MOLECULAR LINE EMISSION FROM MULTIFLUID SHOCK WAVES. I.
NUMERICAL METHODS AND BENCHMARK TESTS

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ABSTRACT

We describe a numerical scheme for studying time-dependent, multifluid, magnetohydrodynamic shock waves in weakly ionized interstellar clouds and cores. Shocks are modeled as propagating perpendicular to the magnetic field and consist of a neutral molecular fluid plus a fluid of ions and electrons. The scheme is based on operator splitting, wherein time integration of the governing equations is split into separate parts. In one part, independent homogeneous Riemann problems for the two fluids are solved using Godunov’s method. In the other, equations containing the source terms for transfer of mass, momentum, and energy between the fluids are integrated using standard numerical techniques. We show that, for the frequent case where the thermal pressures of the ions and electrons are $\ll$ magnetic pressure, the Riemann problems for the neutral and ion–electron fluids have a similar mathematical structure which facilitates numerical coding. Implementation of the scheme is discussed and several benchmark tests confirming its accuracy are presented, including (1) MHD wave packets ranging over orders of magnitude in length- and timescales, (2) early evolution of multifluid shocks caused by two colliding clouds, and (3) a multifluid shock with mass transfer between the fluids by cosmic-ray ionization and ion–electron recombination, demonstrating the effect of ion mass loading on magnetic precursors of MHD shocks. An exact solution to an MHD Riemann problem forming the basis for an approximate numerical solver used in the homogeneous part of our scheme is presented, along with derivations of the analytic benchmark solutions and tests showing the convergence of the numerical algorithm.

Key words: ISM: clouds – ISM: magnetic fields – magnetohydrodynamics (MHD) – methods: numerical – plasmas – shock waves

1. INTRODUCTION AND MOTIVATION

Shock waves in molecular clouds and star-forming regions are energetic events that can process the interstellar gas and dust they entrain. Shock-excited molecular H$_2$O has been detected by Herschel (van Dishoeck et al. 2011) and shocked species such as H$_2$O and CO are among the targets of opportunity for SOFIA (Eisl"offel et al. 2012). Observations of the kinematics and proper motions of protostellar outflows and their associated shocks yield ages as young as $\sim 10^2–10^3$ yr (e.g., Gueth et al. 1998; Hartigan et al. 2001). Because the timescale for these shocks to develop steady flow is typically $\gtrsim 10^4$ yr (see Section 3), time-dependent solutions are generally required to model observations of the shock-excited emission. For instance, Gusdorf et al. (2011) applied approximate time-dependent models to H$_2$, SiO, and H$_2$O lines in the BHR71 bipolar outflow shocks and concluded that the best fit to the observations occurred in models having ages $< 2000$ yr. However, there was a fairly large degeneracy in the parameter space which allowed reasonable fits.

Modeling shocks in weakly ionized interstellar clouds is a complex task due largely to the complex nature of the clouds themselves: molecular clouds are threaded by interstellar magnetic fields of the order of $10–10^3 \mu$G (Crutcher 2004; Rao et al. 2009; Tang et al. 2009) and are only weakly ionized, with ion fractional abundances $\lesssim 10^{-3}$ (Caselli 2002; Miettinen et al. 2011). Because the bulk of the matter is neutral, clouds have a finite electrical conductivity; consequently, the neutral gas and charged particles collectively constitute a nonideal magnetohydrodynamic (MHD) system. Moreover, the neutral particles respond indirectly to magnetic forces via momentum exchange with the (rare) ions and electrons, so that a multifluid treatment of the dynamics is required. Early work on multifluid MHD shocks established that the ions can be accelerated ahead of the neutral shock front in a “magnetic precursor” (Mullan 1971; Draine 1980). Ion collisional drag in the precursor accelerates, compresses, and heats the neutrals, thereby passing information about the disturbance on to the neutral gas upstream from the shock. Depending on the shock speed, the timescale for collisions between the ions and neutrals, and the rate at which the neutral gas can cool, the resulting MHD shock can be a discontinuous flow (J-shock) with a supersonic to subsonic transition at a shock front, a continuous flow (C-shock) which is supersonic everywhere (Draine 1980; Chernoff 1987; Draine & McKee 1993), or a continuous flow with a supersonic to subsonic transition (C*-shock, Roberge & Draine 1990).

The majority of theoretical work has been done on steady multifluid MHD shock waves. The chemical abundances and excitation of certain molecular species in shock waves propagating perpendicular to the ambient magnetic field (“perpendicular shocks”) were studied by Draine et al. (1983). Steady perpendicular shock models with increasingly elaborate physics (including the effects of water emission, extensive chemical networks, etc.) were subsequently presented by Flower et al. (1985, 1986), Pilipp et al. (1990), Tielens et al. (1994), Kaufman & Neufeld (1996a, 1996b), Flower et al. (1996), Schilke et al. (1997), and Guillet et al. (2009). Steady MHD shocks propagating obliquely to the magnetic field have also been studied (e.g., Wardle & Draine 1987; Caselli et al. 1997). However, the applicability of steady models is limited when it comes to interpreting observations of shocks in star-forming regions for the reasons noted above.
A study of time-dependent, multidimensional C shocks using a two-fluid finite-difference formulation, and including the inertia of the ions, was presented by Tóth (1994). Time-dependent simulations of the formation of C-shocks in models that neglected the inertia of the ion–electron fluid were produced by Smith & Mac Low (1997). Mac Low & Smith (1997) investigated the nonlinear development of Wardle instabilities in three-dimensional C-shocks (Wardle 1990) using a two-fluid, time-dependent, finite-difference MHD code, including the inertia of the ions. Ciolek & Roberge (2002) and Ciolek et al. (2004) simulated the formation and evolution of one-dimensional perpendicular shocks, accounting for the inertial effect of charged dust grains (ion inertia was ignored, however). Time-dependent, one-dimensional, perpendicular shock models including the inertia of the ions, heating and cooling, and an extensive network of reactions for 33 different chemical species, were described by LeSaffre et al. (2004a, 2004b). Multidimensional and multifluid (including charged dust grain species, but neglecting ion inertia) evolutionary MHD shock models were discussed by Falle (2003) and Van Loo et al. (2009); the same numerical code was also used to investigate the effect of upstream density perturbations on dusty C-shocks in Ashmore et al. (2010). Finally, a one-dimensional study of the redistribution of mass and magnetic flux and potential protostellar core formation due to ambipolar diffusion (the drift of the charged particles with respect to the neutrals) in transient C-shocks arising from colliding flows was presented by Chen & Ostriker (2012); in their calculation the inertia of the ions (and self-gravity of the system) was not included.

A pseudo-time-dependent (quasi-Lagrangian), one-dimensional MHD shock model was employed in several studies of multifluid shocks (e.g., Chieze et al. 1998; Gusdorf et al. 2008; Flower & Pineau des Forêts 2010, 2012). In these models, time dependence is mimicked by setting the partial derivative \( \partial / \partial t \) equal to zero in the governing equations for conservation of mass, momentum, energy, and magnetic flux (see Equations (1)–(6) below), and then replacing the spatial derivatives \( \partial / \partial x \) in those equations with a “flow derivative” \( (1/v) \partial / \partial t \), where \( v \) is the fluid velocity of either the neutrals or the ions. The pseudo-time-dependent approach has grown to include increasingly more detailed and chemical networks.

To study MHD turbulence in weakly ionized clouds, some have advocated a multifluid MHD method referred to as the “heavy ion approximation” (Li et al. 2006; Oishi & Mac Low 2006; Li et al. 2008). This approximation was introduced to overcome numerical difficulties associated with large variations of the ion Alfvén speed, which are typical in molecular clouds (see Equation (31) below) and limit the size of the time steps a numerical code can take to produce accurate, stable simulations (via the CFL condition; see Equation (29)). In the heavy ion approximation the ions are artificially raised to increase the mass density of the ion fluid so that smaller, more tractable ion Alfvén speeds (and thus, CFL-limited time steps) are attained. To keep the collisional drag terms between the neutral and ion fluids unchanged, the coupling constants that appear in the ion–neutral frictional forces (e.g., see Equations (13)–(15)) are also adjusted so as to offset the effect the inflation of the ion masses has on the forces. However, Tilley & Balsara (2010) showed that while the collisional force terms remain unchanged in the heavy ion approximation, characteristic length scales (which involve different combinations of the ion masses and coupling coefficients) associated with the dissipation range of MHD turbulence are not calculated correctly using this method.

Roberge & Ciolek (2007, hereafter RC07) examined the initial phases of MHD shock formation in weakly ionized clouds, including the effects of ion inertia. They noted that for sufficiently small times the disturbance in the ion–electron fluid is linear, and derived explicit analytic solutions for the time evolution of a perpendicular shock. RC07 found that at very early times the inertia of the ions determines the formation and propagation of the magnetic precursor to a neutral shock. At later times the ions’ inertia becomes negligible, their motion is force free, and the evolution of the precursor is nearly self-similar.

