mu e_z}{F_{p_{ex}}} \\
\kappa e_{\alpha x}
\end{pmatrix} + A \frac{\partial}{\partial x_\alpha}
\begin{pmatrix}
\frac{\rho}{F_{p_{ex}}} \\
\frac{\mu e_x}{F_{p_{ex}}} \\
\frac{\mu e_y}{F_{p_{ex}}} \\
\frac{\mu e_z}{F_{p_{ex}}} \\
\kappa e_{\alpha x}
\end{pmatrix} =
\begin{pmatrix}
0 \\
0 \\
0 \\
0 \\
\Phi
\end{pmatrix}
\]

where

\[
\begin{pmatrix}
\frac{\rho}{F_{p_{ex}}} \\
\frac{\mu e_x}{F_{p_{ex}}} \\
\frac{\mu e_y}{F_{p_{ex}}} \\
\frac{\mu e_z}{F_{p_{ex}}} \\
\kappa e_{\alpha x}
\end{pmatrix} + A \frac{\partial}{\partial x_\alpha}
\begin{pmatrix}
\frac{\rho}{F_{p_{ex}}} \\
\frac{\mu e_x}{F_{p_{ex}}} \\
\frac{\mu e_y}{F_{p_{ex}}} \\
\frac{\mu e_z}{F_{p_{ex}}} \\
\kappa e_{\alpha x}
\end{pmatrix} =
\begin{pmatrix}
0 \\
0 \\
0 \\
0 \\
\Phi
\end{pmatrix}
\]

\[
\begin{pmatrix}
\frac{\rho}{F_{p_{ex}}} \\
\frac{\mu e_x}{F_{p_{ex}}} \\
\frac{\mu e_y}{F_{p_{ex}}} \\
\frac{\mu e_z}{F_{p_{ex}}} \\
\kappa e_{\alpha x}
\end{pmatrix} + A \frac{\partial}{\partial x_\alpha}
\begin{pmatrix}
\frac{\rho}{F_{p_{ex}}} \\
\frac{\mu e_x}{F_{p_{ex}}} \\
\frac{\mu e_y}{F_{p_{ex}}} \\
\frac{\mu e_z}{F_{p_{ex}}} \\
\kappa e_{\alpha x}
\end{pmatrix} =
\begin{pmatrix}
0 \\
0 \\
0 \\
0 \\
\Phi
\end{pmatrix}
\]

\[
\begin{pmatrix}
\frac{\rho}{F_{p_{ex}}} \\
\frac{\mu e_x}{F_{p_{ex}}} \\
\frac{\mu e_y}{F_{p_{ex}}} \\
\frac{\mu e_z}{F_{p_{ex}}} \\
\kappa e_{\alpha x}
\end{pmatrix} + A \frac{\partial}{\partial x_\alpha}
\begin{pmatrix}
\frac{\rho}{F_{p_{ex}}} \\
\frac{\mu e_x}{F_{p_{ex}}} \\
\frac{\mu e_y}{F_{p_{ex}}} \\
\frac{\mu e_z}{F_{p_{ex}}} \\
\kappa e_{\alpha x}
\end{pmatrix} =
\begin{pmatrix}
0 \\
0 \\
0 \\
0 \\
\Phi
\end{pmatrix}
\]

\[
\begin{pmatrix}
\frac{\rho}{F_{p_{ex}}} \\
\frac{\mu e_x}{F_{p_{ex}}} \\
\frac{\mu e_y}{F_{p_{ex}}} \\
\frac{\mu e_z}{F_{p_{ex}}} \\
\kappa e_{\alpha x}
\end{pmatrix} + A \frac{\partial}{\partial x_\alpha}
\begin{pmatrix}
\frac{\rho}{F_{p_{ex}}} \\
\frac{\mu e_x}{F_{p_{ex}}} \\
\frac{\mu e_y}{F_{p_{ex}}} \\
\frac{\mu e_z}{F_{p_{ex}}} \\
\kappa e_{\alpha x}
\end{pmatrix} =
\begin{pmatrix}
0 \\
0 \\
0 \\
0 \\
\Phi
\end{pmatrix}
\]

\[
\begin{pmatrix}
\frac{\rho}{F_{p_{ex}}} \\
\frac{\mu e_x}{F_{p_{ex}}} \\
\frac{\mu e_y}{F_{p_{ex}}} \\
\frac{\mu e_z}{F_{p_{ex}}} \\
\kappa e_{\alpha x}
\end{pmatrix} + A \frac{\partial}{\partial x_\alpha}
\begin{pmatrix}
\frac{\rho}{F_{p_{ex}}} \\
\frac{\mu e_x}{F_{p_{ex}}} \\
\frac{\mu e_y}{F_{p_{ex}}} \\
\frac{\mu e_z}{F_{p_{ex}}} \\
\kappa e_{\alpha x}
\end{pmatrix} =
\begin{pmatrix}
0 \\
0 \\
0 \\
0 \\
\Phi
\end{pmatrix}
\]

\[
\begin{pmatrix}
\frac{\rho}{F_{p_{ex}}} \\
\frac{\mu e_x}{F_{p_{ex}}} \\
\frac{\mu e_y}{F_{p_{ex}}} \\
\frac{\mu e_z}{F_{p_{ex}}} \\
\kappa e_{\alpha x}
\end{pmatrix} + A \frac{\partial}{\partial x_\alpha}
\begin{pmatrix}
\frac{\rho}{F_{p_{ex}}} \\
\frac{\mu e_x}{F_{p_{ex}}} \\
\frac{\mu e_y}{F_{p_{ex}}} \\
\frac{\mu e_z}{F_{p_{ex}}} \\
\kappa e_{\alpha x}
\end{pmatrix} =
\begin{pmatrix}
0 \\
0 \\
0 \\
0 \\
\Phi
\end{pmatrix}
\]

\[
\begin{pmatrix}
\frac{\rho}{F_{p_{ex}}} \\
\frac{\mu e_x}{F_{p_{ex}}} \\
\frac{\mu e_y}{F_{p_{ex}}} \\
\frac{\mu e_z}{F_{p_{ex}}} \\
\kappa e_{\alpha x}
\end{pmatrix} + A \frac{\partial}{\partial x_\alpha}
\begin{pmatrix}
\frac{\rho}{F_{p_{ex}}} \\
\frac{\mu e_x}{F_{p_{ex}}} \\
\frac{\mu e_y}{F_{p_{ex}}} \\
\frac{\mu e_z}{F_{p_{ex}}} \\
\kappa e_{\alpha x}
\end{pmatrix} =
\begin{pmatrix}
0 \\
0 \\
0 \\
0 \\
\Phi
\end{pmatrix}
\]
EULERIAN GODUNOV METHOD FOR SOLID MECHANICS

\[
A = \begin{pmatrix}
    v_\alpha & \rho e_\alpha^T & 0 & 0 & 0 & 0 & 0 & 0 & 0 \\
    0 & v_\alpha I & 0 & 0 & 0 & 0 & 0 & 0 & 0 \\
    0 & -(\sigma e_\alpha)^T / \rho & v_\alpha & 0 & 0 & 0 & 0 & 0 & 0 \\
    0 & \delta_{\alpha\alpha} & v_\alpha I & 0 & 0 & 0 & 0 & 0 & 0 \\
    0 & \delta_{\gamma\alpha} & v_\alpha I & 0 & 0 & 0 & 0 & 0 & 0 \\
    0 & \delta_{\alpha\alpha} & 0 & 0 & 0 & 0 & v_\alpha I & 0 & 0 \\
    0 & 0 & 0 & 0 & 0 & 0 & v_\alpha I & 0 & 0 \\
    0 & 0 & 0 & 0 & 0 & 0 & v_\alpha I & 0 & 0 \\
    0 & -A_{\alpha\alpha} & 0 & 0 & 0 & 0 & 0 & 0 & v_\alpha I
\end{pmatrix}
\]

with
\[
A_{\alpha\beta} = -\frac{\partial \sigma e_\alpha}{\partial e_\beta} g
\]

and
\[
b_\alpha = \frac{\partial \sigma e_\alpha}{\partial F_p} h + \frac{\partial \sigma e_\alpha}{\partial K} K + \frac{\partial \sigma e_\alpha}{\partial \Phi} \Phi.
\]

The eigenvalue decomposition of \( A \) uses the technique of eigenvalue deflation, and hinges upon recognition of the matrices \( A_{\alpha\alpha} \) as being acoustic wave propagation tensors for waves traveling in direction \( e_\alpha \),

\[
\rho \ddot{u}_\beta = (A_{\alpha\alpha})_{\gamma\delta} \frac{\partial^2 u_\delta}{\partial x_\beta \partial x_\gamma},
\]

where \( u \) is the displacement vector. The matrices \( A_{\alpha\alpha} \) are positive definite as a requirement of thermodynamic stability. This is made clear by writing \( A_{\alpha\alpha} \) in terms of gradients of the spatial displacements \( \ddot{u} \) defined relative to the current configuration,

\[
(A_{\alpha\alpha})_{\beta\gamma} = \rho \frac{\partial^2 \Phi}{\partial \ddot{u}_\alpha \partial \ddot{u}_\gamma}.
\]

Here, \( \ddot{u}_\beta \) is related to the deformation tensor \( F_{\beta\alpha} \) with the reference coordinate frame \( \{a\} \) chosen to correspond to the current spatial frame \( \{x\} \):

\[
\ddot{u}_\beta = F_{\beta\alpha} \bigg|_{\{a\} = \{x\}} - \delta_{\beta\alpha}.
\]

\( A_{\alpha\alpha} \) is therefore a component of the Hessian of \( \Phi \), which is positive definite for a thermodynamically stable material, and consequently \( A_{\alpha\alpha} \) has positive real eigenvalues and three linearly-independent eigenvectors.

Recognizing \( A_{\alpha\alpha} \) as being the acoustic wave propagation tensor suggests the wave equation solution

\[
A_{\alpha\alpha} X_{ac} = \rho X_{ac} \Lambda_{ac}^2
\]

where \( \Lambda_{ac} \) is the diagonal matrix of acoustic wave speeds \( c \), \( \Lambda_{ac} = \text{diag}(c_1, c_2, c_3) \), and \( X_{ac} \) are the acoustic displacement vectors.
The linearized 1D matrix $A$ then has eigenvalue decomposition

$$A = X\Lambda X^{-1}$$

with $X$, the matrix of right eigenvectors, given by

$$X = \begin{pmatrix}
1 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & -pe_a^T X_{ac} & -pe_a^T X_{ac} \\
0 & 1 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & X_{ac}\Lambda_{ac} & -X_{ac}\Lambda_{ac} \\
0 & 0 & I & 0 & 0 & 0 & 0 & 0 & 0 & 0 & -gX_{ac}\delta_{xa} & -gX_{ac}\delta_{xa} \\
0 & 0 & 0 & I & 0 & 0 & 0 & 0 & 0 & 0 & -gX_{ac}\delta_{ya} & -gX_{ac}\delta_{ya} \\
0 & 0 & 0 & 0 & I & 0 & 0 & 0 & 0 & 0 & -gX_{ac}\delta_{za} & -gX_{ac}\delta_{za} \\
0 & 0 & 0 & 0 & 0 & I & 0 & 0 & 0 & 0 & 0 & 0 \\
0 & 0 & 0 & 0 & 0 & 0 & I & 0 & 0 & 0 & 0 & 0 \\
0 & 0 & 0 & 0 & 0 & 0 & 0 & I & 0 & 0 & 0 & 0 \\
0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & I & 0 & 0 & 0 \\
0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & I & 0 & 0 \\
0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & X_{ac}\Lambda_{ac}^2 & X_{ac}\Lambda_{ac}^2 \rho
\end{pmatrix}, \quad (20)$$

and $\Lambda$, the diagonal matrix of eigenvalues, given by

$$\Lambda = \begin{pmatrix}
v_a & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 \\
0 & v_a & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 \\
0 & 0 & v_aI & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 \\
0 & 0 & 0 & v_aI & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 \\
0 & 0 & 0 & 0 & v_aI & 0 & 0 & 0 & 0 & 0 & 0 & 0 \\
0 & 0 & 0 & 0 & 0 & v_aI & 0 & 0 & 0 & 0 & 0 & 0 \\
0 & 0 & 0 & 0 & 0 & 0 & v_aI & 0 & 0 & 0 & 0 & 0 \\
0 & 0 & 0 & 0 & 0 & 0 & 0 & v_aI - \Lambda_{ac} & 0 & 0 & 0 \\
0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & v_aI + \Lambda_{ac} & 0 & 0 \\
\end{pmatrix} \quad (21)$$

The wave speeds are Galilean invariant, and properly analogous to the Lagrangian representation, with three $-$ waves with velocities $v_a - c_\gamma$, three $+$ waves with velocities $v_a + c_\gamma$, and 21 material waves with speeds $v_a$.