The essential mathematical difficulty in modeling time-dependent, multifluid shock waves is caused by the transfer of mass, momentum, and energy between the fluids by elastic ion–neutral scattering, ionization/recombination, and numerous other atomic and molecular processes. In the absence of coupling between the fluids, the multifluid shock problem would reduce to independent solutions of Euler’s equations for the neutral gas and the equations of an ideal MHD for the ion–electron fluid. In both cases the governing equations are hyperbolic partial differential equations (PDEs) which can be solved using well-known techniques. Many of the algorithms for Euler’s equations rely on one’s ability to solve the “Riemann problem” (e.g., Courant & Friedrichs 1948), wherein the gas initially is in two uniform but different states separated by a discontinuity. In particular, Godunov (1959) realized that the evolution of gas in two adjacent cells of a finite difference grid is indeed a Riemann problem and exploited this fact to produce an algorithm, Godunov’s method, whose descendants account for a large subset of all algorithms for gas dynamics (e.g., see Toro 2009 and Pirozzoli 2010). Riemann solutions are combinations of the characteristic waves of a fluid—sound waves, rarefactions, and contact discontinuities in the case of a gas—which are orchestrated to satisfy certain matching conditions where the waves intersect (see Appendix A for an example). Riemann–Godunov (RG) algorithms also exist for ideal MHD but are complicated by the fact that seven characteristic waves exist when the flow propagates at an arbitrary angle with respect to the magnetic field (see Figure 2 of Dai & Woodward 1994 or Equation (7) of Torrilhon 2003).

When interfluid coupling is included, the only changes to the hyperbolic PDEs for each fluid are the addition of certain “source terms” which describe mass, momentum, and energy transfer between the fluids (see Equations (1)–(6)). Toro (2009) noted that problems of this nature could be solved by a technique called operator splitting and gave a simple, nonhydrodynamical example. Operator splitting combines separate solutions of the “homogeneous problem” (source terms set to zero) obtained, e.g., with an RG algorithm, and the “inhomogeneous problem” (evolution due only to the source terms) obtained with standard methods for ordinary differential equations (ODEs). Tilley et al. (2012) recently presented an operator-splitting scheme for multifluid MHD that treats general time-dependent flows for a two-fluid system of neutrals and ions traveling at any orientation with respect to the magnetic field. Their algorithm retains the inertia of both the neutrals and the ions. Collisional drag between the two fluids (with a velocity-independence collision rate) is included along with an energy equation with source terms for each fluid. The algorithm of Tilley et al. (2012) does not account for mass exchange between the neutral and ion–electron fluids, presumably because this is not important for the applications of interest to them. Tilley et al. (2012) presented various benchmarks tests, including the development of a Wardle instability, and found that the scheme performed well.
In this paper, the first in a series on shocks in molecular clouds, we present a split-operator method similar to the algorithm of Tilley et al. (2012), but tailored to the one-dimensional geometry appropriate for shock waves. This allows us to exploit the unique property of perpendicular shocks in weakly ionized plasmas: there exists an exact solution to the MHD Riemann problem, which is fundamental to our algorithm. While the assumption of perpendicular shocks is obviously a special case, it does treat the fundamental mathematical problem of coupled hyperbolic PDEs including mass, momentum, and energy transfer. Deferring the case of oblique shocks to future work also makes strategic sense because modeling perpendicular shocks has a particular advantage: for likely cloud and shock parameters, the thermal pressure of the ion–electron fluid can be neglected compared to magnetic pressure (Section 2.1). In this regime the charged fluid acoustic modes play no part in the MHD Riemann problem and the number of characteristic waves in the ion–electron fluid reduces from seven to just three—the same number of waves that occur in the gas dynamic Riemann problem. As a result, it turns out that the Riemann problem for the charged fluid can be solved exactly (Appendix A). From this exact solution an approximate but accurate solver can be constructed for use in the numerical solution of the MHD Riemann problem in the split-operator method. In our algorithm we also include the effect of mass transfer between the neutral and charged fluids, which can have profound effects on multifluid shocks (Flower et al. 1985).

The plan of our paper is as follows: in Section 2 we present a formulation of the model, including the governing equations and assumptions, the respective Riemann problems for the neutral gas and ion–electron fluid, and a description of how the split-operator scheme is implemented to evolve model shocks in time. We assess the accuracy of our numerical solutions with various benchmark tests in Section 3. Our results are summarized in Section 4. In the appendices, we present the exact solution to the MHD Riemann problem for perpendicular flows (Appendix A), an approximate solver based on this solution (Appendix B), a derivation of the analytic solutions used as benchmark tests (Appendix C), and tests confirming the order of spatial and temporal convergence of our numerical algorithm (Appendix D).

2. FORMULATION

Our area of interest is shocks and related flows in interstellar molecular clouds. These clouds contain neutral particles with number density \( n_n \) and particle mass \( m_n \) plus a weakly ionized plasma of singly charged ions and electrons having number densities \( n_i \) and \( n_e \), and particle masses \( m_i \) and \( m_e \), respectively. Since we are not concerned with the problem of gravitational collapse and deal with systems typically having length scales much smaller than the Jeans length, self-gravity of the gas is ignored. We adopt a Cartesian coordinate system \((x, y, z)\) and restrict our attention to models in which all of the physical variables are functions of time and the \( x \)-coordinate only. We assume that there is a magnetic field \( B = B(x, t)\hat{z} \), and that each of the fluids has a velocity \( v_\alpha = v_\alpha(x, t)\hat{x} \), with \( \alpha = n, i, e \). Thus we consider perpendicular shocks only; however, we note that the split-operator method described in this paper can be extended to other geometries.

In the molecular cloud and core environments we study, the magnetic field strengths and neutral gas densities are such that the ion and electron fluids each have Hall parameters (= charged particle gyrofrequency times the mean collision time with the neutral gas) \( \gg 1 \). This means that the ions and electrons gyrate about a magnetic field line many times before suffering a collision with a neutral particle, and can therefore be considered to be attached to the field. The magnetic field is thus “frozen into” the charged fluid, and the electrons and ions move together with \( v_\alpha \approx v_i \).

Finally, we ignore the effects of interstellar dust grains. Grains can become charged and numerical simulations have shown that under certain conditions they can affect the evolution of MHD shocks (Ciolek & Roberge 2002; Ciolek et al. 2004; Van Loo et al. 2009); the influence of dust grains in MHD flows using a split-operator method will be considered at a later time.

2.1. Governing Equations and Assumptions

Comprehensive derivations of the multifluid MHD equations for astrophysical flows are presented in Draine (1986) and Mouschovias (1987). For the geometry adopted here they are

\[
\frac{\partial n_n}{\partial t} + \frac{\partial (n_nv_n)}{\partial x} = S_n
\]

(1)

\[
\frac{\partial (n_nv_n)}{\partial t} + \frac{\partial}{\partial x} \left( n_nv_n^2 + P_n \right) = F_n
\]

(2)

\[
\frac{\partial E_n}{\partial t} + \frac{\partial}{\partial x} \left[ (E_n + P_n)v_n \right] = F_nv_n - \frac{1}{2} S_n v_n^2 + G_n - \Lambda_n
\]

(3)

\[
\frac{\partial \rho_i}{\partial t} + \frac{\partial (\rho_iv_i)}{\partial x} = - S_n
\]

(4)

\[
\frac{\partial (\rho_iv_i)}{\partial t} + \frac{\partial}{\partial x} \left( \rho_i v^2_i + \frac{B^2}{8\pi} \right) = - F_n
\]

(5)

\[
\frac{\partial B}{\partial t} + \frac{\partial (Bv_i)}{\partial x} = 0.
\]

(6)
Equations (1)–(3) express the conservation of mass, momentum, and energy for the neutral gas. Equations (4)–(5) express mass and momentum conservation for the ion–electron fluid and (6) is the induction equation. We also impose macroscopic charge neutrality, 

\[ e(n_i - n_e) = 0. \] (7)

To the set above one should generally add separate energy equations for the ions and electrons, which are needed to calculate the ion and electron temperatures, \( T_i \) and \( T_e \). However, if elastic collisions dominate the transfer of energy between the ions and neutrals, as is usually the case, then

\[ T_i \approx T_n + \frac{m_n}{3k_B} (v_i - v_n)^2 \] (8)

(Chernoff 1987), an approximation we adopt. There is no analogous approximation for \( T_e \) because energy exchange between the electrons and neutrals is dominated by complex inelastic processes such as electron impact excitation and ionization. We temporarily omit the electron energy equation and estimate \( T_e \) when it is needed in one of our benchmark calculations (see Section 3.3).

The quantities \( \rho_n (=m_n n_n) \) and \( \rho_i (=m_i n_i) \) are the mass densities of the neutral gas and ion–electron fluid (we neglect the electrons’ contribution to \( \rho_n \)). The neutral gas has thermal pressure

\[ P_n = \frac{\rho_n k_B T_n}{m_n}, \] (9)

where \( T_n \) is the neutral temperature and \( k_B \) is the Boltzmann constant. Its energy density is

\[ E_n = \frac{1}{2} \rho_n v_n^2 + \rho_n \epsilon_n, \] (10)

where \( \epsilon_n \) is the internal energy per unit mass. In general it is necessary to find \( \epsilon_n \) by a kinetic calculation of the level populations of rotationally and vibrationally excited H\(_2\), CO, H\(_2\)O, etc. (e.g., Flower et al. 2003). Here we temporarily forego this complication and use the internal energy for an ideal gas in thermodynamic equilibrium,

\[ \epsilon_n = \frac{P_n}{\rho_n (\gamma - 1)}, \] (11)

with \( \gamma = 5/3 \) in the calculations presented below. However, it is important to note that the split-operator method does not preclude a kinetic calculation of \( \epsilon_n \) or nonideal equations of state for \( \epsilon_n \).