$X^{-1}$, the inverse of $X$, is given by

$$X^{-1} = \begin{pmatrix}
1 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & -pe_a^T X_{ac} & -pe_a^T X_{ac} \\
0 & 1 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & -(\sigma e_\alpha)^T X_{ac} \Lambda_{ac}^{-2} X_{ac}^{-1} / \rho^2 & -(\sigma e_\alpha)^T X_{ac} \Lambda_{ac}^{-2} X_{ac}^{-1} / \rho \\
0 & 0 & I & 0 & 0 & 0 & 0 & 0 & 0 & 0 & gX_{ac} \Lambda_{ac}^{-2} X_{ac}^{-1} \delta_{xa}/\rho & gX_{ac} \Lambda_{ac}^{-2} X_{ac}^{-1} \delta_{xa}/\rho \\
0 & 0 & 0 & I & 0 & 0 & 0 & 0 & 0 & 0 & gX_{ac} \Lambda_{ac}^{-2} X_{ac}^{-1} \delta_{ya}/\rho & gX_{ac} \Lambda_{ac}^{-2} X_{ac}^{-1} \delta_{ya}/\rho \\
0 & 0 & 0 & 0 & I & 0 & 0 & 0 & 0 & 0 & gX_{ac} \Lambda_{ac}^{-2} X_{ac}^{-1} \delta_{za}/\rho & gX_{ac} \Lambda_{ac}^{-2} X_{ac}^{-1} \delta_{za}/\rho \\
0 & 0 & 0 & 0 & 0 & I & 0 & 0 & 0 & 0 & 0 & 0 \\
0 & 0 & 0 & 0 & 0 & 0 & I & 0 & 0 & 0 & 0 & 0 \\
0 & 0 & 0 & 0 & 0 & 0 & 0 & I & 0 & 0 & 0 & 0 \\
0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & I & 0 & 0 & 0 \\
0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & I & 0 & 0 \\
0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & X_{ac}\Lambda_{ac}^{-2} X_{ac}^{-1} & X_{ac}\Lambda_{ac}^{-2} X_{ac}^{-1} \\
0 & -\frac{1}{2} \Lambda_{ac}^{-1} X_{ac}^{-1} & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & \frac{1}{2} \Lambda_{ac}^{-2} X_{ac} & \frac{1}{2} \Lambda_{ac}^{-2} X_{ac} \\
0 & -\frac{1}{2} \Lambda_{ac}^{-1} X_{ac}^{-1} & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & \frac{1}{2} \Lambda_{ac}^{-2} X_{ac} & \frac{1}{2} \Lambda_{ac}^{-2} X_{ac} \\
\end{pmatrix} \quad (22)$$

and

$$X^{-1} = \begin{pmatrix}
1 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & -pe_a^T X_{ac} \Lambda_{ac}^{-2} X_{ac}^{-1} & -pe_a^T X_{ac} \Lambda_{ac}^{-2} X_{ac}^{-1} \\
0 & 1 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & -(\sigma e_\alpha)^T X_{ac} \Lambda_{ac}^{-2} X_{ac}^{-1} / \rho^2 & -(\sigma e_\alpha)^T X_{ac} \Lambda_{ac}^{-2} X_{ac}^{-1} / \rho \\
0 & 0 & I & 0 & 0 & 0 & 0 & 0 & 0 & 0 & gX_{ac} \Lambda_{ac}^{-2} X_{ac}^{-1} \delta_{xa}/\rho & gX_{ac} \Lambda_{ac}^{-2} X_{ac}^{-1} \delta_{xa}/\rho \\
0 & 0 & 0 & I & 0 & 0 & 0 & 0 & 0 & 0 & gX_{ac} \Lambda_{ac}^{-2} X_{ac}^{-1} \delta_{ya}/\rho & gX_{ac} \Lambda_{ac}^{-2} X_{ac}^{-1} \delta_{ya}/\rho \\
0 & 0 & 0 & 0 & I & 0 & 0 & 0 & 0 & 0 & gX_{ac} \Lambda_{ac}^{-2} X_{ac}^{-1} \delta_{za}/\rho & gX_{ac} \Lambda_{ac}^{-2} X_{ac}^{-1} \delta_{za}/\rho \\
0 & 0 & 0 & 0 & 0 & I & 0 & 0 & 0 & 0 & 0 & 0 \\
0 & 0 & 0 & 0 & 0 & 0 & I & 0 & 0 & 0 & 0 & 0 \\
0 & 0 & 0 & 0 & 0 & 0 & 0 & I & 0 & 0 & 0 & 0 \\
0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & I & 0 & 0 & 0 \\
0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & I & 0 & 0 \\
0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & X_{ac}\Lambda_{ac}^{-2} X_{ac} & X_{ac}\Lambda_{ac}^{-2} X_{ac} \\
0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & X_{ac}\Lambda_{ac}^{-2} X_{ac} \\
0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & X_{ac}\Lambda_{ac}^{-2} X_{ac} \\
\end{pmatrix} \quad (23)$$
3. NUMERICAL METHOD: 1D

In 1D we discretize space into cells, indexed with subscript \(i\), with width \(\Delta x_i\). Time is discretized in steps of \(\Delta t\) with integer superscript \(n\); \(t^{n+1} - t^n = \Delta t\). The generalization to 2D and 3D is similar, with indices \(j\) and \(k\) used for the second and third dimensions respectively. Half-integral subscript indices represent edge-centered quantities. Lower case Greek subscripts are used to denote vector and tensor indices.

We begin by evaluation of the equation of state in each cell to determine the Cauchy stress \(\sigma\), the acoustic wave propagation tensor \(A_{\alpha\alpha}\), and the thermodynamic derivatives \(\partial_\sigma/\partial E|_{\epsilon,J,Y,\Omega,\kappa,\partial\sigma/\partial F|_{\epsilon,J,Y,\Omega,\kappa,\partial \sigma/\partial K|_{\epsilon,J,Y,\Omega}}\).

Next, we evaluate the 1D slopes \(dq/dx\) of the 27 primitive cell-centered variables \(q\),

\[
q = (\rho, v, E, g, F, \kappa, \sigma e_{\alpha}).
\]  

We construct these slopes beginning with the van Leer slope in cell \(i\) which uses the monotinized limiter \([25]\):

\[
\left(\frac{\partial q}{\partial x}\right)_i^{UL} = \text{sign}(q_{i+1} - q_{i-1}) \min \left(\frac{2|q_{i+1} - q_{i-1}|}{\Delta x_{i-1} + 2\Delta x_i + \Delta x_{i+1}}, \frac{2|q_i - q_{i-1}|}{\Delta x_i}, \frac{2|q_{i+1} - q_i|}{\Delta x_i}\right).
\]  

A 4th-order accurate slope is then constructed as \([5]\):

\[
\left(\frac{\partial q}{\partial x}\right)_i^{4th} = \left[\frac{q_{i+1} - \frac{1}{4}\Delta x_{i+1} \left(\frac{\partial q}{\partial x}\right)_{i+1}^{UL}}{\frac{1}{4}\Delta x_{i-1} + \Delta x_i + \frac{1}{4}\Delta x_{i+1}}\right] - \left[\frac{q_{i-1} + \frac{1}{4}\Delta x_{i-1} \left(\frac{\partial q}{\partial x}\right)_{i-1}^{UL}}{\frac{1}{4}\Delta x_{i-1} + \Delta x_i + \frac{1}{4}\Delta x_{i+1}}\right].
\]  

To prevent overshoot and ringing, dissipation at strong shocks may be introduced via a "flattening parameter" \(\chi\), \(0 \leq \chi \leq 1\), whence \([5, 7, 6]\)

\[
\left(\frac{\partial q}{\partial x}\right)_i = \chi \left(\frac{\partial q}{\partial x}\right)_i^{4th}.
\]  

The determination of this flattening parameter is described in a later section.

These limited slopes are used to construct time-centered edge-valued estimates of the primitive variables. The exact solution of the linearized equations, which we abbreviate as

\[
\frac{\partial q}{\partial t} + A \frac{\partial q}{\partial x} = s,
\]

gives time-centered edge values

\[
q_{R,i+1/2}^{n+1/2} = q_i^n - \frac{\Delta x_i}{2} X_i \left(\frac{\Delta t}{\Delta x_i} \Lambda_i + I\right) X_i^{-1} \left(\frac{\partial q}{\partial x}\right)_i + \frac{\Delta t}{2} s_i
\]  

\[
q_{L,i+1/2}^{n+1/2} = q_i^n - \frac{\Delta x_i}{2} X_i \left(\frac{\Delta t}{\Delta x_i} \Lambda_i - I\right) X_i^{-1} \left(\frac{\partial q}{\partial x}\right)_i + \frac{\Delta t}{2} s_i.
\]

However, this construction uses both upwind and downwind characteristics. We make the method strictly upwind by filtering out the downwind characteristics:

\[
q_{R,i-1/2}^{n+1/2} = q_i^n - \frac{\Delta x_i}{2} X_i \left(\frac{\Delta t}{\Delta x_i} \Lambda_i + I\right) X_i^{-1} \left(\frac{\partial q}{\partial x}\right)_i + \frac{\Delta t}{2} s_i
\]  

\[
q_{L,i+1/2}^{n+1/2} = q_i^n - \frac{\Delta x_i}{2} X_i \left(\frac{\Delta t}{\Delta x_i} \Lambda_i - I\right) X_i^{-1} \left(\frac{\partial q}{\partial x}\right)_i + \frac{\Delta t}{2} s_i.
\]  

\[
q_{R,i-1/2}^{n+1/2} = q_i^n - \frac{\Delta x_i}{2} X_i \left(\frac{\Delta t}{\Delta x_i} \Lambda_i + I\right) X_i^{-1} \left(\frac{\partial q}{\partial x}\right)_i + \frac{\Delta t}{2} s_i
\]  

\[
q_{L,i+1/2}^{n+1/2} = q_i^n - \frac{\Delta x_i}{2} X_i \left(\frac{\Delta t}{\Delta x_i} \Lambda_i - I\right) X_i^{-1} \left(\frac{\partial q}{\partial x}\right)_i + \frac{\Delta t}{2} s_i.
\]  

\[
q_{R,i-1/2}^{n+1/2} = q_i^n - \frac{\Delta x_i}{2} X_i \left(\frac{\Delta t}{\Delta x_i} \Lambda_i + I\right) X_i^{-1} \left(\frac{\partial q}{\partial x}\right)_i + \frac{\Delta t}{2} s_i
\]  

\[
q_{L,i+1/2}^{n+1/2} = q_i^n - \frac{\Delta x_i}{2} X_i \left(\frac{\Delta t}{\Delta x_i} \Lambda_i - I\right) X_i^{-1} \left(\frac{\partial q}{\partial x}\right)_i + \frac{\Delta t}{2} s_i.
\]
\[ q_{L,i+1/2}^{n+1/2} = q_{L,i}^{n} - \frac{\Delta x_i}{2} X_i \mathcal{P}^+ \left( \frac{\Delta t}{\Delta x_i} \Lambda_i - I \right) X_i^{-1} \left( \frac{\partial q}{\partial x} \right)_i + \frac{\Delta t}{2} s_i \]  

with projection operators \( \mathcal{P}_\pm \) defined as:

\[
\begin{align*}
(\mathcal{P}_- \left( \frac{\Delta t}{\Delta x} \Lambda + I \right))_{\alpha\beta} &= \left\{ \begin{array}{ll}
(\frac{\Delta t}{\Delta x} \Lambda_{\alpha\alpha} + 1) \delta_{\alpha\beta} & \Lambda_{\alpha\alpha} \leq 0 \\
0 & \Lambda_{\alpha\alpha} > 0
\end{array} \right. \\
(\mathcal{P}_+ \left( \frac{\Delta t}{\Delta x} \Lambda - I \right))_{\alpha\beta} &= \left\{ \begin{array}{ll}
(\frac{\Delta t}{\Delta x} \Lambda_{\alpha\alpha} - 1) \delta_{\alpha\beta} & \Lambda_{\alpha\alpha} \geq 0 \\
0 & \Lambda_{\alpha\alpha} < 0
\end{array} \right.
\end{align*}
\]

At each cell edge \((i+1/2)\), time-centered values are thus obtained from the left \((i)\) and right \((i+1)\) neighboring cells. These edge values are then used to pose a Riemann problem: an initial value problem with constant left and right initial states given by \(q_{L,i+1/2}^{n+1/2}\) and \(q_{R,i+1/2}^{n+1/2}\) respectively. We approximate the solution to the Riemann problem by decomposing the jump \(q_{R,i+1/2}^{n+1/2} - q_{L,i+1/2}^{n+1/2}\) in terms of the eigenvectors \(X\) of the linearized coefficients \(\Lambda\). Specifically,

\[
q_{R,i+1/2}^{n+1/2} - q_{L,i+1/2}^{n+1/2} = \sum_{\gamma=1}^{27} \varphi_{\gamma} X_{\gamma,i+1/2}
\]

where eigenvector column \(X_{\gamma,i+1/2}\) is evaluated with certain \(L\) or cell-\(i\) properties if \(\Lambda_{\gamma\gamma}\) is a member of the \(-\) family (i.e., of the form \(ve_{\alpha} - c\)), or with certain \(R\) or cell-(\(i+1\)) properties if \(\Lambda_{\gamma\gamma}\) is a member of the \(+\) family (of the form \(ve_{\alpha} + c\), as given by the discretization:

\[
X_{i+1/2} = \begin{pmatrix}
1 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & -\rho_i e_{\alpha}^T X_{ac,i} & -\rho_{i+1} e_{\alpha}^T X_{ac,i+1} \\
0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & X_{ac,i} \Lambda_{ac,i} & -X_{ac,i+1} \Lambda_{ac,i+1} \\
0 & 1 & 0 & 0 & 0 & 0 & 0 & 0 & (\varepsilon_{i+1/2} e_{\alpha})^T X_{ac,i} & (\varepsilon_{i+1/2} e_{\alpha})^T X_{ac,i+1} \\
0 & 0 & I & 0 & 0 & 0 & 0 & 0 & -g_{i+1/2,L} X_{ac,i} \delta_{2\alpha} & -g_{i+1+1/2,R} X_{ac,i} \delta_{2\alpha} \\
0 & 0 & 0 & I & 0 & 0 & 0 & 0 & -g_{i+1/2,L} X_{ac,i} \delta_{2\alpha} & -g_{i+1+1/2,R} X_{ac,i} \delta_{2\alpha} \\
0 & 0 & 0 & 0 & I & 0 & 0 & 0 & 0 & 0 \\
0 & 0 & 0 & 0 & 0 & I & 0 & 0 & 0 & 0 \\
0 & 0 & 0 & 0 & 0 & 0 & I & 0 & 0 & 0 \\
0 & 0 & 0 & 0 & 0 & 0 & 0 & I & 0 & 0 \\
0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & X_{ac,i} \Lambda_{ac,i} \rho_i & X_{ac,i+1} \Lambda_{ac,i+1} \rho_{i+1}
\end{pmatrix}
\]

In this expression, the density \(\rho\), and the components \((X_{ac}, \Lambda_{ac})\) of the acoustic propagation tensor, are evaluated at the cell centers in order to avoid multiple evaluations of the equation of state.

From the coefficients \(\varphi_{\gamma}\) of the jump decomposition, the material velocity \(v^* \cdot e_{\alpha}\) at the cell edge is determined by adding to the \(L\) state the contributions of the \(-\) family, or by subtracting from the \(R\) state the contributions of the \(+\) family:

\[
v_{i+1/2}^* \cdot e_{\alpha} = v^* L \equiv v_{i+1/2,L} \cdot e_{\alpha} + \varphi_6 X_{6\beta,i+1/2} + \varphi_7 X_{7\beta,i+1/2} + \varphi_8 X_{8\beta,i+1/2}
\]
or

\[ v_{i+1/2}^{* \cdot e_{\alpha}} = R \equiv v_{i+1/2,R \cdot e_{\alpha}} - \varphi_{9} X_{9, \beta, i+1/2} - \varphi_{10} X_{10, \beta, i+1/2} - \varphi_{11} X_{11, \beta, i+1/2} \]

where \( \beta = 2, 3, 4 \) for directions \( e_{\alpha} \) equal to \( e_{x}, e_{y}, e_{z} \) respectively. We average the results of these calculations to determine the normal-direction edge velocity \( v_{i+1/2}^{* \cdot e_{\alpha}} \).

\[ v_{i+1/2}^{* \cdot e_{\alpha}} = \frac{1}{2} \left( v^{L} + v^{R} \right) . \]

For other properties to be evaluated at the cell edge as solutions to the Riemann problem we do not average the values evaluated from the \( L \) and the \( R \) states as above. Instead, we evaluate from the \( L \) state if \( v_{i+1/2}^{* \cdot e_{\alpha}} \) is positive, or from the \( R \) state if \( v_{i+1/2}^{* \cdot e_{\alpha}} \) is negative. Only if \( v_{i+1/2}^{* \cdot e_{\alpha}} \) is approximately zero do we average these estimates. The evaluations include only upwind characteristics by writing

\[
q_{i+1/2}^{*} = \begin{cases} 
q_{i+1/2,R} + \sum_{\gamma} w_{L, \gamma} \varphi_{7} X_{\gamma} & v_{i+1/2}^{* \cdot e_{\alpha}} > \epsilon \\
q_{i+1/2,L} + \sum_{\gamma} w_{R, \gamma} \varphi_{7} X_{\gamma} & v_{i+1/2}^{* \cdot e_{\alpha}} < -\epsilon \\
\frac{1}{2} \left( q_{i+1/2,L} + q_{i+1/2,R} + \sum_{\gamma} (w_{L, \gamma} + w_{R, \gamma}) \varphi_{7} X_{\gamma} \right) & |v_{i+1/2}^{* \cdot e_{\alpha}}| \leq \epsilon
\end{cases}
\]

with

\[
w_{L, \gamma} = \begin{cases} 
1 & \Lambda_{\gamma,i} - v_{i} \cdot e_{\alpha} + v_{i+1/2}^{* \cdot e_{\alpha}} < -\epsilon \\
0 & \text{otherwise}
\end{cases}
\]

\[
w_{R, \gamma} = \begin{cases} 
1 & \Lambda_{\gamma,i+1} - v_{i+1} \cdot e_{\alpha} + v_{i+1/2}^{* \cdot e_{\alpha}} > \epsilon \\
0 & \text{otherwise}
\end{cases}
\]

\( w_{L, \gamma} \) is 1 when eigenvalue \( \gamma \), estimated using the \( * \) value of the material velocity together with the \( i \) cell-centered acoustic wave speeds, is negative; and 0 otherwise. \( w_{R, \gamma} \) is 1 when the approximated value of eigenvalue \( \gamma \) is positive; and 0 otherwise. In our computations presented below, we use a value of \( \epsilon = 10^{-9} \).

By this procedure, we obtain the edge \( * \) value solutions of the Riemann problem, \( \rho^{*}, v^{*}, E^{*}, \sigma^{*}, F^{*}, \kappa^{*}, \) and \( (\sigma e)^{*} \). These are then used to compute edge-valued fluxes (cf. (9,10)). For example, in direction \( e_{z} \),

\[
F_{x,i+1/2}^{*} = \begin{pmatrix}
\rho v_{z} \\
\rho v_{z}^{2} - \sigma_{zz} \\
\rho v_{y} v_{z} - \sigma_{yz} \\
\rho v_{z} u_{z} - \sigma_{zz} \\
\rho \dot{E} v_{z} - v_{z} \sigma_{zx} - v_{y} \sigma_{yz} - v_{z} \sigma_{xz} \\
v_{z} g e_{z} + v_{y} g e_{y} + v_{z} g e_{z} \\
0 \\
0 \\
\rho v_{z} F^{*} e_{z} \\
\rho v_{z} F^{*} e_{y} \\
\rho v_{z} F^{*} e_{z} \\
\rho v_{z} \kappa
\end{pmatrix}_{i+1/2}
\]
In 1D we obtain a preliminary update $\bar{U}$ of the variables $U$ by conservatively differencing the fluxes:

$$\bar{U}_i^{n+1} = U_i^n - \frac{\Delta t}{\Delta x_i} (F_i^{n+1/2} - F_i^{n-1/2})$$

(40)

the final time-$(n+1)$ value of the variables $U$ is obtained from the preliminary values $\bar{U}$ by addition of the source terms $S$

$$U_i^{n+1} = \bar{U}_i^{n+1} + \Delta t S_i$$

(41)

We discretize this in the general 3D (Cartesian) case as

$$U_i^{n+1} = \bar{U}_i^{n+1} + \Delta t$$

(42)

In 1D we use the 3D discretization above, but retain only terms in $\partial / \partial x$ and $\partial^2 / \partial x^2$, and omit derivatives in all transverse directions.

In the above expression, time-centered terms (e.g., $\rho_i^{n+1/2}$) are estimated with

$$q_{ijk}^{n+1/2} \approx \frac{1}{2} (q_{ijk}^n + q_{ijk}^{n+1}),$$

(43)

except for the $g^*$s appearing in the $(\mathbf{v} \times \mathbf{G})$ terms. These are obtained at the half time step and cell edges as components of the Riemann problem solutions.

4. NUMERICAL METHOD: 2D AND 3D

To extend the 1D method described above to multiple spatial dimensions, we use a spatially-unsplit fully corner-coupled 2nd-order accurate scheme after [6] and [21]. In 2D, this predictor-corrector approach begins by estimating the 1D $x-$ and $y-$ fluxes at each cell edge, using the higher-order 1D approach described in the previous section. These predictor fluxes, $\bar{F}_x$ and $\bar{F}_y$, are given schematically as solutions to the Riemann problem
The predictor fluxes are used to pose a corrector problem, wherein the edge values are augmented by transverse predictor fluxes. Schematically,

\[
\begin{align*}
F_{x,i+1/2,j} &= F_x(R(q^{n+1/2}_{xL,i+1/2,j}; q^{n+1/2}_{xR,i+1/2,j})) \\
F_{y,i,j+1/2} &= F_y(R(q^{n+1/2}_{yL,i,j+1/2}; q^{n+1/2}_{yR,i,j+1/2}))
\end{align*}
\]  

with, for example,

\[
\begin{align*}
q^{n+1/2}_{xL,i+1/2,j} &= q^{n+1/2}_{xL,i+1/2,j} - \frac{\Delta t}{2\Delta y_j} \left( F^y_{i,j+1/2} - F^y_{i,j-1/2} \right) \\
&\quad + \frac{\Delta t}{2\Delta y_j} \sum_{\gamma \delta} \left( u_{i,j} \times e_y \times \left( g^y_{i,j+1/2} - g^y_{i,j-1/2} \right) \right)_{T, \gamma \delta} \Gamma(\gamma, \delta).
\end{align*}
\]

\(\Gamma(\gamma, \delta)\) is a vector introduced to align the elements of the matrix \((v \times \nabla \times g^T)^T\) with the appropriate elements of the vector \(q\):

\[
\Gamma(\gamma, \delta) = (0, 0, 0, 0, 0, \delta_1 \delta_8, 0, 0, \delta_2 \delta_1, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0)^T.
\]

In 2D there are therefore 4 Riemann problems solved per cell: 2 in the predictor and 2 in corrector steps.

In setting up the corrector step, the components \(p, v, \mathcal{E}, g, \mathcal{F}^p, \) and \(\kappa\) of the vectors \(q'\) are updated as indicated above (Eq. (46)). Our 1D Riemann solver also requires time-centered edge values of the stresses, \((\sigma e_\alpha)_{L/R}\) in direction \(e_\alpha\), and these components of \(q'\) are calculated by updating the \((\sigma e_\alpha)_{L/R}\) components of \(q_{L/R}\) with the change in stress accompanying the changes \(q'_{L/R} - q_{L/R}\) in \(\mathcal{E}, g, \mathcal{F}^p, \) and \(\kappa\) using cell-centered thermodynamic derivatives. For example,

\[
\begin{align*}
(\sigma' e_\alpha)_{L,i+1/2,j} &= (\sigma e_\alpha)_{L,i+1/2,j} + \left( \frac{\partial \sigma e_\alpha}{\partial \mathcal{E}} \bigg|_{y, \mathcal{F}^p, \kappa} \right)_{ij}^{n} \left( \mathcal{E}^i_{L,i+1/2,j} - \mathcal{E}^i_{L,i+1/2,j} \right) \\
&\quad + \left( \frac{\partial \sigma e_\alpha}{\partial \mathcal{E}} \bigg|_{y, \mathcal{F}^p, \kappa} \right)_{ij}^{n} \left( \kappa_{L,i+1/2,j}^i - \kappa_{L,i+1/2,j}^i \right) \\
&\quad + \sum_{\gamma \delta} \left( \frac{\partial \sigma e_\alpha}{\partial \mathcal{F}^p} \bigg|_{y, \mathcal{F}^p, \kappa, \mathcal{E}} \right)_{ij}^{n} \left( (g_{\gamma \delta})_{L,i+1/2,j}^i - (g_{\gamma \delta})_{L,i+1/2,j}^i \right) \\
&\quad + \sum_{\gamma \delta} \left( \frac{\partial \sigma e_\alpha}{\partial \mathcal{F}^p} \bigg|_{y, \mathcal{F}^p, \kappa, \mathcal{E}} \right)_{ij}^{n} \left( (\mathcal{F}^p_{\gamma \delta})_{L,i+1/2,j}^i - (\mathcal{F}^p_{\gamma \delta})_{L,i+1/2,j}^i \right)
\end{align*}
\]

By employing this approximation we require only one equation of state evaluation per time step per cell for problems involving only elasticity. In problems that also include plasticity,
additional equation of state evaluations are required for the computation of plastic source terms.

In 3D there are 2 corrector steps: first,

\[
\tilde{F}_{i+1/2,j,k} = F_x\left(\mathcal{R}\left(q^{n+1/2}_{xL,i+1/2,j,k}, q^{n+1/2}_{xR,i+1/2,j,k}\right)\right)
\]

\[
\tilde{F}_{i+1/2,j,k} = F_y\left(\mathcal{R}\left(q^{n+1/2}_{yL,i,j+1/2,k}, q^{n+1/2}_{yR,i,j+1/2,k}\right)\right)
\]

\[
\tilde{F}_{i,j,k+1/2} = F_z\left(\mathcal{R}\left(q^{n+1/2}_{zL,i,j,k+1/2}, q^{n+1/2}_{zR,i,j,k+1/2}\right)\right)
\]

then

\[
\tilde{F}_{i+1/2,j,k} = F_x\left(\mathcal{R}\left(q^{n+1/2}_{xL,i+1/2,j,k}, q^{n+1/2}_{xR,i+1/2,j,k}\right)\right)
\]

\[
\tilde{F}_{i+1/2,j,k} = F_y\left(\mathcal{R}\left(q^{n+1/2}_{yL,i,j+1/2,k}, q^{n+1/2}_{yR,i,j+1/2,k}\right)\right)
\]

\[
\tilde{F}_{i,j,k+1/2} = F_z\left(\mathcal{R}\left(q^{n+1/2}_{zL,i,j,k+1/2}, q^{n+1/2}_{zR,i,j,k+1/2}\right)\right)
\]

with, e.g.,

\[
q^{n+1/2}_{xL,i+1/2,j,k} = q^{n+1/2}_{xL,i+1/2,j,k} - \Delta t \frac{\Delta y_j}{2} \left( \tilde{F}_{i,j+1/2,k} - \tilde{F}_{i,j-1/2,k} \right)
\]

\[
+ \Delta t \frac{\Delta y_j}{3 \Delta y_j} \sum_{\gamma \delta} \left( v_{i,j,k} \times e_y \times \left( \tilde{g}^{y}_{i,j+1/2,k} - \tilde{g}^{y}_{i,j-1/2,k} \right) \right)^T \Gamma(\gamma, \delta).
\]

The final fluxes, which enter the conservative differencing step of the integration, are then computed as:

\[
F_{x,i+1/2,j,k} = F_x\left(\mathcal{R}\left(q^{n+1/2}_{xL,i+1/2,j,k}, q^{n+1/2}_{xR,i+1/2,j,k}\right)\right)
\]

\[
F_{y,i,j+1/2,k} = F_y\left(\mathcal{R}\left(q^{n+1/2}_{yL,i,j+1/2,k}, q^{n+1/2}_{yR,i,j+1/2,k}\right)\right)
\]

\[
F_{z,i,j,k+1/2} = F_z\left(\mathcal{R}\left(q^{n+1/2}_{zL,i,j,k+1/2}, q^{n+1/2}_{zR,i,j,k+1/2}\right)\right)
\]

with, e.g.,

\[
q^{n+1/2}_{zL,i+1/2,j,k} = q^{n+1/2}_{zL,i+1/2,j,k} - \Delta t \frac{\Delta z_k}{2} \left( \tilde{F}_{i,j+1/2,k} - \tilde{F}_{i,j-1/2,k} \right)
\]

\[
+ \Delta t \frac{\Delta z_k}{2 \Delta y_j} \sum_{\gamma \delta} \left( v_{i,j,k} \times e_z \times \left( \tilde{g}^{z}_{i,j+1/2,k} - \tilde{g}^{z}_{i,j-1/2,k} \right) \right)^T \Gamma(\gamma, \delta)
\]

\[
+ \Delta t \frac{\Delta z_k}{2 \Delta y_j} \sum_{\gamma \delta} \left( v_{i,j,k} \times e_z \times \left( \tilde{g}^{z}_{i,j+1/2,k} - \tilde{g}^{z}_{i,j-1/2,k} \right) \right)^T \Gamma(\gamma, \delta)
\]
There are a total of 12 Riemann solves per unit cell in 3D: 9 in the predictor steps, and 3 in the corrector step. The $\sigma$ components of the vectors $q$ and $q''$ are computed as in the 2D case.

5. PLASTIC SOURCE TERMS

We present here an associated plasticity evolution equation for the rate of change of the plastic deformation tensor $\mathcal{F}^p$ with time. The more common approach (e.g., [22, 18]) is to consider evolution equations for the plastic strain $\eta^p = \frac{1}{2} (\mathcal{F}^p T \mathcal{F}^p - I)$, the plastic Green tensor $C^p = \mathcal{F}^p T \mathcal{F}^p$, or the plastic Finger tensor $b^p = \mathcal{F}^p T \mathcal{F}^p T$. We choose instead to evolve the full 9-component plastic deformation tensor $\mathcal{F}^p$. This choice is necessary to be capable of modeling arbitrary crystal systems (see e.g., [23]). For example, the elastic response of the lowest symmetry crystal system (triclinic) depends upon all 6 components of the elastic Green tensor. If one were to specify the total inverse deformation $g$, and either $\eta^p$, $C^p$, or $b^p$, then all 6 components of $C^\varepsilon$ could not be determined. Although our examples will make use of isotropic equation of state models (whose elastic invariants may be determined using $g$ and $C^p$), our goal is to construct a framework of more general applicability.