In the ion–electron momentum equation (5) we have dropped the thermal-pressure force because it is typically much smaller than the magnetic-pressure force. The ratio of the ion+electron to magnetic pressure is

\[ \frac{P_n}{\rho_n |v_n - v_i|} = \frac{n_i k_B (T_i + T_e)}{(B^2/8\pi)} = 1.39 \times 10^{-6} \left( \frac{50 \mu G}{B} \right)^2 \left( \frac{n_n}{10^4 \, \text{cm}^{-3}} \right) \left( \frac{x_i}{10^{-7}} \right) \left( \frac{T_i + T_e}{10^3 \, \text{K}} \right), \] (12)

where \( x_i = n_i/n_n \) is the fractional ionization and we have normalized quantities to representative values for a shock wave. We conclude that thermal pressure of the ions and electrons is indeed negligible for the conditions of interest here.

The source terms \( S_n, F_n, \) and \( G_n - \Lambda_n \) are the net rates per unit volume at which mass, momentum, and thermal energy are added to the neutral gas, where \( \Lambda_n \) is rate of radiative cooling and \( G_n \) is the net rate of heating by all other processes. In Section 3, we describe benchmark calculations with various assumptions about the source terms. In all cases momentum exchange is dominated by elastic ion–neutral scattering with

\[ F_n \approx F_{n,el} = -\frac{\rho_n}{\tau_m} (v_n - v_i). \] (13)

The timescale for drag to accelerate the ion–electron fluid is

\[ \tau_m = \frac{(m_i + m_n)}{\rho_n \langle \sigma w \rangle_{in}} \left( 1 + \left[ \frac{\sigma_{geo}|v_n - v_i|}{\langle \sigma w \rangle_{in}} \right]^2 \right)^{-1/2}, \] (14)

where \( \langle \sigma w \rangle_{in} \) is the Langevin collision rate (Gioumousis & Stevenson 1958; Flower 2000) and \( \sigma_{geo} \) is the geometric cross section. It follows from Newton’s 3rd Law that the time to accelerate the neutral gas is

\[ \tau_{ni} = (\rho_n/\rho_i) \tau_m \gg \tau_m. \] (15)

The ion–neutral and neutral–ion drag times, \( \tau_m \) and \( \tau_{ni} \) respectively, are fundamental timescales for the flow.
2.2. The Split Operator Method

The system of governing PDEs (1)–(6) has the conservative form

$$\frac{\partial U}{\partial t} + \frac{\partial F}{\partial x} = S,$$  \hspace{1cm} (16)

where $U \equiv [U_n, U_i]^T$ is the array of conserved dependent variables, $F \equiv [F_n, F_i]^T$ is the corresponding array of fluxes, and $S \equiv [S_n, S_i]^T$ is the array of source terms, with

$$U_n \equiv [\rho_n, \rho_n v_n, E_n]^T,$$  \hspace{1cm} (17)
$$U_i \equiv [\rho_i, \rho_i v_i, B]^T,$$  \hspace{1cm} (18)
$$F_n \equiv [\rho_n v_n, \rho_n v_n^2 + P_n, (E_n + P_n) v_n]^T,$$  \hspace{1cm} (19)
$$F_i \equiv [\rho_i v_i, \rho_i v_i^2 + B^2/8\pi, B v_i]^T,$$  \hspace{1cm} (20)
$$S_n \equiv \left[ S_n, F_n, F_n v_n - \frac{1}{2} S_n v_n^2 + G_n - A_n \right]^T,$$  \hspace{1cm} (21)
$$S_i \equiv [-S_n, -F_n, 0]^T.$$  \hspace{1cm} (22)

There are several different ways to solve system (16). In this paper we apply a scheme known as operator splitting, an authoritative discussion of which can be found in Toro (2009, see chapter 15). The basic idea is to split the full problem into two parts, called the homogeneous and inhomogeneous sub-problems. In the former solves the hyperbolic system

$$\frac{\partial U}{\partial t} + \frac{\partial F}{\partial x} = 0$$  \hspace{1cm} (23)

on a discretized computational domain using, say, an RG solver (see Section 2.3). In the latter solves

$$\frac{dU}{dt} = S$$  \hspace{1cm} (24)

using, say, a Runge–Kutta algorithm for ODEs.

The accuracy of the resulting solution depends on the manner in which the sub-problems are coordinated. For example, $U(x, t)$ can be advanced from time $t$ to $t + \Delta t$ with first-order accuracy in $\Delta t$ by setting

$$U(x, t + \Delta t) = T_S^{(\Delta t)} T_{RG}^{(\Delta t)} [U(x, t)],$$  \hspace{1cm} (25)

where the operator $T_{RG}^{(\Delta t)}$ advances the solution from time $t$ to $t + \Delta t$ by solving (23). The result is then fed to $T_S^{(\Delta t)}$, which advances the solution from $t$ to $t + \Delta t$ by solving (24). Second-order accuracy can be attained by using what is sometimes referred to as “Strang splitting” (Strang 1968; Toro 2009),

$$U(x, t + \Delta t) = T_S^{(\Delta t/2)} T_{RG}^{(\Delta t)} T_S^{(\Delta t/2)} [U(x, t)],$$  \hspace{1cm} (26)

where the inhomogeneous sub-problem is advanced by a “predictor” half step $\Delta t/2$, the homogeneous sub-problem is advanced by a full step $\Delta t$, and the inhomogeneous sub-problem is integrated again over another half step. We use Strang splitting in our algorithm (see Section 2.3).

In the operator-splitting method, the homogeneous subproblem (23) can be further separated into independent RG problems for each fluid:

$$\frac{\partial U_n}{\partial t} + \frac{\partial F_n(U_n)}{\partial x} = 0,$$  \hspace{1cm} (27)
$$\frac{\partial U_i}{\partial t} + \frac{\partial F_i(U_i)}{\partial x} = 0.$$  \hspace{1cm} (28)

This means that the neutral gas and ion–electron fluid do not directly influence one another in either of their separate RG problems. The two fluids are therefore dynamically uncoupled during this stage of the calculation. Operationally, this has the advantage that existing well-developed numerical techniques (such as approximate Riemann solvers for gas dynamics) can be applied readily to RG problem (27) for the neutral gas. In Appendix B, we describe an accurate MHD Riemann solver for RG problem (28) for the ion–electron fluid. Thus there is an overall symmetry to the solution of the two separate RG problems when the operator splitting scheme is used. Exploiting this symmetry greatly enhances the efficiency, and simplifies the writing, of a numerical code designed to model multifluid shock waves. Another advantage of our scheme is that by having RG problem (28) as one of the governing equations, the inherent MHD hyperbolic (i.e., wave) structure is unambiguously built into the time evolution. That is, the inertia of the ion–electron fluid is always included. Of course the neutral gas and ion–electron fluid do influence one another. Dynamical recoupling of the fluids occurs during the source integration stage (24) of the split-operator method. As we show in our benchmark tests (Section 3), the interaction between the fluids is indeed accurately accounted for during this stage of the calculation.
2.3. Outline of the Algorithm

Models are calculated on a spatial domain consisting of a set of \( N \) fixed mesh points \( \{x_j, j = 1, \ldots, N\} \) with uniform spacing \( \Delta x = x_j - x_{j-1} \). Variables are calculated at \( x_j \). Located midway between cell \( j \) (centered about \( x_j \)) and cell \( j + 1 \) (centered about \( x_{j+1} \)) is the cell face \( x_{j+1/2} = x_j + \Delta x/2 \).

To advance a model in time, a stable numerical time step \( \Delta t \) has to be chosen. One possibility is to impose the Courant–Friedrichs–Lewy (CFL; e.g., Toro 2009) condition at each mesh point \( j \),

\[
\Delta t_{\text{CFL},j} = \frac{\nu \Delta x}{\lambda_{\text{max},j}},
\]

where \( \lambda_{\text{max},j} \) is the maximum (for all wave modes including both fluids) wave speed at grid point \( x_j \) and \( \nu \) is a number such that \( 0 < \nu < 1 \). In our application it is typically the case that the MHD wave speeds far exceed those of the neutral gas. Therefore we set

\[
\lambda_{\text{max},j} = (|v| + V_{\text{ims}})_j,
\]

where \( V_{\text{ims}} \) is the ion magnetosound speed. When ion–electron pressure is neglected, \( V_{\text{ims}} = V_{\text{IA}} \), where the ion Alfvén speed

\[
V_{\text{IA}} \equiv \frac{B}{\sqrt{4\pi \rho_i}} = 6.90 \times 10^2 \left( \frac{B}{50 \, \mu G} \right) \left( \frac{25 \, m_p}{m_i} \right)^{1/2} \left( \frac{10^{-7} \, \Omega \cdot m}{x_i} \right)^{1/2} \left( \frac{10^4 \, \text{cm}^{-3}}{n_n} \right)^{1/2} \text{ km s}^{-1}
\]

(\( m_p \) is the proton mass). However, the source terms \( \mathbf{S}_n \) and \( \mathbf{S}_i \) also contain important timescales, including the collisional drag times, \( \tau_m \) and \( \tau_{\text{in}} \), as well as scales related, e.g., to the heating and cooling of the gas. Let \( \Delta t_{S_j} \) be some fraction of the smallest timescale associated with the source terms at \( x_j \). To have a stable numerical time integration throughout the entire computational mesh, it is then necessary that the time step be such that

\[
\Delta t \leq \min[\Delta t_{S_j}, \Delta t_{\text{CFL},j}].
\]

Once a stable step size has been determined, integration proceeds as given by Equation (26) in the fashion described in the subsections below.