To motivate our choice of evolution equations for $\mathcal{F}^p$ we begin by postulating the existence of a hyperelastic equation of state,

$$e = \dot{\mathcal{E}}(g, \mathcal{F}^p, \kappa, S),$$

where $S$ is the specific entropy. The material derivative of $\mathcal{E}$ is

$$\dot{\mathcal{E}} = \frac{\partial \mathcal{E}}{\partial g_{\alpha\beta}} \dot{g}_{\alpha\beta} + \frac{\partial \mathcal{E}}{\partial \mathcal{F}^p_{\alpha\beta}} \dot{\mathcal{F}^p}_{\alpha\beta} + \frac{\partial \mathcal{E}}{\partial \kappa} \dot{\kappa} + \frac{\partial \mathcal{E}}{\partial S} \dot{S} = -\frac{1}{\rho} \sigma_{\rho\gamma} \mathcal{F}^p_{\gamma\alpha} \dot{g}_{\alpha\beta} + \Phi,$$

where the second equality equates energy change with the sum of work and heat. Solving for entropy production (dissipation) we have

$$0 \leq \dot{S} = -\frac{1}{T} \left( \frac{\sigma_{\rho\gamma} \mathcal{F}^p_{\gamma\alpha}}{\rho} + \frac{\partial \mathcal{E}}{\partial g_{\alpha\beta}} \right) \dot{g}_{\alpha\beta} - \frac{1}{T} \frac{\partial \mathcal{E}}{\partial \mathcal{F}^p_{\alpha\beta}} \dot{\mathcal{F}^p}_{\alpha\beta} - \frac{1}{T} \frac{\partial \mathcal{E}}{\partial \kappa} \dot{\kappa} + \frac{\partial \mathcal{E}}{\partial S} \dot{S}$$

$$= \frac{1}{\rho T} (\Psi_{\text{plast}} + \Psi_{\text{therm}}).$$

Here $\partial \mathcal{E} / \partial S = T$ is the temperature, and we have introduced the specific power of thermal dissipation,

$$\Psi_{\text{therm}} = \rho \Phi,$$

and the specific power of plastic dissipation,

$$\Psi_{\text{plast}} = -\rho \frac{\partial \mathcal{E}}{\partial \mathcal{F}^p_{\alpha\beta}} \dot{\mathcal{F}^p}_{\alpha\beta} - \rho \frac{\partial \mathcal{E}}{\partial \kappa} \dot{\kappa}$$

$$= g_{\beta\gamma} \sigma_{\gamma\alpha} \mathcal{F}^p_{\delta\alpha} \dot{\mathcal{F}^p}_{\delta\beta} - \tilde{\vartheta} \kappa$$

$$= g_{\beta\gamma} \sigma_{\gamma\alpha} \mathcal{F}^p_{\delta\alpha} \mathcal{F}^p_{\delta\beta} - \tilde{\vartheta} \kappa$$

$$= \Sigma : L^p - \tilde{\vartheta} \kappa.$$
with $\vartheta = \rho \partial \mathcal{E} / \partial \kappa$ being the work hardening modulus. $L_{\nu\beta}^p = \frac{g_{\nu\beta}^p \dot{F}_{\alpha\beta}^p}{(F^p)^{-1}}$ is the plastic distortion rate [13], and $\Sigma_{\gamma\delta} = g_{\beta\gamma} \sigma_{\gamma\delta} F_{\delta\nu}$ is the thermodynamic force conjugate to $L^p$. The dependence of $\dot{S}$ on $\dot{g}_{\alpha\beta}$ vanishes because

$$\sigma_{\alpha\beta} = -\rho \frac{\partial \mathcal{E}}{\partial g_{\alpha\beta}} g_{\gamma\alpha}. \hspace{1cm} (59)$$

In evaluating (58) we have assumed that $\mathcal{E}$ depends on $g$ and $F^p$ only through the elastic deformation $F^e = F^p g = (F^p g)^{-1}$, e.g.,

$$\mathcal{E} = \dot{\mathcal{E}}(F^e, S, \kappa), \hspace{1cm} (60)$$

whence

$$\frac{\partial \mathcal{E}}{\partial F^p_{\alpha\beta}} \bigg|_{S, \kappa} = -\frac{1}{\rho} g_{\gamma\beta} \sigma_{\gamma\delta} F^e_{\delta\alpha}. \hspace{1cm} (61)$$

Thermodynamics requires that the internal energy depends upon the volume, and we assume by (60) that this energy dependence is carried by the tensor $F^e$, e.g., $V = V_0 \det F^e$. For this to be true, it is necessary that $\det F^p = 1$ at all times (i.e., $V = V_0 \det F = V_0 \det F^e \det F^p$; $V = V_0 \det F^e$ iff $\det F^p = 1$.) Therefore, (60) assumes that plastic flow is volume-preserving.

We postulate a plastic yield surface $f = 0$, which we represent for illustrative purposes with a Mises-Huber constitutive model written in terms of the Cauchy stress $\sigma$, a constant yield stress parameter $\sigma_Y$, and the work hardening modulus $\vartheta$:

$$f(\sigma, \vartheta) = \|\text{dev}\sigma\| - \sqrt{\frac{2}{3}} (\sigma_Y + \vartheta). \hspace{1cm} (62)$$

Here, $\text{dev\sigma} = \sigma - \frac{1}{3} (tr\sigma) I$ is the stress deviator, and $\|A\|$ is the Schur norm of $A$, $\|A\|^2 = A_{\alpha\beta} A_{\alpha\beta} = \text{tr}(A^T A)$.

The flow model we adopt is derived from (62) by the postulate of maximum plastic dissipation [10, 12]. The plastic dissipation (58) is considered as a function of the variables $\Sigma$ and $\vartheta$, with fixed parameters $L^p$ and $\kappa$; $\Psi_{\text{plast}} = \Psi_{\text{plast}}(\Sigma, \vartheta; L^p, \kappa)$. The plastic dissipation is then maximized with respect to $\Sigma$ and $\vartheta$, subject to the constraint that $f = 0$ during plastic flow. The resulting flow laws are:

$$\dot{\mathcal{F}}^p = \zeta F^p g \frac{\text{dev}(\sigma)}{\|\text{dev}(\sigma)\|} \mathcal{F} \hspace{1cm} (63)$$

$$\dot{\kappa} = \zeta \sqrt{\frac{2}{3}} \hspace{1cm} (64)$$

with $\zeta$ a parameter chosen to satisfy the Kuhn-Tucker complementarity conditions and the "consistency condition" [22]

$$f = 0 \hspace{1cm} (65)$$

$$\zeta \geq 0 \hspace{1cm} (66)$$

$$\zeta f = 0 \hspace{1cm} (67)$$

$$\zeta f = 0 \hspace{1cm} \text{(if } f = 0) \hspace{1cm} (68)$$
The flow model (63) is consistent with the assumption that plastic flow is volume-preserving,

\[(\det \mathcal{F}^p) = (\det \mathcal{F}^p) g^p_{\alpha\beta} \frac{\mathcal{F}^p_{\gamma\nu}}{||\text{dev}\sigma||} = 0 \quad \text{because } \text{tr}(\text{dev}\sigma) = 0, \]

and is therefore compatible with the assumption made in evaluating $\Psi_{\text{plast}} (58)$.

As an example, we use a modified Mooney-Rivlin equation of state:

\[
\rho_0 \varepsilon(C^e, S) = \frac{\lambda(S)}{2} (\ln \sqrt{\det C^e})^2 + \frac{\mu(S)}{2} \text{tr} C^e - \frac{\mu(S)}{2} \log \det C^e + \frac{\rho_0 \vartheta_0}{\rho} \left( \kappa + \frac{1}{\vartheta_1} e^{-\vartheta_1 \kappa} \right)
\]

This equation of state gives a work hardening modulus,

\[
\vartheta(\kappa) = \rho \frac{\partial \varepsilon}{\partial \kappa} = \vartheta_0 \left( 1 - e^{-\vartheta_1 \kappa} \right)
\]

in terms of two parameters: $\vartheta_0$ is the ultimate, asymptotic value of the work hardening modulus, and $\vartheta_1$ dictates the rate of approach of the asymptotic limit.

The combined elastic-plastic evolution problem is solved with a predictor-corrector strategy. The inverse total deformation $g$ is advanced in accordance with the equations of motion, with the plastic deformation $\mathcal{F}^p$ being conservatively advected. This step may predict a coordinate in state space that lies outside the convex manifold of permissible states $f(\sigma, \vartheta) \leq 0$, in which case a plastic corrector step is used to bring state back to the yield surface. The algorithmic approach is a return mapping algorithm [22], modified to require only one equation of state evaluation.

Begin the iteration sequence with iteration index $m = 0$,

\[
\mathcal{F}^p(0) = \mathcal{F}^{p,n+1} \\
\kappa(0) = \kappa^{n+1} \\
\sigma(0) = \sigma(g^{n+1}, \mathcal{F}^{p,n+1}, \kappa^{n+1}) \\
\vartheta(0) = \vartheta(\kappa^{n+1})
\]

There is one equation of state evaluation at the beginning of the iteration in which the Cauchy stress $\sigma$, work hardening modulus $\vartheta$, and the derivatives $\partial \sigma / \partial \mathcal{F}^p |_{\varepsilon, g, \kappa}$, $\partial \sigma / \partial \kappa |_{\varepsilon, g, \mathcal{F}^p}$, and $\partial \vartheta / \partial \kappa$, are calculated. Next, evaluate the yield criterion

\[
f(m) = f(\sigma(m), \vartheta(m)).
\]
If $m = 0$ and $f^{(m)} \leq \epsilon$, then the state point is interior to the yield surface, and no plastic flow occurs. If $f^{(0)} > \epsilon$ and $|f^{(m)}| \leq \epsilon$, then

$$\mathcal{F}^{p,n+1} \leftarrow \mathcal{F}^{p}(m)$$
$$\kappa^{n+1} \leftarrow \kappa^{(m)}$$

and stop. Otherwise, calculate $\Delta \zeta^{(m)} = \zeta^{(m+1)} - \zeta^{(m)}$ using Newton’s method

$$\Delta \zeta^{(m)} = - f^{(m)} \left( \frac{df^{(m)}}{d\zeta} \right)^{-1}$$

with $df/d\zeta$ estimated from

$$\frac{df^{(m)}}{d\zeta} \approx \left( \frac{\partial f}{\partial \sigma} \right)^{(m)} \left[ \left( \frac{\partial \sigma}{\partial \mathcal{F}^{p}} \right)_{\varepsilon,\kappa} \left( \frac{d\mathcal{F}^{p}}{d\zeta} \right)^{(m)} + \left( \frac{\partial \sigma}{\partial \kappa} \right)_{\mathcal{F}^{p},\varepsilon,\kappa} \left( \frac{d\kappa}{d\zeta} \right)^{(m)} \right] +$$

$$\left( \frac{\partial f}{\partial \kappa} \right)^{(m)} \left( \frac{\partial \kappa}{\partial \mathcal{F}^{p}} \right)_{\varepsilon,\kappa} \left( \frac{d\mathcal{F}^{p}}{d\zeta} \right)^{(m)}. \quad (77)$$

Next, calculate revised estimates

$$\mathcal{F}^{p} = \mathcal{F}^{p}(m) + \left( \frac{\partial \mathcal{F}^{p}}{\partial \zeta} \right)^{(m)} \Delta \zeta^{(m)}$$
$$\mathcal{F}^{p}(m+1) = (\det. \mathcal{F}^{p})^{-1/3} \mathcal{F}^{p}$$
$$\kappa^{(m+1)} = \kappa^{(m)} + \left( \frac{\partial \kappa}{\partial \zeta} \right)^{(m)} \Delta \zeta^{(m)}$$
$$\sigma^{(m+1)} = \sigma^{(0)} + \left( \frac{\partial \sigma}{\partial \mathcal{F}^{p}} \right)_{\varepsilon,\kappa} \left( \mathcal{F}^{p}(m+1) - \mathcal{F}^{p}(0) \right)$$

$$+ \left( \frac{\partial \sigma}{\partial \kappa} \right)_{\varepsilon,\kappa,\mathcal{F}^{p}} \left( \kappa^{(m+1)} - \kappa^{(0)} \right)$$
$$\vartheta^{(m+1)} = \vartheta(\kappa^{(m+1)}),$$

set $m \leftarrow m + 1$, and retest the stopping criterion.

In this procedure we evaluate the equation of state once to determine $\sigma$ and the thermodynamic derivatives $\partial \sigma / \partial \mathcal{F}^{p}$ and $\partial \sigma / \partial \kappa$. The stress $\sigma^{(m)}$, for $m > 0$, is approximated by first-order Taylor expansion about the initial $m = 0$ value. The method converges in 1 or 2 iterations, with $\epsilon = 10^{-6}$, in each of the test problems involving plasticity described below.

The framework described by Eq. (9) calls for rates of plastic deformation $h$ and rates of work hardening $K$. In the example above, which is rate-independent, we use

$$\tau h = \dot{\mathcal{F}}^{p}$$
$$\tau K = \dot{\kappa}$$

where $\tau = \Delta t/2$ in the predictor step of the method (Eqs. (30a,30b), and $\tau = \Delta t$ in the corrector (Eq. (41)). A generalization of this approach to rate-dependent plasticity is described in [18].
6. DISSIPATION

In certain problems in hydrodynamics it has been found that the higher-order Godunov strategy we adapted here will give rise to spurious post-shock oscillations (e.g., [5]). A solution that rectifies this problem is the addition of a small amount of additional dissipation at strong shocks. This dissipation is added by introducing an additional slope limiter via a "flattening" parameter $\chi$ (see Eq. (27)).

A variety of flattening strategies have been proposed. Perhaps the simplest variant, employed by Miller and Puckett [14], uses the divergence of the velocity to detect potential shocks, and uses a simple measure of shock strength, the ratio of pressure jump across a cell to the isentropic bulk modulus, $|\Delta P|/K_S$ where $K_S = \partial P/\partial \log \rho|_S$, to compute a flattening measure. This introduces additional dissipation in regions where the pressure change is large compared to the bulk modulus—where linearization of the equation of state is expected to become error-prone. This strategy may introduce extra dissipation in regions that do not require it, however, as when a shock is spread over a large (> 5 or 6) number of grid cells. It is therefore desirable to also include measures of the shock structure in order to minimize application of this dissipation mechanism.