As discussed early on by Paleologou & Mouschovias (1983; see their Appendix A), large disparities exist between the magnitudes of fluid timescales and the multifluid MHD equations rendering the governing equations mathematically "stiff." This means that \( \Delta t \) will often have to be less than the smallest physical timescale in a system containing other timescales which are much greater in size. In this circumstance a large number of small time steps will then be required to follow a state that is evolving on a much longer natural timescale. An extreme example occurs in the third benchmark test presented in Section 3.1.3 which is followed in its development until several neutral–ion collision times \( \tau_m \) have elapsed, up to a time \( \approx 2.6 \times 10^5 \text{ yr} \) (\( \sim 10–100 \) times greater than the ages of shocks in star-forming regions that we intend to study; see Section 1). To stably integrate that model to that time using an explicit method, \( \Delta t \) had to be kept below the ion–neutral timescale \( \tau_{\text{in}} \approx 1.2 \times 10^{-2} \text{ yr} \); \( \Delta t = 0.4 \tau_{\text{in}} \) was actually used for that model. Thus, \( \sim 8 \times 10^7 \) time steps were taken to reach completion. While this is indeed a large number of computational steps, it is not prohibitively so with modern computational methods.

2.3.1. Step 1: First Source Integration

Given some initial values \( \mathbf{U}(x, t) \), system (24) is advanced in time by \( \Delta t/2 \). This can be carried out using well-known and tested numerical integration methods such as Runge–Kutta schemes, or explicit multistep schemes, etc. (e.g., see chap. 6 of Atkinson 1989; or chapter 16 of Press et al. 1996). For the benchmark calculations presented in Section 3, we used a second-order Runge–Kutta integrator. The result of Step 1 is an updated state vector \( \mathbf{U}^1 = [\mathbf{U}_n^1, \mathbf{U}_l^1]^T \) which becomes the input for Step 2.

2.3.2. Step 2: Riemann–Godunov Integration

In Step 2 the dynamically decoupled RG problems (27) and (28) are advanced a full time step. The solution for the conserved dependent variables at the end of this integration step, \( \mathbf{U}^2 = [\mathbf{U}_n^2, \mathbf{U}_l^2]^T \), is given by

\[
\mathbf{U}_j^2 = \mathbf{U}_j^1 - \frac{\Delta t}{\Delta x} (\mathbf{F}_j(x_{j+1/2}) - \mathbf{F}_j(x_{j-1/2})),
\]

with \( j = n, l \).

The variables are known at the grid points, \( x_j \), but the fluxes in Equation (33) are required at the cell faces. Second-order accuracy in space and time is attained by first "reconstructing" the data and then interpolating the variables from the grid points to the cell faces using total variation diminishing (TVD) slope limiter methods such as those employed in the MUSCL–Hancock scheme (Toro 2009). For any dependent variable \( q \) in the cell centered about \( x_j \), reconstruction and interpolation to the cell faces are carried out by setting

\[
q_j^R = q(x_j) + \frac{1}{2} \Delta_j, \quad q_j^L = q(x_j) - \frac{1}{2} \Delta_j,
\]

with \( \Delta_j \) being the "limited average slope value" of \( q \). There are many different kinds of slope limiter. For example, the MINBEE slope limiter is

\[
\Delta_j = \begin{cases} 
\max[0, \min(\Delta_j/2, \Delta_{j+1/2})] & \text{if } \Delta_{j+1/2} > 0, \\
\min[0, \max(\Delta_j/2, \Delta_{j+1/2})] & \text{if } \Delta_{j+1/2} < 0,
\end{cases}
\]
Based on Colella (1997), the latter solver is similar to the approximate Riemann solver presented in Toro (2009, chapter 9), but has been extended to RG problems that also include nonideal gases (e.g., Colella & Glaz, 1985).

We normally use the van Leer limiter but both are easy to implement and yield good results. We also impose the positivity conditions of Waagan (2009) to the slope limiting and interpolation procedure for the neutral gas, which ensures that \( \rho_n, P_n, T_n \) and \( \epsilon_n \) are always \( >0 \) at all \( x_j \), even for very hypersonic flows.

Another TVD-limiter which can be used for the data reconstruction of the charged fluid and magnetic field variables is the one derived by \( \dot{\text{C}} \)ada & Torrilhon (2009). It has the form:

\[
\Delta j_{+1/2} = q(x_{j+1}) - q(x_j), \quad \Delta j_{-1/2} = q(x_j) - q(x_{j-1})
\]

(Toro 2009), and a generalized version of the monotonic slope limiter of van Leer (1979) is

\[
\bar{\Delta}_j = \sigma_j \min \left[ \Theta |\Delta j_{+1/2}|, \frac{1}{2} |\Delta j_{+1/2} + \Delta j_{-1/2}|, \Theta |\Delta j_{-1/2}| \right], \quad 1 \leq \Theta \leq 2,
\]

(37)

\[
\sigma_j = \begin{cases} 
1 & \text{if } \Delta j_{-1/2} > 0, \quad \Delta j_{+1/2} > 0, \\
-1 & \text{if } \Delta j_{-1/2} < 0, \quad \Delta j_{+1/2} < 0, \\
0 & \text{otherwise}.
\end{cases}
\]

(38)

We integrate Equations (27) and (28) over a virtual (predictor) time increment \( \Delta t/2 \) further improves the left and right cell face values in each cell:

\[
\bar{U}_{\beta,i,j}^R = U_{\beta,i,j}^R - \frac{\Delta t}{2\Delta x} (\mathcal{F}_\beta(U_{\beta,i,j}^R) - \mathcal{F}_\beta(U_{\beta,i,j}^L)),
\]

(42a)

\[
\bar{U}_{\beta,i,j}^L = U_{\beta,i,j}^L - \frac{\Delta t}{2\Delta x} (\mathcal{F}_\beta(U_{\beta,i,j}^R) - \mathcal{F}_\beta(U_{\beta,i,j}^L)).
\]

(42b)

(Toro 2009).

Data reconstruction and interpolation yields variable pairs \( \{\bar{U}_{\beta,j}^R, \bar{U}_{\beta,j+1/2}\} \) at \( x_{j+1/2} \) and \( \{\bar{U}_{\beta,j-1}, \bar{U}_{\beta,j}^L\} \) at \( x_{j-1/2} \), which can be used to calculate the fluxes \( \mathcal{F}_\beta(x_{j+1/2}) \) and \( \mathcal{F}_\beta(x_{j-1/2}) \) needed in Equation (33). Calculation of the fluxes is accomplished by solving the Riemann problem at the cell faces using the method of Godunov (1959; for an extremely comprehensive discussion, see Toro 2009). Solution of the RG problems at each cell interface is carried out using approximate Riemann solvers, one for the neutral gas and another for the ion–electron fluid. The approximate MHD Riemann solver used for the charged fluid is derived in Appendix B. For the neutral gas we use the approximate gas dynamic Riemann solver described in Section 5 of Almgren et al. (2010; reportedly based on Colella 1997). The latter solver is similar to the approximate Riemann solver presented in Toro (2009, chapter 9), but has been extended to RG problems that also include nonideal gases (e.g., Colella & Glaz, 1985). Using the solver of Almgren et al. allows the modeling of fluids having nonideal equations of state, or those that are not in LTE. Our approximate gas dynamic Riemann solver is very similar in structure to our approximate MHD Riemann solver. (Hence the general symmetry of the problem when using the split-operator method, as noted above.)

The approximate gas dynamic and MHD Riemann solvers are used to obtain the (Godunov) fluxes at the faces of each cell, \( \mathcal{F}_\beta(x_{j+1/2}) \) and \( \mathcal{F}_\beta(x_{j-1/2}) \). The details of how this is done for the ion–electron fluid MHD problem are presented in Appendix B, and the details for the neutral gas problem follow by analogy (or see Section 5 of Almgren et al. 2010). These values are then inserted into Equation (33), thereby giving \( U_{\beta}^2 \) and \( U_{\beta}^3 \). The updated variables \( U_{\beta}^2 \) are the input for Step 3.