Elaborate strategies for computing $\chi$ are described by Colella and Woodward [7]. One of their strategies is to restrict the use of this dissipative mechanism to regions where the detected shock is steep. In our solid mechanics computations we found this strategy to be useful, and in conjunction with a measure of shock strength provides judicious, adequate additional dissipation.

We detect a strong shock by measuring in 1D the divergence of the velocity field, and calculating a normalized jump in stress. We define

$$z_i = \frac{\|[(\sigma e_\alpha)_{i+1} - (\sigma e_\alpha)_{i-1}]\|_\infty}{(\det A_{\alpha\alpha,i})^{1/3}}$$

(80)

as a measure of shock strength in the neighborhood of cell $i$ in direction $e_\alpha$. The numerator is the maximum of the absolute value of the jump in those stress components that may change in direction $e_\alpha$ 1D purely elastic flow, and the denominator is a mean modulus of the acoustic propagation tensor in direction $e_\alpha$.

Following Colella and Woodward, we discriminate between steep and broad shocks by the ratio

$$\beta_i = \frac{\|[(\sigma e_\alpha)_{i+1} - (\sigma e_\alpha)_{i-1}]\|_\infty}{\|[(\sigma e_\alpha)_{i+2} - (\sigma e_\alpha)_{i-2}]\|_\infty}$$

(81)

In the limit $\beta_i = \frac{1}{2}$, stress is approximately linear across 5 grid cells, and so a shock discontinuity is not being captured. When $\beta_i \approx 1$ the discontinuity is captured in 3 cells: the shock may be overly steep, and post-shock oscillations are expected. Accordingly, the minimum value that our flattening parameter $\chi$ should have, based upon shock steepness, is

$$\chi_{\min\ i} = \max \left(0, \min \left(1, \frac{a_1 - \beta_i}{a_1 - a_0} \right) \right)$$

(82)

where $a_0$ and $a_1$ are numerical constants. We use the values $a_0 = 0.75$ and $a_1 = 0.85$ in the computations presented here.
A local shock-strength-sensitive flattening parameter $\bar{\chi}_i$, $\chi_{\text{min}} \leq \bar{\chi}_i \leq 1$, is thus

$$
\bar{\chi}_i = \begin{cases} 
\min \left( 1, \max \left( \frac{z_i - z_{i-1}}{z_i}, \chi_{\text{min}} \right) \right) & (v \cdot e_i)_{i+1} < (v \cdot e_i)_{i-1} \\
\chi_{\text{min}} & \text{otherwise}
\end{cases}
$$

(83)

In our example calculations we use the numerical values $z_0 = 0.25$ and $z_1 = 0.75$.

In 1D we limit the slopes by the minimum over nearest neighbor cells of the local flattening parameter,

$$
\chi_i = \min(\bar{\chi}_i, \bar{\chi}_{i-1}, \bar{\chi}_{i+1}).
$$

(84)

In higher dimensions, we employ the same 1D local flattening parameters — measured separately in each direction. All slopes ($\partial q/\partial x$, $\partial q/\partial y$, and $\partial q/\partial z$) are limited by the same cell-valued flattening parameter, which is given by the minimum of the directional local measures. In 2D,

$$
\chi_{ij} = \min(\bar{\chi}_{x,i-1,j}, \bar{\chi}_{x,i,j}, \bar{\chi}_{x,i+1,j}, \bar{\chi}_{y,i,j-1}, \bar{\chi}_{y,i,j}, \bar{\chi}_{y,i,j+1}),
$$

(85)

and in 3D,

$$
\chi_{ijk} = \min(\bar{\chi}_{x,i-1,j,k}, \bar{\chi}_{x,i,j,k}, \bar{\chi}_{x,i+1,j,k}, \bar{\chi}_{y,i,j-1,k}, \bar{\chi}_{y,i,j,k}, \bar{\chi}_{y,i,j+1,k}, \bar{\chi}_{z,i,j,k}, \bar{\chi}_{z,i,j,k+1}).
$$

(86)

7. ACCURACY

The term $(v \times \nabla \times g^T)^T$ was introduced to the evolution equations of the inverse deformation gradient $g$ to make the system of equations stable and well-posed when the gauge constraint $\nabla \cdot g = 0$ fails to be satisfied. Although the partial differential equations show that when satisfied initially, it will be satisfied for all times, numerical errors cause the constraint to be violated to some degree.

We propose a modification of (9) to control inaccuracy that may arise from violation of the gauge constraint. The conservation law (11) indicates that $\mathcal{G}$ will be created by numerical errors as dipoles. Thus, a numerical strategy that will control this truncation error is to diffuse $\mathcal{G}$,

$$
\frac{\partial \mathcal{G}}{\partial t} + \nabla \cdot (v \mathcal{G} - \mathcal{G}v) = \mathcal{D}(\nabla^2 \mathcal{G})
$$

(87)

or, equivalently,

$$
\frac{\partial g e_\alpha}{\partial t} + \frac{\partial}{\partial x_\alpha} (g v) = (v \times g)^T e_\alpha - \mathcal{D}(\nabla \times g)^T e_\alpha.
$$

(88)

g is also related to the density via

$$
\rho = \rho_0 \det g.
$$

(89)

where $\rho_0$ is the mass density in the reference state $\mathcal{F} = g = I$. Multiplying the $g$ equations by $\rho_0 \det(g) g^{-T}$, and summing over the 9 components of $g$, gives a conservation law for
\( \hat{p} \equiv \rho_0 \det(g) \):

\[
\frac{\partial \hat{p}}{\partial t} + \nabla \cdot (\hat{p} \nu) = 0,
\]  

(90)

thus the continuity equation is embodied in the \( g \) equations as well. However, because of discretization errors the equivalence of \( \hat{p} \) and the mass density \( \rho \) cannot be assured. To make the method strictly conservative, we keep \( \rho \) as a redundant variable, and we invoke a relaxation mechanism on \( g \) to enforce the condition \( \hat{p} = \rho \). This relaxation alone (not including the diffusion modification) is accomplished by writing

\[
\frac{\partial g e_\alpha}{\partial t} + \frac{\partial}{\partial x_\alpha} (g \nu) = (v \times \mathcal{G})^T e_\alpha + \eta \left( \frac{\rho}{\rho} - 1 \right) g e_\alpha.
\]  

(91)

The "continuity" equation for \( \hat{p} \) is then

\[
\frac{D\hat{p}}{Dt} = \hat{p} \mathcal{T} \alpha \frac{Dg_\alpha}{Dt} = \hat{p} \mathcal{T} \alpha \eta (v \times \mathcal{G})_\beta \alpha + 3\eta (\rho - \hat{p})
\]  

(92)

\[
\frac{\partial \hat{p}}{\partial t} + \nabla \cdot (\hat{p} \nu) = 3\eta (\rho - \hat{p}) \quad \text{when} \; \mathcal{G} = 0.
\]  

(93)

Including the diffusion and relaxation terms, the system of equations we will solve is:

\[
\begin{pmatrix}
\rho \\
\rho v \\
\rho E \\
ge_x \\
ge_y \\
ge_z \\
\rho \mathcal{T} \nu e_x \\
\rho \mathcal{T} \nu e_y \\
\rho \mathcal{T} \nu e_z \\
\rho K
\end{pmatrix}
+ \nabla \frac{\partial}{\partial x_\alpha}
\begin{pmatrix}
\rho v_\alpha \\
\rho v_\nu - \sigma e_\alpha \\
\rho E v_\alpha - v_\beta \sigma_\beta \alpha \\
ge_\nu v_\alpha \\
ge_\nu - \nu_\nu \nu_\nu \\
\rho \mathcal{T} \nu e_\nu v_\alpha \\
\rho \mathcal{T} \nu e_\nu v_\alpha \\
\rho \mathcal{T} \nu e_\nu v_\alpha \\
\rho K v_\alpha
\end{pmatrix}
= \begin{pmatrix}
0 \\
\rho f \\
\rho (\mathcal{F} + v \cdot f) \\
(v \times \mathcal{G})^T e_1 \\
(v \times \mathcal{G})^T e_2 \\
(v \times \mathcal{G})^T e_3 \\
\rho e_x \\
\rho e_y \\
\rho e_z
\end{pmatrix}
\]  

(94)

\[
-D(\nabla \times \mathcal{G})^T e_x + \eta \left( \frac{\rho}{\rho_0 \det g} - 1 \right) g e_x
\]

\[
-D(\nabla \times \mathcal{G})^T e_y + \eta \left( \frac{\rho}{\rho_0 \det g} - 1 \right) g e_y
\]

\[
-D(\nabla \times \mathcal{G})^T e_z + \eta \left( \frac{\rho}{\rho_0 \det g} - 1 \right) g e_z
\]
Our discretization of the diffusion and relaxation terms takes the form

\[
\begin{pmatrix}
g_{e_x}^{n+1} \\
g_{e_y}^{n+1} \\
g_{e_z}^{n+1}
\end{pmatrix}_{ijk} = \begin{pmatrix}
g_{e_x}^{n} \\
g_{e_y}^{n} \\
g_{e_z}^{n}
\end{pmatrix}_{ijk} + \Delta t \begin{pmatrix}
-D(\nabla \times G)^T n \cdot e_x + \eta \left( \frac{\tilde{g}_{n+1}^{+1}}{\rho_0 \det \tilde{g}^{+1}} - 1 \right) \tilde{g}_{n+1}^{+1} e_x \\
-D(\nabla \times G)^T n \cdot e_y + \eta \left( \frac{\tilde{g}_{n+1}^{+1}}{\rho_0 \det \tilde{g}^{+1}} - 1 \right) \tilde{g}_{n+1}^{+1} e_y \\
-D(\nabla \times G)^T n \cdot e_z + \eta \left( \frac{\tilde{g}_{n+1}^{+1}}{\rho_0 \det \tilde{g}^{+1}} - 1 \right) \tilde{g}_{n+1}^{+1} e_z
\end{pmatrix}_{ijk}
\] (95)

where \( \tilde{g}_{n+1}^{+1} \) denotes \( g \) after flux differencing and evaluation of source terms in (10) (cf. (40)).

The second derivatives of \( g \) appearing in the \( G \) diffusion term,

\[
(\nabla \times G)^T = \begin{pmatrix}
\frac{\partial^2 g_{e_x}}{\partial x^2} & \frac{\partial^2 g_{e_x}}{\partial y \partial z} & \frac{\partial^2 g_{e_x}}{\partial z^2} \\
\frac{\partial^2 g_{e_y}}{\partial x \partial z} & \frac{\partial^2 g_{e_y}}{\partial y^2} & \frac{\partial^2 g_{e_y}}{\partial z^2} \\
\frac{\partial^2 g_{e_z}}{\partial x \partial z} & \frac{\partial^2 g_{e_z}}{\partial y \partial z} & \frac{\partial^2 g_{e_z}}{\partial z^2}
\end{pmatrix}
\] (96)

are computed using time-\( n \) cell-centered values of \( g \), with a standard 3 point stencil for homogeneous second derivatives, e.g.,

\[
\left( \frac{\partial^2 g}{\partial x^2} \right)_{ijk} = \frac{1}{\Delta x_i} \left( \frac{2(g_{i+1,j,k}^{n} - g_{ij,k}^{n})}{\Delta x_{i+1} + \Delta x_i} - \frac{2(g_{ij+1,k}^{n} - g_{ij-1,k}^{n})}{\Delta x_i + \Delta x_{i-1}} \right),
\] (97)

and heterogeneous derivatives are computed with a 4 point stencil, e.g.,

\[
\left( \frac{\partial^2 g}{\partial x \partial y} \right)_{ijk} = \frac{4}{(\Delta x_i - 1 + 2\Delta x_i + \Delta x_{i+1})(\Delta y_{j-1} + 2\Delta y_j + \Delta y_{j+1})} \left( g_{i+1,j+1,k}^{n} - g_{i-1,j-1,k}^{n} - g_{i+1,j-1,k}^{n} + g_{i-1,j+1,k}^{n} \right),
\] (98)

A von Neumann stability analysis of the diffusion update in (94), considered independently of other source terms or the basic solid mechanics equations, gives a bound on the diffusion coefficient:

\[
D \leq \begin{cases} 
\frac{\Delta^2}{2\Delta x^2} & \text{in 1D,} \\
\frac{\Delta^2}{4\Delta x^2} & \text{in 2D or 3D.}
\end{cases}
\] (99)

This suggest an approximate overall Courant-Friedrichs-Lewy stability criterion of

\[
\text{CFL} = \begin{cases} 
\frac{\Delta t}{\Delta x} + \frac{2D\Delta t}{\Delta x^2} & \text{in 1D,} \\
\frac{\Delta t}{\Delta x} + \frac{4D\Delta t}{\Delta x^2} & \text{in 2D or 3D,}
\end{cases}
\] (100)

\[
\text{CFL} < 1,
\] (101)

where here it is assumed that \( \Delta x = \Delta y = \Delta z \), a constant. The more rigorous CFL condition \( \text{CFL} = \max \left( \frac{\Delta t (|v| + c_{\text{max}})}{\Delta x}, \frac{4D\Delta t}{\Delta x^2} \right) \) (in 2D or 3D) would hold if the mechanics equations and \( G \) diffusion steps were performed sequentially.
The optimal damping conditions for the diffusion of $\mathcal{G}$ is obtained by choosing the empirical diffusion constant $D$ to satisfy

$$D = \frac{(\Delta x)^2}{4d\Delta t}$$  \hspace{1cm} (102)$$

where $d = 1, 2, 3$ is the dimensionality of the problem. Similarly, optimal relaxation is obtained by choosing the empirical relaxation parameter $\eta$ to satisfy

$$\eta = \frac{1}{6\Delta t}$$  \hspace{1cm} (103)$$

for all dimensions. According to the approximate CFL condition (101), the optimal value of $D$ will contribute $1/2$ to the CFL value in 1D and 2D, and $1/3$ in 3D; limiting the overall step size $\Delta t$ by factors of $1/2$ and $2/3$ (respectively) relative to the $D = 0$ value. Thus, in some of the examples presented below use smaller values of $D$ than indicated by (102). In some cases, however, we find that values of CFL $> 1$ provide stable solutions (consistent with (101) being only an approximation).