2.3.3. Step 3: Final Source Integration

Another source integration of size \( \Delta t/2 \) is performed as in Step 1. This yields the last update of the physical variables, \( U^3 = [U_{\beta}^2, U_{\beta}^3]^T \). The split-operator integration scheme is thus completed with the second-order accurate (in space and time) final result \( U(x, t + \Delta t) \approx U^3 \).
3. BENCHMARK TESTS

In this section we present some numerical calculations that establish the validity of the split-operator method for multifluid MHD shocks and other flows in weakly ionized interstellar clouds. The accuracy of the algorithm is tested by comparing our numerical results to accurate analytic solutions. Data regarding the convergence and second-order scaling of our algorithm are presented in Appendix D. In all of the test models the neutrals are assumed to be an ideal gas of H$_2$ molecules having a ratio of specific heats $\gamma = 5/3$, and the ion mass $m_i$ is set to a generic value of $25m_p$, which could represent ions such as HCO$^+$ and Na$^+$. For momentum transfer by elastic ion–neutral collisions we use a Langevin collision rate $\langle \sigma w \rangle_{in} = 1.7 \times 10^{-9}$ cm$^3$ s$^{-1}$ (McDaniel & Mason 1973) and geometric cross section $\sigma_{geo} = \pi (r_{H_2} + r_i)^2 = 2.86 \times 10^{-15}$ cm$^2$ ($r_{H_2}$ and $r_i$ are the molecular and ionic radii, respectively).

3.1. Linear Wave Packets

In this test we set the source terms $S_n$, $G_n$, and $\Lambda_n$ to zero in Equations (1)–(6) and assume that momentum transfer is entirely due to elastic ion–neutral scattering. We initiate small-amplitude disturbances in the plasma and follow the subsequent evolution of the charged and neutral fluids. As discussed in Ciolek & Roberge (2002, Section 2.3.2) and Mouschovias et al. (2011, Section 3.2), there exist two important length scales relevant to MHD waves propagating perpendicular to the magnetic field. The first length scale is the ion magnetosound wave upper cutoff $L_{ims} = 4\pi V_{ims} \tau_{in}$. Ion magnetosound waves with wave speed $V_{ims} = V_{iA}$ can propagate when their wavelength is less than $L_{ims}$. The second length scale is the neutral magnetosound wave lower cutoff length scale $L_{nms} = \pi V_{nA} \tau_{ni}/V_{nms}$. Magnetosound waves in the neutral fluid travel at the neutral magnetosound speed,

$$V_{nms} = \left( C_n^2 + V_{nA}^2 \right)^{1/2},$$

where

$$C_n \equiv \left( \frac{\partial P_n}{\partial \rho_n} \right)_{s_n}^{1/2} \left( \frac{\gamma P_n}{\rho_n} \right)^{1/2} = 0.262 \left( \frac{\gamma}{5/3} \right)^{1/2} \left( \frac{T}{10 \text{K}} \right)^{1/2} \left( \frac{2m_p}{m_n} \right)^{1/2} \text{km s}^{-1},$$

and $U(x,t+\Delta t) = U^3$. 

Figure 1. Schematic diagram of our split-operator algorithm to simulate multifluid MHD shocks and related flows in interstellar clouds and cores.
is the adiabatic sound speed (the derivative is taken at constant neutral entropy \( s_n \)), and

\[
V_{nA} = \frac{B}{\sqrt{4\pi \rho_n}} = 0.771 \left( \frac{B}{50 \mu G} \right) \left( \frac{2m_p}{m_n} \right)^{1/2} \left( \frac{10^4 \text{ cm}^{-3}}{n_n} \right)^{1/2} \text{ km s}^{-1}
\]  

is the neutral Alfvén speed. Neutral magnetosound waves can propagate for wavelengths greater than \( \lambda_{ms} \). At wavelengths between \( \lambda_{ms} \) and \( \lambda_{mm} \), there is no MHD wave propagation; any mode excited on these scales produces ambipolar diffusion of the ions, electrons, and magnetic field through the neutral gas.

To selectively excite different wave modes we impose Gaussian initial perturbations in the ion density and magnetic field of the form

\[
B(x, 0) = B_0 \left[ 1 + A_p \exp \left( -\frac{x^2}{L_G^2} \right) \right]
\]

\[
n_i(x, 0) = n_0 \frac{B(x, 0)}{B_0}.
\]

The charged fluid is initially taken to be stationary with \( n_p(x, 0) = n_{\text{eq}}, \) \( P_n(x, 0) = P_{\text{eq}}, \) and \( v_n(x, 0) = 0. \) Quantities with a ‘0’ subscript are constant. \( A_p \) is the dimensionless amplitude of the perturbation, and \( L_G \) is its width; wave modes excited by this perturbation will have wavelengths clustered about \( L_G.\)

Below we present three benchmark tests having different packet widths, \( L_G, \) but the same perturbation amplitude \( A_p = 0.01. \) All three tests have \( n_0 = 2 \times 10^4 \text{ cm}^{-3} \) and \( x_0 = n_\text{eq}/n_0 = 3.16 \times 10^{-8}. \) All particle species are assigned the same reasonable but somewhat arbitrary temperature: \( T_\text{eq} = T_p = T_0 = 10 \text{ K}. \) The unperturbed magnetic field strength is \( B_0 = 50 \mu G, \) yielding an ion magnetosound speed \( V_{ims} = V_{IA} = 868 \text{ km s}^{-1}, \) and a neutral Alfvén speed \( V_{nA} = 0.545 \text{ km s}^{-1}. \) The sound speed \( C_n = 0.262 \text{ km s}^{-1} \) and the neutral magnetosound speed \( V_{nms} = 0.605 \text{ km s}^{-1}. \) It follows that \( \tau_{\text{in}} = 1.26 \times 10^{-2} \text{ yr}, \) \( \tau_{\text{ni}} = 3.19 \times 10^4 \text{ yr}, \)

\[
L_{ims} = 4.33 \times 10^{14} \text{ cm}, \quad \text{and} \quad L_{nms} = 1.55 \times 10^{17} \text{ cm}.
\]

### 3.1.1. Wave Packet 1: \( L_G = 9.35 \times 10^{11} \text{ cm} \)

Because this wave packet has \( L_G \ll L_{ims}, \) we expect that the initial perturbation will generate two traveling wave packets in the ions and the magnetic field, one traveling leftward from the origin and the other rightward. Although acoustic waves in the neutrals can exist on these scales (see, e.g., Figure 1 of Ciolek et al. 2004 or Figure 6 of Mouschovias et al. 2011), they will not be excited by this perturbation: the neutrals are initially uniform and motionless and the neutral–ion drag time, \( \tau_{\text{ni}}, \) is much longer than the decay timescale \( \tau_{\text{dec}} \) for the waves (see below). We expect that in this model the ion and magnetic field wave packets will propagate through a stationary neutral fluid.

Given the assumption of stationary neutrals, analytic expressions for the small-amplitude wave packets in the ion–electron fluid, which are expected to occur in this model, can be derived from the governing equations (1)–(6) in the limit where the wavelengths \( (\sim L_G) \) of the Fourier components comprising the packets are \( \ll L_{ims}. \) For the initial conditions (46), (47), and \( v_i(x, 0) = 0, \) the solutions are

\[
B_{ims}(x, t) = B_0 \left[ 1 + \frac{A_p}{2} \exp \left( -\frac{t}{2\tau_{\text{in}}} \right) \right] \left\{ \exp \left( -\frac{(x - V_{ims}t)^2}{L_G^2} \right) + \exp \left( -\frac{(x + V_{ims}t)^2}{L_G^2} \right) \right\}
\]

\[
+ \frac{\sqrt{\pi}A_pL_G}{8V_{ims}\tau_{\text{in}}} \left[ \exp \left( -\frac{t}{2\tau_{\text{in}}} \right) \right] \left\{ \text{erf} \left( \frac{x + V_{ims}t}{L_G} \right) - \text{erf} \left( \frac{x - V_{ims}t}{L_G} \right) \right\},
\]

\[
n_{i,ims}(x, t) = n_0 \frac{B_{ims}(x, t)}{B_0},
\]

\[
v_{i,ims}(x, t) = \frac{A_p}{2} V_{ims} \exp \left( -\frac{t}{2\tau_{\text{in}}} \right) \left\{ \exp \left( -\frac{(x - V_{ims}t)^2}{L_G^2} \right) - \exp \left( -\frac{(x + V_{ims}t)^2}{L_G^2} \right) \right\}
\]

(see Appendix C.1). As expected, the solution consists of two traveling Gaussian packets, with propagation velocities \( \pm V_{ims} \) and amplitudes \( A_p/2. \) Momentum exchange (i.e., friction) with the fixed neutrals causes the pulses to decay exponentially on a characteristic timescale \( \tau_{\text{dec}} = 2\tau_{\text{in}}. \) The decay time exceeds \( \tau_{\text{in}} \) by a factor of two because of the equipartition of magnetic and kinetic energy in the ion magnetosound waves which make up the wave packets. Elastic collisions with the neutrals directly reduce the kinetic energy of the ion–electron fluid; the energy stored in the magnetic field is reduced indirectly as it is gradually converted into ion–electron kinetic energy.