Assumptions underlying our plastic yield model require that $\det \mathcal{F}^p$ be constant. The differential equations describing our plastic flow model $\dot{\mathcal{F}}^p$ preserves $\det \mathcal{F}^p$, but again numerical errors will lead to some violation of this constraint. To remedy this problem we renormalize the plastic deformation tensor at the end of each time step,

$$\mathcal{F}^p \leftarrow (\det \mathcal{F}^p)^{-1/3} \mathcal{F}^p.$$

### 8. EXAMPLES

#### 8.1. Convergence: elasticity

To demonstrate the convergence properties of the algorithm we model in 1D the smooth flow resulting from an initial Gaussian-shaped disturbance. For these computations we use a hyperthermoelastic model of the Mooney-Rivlin variety,

$$\rho_0 \mathcal{E}(C^e, S) = \frac{\lambda(S)}{2} \left( \log \sqrt{\det C^e} \right)^2 + \frac{\mu(S)}{2} \text{tr} C^e - \frac{\mu(S)}{2} \log \det C^e + \frac{\rho_0 \theta_0}{\rho} \left( \kappa + \frac{1}{\theta_1} e^{-\phi_1 \kappa} \right)$$  \hspace{1cm} (105)$$

where $S$ is the entropy. Entropy dependence is introduced by supposing

$$\lambda(S) = \lambda_0 + \lambda_S f(S)$$  \hspace{1cm} (106a)$$
$$\mu(S) = \mu_0 + \mu_S f(S)$$  \hspace{1cm} (106b)$$

where $f(S)$ is an unspecified function of the entropy. From this equation we evaluate $\sigma(\mathcal{E}, g, \mathcal{F}^p, \kappa)$, and other derivatives including the acoustic propagation tensors, by first solving this equation of state for $f(S)$, then differentiating $\mathcal{E}$ with respect to the elements of $C^e$ while holding $f(S)$ constant. We use values $\rho_0 = 1$, $\mu_0 = \lambda_0 = 0.6$, and $\mu_S = \lambda_S = 0.01$, with initial values $g_0 = 1.1I$, $\mathcal{F}_0^p = I$, $\kappa = 0$, and $v = 0$. The initial disturbance is generated by distributing internal energy from $\mathcal{E}'$ to $\mathcal{E}''$,

$$\mathcal{E}' = \mathcal{E}(g_0, \mathcal{F}_0^p, f(S) = 0)$$  \hspace{1cm} (107a)$$
$$\mathcal{E}'' = 10\mathcal{E}_0.$$  \hspace{1cm} (107b)$$
These limiting values are used to construct a Gaussian initial profile, via

\[ E_i = E' \omega_i + (1 - \omega_i)E'' \]  

with

\[ \omega_i = \frac{1}{a\sqrt{2\pi}} \exp \left( -\frac{r_i^2}{2a^2} \right) \]

where \( a^2 = 100 \) is the variance of the distribution, and where \( r_i \) is the coordinate of the center of cell \( i \), in the domain \([0,40]\). Boundary conditions are reflecting at \( r = 0 \) and \( r = 40 \). We pick the time step \( \Delta t \) to satisfy the Courant-Friedrichs-Lewy constraint (101), with CFL = 0.8.

This problem was chosen to give a nontrivial shockless flow, with initial conditions that strictly obey \( G = 0 \). Plasticity is not incorporated into this test problem, and no flattening is required.

Figure 1 shows the initial and final conditions of this test problem in Cartesian geometry. At this scale, the difference between results at 40, 80, and 160 Cartesian points is not resolvable.

A comparison of results using 40, 80, and 160 grid points is used to estimate the \( L_1 \), \( L_2 \), and \( L_\infty \) (max) norm rates of convergence using the volume-weighted variables (Table 1). In Cartesian geometry the method exhibits approximately third-order convergence: as the number of grid cells is doubled, the error diminishes by a factor of 2.3. Slightly lower rates of convergence are seen in cylindrical and spherical geometries, but in all cases the order exceeds 2.

8.2. Convergence: plasticity

To assess the rate of convergence in a plasticity-dominated flow we pose a model problem similar to the purely elastic problem presented above. A Gaussian distribution with width 5 is used to vary the \( g_{yy} \) and \( g_{zz} \) as functions of coordinate \( x \) according to

\[ g_{xx,i} = 1.1 \]
\[ g_{yy,i} = (1 + 9\omega_i)1.1 \]
\[ g_{zz,i} = 1.1/(1 + 9\omega_i) \]

with homogeneous initial density, internal energy, and zero velocity. We use the equation of state (105) with yield model (62) and flow rates (63,64). The equation of state parameters are as used in the purely elastic convergence test, and the plastic constitutive parameters are \( \sigma_Y = 0.1 \), \( \phi_0 = 0.1 \), and \( \phi_1 = 10.0 \).

The flow field in this problem is \( C^\omega \), which lowers the overall order of convergence. Density converges at greater than 2nd order (Table 2), but the tangential stress components converge only at 1st order.

8.3. Blake’s problem

Blake [3] presented an analytical solution to the problem of an unbounded solid medium characterized by an isotropic linear elastic equation of state,

\[ \rho_0 \mathcal{E} = \frac{1}{8} \lambda [\text{tr}(C_e - I)]^2 + \frac{1}{4} \mu \text{tr} \left( C_e^TC_e - 2C^e - I \right), \]
loaded by a prescribed pressure boundary condition on the interior of a spherical cavity of initial radius \(a\). We present a numerical solution to this problem in 1D spherical coordinates (see Appendix), with slight modification of the code to accommodate the moving boundary with prescribed flux (Neumann) boundary conditions. This problem is selected to verify the behavior of the elastic algorithm in the weak shock limit.

The cavity wall represents a material interface across which the mass flux will be zero. Accordingly, the flux at this boundary is given by \(F(U^B) - sU^B\) where \(U^B\) is the vector of conserved quantities at the boundary, \(s\) is the velocity of the boundary, and \(F(U^B)\) is the radial flux vector evaluated at the boundary. Blake’s solution provides \(u(r, t)\), the displacement of a mass element in the radial direction. In spherical coordinates, this gives rise to an inverse deformation tensor

\[
g(r, t) = \begin{pmatrix}
(1 + \partial u / \partial r)^{-1} & 0 & 0 \\
0 & (1 + u/r)^{-1} & 0 \\
0 & 0 & (1 + u/r)^{-1)
\end{pmatrix}.
\]

(112)

The velocity of the material interface is \(s = \partial u / \partial t\) at \(r = a\).

Cell 1, whose left boundary is \(r = a\) at \(t = 0\), and whose right boundary is fixed at \(a + \Delta r\), has a volume which varies with time. Applying Gauss’s divergence theorem to this cell gives

\[
V_{1}^{n+1}U_{1}^{n+1} = V_{1}^{n}U_{1}^{n} + \Delta t \left( A_{1/2}F_{1/2}^{n+1/2} - A_{3/2}F_{3/2}^{n+1/2} \\
+ \bar{A}_{1}(H_{1/2}^{n+1/2} - H_{3/2}^{n+1/2}) + \Delta t\bar{V}_{1}G_{1}^{n+1/2},
\]

(113)

where \(F\) denotes the radial flux component that enters as \((1/r^2)\partial(r^2F)/\partial r\), \(H\) denotes the radial flux component that enters as \(\partial H/\partial r\) (see Appendix), \(\bar{A}_{1}\) is the average area \((r^2)\) over \(r\) in \([a - u(a, t), a + \Delta r]\), \(\bar{V}_{1}\) is the time-averaged cell volume, and \(G_{1}^{n+1/2}\) is the cell-centered vector of (geometric) source terms, which we time-center with a predictor-corrector strategy.

In general (see Wilkins’ problem below), algebraic solution of this discretization is unstable. In the particular case of our discretization of Blake’s problem, however, \([u(a, t)] \ll \Delta r\) and so \(V_{1}\) does not vary appreciably with time and in particular is of order \(a^2 \Delta r\). Our solution of Blake’s problem therefore uses (113) as written. It is also necessary to modify the algorithm to account for the absence of cell values at \(i - 1\) and \(i - 2\). The gradient \(\partial q/\partial r\) at \(i = 1\) is obtained by 1st-order forward finite difference with a van Leer limiter. The flattening parameter \(\chi\) operates on a stencil that requires cell values at 0 and \(-1\). However, for this weak problem additional flattening is never required, so the algorithm is modified by omission of the flattening computation (\(\chi = 1\)).

Following Trangenstein and Colella [24] we use parameters \((a = 0.1\ m, \rho_0 = 3000\ kg/m^3, \lambda = 2.36 \times 10^{10}\ Pa, \mu = 2.78 \times 10^{10}\ Pa)\). The pressure inside the spherical cavity is \(10^6\ Pa\), and the solution is plotted at time \(1.6 \times 10^4\ s\).

We compare in Figs (3-6) our computed results for radial stress,

\[
\sigma_{rr} = (\lambda + 2\mu)(\partial u / \partial r) + 2\lambda(u/r),
\]

(114)

and hoop stress

\[
\sigma_{\theta\theta} = \sigma_{\phi\phi} = \lambda(\partial u / \partial r) + 2(\lambda + \mu)(u/r),
\]

(115)
pressure
\[ P = -\frac{1}{3} \sigma_{ii} = -\left( \lambda + \frac{2}{3} \mu \right) \left[ (\partial u/\partial r) + 2(u/r) \right], \] (116)

and radial velocity \( v_r \) against Blake's analytical results.

These results verify the method in the case of weak (linear) waves. The leading shock is captured in approximately 5 grid cells. A single stress undershoot precedes the shock, and a corresponding overshoot follows it, but the wave speed and amplitude are correctly modeled.

8.4. Wilkins' Problem

Wilkins' flying plate problem [26] involves a 5mm-thick aluminum plate impacting an initially-stationary aluminum halfspace. The rear (left) surface of the flying plate is a free surface (vacuum). Initially, left- and right-traveling shocks propagate outward from the point of contact of the plate with the halfspace. When the left-traveling shock reaches the free surface, a right-traveling rarefaction is created, which ultimately overtakes the right-traveling shock. This problem incorporates plasticity.

To model this problem, we modify our 1D algorithm to allow for the moving free-surface boundary. This is an example of volume-of-fluid front reconstruction applied to multi-fluid modeling, and details will be described in a future correspondence. Briefly, we modify the approach adopted for Blake's problem using the flux redistribution ideas of Chern and Colella [4]. Application of this approach to stationary incompressible boundaries is described in [15], and to reaction front tracking in [1, 17]. Our implementation is similar, but the free-surface boundary moves at a velocity determined by the solid-vacuum Riemann problem. This problem is solved as described above for the solid-solid case, but uses only the \( 3 \times 3 \) stress component of the eigenvectors. This interface velocity, and the surrounding material velocities, are used with a volume-pushing algorithm (after [2]) to update the fractional occupancy of the interface cells.

We construct a hyperelastic model of aluminum in close correspondence to Wilkins' (rate model) description, with

\[ \mathcal{E}(g, T^p) = \left( \int \frac{P(\rho)}{\rho^2} \, d\rho \right) + \frac{\mu_0}{2\rho_0} \left( \text{tr} \mathbf{C}^e - 3 \left( \frac{\rho_0}{\rho} \right)^{2/3} \right), \] (117)

where \( P(\rho) \) is the hydrostatic pressure (in GPa)

\[ P(\rho) = 72(\rho/\rho_0 - 1) + 172(\rho/\rho_0 - 1)^2 + 40(\rho/\rho_0 - 1)^3, \] (118)

with \( \rho_0 = 2.7 \text{ kg/m}^3 \). The shear modulus is \( \mu_0 = 24.8 \text{ GPa} \). The problem is perfectly plastic (no work hardening), and uses the von Mises yield surface function

\[ f(\sigma) = ||\text{dev}\sigma|| - \sqrt{\frac{2}{3}} \sigma_Y \] (119)

with constant flow stress \( \sigma_Y = 0.2976 \text{ GPa} \).

 Computations with impact velocities of 0.8 km/s and 2.0 km/s were obtained with CFL = 0.80 and 500 Cartesian grid points. At 0.8 km/s, a plastic shock trails a leading
elastic shock precursor. When the left-facing shocks reach the free surface, right-traveling elastic and trailing plastic rarefaction waves begin to overtake the initial right-facing shocks. The shock stress at 2.0 km/s is above the elastic limit, so only plastic shocks are formed. On rarefaction from the left free surface, a leading right-facing elastic rarefaction is formed, followed by the plastic wave. These results are in good quantitative agreement with Wilkins'.

8.5. A test in 2D

This test problem compares a 1D cylindrical coordinate computation against a 2D Cartesian result, for a problem with cylindrical symmetry. We use the modified Mooney-Rivlin model presented in Eq. (70) with initial conditions \( \rho_0 = 1, \ g = 1.1I, \ F^p = I, \) and \( \kappa = 0. \) The plasticity parameters are \( \sigma_Y = 0.1, \ \theta_0 = 0.1, \) and \( \theta_1 = 10. \) All boundary conditions are reflecting. The material is initially at rest, except for a cylindrical shell \( r \in [5, 15] \) which moves toward the axis with a velocity of \(-1\). This generates a diverging rarefaction, and a convergent shock, which reflects off the axis of symmetry.

In Figures (9-12) we compare results from a 1D cylindrical calculation (500 cells, \( \text{CFL} = 0.8 \)), and an equivalent 2D Cartesian calculation using \( 250 \times 250 \) cells, also at \( \text{CFL} = 0.8 \). The 2D results are presented as 1D scatter plots in order to demonstrate the accurate preservation of cylindrical symmetry obtained with the spatially-unsplit 2D method. The high-resolution 1D results and lower-resolution 2D results are in good agreement, although there is some discrepancy in \( \kappa \) and \( \sigma_{rr} \) near the axis.

Using this same 2D test we demonstrate the errors associated with the gauge constraints

\[
\rho - \rho_0 \det(g) = 0 \quad (120)
\]

and

\[
\mathcal{G} = \nabla \times g^T = 0. \quad (121)
\]

These conditions are enforced in the computation by way of a relaxation term to satisfy (120) and a diffusion-like term to satisfy (121). In Figure 13 we plot the left hand side of (120) comparing results from the computation presented above (in Figs. 9-12), and results from a similar computation but in which neither a relaxation nor a diffusion correction was applied. In Figure 14 we plot the \( L_2 \) norm of the tensor \( \nabla \times g^T \), comparing results from the computation with relaxation and diffusion with results from a computation using neither correction. These figures demonstrate over an order of magnitude reduction in density error is achieved by the relaxation mechanism. Approximately a factor of 2 reduction of \( ||\mathcal{G}||_2 \) is achieved by the diffusion mechanism.