Figure 2 shows the results for this model at \( t = 5.42 \times 10^{-3} \text{ yr} \) using our fully nonlinear split-operator method code on a mesh with 5000 grid points. Also shown in each panel are the analytic wave solutions (48a)–(48c) evaluated at the same time. The split-operator code results are in excellent agreement with the analytic solutions, with the relative error in the ion density having a maximum value of \( 2.5 \times 10^{-4}. \) The split-operator scheme we have employed has accurately reproduced all of the fundamental aspects of the physical evolution for this model. This includes the propagation of the magnetosound pulses, the dependence of their speed on the magnetic field strength and inertia of the ions (inherent to the solution from the RG step of the integration), as well as the decay of the two pulse peaks caused by ion–neutral friction (which can only result from the source integration steps). Although the ions and neutrals are dynamically decoupled during the RG stage of the integration, recoupling and realistic interaction of the two fluids are faithfully reproduced during the source integration stages of the algorithm.
3.1.2. Wave Packet 2: \( L_G = 3.74 \times 10^{15} \) cm

This wave packet has \( L_{\text{ims}} \ll L_G \ll L_{\text{ims}} \). As a result, there will be no propagating MHD waves in this case. Instead, the initial perturbation will give rise to an ambipolar diffusion mode, in which the ions, electrons, and magnetic field diffuse outward from the origin through a neutral fluid which is still effectively stationary on the length and timescales characterizing this particular model. The motion of the ions is essentially inertialess or force-free; that is, the driving magnetic-pressure force is almost exactly balanced by the retarding ion–neutral friction (5). The linearized ambipolar diffusion mode is found to have the analytic solutions

\[
B_{\text{ad}}(x,t) = B_0 \left[ 1 + \frac{A_p}{(1 + 4D_{\text{ad}}/L_G^2)^{1/2}} \exp \left( \frac{-x^2/L_G^2}{1 + 4D_{\text{ad}}t/L_G^2} \right) \right] \\
+ \frac{2B_0A_pD_{\text{ad}}\tau_{\text{in}}/L_G^2}{(1 + 4D_{\text{ad}}/L_G^2)^{3/2}} \left[ 1 - \frac{2x^2/L_G^2}{1 + 4D_{\text{ad}}t/L_G^2} \right] \exp \left( \frac{-x^2/L_G^2}{1 + 4D_{\text{ad}}t/L_G^2} \right) \\
- \frac{2B_0A_pD_{\text{ad}}\tau_{\text{in}}/L_G^2}{(1 - 4D_{\text{ad}}t/L_G^2)^{3/2}} \left[ 1 - \frac{2x^2/L_G^2}{1 - 4D_{\text{ad}}t/L_G^2} \right] \exp \left( \frac{-t}{\tau_{\text{in}}} \right) \exp \left( \frac{-x^2/L_G^2}{1 - 4D_{\text{ad}}t/L_G^2} \right), \tag{49a}
\]

\[
n_{i,\text{ad}}(x,t) = n_0 \frac{B_{\text{ad}}(x,t)}{B_0}, \tag{49b}
\]

\[
v_{i,\text{ad}}(x,t) = \frac{2D_{\text{ad}}A_p x}{L_G^2(1 + 4D_{\text{ad}}t/L_G^2)^{3/2}} \exp \left( \frac{-x^2/L_G^2}{1 + 4D_{\text{ad}}t/L_G^2} \right) \\
- \frac{2A_pD_{\text{ad}} x}{L_G^2(1 - 4D_{\text{ad}}t/L_G^2)^{3/2}} \exp \left( \frac{-t}{\tau_{\text{in}}} \right) \exp \left( \frac{-x^2/L_G^2}{1 - 4D_{\text{ad}}t/L_G^2} \right), \tag{49c}
\]

(a derivation is given in Appendix C.2). The quantity

\[
D_{\text{ad}} \equiv V_{\text{ims}}^2 \tau_{\text{in}} \tag{49d}
\]
3.1.3. Wave Packet 3: $L_G = 2.24 \times 10^{17}$ cm

This wave packet has $L_G > L_{\text{nms}}$, which means it will excite low-frequency ($< 1/\tau_{ni}$) neutral magnetosound waves over sufficiently large length and timescales. In these waves, magnetic forces on the ions are transmitted to the neutrals by collisional drag, allowing the neutrals, ions, and magnetic field to participate in collective wave behavior. Because $\rho_n \gg \rho_i$, the overall inertia of the collective motion is due to the neutral fluid. The wave speed is therefore the neutral magnetosound speed $V_{\text{nms}}$ (Equation (43)), with the bulk neutral fluid acting as if it is responding directly to magnetic forces on these scales.

For wavelengths $\geq L_{\text{nms}}$ the ions can still be described as being in force-free motion because $L_{\text{nms}} \gg L_{\text{ims}}$. Using this fact, we find that the Fourier modes for perturbations with $L_G \gg L_{\text{nms}}$ yield the analytic solutions

$$B_{\text{nms}}(x, t) = B_0 \left[ 1 + \frac{A_p G_3}{2} \left( 1 - \frac{2(x + V_{\text{nms}} t)D_{\text{th}}/L_G^2}{V_{\text{nms}}[1 + (2D_{\text{ad}} t/L_G^2)]} \right) \exp \left( \frac{-(x + V_{\text{nms}} t)^2}{L_G^2[1 + (2D_{\text{ad}} t/L_G^2)]} \right) 
+ \left( 1 + \frac{2(x - V_{\text{nms}} t)D_{\text{th}}/L_G^2}{V_{\text{nms}}[1 + (2D_{\text{ad}} t/L_G^2)]} \right) \exp \left( \frac{-(x - V_{\text{nms}} t)^2}{L_G^2[1 + (2D_{\text{ad}} t/L_G^2)]} \right) 
+ A_p G_2 \left( \frac{C_n}{V_{\text{nms}}} \right)^2 \exp \left( \frac{-x^2}{L_G^2[1 + (4D_{\text{ad}} t/L_G^2)]} \right) \right], \tag{50a}$$
\[ n_{1,nms}(x, t) = n_{i0} \frac{B_{nms}(x, t)}{B_0}, \]  

\[ v_{n,nms}(x, t) = \frac{A_p V_{nms} G_1}{2} \left[ \left( \frac{2(x + V_{nms} t) D_{lh}/L_G^2}{V_{nms}[1 + (2 D_{ad}/L_G^2)]} - 1 \right) \exp \left( \frac{-(x + V_{nms} t)^2}{L_G^2[1 + (2 D_{ad}/L_G^2)]} \right) \right. \]
\[ + \left. \left( \frac{1 + 2(x - V_{nms} t) D_{lh}/L_G^2}{V_{nms}[1 + (2 D_{ad}/L_G^2)]} \exp \left( \frac{-(x - V_{nms} t)^2}{L_G^2[1 + (2 D_{ad}/L_G^2)]} \right) \right) \right] \]
\[ - \frac{2 A_p G_2 D_{ih} x}{L_G^2[1 + (4 D_{ad}/L_G^2)]} \exp \left( \frac{-x^2}{L_G^2[1 + (4 D_{ad}/L_G^2)]} \right), \]

\[ n_{n,nms}(x, t) = n_{i0} \left[ 1 + \frac{A_p G_1}{2} \left( \frac{1 - 2(x + V_{nms} t) D_{lh}/L_G^2}{V_{nms}[1 + (2 D_{ad}/L_G^2)]} \right) \exp \left( \frac{-(x + V_{nms} t)^2}{L_G^2[1 + (2 D_{ad}/L_G^2)]} \right) \right. \]
\[ + \left. \left( \frac{1 + 2(x - V_{nms} t) D_{lh}/L_G^2}{V_{nms}[1 + (2 D_{ad}/L_G^2)]} \exp \left( \frac{-(x - V_{nms} t)^2}{L_G^2[1 + (2 D_{ad}/L_G^2)]} \right) \right) \right] \]
\[ - A_p G_2 \exp \left( \frac{-x^2}{L_G^2[1 + (4 D_{ad}/L_G^2)]} \right), \]

\[ v_{1,nms}(x, t) = v_{n,nms}(x, t) \]
\[ + \frac{A_p D_{ad} G_1}{L_G^2[1 + (2 D_{ad}/L_G^2)]} \left[ \left( \frac{1 + 2(x + V_{nms} t) D_{lh}/L_G^2}{V_{nms}[1 + (2 D_{ad}/L_G^2)]} \right)(x + V_{nms} t) - \frac{D_{lh}}{V_{nms}} \exp \left( \frac{-(x + V_{nms} t)^2}{L_G^2[1 + (2 D_{ad}/L_G^2)]} \right) \right] \]
\[ + \frac{A_p G_1 D_{ad}}{L_G^2[1 + (2 D_{ad}/L_G^2)]} \left[ \frac{D_{lh}}{V_{nms}} + \left( \frac{1 - 2(x - V_{nms} t) D_{lh}/L_G^2}{V_{nms}[1 + (2 D_{ad}/L_G^2)]} \right)(x - V_{nms} t) \right) \exp \left( \frac{-(x - V_{nms} t)^2}{L_G^2[1 + (2 D_{ad}/L_G^2)]} \right) \]
\[ + \frac{2 A_p G_2 D_{ih} x}{L_G^2[1 + (4 D_{ad}/L_G^2)]} \exp \left( \frac{-x^2}{L_G^2[1 + (4 D_{ad}/L_G^2)]} \right), \]

where

\[ G_1 = \left[ 1 + \left( \frac{C_n/V_{nms}}{2} \right)^2 \right]^{-1} \left[ 1 + \left( 2 D_{ad}/L_G^2 \right) \right]^{-1/2}, \]  

\[ G_2 = \left[ 1 + \left( \frac{C_n/V_{nms}}{2} \right)^2 \right]^{-1} \left[ 1 + \left( 4 D_{ad}/L_G^2 \right) \right]^{-1/2}, \]

(for a derivation, see Appendix C.3). Appearing in these expressions is the neutral thermal diffusion coefficient

\[ D_{lh} = \left( \frac{C_n}{V_{nms}} \right)^2 D_{ad} = \frac{C_n^2 \tau_{ni}}{1 + (C_n/V_{nms})^2}. \]

The last equality follows from Equations (49d), (15), (31), (45), and (43). The analytic solutions (50a)–(50d) show that there will be Gaussian wave packets propagating with velocities ±\(V_{nms}\), verifying that it is the inertia of the neutral fluid that determines the rate at which the packets propagate. The solutions also reveal that the neutral magnetosol pulse decay by ambipolar diffusion on a timescale \(\tau_{dec} = 2\tau_{ad} = L_G^2/2D_{ad} = 2.66 \times 10^5\) yr; the decay time is a factor of two longer than the diffusion timescale because, once again, there is equipartition of magnetic and kinetic energy.

Figure 4 shows the results of the split-operator algorithm at time \(t = 2.55 \times 10^5\) yr calculated on a mesh with 2000 points. The figure shows that there are indeed two oppositely directed wave packets emerging from the initial disturbance about the origin. There is a decrease in the neutral gas density from the central region as the matter once located there is set into motion and becomes “scooped up” by collisions with the ions, which are being driven outward by magnetic field pressure gradients. The transported neutral gas piles up just outside the depleted central region and increases the density in the two adjacent pulses moving away from the origin. Note that the ion fluid velocity slightly leads that of the neutral fluid. This effect is caused by the inertia of the neutrals, which delays the acceleration of the neutrals from rest up to the same velocity as the ions. Figure 4 also shows the analytic solutions (50a)–(50d) at the same output time. The numerical results are in good agreement with the analytic solutions, with a relative neutral density difference \(\leq 3.1 \times 10^{-4}\) and magnetic field relative difference \(\leq 2.2 \times 10^{-4}\).
Figure 4. As in Figure 2 but for $L_G = 2.24 \times 10^{17}$ cm and time $t = 2.55 \times 10^5$ yr. Top: neutral and ion densities. Also shown are the analytic solutions (50b) and (50d) for the ions (dash-dotted) and the neutrals ($\times$). Middle: neutral and ion velocities. Included are the solutions (50c) and (50e) for the neutrals ($\times$) and the ions (dash-dotted). Bottom: magnetic field. The solution (50a) is also displayed (dash-dotted).

Because the initial state of the neutral fluid is static and completely lacks density or thermal-pressure gradients, the RG integration step for the neutrals does not initiate the wave motion in the neutrals in this example. Rather it is the collisional drag of the magnetically driven ions during the source integration steps which starts the neutrals moving after a time $\gtrsim \tau_{ni}$ has passed for regions near the origin. Once this happens, as we noted above, gradients in the neutral density and pressure form, and thermal-pressure gradient forces start to act on the neutrals during the RG integration step for that fluid. From that point on, there is a combination of wave-driving thermal pressure and magnetic forces (through ion collisions) on the neutrals during all integration stages of the split-operator method scheme. (It is this combination of thermal pressure and magnetic forces acting on the neutrals that explains the appearance of both the sound and Alfvén speeds in the expression for the neutral magnetosound speed $V_{nms}$, Equation (43).)

To summarize: the benchmark tests for wave packets show that our split-operator method successfully incorporates the relevant physics and accurately follows the evolution of both propagating and diffusive flows traveling perpendicular to the magnetic field, over many orders of magnitude in both the temporal and spatial scales.

3.2. Early Evolution of Shocks Caused by a Cloud–Cloud Collision

We now consider a more dynamic test which shows that the split-operator method is also suitable for multifluid shock waves. RC07 considered the collision of two identical clouds and found analytic solutions (using Green functions) which describe the early-time behavior of the ensuing disturbances in the ion–electron fluid. For the parameters RC07 considered, the collision produced forward- and reverse J-shocks in the neutral gas. They referred to the disturbances in the ion–electron fluid as “driven waves” because they are driven by frictional coupling to the neutral flow. Their results describe initial stages in the formation of magnetic precursors on the J-shocks.

The flow geometry in RC07 is the same as the geometry in this paper, with fluid velocities perpendicular to the magnetic field. The clouds are identical and semi-infinite; the collision takes place when their free surfaces coincide at the plane $x = 0$ at $t = 0$. Prior to the collision the charged and neutral fluids in each cloud move together with $v_n(x, 0) = v_i(x, 0) = v(x)$, where

$$v(x) = \begin{cases} 
-10 \text{ km s}^{-1} & \text{for } x > 0, \\
+10 \text{ km s}^{-1} & \text{for } x < 0. 
\end{cases}$$

Each cloud has an initial neutral density $n_n(x, 0) = n_{n0} = 2 \times 10^4$ cm$^{-3}$, ion density $n_i(x, 0) = n_{i0} = 5.72 \times 10^{-4}$ cm$^{-3}$, fractional ionization $x_i(x, 0) = n_{i0}/n_{n0} = 2.86 \times 10^{-3}$, temperatures $T_n(x, 0) = T_i(x, 0) = 10$ K, and magnetic field $B(x, 0) = B_0 = 50 \mu$G. The ion magnetosound speed is $V_{nms} = V_{iA} = 912$ km s$^{-1}$ in the undisturbed charged fluid.
We adopt the same source terms as RC07, who set $S_n = G_n = \Lambda_n = 0$ and assumed that momentum transfer is due solely to elastic ion–neutral scattering. RC07 assumed further that $\tau_n$ is independent of the relative velocity between the ions and neutrals. To make a legitimate comparison with the RC07 solution, we drop the term $\propto |v_n - v_i|^2$ on the right-hand side (rhs) of Equation (14) when calculating $F_n$. The ion–neutral mean collision time in the clouds is then $\tau_n = 1.26 \times 10^{-2}$ yr, and the neutral–ion drag time is $\tau_n = 3.52 \times 10^4$ yr. Because $\tau_n$ is much longer than the times considered by RC07, the motion of the neutral gas is unaffected by coupling to the ions and magnetic field.

Figure 5 shows the multfluid MHD split-operator code results for the colliding clouds test at $t = 0.1\tau_n = 1.26 \times 10^{-3}$ yr. At this very early time ($\ll \tau_n$), the ions and magnetic field have yet to be affected by friction from the neutrals. Two discontinuities in the magnetic field and charged fluid, symmetric about the origin, travel outward at the ion magnetosound speed $V_{ims}$, with their fronts located at $|x_{idis}| = 3.63 \times 10^{12}$ cm. There are also two symmetric neutral shocks propagating away from $x = 0$ at a speed $v_{nshk} = 3.34$ km s$^{-1}$. The neutral shock fronts are located at $|x_{nshk}| = 1.33 \times 10^{10}$ cm so they are not visible on the scale of Figure 5. Also shown (dash-dotted curves) in each panel of the figure are the analytic solutions of RC07. The split-operator results agree very well with the RC07 solution, with a relative difference in the magnetic field that is $\ll 0.015$ behind the discontinuities.

Results for the colliding clouds test at $t = 5\tau_n = 6.29 \times 10^{-2}$ yr are presented in Figure 6. The shock fronts in the neutral fluid continue to travel away from the origin at $3.34$ km s$^{-1}$, and are located at $|x_{nshk}| = 6.63 \times 10^{14}$ cm. At this stage of the evolution, pronounced precursors in the ion velocity and the magnetic field extend to much greater distances beyond the neutral shock fronts. At the time shown ($\gg \tau_n$), collisional drag from the inflowing neutrals has a pronounced effect on the precursors. Notably, they continue to be led by discontinuities in the velocity and magnetic field heading outward at the ion magnetosound speed $V_{ims}$, but the jumps at the discontinuities are significantly diminished in magnitude and strength compared to the earlier time (Figure 5). The RC07 solutions are also plotted in Figure 6. Once again the split-operator results agree extremely well with the analytic solutions.

The solutions at time $t = 50\tau_n = 0.629$ yr are shown in Figure 7 along with the RC07 solutions. The agreement between the split-operator results and the analytic solutions is again very good. At this time ($\gg \tau_n$) the ion and magnetic field discontinuities have virtually vanished, with the precursors ahead of the neutral shocks having profiles that are smooth and continuous. Inside the precursors the motion of the ions is dictated by a balance between the collisional drag from the neutrals flowing toward the origin and the outwardly directed magnetic-pressure gradient. This is consistent with the RC07 solution, which found that on these timescales the solution for the charged fluid motion tended toward a force-free diffusion mode.