8.6. A test in 3D

This test problem compares a 1D spherical coordinate computation against a 3D Cartesian result, for a problem with spherical symmetry. The equation of state is identical to the 2D test above, and the initial conditions are similar: a spherical shell \( r \in [5, 15] \) is given an initial velocity of \(-1\). This computation, with \( 100 \times 100 \times 100 \) cells at \( \text{CFL} = 0.8 \) is underresolved. Nevertheless, there is good agreement between the 3D Cartesian results and the 1D spherical calculation, and excellent preservation of spherical symmetry (Figs. 15,16).
9. CONCLUSIONS

We have presented a new method for the solution to equations of solid mechanics in 1, 2, and 3 spatial dimensions on Eulerian grids. Our method addresses the problem of gauge constraints ($\nabla \times g^T = 0$) by adopting a non-conservation approach first proposed by [8] for the equations of magnetohydrodynamics. We write the partial differential equations of solid mechanics in such a way that the constraint, if applicable initially, holds true for all time. The constraint is violated by the truncation error of the method, and reinforced with an explicit diffusion term which annihilates the dipolar field of $\nabla \times g^T$. Another constraint of the system, a correspondence between density variation and the deformation field ($\rho = \rho_0 \det g$) is also satisfied for all times by the PDEs, if satisfied in the initial conditions. Truncation errors of the method are compensated with an explicit relaxation term.

The method presented here does not incorporate artificial viscosity, but its solutions are sensitive to 6 adjustable parameters: $D$ and $\eta$ control accuracy of the gauge constraints, and $a_0$, $a_1$, $z_0$, and $z_1$ in Eqs. (82,83)) govern the introduction of dissipation near strong shocks to prevent overshoot and ringing by locally reducing the high-order Godunov method to first-order.

Our strategy for damping modes violating the curl gauge constraint,

$$g^T := g^T - \lambda \nabla \times \nabla \times g^T,$$

$$= g^T + \lambda (\nabla^2 g^T - \nabla(\nabla \cdot g^T)),$$

$$= g^T + \lambda (\nabla^2 g^T - \nabla^2 Q(g^T)),$$

($\lambda = \Delta t D$) uses a single central difference operator acting on cell-centered variables. Here, $Q(g^T) = \nabla^{-2} \nabla(\nabla \cdot g^T)$ is the projection onto the curl-free part of $g^T$. Defining $P(x) = 1 - Q(x)$ as the projection onto the divergence-free part of $x$, and noting $PQ = QP = 0$, we have

$$P(g^T) := P(g^T) + \lambda \nabla^2 P(g^T),$$

(123)

thus we are diffusing the divergence-free part of $g^T$ without modifying the curl-free part. A similar scheme may be used to modify a vector field $B$ subject to a divergence-free constraint:

$$B := B + \lambda \nabla (\nabla \cdot B),$$

$$Q(B) := Q(B) + \lambda \nabla^2 Q(B)$$

(124)

with a single matrix-valued central difference operator for the projection $\nabla(\nabla \cdot B)$. This will directly target odd-even and checkerboard short-wavelength modes of $\nabla \cdot B$ by diffusing the curl-free part of $B$. The application of this extension to magnetohydrodynamics, where $B$ is the magnetic field subject to gauge constraint div $B = 0$, is currently being investigated (R. Crockett, personal communication).

APPENDIX: CYLINDRICAL AND SPHERICAL COORDINATES

The equations of solid mechanics in cylindrical and in spherical coordinates (like those of gas dynamics) differ from the Cartesian equations by the existence of both spatial and volumetric spatial derivatives, and by the introduction of "geometric source terms". The
coordinate transformation is accomplished by rotating the Cartesian basis vectors into the curved coordinate frame via the rotations matrices

\[ R_{\text{cyl}} = \begin{pmatrix} 
\cos \theta & -\sin \theta & 0 \\
\sin \theta & \cos \theta & 0 \\
0 & 0 & 1
\end{pmatrix} \quad (A.1) \]

\[ R_{\text{sph}} = \begin{pmatrix} 
\sin \theta \cos \phi & \cos \theta \cos \phi & -\sin \phi \\
\sin \theta \sin \phi & \cos \theta \sin \phi & \cos \phi \\
\cos \theta & -\sin \theta & 0
\end{pmatrix} \quad (A.2) \]

where we adopt the standard curved coordinate notation

\[ x = r \cos \theta \quad (A.3) \]
\[ y = r \sin \theta \]
\[ z = z \]

in cylindrical coordinates, and

\[ x = r \sin \theta \cos \phi \quad (A.4) \]
\[ y = r \sin \theta \sin \phi \]
\[ z = r \cos \theta \]

in spherical coordinates. \( R \) is the matrix of inner products of unit vectors in the curved coordinate system, \( e_{\alpha}' \), and the Cartesian system \( e_{\beta}; R_{\alpha\beta} = e_{\alpha}' \cdot e_{\beta}. \) These rotation matrices transform the Cartesian tensors \( F, g, \) and \( \sigma \) transform as \( \sigma_{\text{cyl}} = R^T \sigma_{\text{Cart}} R, \) etc, and transform the velocity vector \( v \) as \( v_{\text{cyl}} = R^T v_{\text{Cart}}. \)

In cylindrical coordinates, the system of transformed equations may be written (cf., Eq. 9):

\[
\begin{align*}
\frac{\partial}{\partial t} & + \frac{1}{r} \frac{\partial}{\partial r} \left( r \rho v_r \right) + \frac{\partial}{\partial \theta} \left( \rho v_\theta \right) + \frac{\partial}{\partial z} \left( \rho v_z \right) = 0 \\
\rho v_r \frac{\partial v_r}{\partial r} & + \rho v_\theta \frac{\partial v_r}{\partial \theta} + \rho v_z \frac{\partial v_r}{\partial z} + \rho F^r e_r v_r & = \frac{\sigma_{rr}}{r} \\
\rho v_\theta \frac{\partial v_r}{\partial r} & + \rho v_\theta \frac{\partial v_\theta}{\partial \theta} + \rho v_z \frac{\partial v_\theta}{\partial z} + \rho F^\theta e_\theta v_r & = \frac{\sigma_{r\theta}}{r} \\
\rho v_z \frac{\partial v_r}{\partial r} & + \rho v_\theta \frac{\partial v_z}{\partial \theta} + \rho v_z \frac{\partial v_z}{\partial z} + \rho F^z e_z v_r & = \frac{\sigma_{rz}}{r}
\end{align*}
\]

(\( A.5 \))
This equation does not include the det $g$ relaxation term, whose representation is unaffected by the change in variables, nor does it include the $G$ diffusion correction, which will be described separately below.

There is some latitude in the partitioning of terms between the LHS and the RHS geometric source vector. This is particularly evident in the stress terms appearing in the momentum equations. The choice of representations described here was chosen in order that some cancellation between $\sigma_{rr}$ and $\sigma_{r\theta}$ occur in the $r$-momentum source term.

The linearized equations of solid mechanics (cf., (28)), used in the construction of $L$ and $R$ edge states, also has a geometric source vector. Expressed in terms of the primitive,
variables \( q \), but omitting stress components which are described later,

\[
\begin{pmatrix}
-\rho \dot{v}_r \\
\frac{\partial}{\partial r} (\sigma_{rr} / \rho) + \frac{\partial}{\partial \theta} (\sigma_{r\theta} / \rho) + \frac{\partial}{\partial z} (\sigma_{rz} / \rho) \\
-\rho \ddot{v}_r + 2\sigma_{rr} / \rho + \sigma_{rz} / \rho \\
\sigma_{rz} / \rho \\
(v_r \sigma_{\theta \theta} - v_\theta \sigma_{r \theta}) / \rho \\
0 \\
0 \\
0 \\
v_r g_{rr} + v_\theta g_{r \theta} + v_z g_{rz} \\
-\rho \ddot{v}_r + \rho v_r g_{r \theta} - v_\theta g_{rz}
\end{pmatrix}
\]

\( s = \frac{1}{r} \)  

(A.6)

The stress evolution equations, used in the predictor steps of the method, are (plastic
source terms omitted):

\[
\begin{aligned}
\frac{\partial}{\partial t} (\sigma_{rr} / \rho) + v_r \frac{\partial}{\partial r} (\sigma_{rr} / \rho) + v_\theta \frac{\partial}{\partial \theta} (\sigma_{r\theta} / \rho) + v_z \frac{\partial}{\partial z} (\sigma_{rz} / \rho) &= 0 \\
(A_{rr} / \rho) + \cdots + (\sigma_{r\theta} / \rho) + v_\theta \frac{\partial}{\partial \theta} (\sigma_{rz} / \rho) &= 0 \\
A_{zr} &+ \cdots + (\sigma_{rz} / \rho) + v_z \frac{\partial}{\partial z} (\sigma_{rz} / \rho) &= 0
\end{aligned}
\]

(A.7)

where the tensors \( A \) are defined by (14).

The \( g \) relaxation term, \(-D(\nabla \times \nabla \times g^T)^T\), transforms in cylindrical coordinates as

\[
-D \times
\]

(A.8)
The transformed system of equations in spherical coordinates may be written:

\[
\frac{\partial}{\partial t} + \frac{\partial}{\partial r} r^2 + \frac{\partial}{\partial \theta} \sin \theta + \frac{\partial}{\partial \phi} = \rho \nabla \cdot \mathbf{v} = \rho
\]

\[
\rho \mathbf{v} \\
\rho u \mathbf{v} \\
\rho E \\
\ge_r \\
\ge_\theta \\
\rho F_{\rho r} \\
\rho F_{\rho \theta} \\
\rho F_{\rho \phi}
\]

\[
\frac{\partial}{\partial t} + \frac{\partial}{\partial r} r + \frac{\partial}{\partial \theta} \sin \theta = \rho \nabla \cdot \mathbf{v} = \rho
\]

\[
\rho v_r \\
\rho v_\theta \\
\rho v_\phi \\
\rho F_{vr_\theta} \\
\rho F_{vr_\phi} \\
\rho F_{v_\theta} \\
\rho F_{v_\phi}
\]

\[
\frac{\partial}{\partial t} + \frac{\partial}{\partial r} r^2 + \frac{\partial}{\partial \theta} \sin \theta + \frac{\partial}{\partial \phi} = \rho \nabla \cdot \mathbf{v} = \rho
\]

\[
\frac{\partial}{\partial t} + \frac{\partial}{\partial r} r^2 + \frac{\partial}{\partial \theta} \sin \theta + \frac{\partial}{\partial \phi} = \rho \nabla \cdot \mathbf{v} = \rho
\]

\[
\frac{\partial}{\partial t} + \frac{\partial}{\partial r} r + \frac{\partial}{\partial \theta} \sin \theta = \rho \nabla \cdot \mathbf{v} = \rho
\]
Again, the det $g$ source term, not included above, is unaffected by the transformation of variables. The $G$ diffusion term is described separately below.
The geometric source terms in the vector $s$ (cf. (28)) corresponding to the primitive variables $q$, but omitting the direction-dependent stress terms are:

\[
\begin{pmatrix}
-\rho(2v_r + \cot \theta v_\theta) \\
(\rho + 2\sigma_{rr} - \sigma_{\theta\theta} - \sigma_{\phi\phi})/\rho \\
(-v_r v_\theta + \rho_2)/\rho \\
(-v_r v_\phi - \rho_3)/\rho \\
(v_r (\sigma_{\theta\theta} + \sigma_{\phi\phi}) + v_\phi (-\sigma_{\phi\phi} + \rho_4))/\rho \\
0 \\
0 \\
0 
\end{pmatrix}
\]

\[
s = \frac{1}{r}
\begin{pmatrix}
v_r g_{\theta\theta} + v_\phi g_{\phi\phi} \\
-v_r g_{\phi\phi} - v_\psi g_{\phi\phi} \\
0 \\
v_r g_{\phi\phi} + v_\phi g_{\phi\phi} \\
\rho (v_r (F_{\theta\theta} + F_{\phi\phi}) + v_\phi (F_{\theta\phi} + F_{\phi\theta})) \\
(v_r (F_{\theta\theta} - F_{\phi\phi}) + v_\phi (F_{\theta\phi} - F_{\phi\theta})) \\
\rho (v_r (F_{\phi\phi} + F_{\phi\phi}) + v_\phi (F_{\theta\phi} + F_{\phi\theta})) \\
\rho (v_r (F_{\phi\phi} - F_{\phi\phi}) + v_\phi (F_{\theta\phi} - F_{\phi\theta})) \\
\rho (v_r (F_{\phi\phi} + F_{\phi\phi}) + v_\phi (F_{\theta\phi} - F_{\phi\theta})) \\
0 
\end{pmatrix}
\]

(A.10)

The stress evolution equations, used in the predictor steps of the method, are (non-geometric source terms omitted):

\[
\begin{pmatrix}
\frac{\partial}{\partial t} (\sigma_{rr} + \rho_1) \\
\frac{\partial}{\partial r} (\sigma_{\theta\theta}) \\
\frac{\partial}{\partial \theta} (\sigma_{\phi\phi}) \\
\frac{\partial}{\partial \phi} (\sigma_{\phi\phi}) \\
\rho (\sigma_{rr} + \rho_2) \\
\rho (\sigma_{\theta\theta} + \rho_3) \\
\rho (\sigma_{\phi\phi} + \rho_4) \\
\rho (\sigma_{rr} + \rho_5) \\
\rho (\sigma_{\theta\theta} + \rho_6) \\
\rho (\sigma_{\phi\phi} + \rho_7) \\
0 
\end{pmatrix}
= \begin{pmatrix}
A_{rr} \\
A_{\theta\theta} \\
A_{\phi\phi} \\
A_{\phi\phi} \\
A_{rr} \\
A_{\theta\theta} \\
A_{\phi\phi} \\
A_{rr} \\
A_{\theta\theta} \\
A_{\phi\phi} \\
0 
\end{pmatrix}
\begin{pmatrix}
v_r \\
v_\theta \\
v_\phi \\
v_r \\
v_\theta \\
v_\phi \\
v_r \\
v_\theta \\
v_\phi \\
v_\theta \\
v_\phi 
\end{pmatrix}
\]
The $g$ relaxation term $-D (\nabla \times \nabla \times g T)^T$ transforms in spherical coordinates as:

\[
((\nabla \times \nabla \times g T)^T)_{rr} = \frac{\partial^2 g_{rr}}{r^2} - \frac{\partial^2 g_{\theta\theta}}{r^2} + \frac{\partial^2 g_{\phi\phi}}{r^2} + \frac{\partial^2 g_{\theta\phi}}{r^2} + \frac{\partial^2 g_{\phi r}}{r^2} + \frac{\partial^2 g_{\phi \theta}}{r^2} \]

\[
((\nabla \times \nabla \times g T)^T)_{r\theta} = \frac{\partial^2 g_{rr}}{r^2} + \frac{\partial^2 g_{\theta\theta}}{r^2} + \frac{\partial^2 g_{\phi\phi}}{r^2} - \frac{\partial^2 g_{\theta\phi}}{r^2} + \frac{\partial^2 g_{\phi r}}{r^2} + \frac{\partial^2 g_{\phi \theta}}{r^2} \]

\[
((\nabla \times \nabla \times g T)^T)_{r\phi} = \frac{\partial^2 g_{rr}}{r^2} + \frac{\partial^2 g_{\theta\theta}}{r^2} + \frac{\partial^2 g_{\phi\phi}}{r^2} - \frac{\partial^2 g_{\theta\phi}}{r^2} + \frac{\partial^2 g_{\phi r}}{r^2} + \frac{\partial^2 g_{\phi \theta}}{r^2} \]
\((\nabla \times \nabla \times g^T)^{T}_{\theta \theta} = \frac{\partial^2 g_{\theta \theta}}{\partial r^2} + \frac{\partial^2 g_{\theta \phi}}{\partial \phi^2} + \frac{\partial^2 g_{\phi \phi}}{\partial \phi^2} - \frac{\partial^2 g_{\theta \phi}}{\partial r \partial \phi} + \frac{\partial^2 g_{\phi \phi}}{\partial \phi \partial r} - \frac{\partial g_{\phi \phi}}{r^2 \sin \theta} + \frac{\cot \theta g_{\phi \phi}}{r^2} + \frac{\partial g_{\theta \phi}}{r} + \frac{\partial g_{\phi \theta}}{r} + \frac{\cot \theta g_{\theta \phi}}{r^2} + \frac{\partial g_{r \phi}}{r} + \frac{\partial g_{\phi r}}{r} + \frac{\cot \theta g_{r \phi}}{r^2} \) (A.12c)

\((\nabla \times \nabla \times g^T)^{T}_{\theta \phi} = \frac{\partial^2 g_{\theta \phi}}{\partial r^2} + \frac{\partial^2 g_{\theta \phi}}{\partial \phi^2} + \frac{\partial^2 g_{\phi \phi}}{\partial \phi^2} - \frac{\partial^2 g_{\phi \phi}}{\partial r \partial \phi} + \frac{\partial^2 g_{\phi \phi}}{\partial \phi \partial r} - \frac{\partial g_{\phi \phi}}{r^2 \sin \theta} + \frac{\cot \theta g_{\phi \phi}}{r^2} + \frac{\partial g_{\phi \theta}}{r} + \frac{\partial g_{r \phi}}{r} + \frac{\cot \theta g_{r \phi}}{r^2} \) (A.12d)

\((\nabla \times \nabla \times g^T)^{T}_{\theta r} = \frac{\partial^2 g_{\theta r}}{\partial r^2} + \frac{\partial^2 g_{\theta r}}{\partial \phi^2} + \frac{\partial^2 g_{\phi \phi}}{\partial \phi^2} - \frac{\partial^2 g_{\phi \phi}}{\partial r \partial \phi} + \frac{\partial^2 g_{\phi \phi}}{\partial \phi \partial r} - \frac{\partial g_{\phi \phi}}{r^2 \sin \theta} + \frac{\cot \theta g_{\phi \phi}}{r^2} + \frac{\partial g_{\theta \phi}}{r} + \frac{\partial g_{r \phi}}{r} + \frac{\cot \theta g_{r \phi}}{r^2} \) (A.12e)

\((\nabla \times \nabla \times g^T)^{T}_{\phi r} = \frac{\partial^2 g_{\phi r}}{\partial r^2} + \frac{\partial^2 g_{\phi r}}{\partial \phi^2} + \frac{\partial^2 g_{\phi \phi}}{\partial \phi^2} - \frac{\partial^2 g_{\phi \phi}}{\partial r \partial \phi} + \frac{\partial^2 g_{\phi \phi}}{\partial \phi \partial r} - \frac{\partial g_{\phi \phi}}{r^2 \sin \theta} + \frac{\cot \theta g_{\phi \phi}}{r^2} + \frac{\partial g_{\phi \theta}}{r} + \frac{\partial g_{r \phi}}{r} + \frac{\cot \theta g_{r \phi}}{r^2} \) (A.12f)

\((\nabla \times \nabla \times g^T)^{T}_{\phi \phi} = \frac{\partial^2 g_{\phi \phi}}{\partial r^2} + \frac{\partial^2 g_{\phi \phi}}{\partial \phi^2} + \frac{\partial^2 g_{\phi \phi}}{\partial \phi^2} - \frac{\partial^2 g_{\phi \phi}}{\partial r \partial \phi} + \frac{\partial^2 g_{\phi \phi}}{\partial \phi \partial r} - \frac{\partial g_{\phi \phi}}{r^2 \sin \theta} + \frac{\cot \theta g_{\phi \phi}}{r^2} + \frac{\partial g_{\phi \theta}}{r} + \frac{\partial g_{r \phi}}{r} + \frac{\cot \theta g_{r \phi}}{r^2} \) (A.12g)

\((\nabla \times \nabla \times g^T)^{T}_{\phi \theta} = \frac{\partial^2 g_{\phi \theta}}{\partial r^2} + \frac{\partial^2 g_{\phi \theta}}{\partial \phi^2} + \frac{\partial^2 g_{\phi \phi}}{\partial \phi^2} - \frac{\partial^2 g_{\phi \phi}}{\partial r \partial \phi} + \frac{\partial^2 g_{\phi \phi}}{\partial \phi \partial r} - \frac{\partial g_{\phi \phi}}{r^2 \sin \theta} + \frac{\cot \theta g_{\phi \phi}}{r^2} + \frac{\partial g_{\phi \theta}}{r} + \frac{\partial g_{r \phi}}{r} + \frac{\cot \theta g_{r \phi}}{r^2} \) (A.12h)
We have implemented these equations in 1D, direction $r$, with only slight modifications to the strategy described for Cartesian geometry. Schematically, we represent the overall system of equations in the form

$$\frac{\partial U}{\partial t} + \frac{\partial AF(U)}{\partial V} + \frac{\partial H(U)}{\partial r} = G(q, r) + S(q)$$  \hspace{1cm} (A.13)

Here we distinguish between the area-weighted volumetric flux terms, $AF$, and the spatial flux terms $H$. The geometric source terms are represented by $G(q, r)$, and the plastic source terms are $S(q)$. Note that strict 1D-$r$ flow, there is no angular dependence to any flow variable, and therefore terms proportional to $\cot \theta$ (for example) vanish identically.

As in the Cartesian case, we solve the time-centered edge Riemann problems to deduce single-valued time-centered edge states $U_{i-1/2}^{n+1}$. These edge states are then used to construct the flux terms $F$ and $H$, which are used to compute a preliminary update $\tilde{U}_{i}^{n+1}$ via the difference scheme

$$\tilde{U}_{i}^{n+1} = U_{i}^{n} - \frac{\Delta t}{V_{i}} \left( A_{i+1/2}^{*} F_{i+1/2}^{n} - A_{i-1/2}^{*} F_{i-1/2}^{n} \right) - \frac{\Delta t}{\Delta r_{i}} \left( H_{i+1/2}^{*} - H_{i-1/2}^{*} \right)$$  \hspace{1cm} (A.14)

Next, we modify the preliminary update by inclusion of the geometric source terms. This is made second-order using a predictor-corrector strategy,

$$\tilde{U}_{i}^{n+1} = \tilde{U}_{i}^{n} + \Delta t G(q_{i}^{n}, \sigma_{i}^{n}) \hspace{1cm} (A.15)$$

$$\sigma' = \sigma + \left( \frac{\partial \sigma}{\partial q} \right)^{n} (q'_{i} - q_{i}^{n})$$

$$\tilde{U}_{i}'' = \tilde{U}_{i}^{n+1} + \frac{\Delta t}{2} (G(q_{i}^{n}, \sigma_{i}^{n}) + G(q_{i}^{n}, \sigma'_{i}))$$

The plastic source terms are then evaluated at the half time step, giving the final result

$$\tilde{q}_{i} = \frac{1}{2} (q_{i}^{n} + \tilde{q}_{i}^{n})$$

$$U_{i}^{n+1} = \tilde{U}_{i}^{n} + \Delta t S(\tilde{q}_{i})$$  \hspace{1cm} (A.16)

ACKNOWLEDGMENT

GM benefitted from correspondences with B. Plohr (Los Alamos National Lab), V. Barcilon (University of Chicago), and M. Ortiz (Caltech).

REFERENCES


FIGURES

FIG. 1. Initial conditions and computed results using 160 cells in Cartesian geometry for (a) density, $\rho$; and (b) stress, $\sigma_{xx}$.

FIG. 2. Initial conditions and computed results using 160 cells in Cartesian geometry for (a) density, $\rho$; (b) stress, $\sigma_{xx}$; (c) plastic deformation $\varepsilon_{xx}^p$; and (d) work hardening parameter $\kappa$. 
FIG. 3. $\sigma_{rr}$ calculated from Blake's analytical solution at time $1.6 \times 10^4$, compared with values computed using 200 grid cells.

FIG. 4. Hoop stress, $\sigma_{\theta\theta} = \sigma_{\phi\phi}$. 
FIG. 5. Pressure, \( \sigma_{rr} + \sigma_{\theta \theta} + \sigma_{\phi \phi} \)/3.

FIG. 6. Material velocity \( v_r \) for Blake's problem.
FIG. 7. Longitudinal stress $\sigma_{xx}$ for Wilkins' problem with impact velocity 0.8 km/s. Time in $\mu$s.
FIG. 8. Longitudinal stress $\sigma_{xx}$ for Wilkins' problem with impact velocity 2.0 km/s. Time in $\mu$s.
FIG. 9. Radial velocity at time 4.

FIG. 10. $\sigma_{rr}$ at time 4.
FIG. 11. $\mathcal{F}_{rr}$ component of plastic deformation tensor.

FIG. 12. Work hardening parameter $\kappa$. 
FIG. 13. Error in density constraint, $\rho - \rho_0 \det(g)$.

FIG. 14. Error in curl constraint, $\sqrt{\partial_\alpha \partial_\beta g_{\alpha\beta}}$. 
FIG. 15. Radial velocity at time 4.

FIG. 16. $\sigma_{rr}$ at time 4.
### TABLE 1
Convergence test: pure elasticity

<table>
<thead>
<tr>
<th>geometry</th>
<th>field</th>
<th>$L_1$</th>
<th>$L_2$</th>
<th>$L_{\infty}$</th>
</tr>
</thead>
<tbody>
<tr>
<td>Cartesian</td>
<td>$\rho$</td>
<td>3.33</td>
<td>3.24</td>
<td>3.06</td>
</tr>
<tr>
<td></td>
<td>$u_x$</td>
<td>3.02</td>
<td>2.97</td>
<td>2.89</td>
</tr>
<tr>
<td></td>
<td>$\sigma_{xx}$</td>
<td>3.30</td>
<td>3.47</td>
<td>3.80</td>
</tr>
<tr>
<td></td>
<td>$\sigma_{yy}, \sigma_{zz}$</td>
<td>2.95</td>
<td>2.91</td>
<td>2.69</td>
</tr>
<tr>
<td>cylindrical</td>
<td>$\rho$</td>
<td>2.51</td>
<td>2.68</td>
<td>2.80</td>
</tr>
<tr>
<td></td>
<td>$u_r$</td>
<td>2.84</td>
<td>2.70</td>
<td>2.53</td>
</tr>
<tr>
<td></td>
<td>$\sigma_{rr}$</td>
<td>2.82</td>
<td>2.78</td>
<td>2.68</td>
</tr>
<tr>
<td></td>
<td>$\sigma_{\theta\theta}$</td>
<td>3.12</td>
<td>3.24</td>
<td>3.30</td>
</tr>
<tr>
<td></td>
<td>$\sigma_{zz}$</td>
<td>3.22</td>
<td>3.17</td>
<td>2.97</td>
</tr>
<tr>
<td>spherical</td>
<td>$\rho$</td>
<td>2.42</td>
<td>2.53</td>
<td>2.66</td>
</tr>
<tr>
<td></td>
<td>$u_r$</td>
<td>2.77</td>
<td>2.64</td>
<td>2.55</td>
</tr>
<tr>
<td></td>
<td>$\sigma_{rr}$</td>
<td>2.85</td>
<td>2.78</td>
<td>2.79</td>
</tr>
<tr>
<td></td>
<td>$\sigma_{\theta\phi}, \sigma_{\phi\phi}$</td>
<td>3.31</td>
<td>3.34</td>
<td>3.37</td>
</tr>
</tbody>
</table>

### TABLE 2
Convergence test: elastic-plastic flow

<table>
<thead>
<tr>
<th>geometry</th>
<th>field</th>
<th>$L_1$</th>
<th>$L_2$</th>
<th>$L_{\infty}$</th>
</tr>
</thead>
<tbody>
<tr>
<td>Cartesian</td>
<td>$\rho$</td>
<td>2.31</td>
<td>2.26</td>
<td>2.08</td>
</tr>
<tr>
<td></td>
<td>$u_x$</td>
<td>2.61</td>
<td>2.43</td>
<td>2.20</td>
</tr>
<tr>
<td></td>
<td>$\sigma_{xx}$</td>
<td>2.39</td>
<td>2.40</td>
<td>2.28</td>
</tr>
<tr>
<td></td>
<td>$\sigma_{yy}$</td>
<td>1.31</td>
<td>1.13</td>
<td>1.04</td>
</tr>
<tr>
<td></td>
<td>$\sigma_{zz}$</td>
<td>1.75</td>
<td>1.63</td>
<td>1.51</td>
</tr>
<tr>
<td></td>
<td>$\mathcal{F}_{xx}^p$</td>
<td>3.34</td>
<td>2.86</td>
<td>2.25</td>
</tr>
<tr>
<td></td>
<td>$\mathcal{F}_{yy}^p$</td>
<td>2.52</td>
<td>2.31</td>
<td>2.34</td>
</tr>
<tr>
<td></td>
<td>$\mathcal{F}_{zz}^p$</td>
<td>2.62</td>
<td>2.46</td>
<td>2.56</td>
</tr>
<tr>
<td></td>
<td>$\kappa$</td>
<td>2.59</td>
<td>2.47</td>
<td>2.59</td>
</tr>
</tbody>
</table>