Figure 7 also contains insets showing the fluid velocities and magnetic field on an expanded scale which resolves the region between the neutral J-shocks. The vertical dashed lines in the insets mark the location of the forward- and reverse shocks at that time. The RC07 solutions agree well with the split-operator numerical simulation, even on these greatly magnified scales, with the numerical solution for the magnetic field differing from the RC07 solution by less than 0.3% within the inset region. This is an
Figure 6. Colliding model clouds at time $t = 5\tau_{\text{in}} = 6.29 \times 10^{-2}$ yr. All normalizations and symbols have the same meaning as in Figure 5.

Figure 7. Colliding model clouds at time $t = 50\tau_{\text{in}} = 0.629$ yr. All normalizations and symbols have the same meaning as in Figure 5. The insets show values about the collision point at $x = 0$, including the region between the leftward and rightward traveling neutral shock fronts, each indicated by the vertical dashed lines.
especially interesting result: to be able to use the method of Fourier transforms to obtain their analytic solutions, RC07 were forced to assume that the neutral gas density is uniform—even in the region between the two neutral shock fronts—despite the fact that for strong (i.e., large Mach-number) adiabatic shocks the density behind a shock front increases by a factor of four (for see, e.g., Zeldovich & Raizer1966 or Spitzer1978). RC07 argued on physical grounds that their analytic solution would nevertheless be correct. The colliding clouds test results and their agreement with the analytic solutions of RC07 show that our operator-splitting scheme faithfully reproduces the physics of dynamic MHD multifluid astrophysical flows. The split-operator code successfully models the evolution of shocks and discontinuities in both the neutral gas and the charged fluid, having no difficulty in handling the interaction between the fluids which leads to the formation of a magnetic precursor.

3.3. Shocks with Mass Transfer

As a final test we present a model which allows for the transfer of mass between the neutral and charged fluids in shocks. We introduce a nonzero source term for the conversion of ion–electron mass to neutral mass, with

\[
S_n = m_i(\alpha_{\text{DR}} n_i^2 - \zeta_{\text{CR}} n_n),
\]

(55)

where \(\zeta_{\text{CR}}\) is the cosmic-ray ionization rate and \(\alpha_{\text{DR}}\) is the rate coefficient for dissociative recombination of molecular HCO\(^+\) with electrons; note that in this expression we have used \(n_e = n_i\), which follows from (7). In this test we use the representative ionization rate \(\zeta_{\text{CR}} = 5 \times 10^{-17} \text{ s}^{-1}\) (Dalgarno 2006) and take \(\alpha_{\text{DR}}\) from the UMIST Astrochemistry Database (http://www.udfa.net; Millar et al. 1997),

\[
\alpha_{\text{DR}} = 2.4 \times 10^{-7} \left(\frac{300 \text{ K}}{T_e}\right)^{0.69} \text{ cm}^3 \text{s}^{-1}.
\]

(56)

Mass exchange also adds to the momentum source term, so that

\[
F_n = F_{n,\text{el}} + F_{n,\text{inel}},
\]

(57)

where

\[
F_{n,\text{inel}} = m_i(\alpha_{\text{DR}} n_i^2 v_i - \zeta_{\text{CR}} n_n v_n).
\]

(58)

In this test we again take \(G_i = \Lambda_n = 0\). Neglecting radiative cooling means that the temperatures in this test will be much larger than a realistic model with cooling included. However, it allows us to isolate the effects of mass-transfer between the neutrals and the ions as a test of the split-operator code. The rate coefficient for dissociative recombination depends on the electron temperature, which is not calculated in the current version of our code. For expediency we set

\[
T_e = \max(T_n, 0.15 T_i),
\]

(59)

in rough agreement with realistic simulations of steady multifluid shocks (e.g., see Figures 1–3 of Draine et al. 1983).

Initially this model has a discontinuity at \(x_0 = 1.12 \times 10^{16} \text{ cm}\). For \(x < x_0\) the neutral gas and ions are uniform and flowing in the \(+x\)-direction with a common supersonic velocity, while for \(x > x_0\) the matter is uniform and stationary. The other initial conditions are listed in Table 1. The initial ion density was calculated from the expression

\[
n_i = \left(\frac{\zeta_{\text{CR}} n_n}{\alpha_{\text{DR}}}\right)^{1/2},
\]

(60)

which assumes that the creation and destruction rates of ions in Equation (55) are equal.

Figure 8 presents the mass transfer (MT) test solution at \(t = 632 \text{ yr}\). For contrast we also show the results for a model with no mass transfer (NMT) which is identical otherwise. At the relatively early time shown in the figure (\(\ll \tau_{\text{el}} \sim 3 \times 10^4 \text{ yr}\)) the neutral gas is virtually unaffected by ion–electron drag. The effect of mass transfer on the neutral gas is also imperceptible in the MT model for the following two reasons: (1) even if all of the ions and electrons were to recombine, the added mass to the neutrals would be negligible because \(\rho_i \ll 10^{-6} \rho_n\); (2) the rate of change of the neutral density caused by cosmic-ray ionization is orders of magnitude smaller than the density change caused by advection.

In Figure 8(a) it is evident that the collision of the inflowing and stationary material at \(x_0\) has resulted in two neutral J-shocks: a left shock located at \(x_{L,\text{shk}} = 2.60 \times 10^{16} \text{ cm}\) traveling with a velocity \(v_{L,\text{shk}} = +7.41 \text{ km s}^{-1}\), plus a right shock at \(x_{R,\text{shk}} = 3.93 \times 10^{16} \text{ cm}\) with velocity \(v_{R,\text{shk}} = +14.1 \text{ km s}^{-1}\). The figure also reveals a contact discontinuity in the neutral gas at \(x_{n,\text{con}} = 3.22 \times 10^{16} \text{ cm}\). The neutral density profile is plotted for the MT model; the neutral density in the NMT is indistinguishable from that of the MT model.

| Table 1 | Initial Conditions for the Mass Transfer Test |
|---------|-----------------------------------------------|
| \(x < x_0\) | \(20 \text{ km s}^{-1}\) | \(2.5 \times 10^4 \text{ cm}^3\) | \(8.12 \times 10^4 \text{ cm}^3\) | \(15 \text{ K}\) | \(50 \text{ \mu G}\) | \(765 \text{ km s}^{-1}\) |
| \(x > x_0\) | \(0\) | \(2 \times 10^4 \text{ cm}^3\) | \(6.31 \times 10^4 \text{ cm}^3\) | \(10 \text{ K}\) | \(25 \text{ \mu G}\) | \(434 \text{ km s}^{-1}\) |
The effect of mass transfer on the ion fluid in the MT model can be seen in Figure 8(b). Between the shock fronts the neutral gas is heated, with a corresponding increase in the temperatures of the ions and electrons. This results in a decrease in the number of ion–electron recombinations there (see Equation (56)). At the same time, shock compression of the neutral gas leads to a greater number of cosmic-ray created ions (see Equation (55)). The interplay of these two effects leads to ion densities between the shock fronts that are a factor $\approx 5$ greater compared to the NMT model. Not only does the NMT model have a significantly smaller ion density in the shocked region, it also has an ion contact discontinuity located at $x_{i,\text{con}} = 3.32 \times 10^{16}$ cm. There is no ion contact discontinuity in the MT model because the source integration steps in the algorithm completely overwhelm the RG step (Figure 1). That is, the ion density profile is determined almost entirely by chemistry rather than advection. There is a small jump in the ion density at the neutral contact discontinuity caused by the jump in the neutral density, which is proportional to the rate of ionizations per unit volume. Increased $T_i$ and $T_e$ (due to significant drift speeds between the ions and the neutrals (see Equations (8) and (59)) inside the magnetic precursors upstream from the shock fronts reduces the rate of ion–electron recombinations there. This raises the ion density inside the precursors to values noticeably larger than those in the NMT model.

The width of the magnetic precursors depends on the amount of ion mass loaded onto the magnetic field lines there, as a comparison of the MT and NMT models shows (Figure 8(c)). The width of a magnetic precursor scales as $L_{\text{pre}} \sim V_{\text{ims}}^2 \tau_{\text{m}}/v_{\text{shk}}$, where $v_{\text{shk}}$ is the shock speed; this result can be derived by balancing magnetic pressure and ion–neutral drag in the precursor (Draine 1980) or by

---

**Figure 8.** Mass transfer test at $t = 632$ yr. In each panel the initial state is shown as black dashed lines. Also displayed are the locations of a neutral contact discontinuity (vertical line with squares) in the MT model and an ion contact discontinuity (vertical line with triangles) in the NMT model. (a) Neutral density. (b) Ion density (curve with circles) in the MT model and NMT model (curve with diamonds). The quasi-mass-equilibrium relation (60) is the dash-dotted curve. (c) Velocities of the neutrals (solid curve) and ions (curve with circles) in the MT model, and the ions in the NMT model (curve with diamonds). (d) Magnetic field in the MT model (curve with circles) and the NMT model (curve with diamonds).
setting the precursor crossing time \( t_{\text{pre}} = L_{\text{pre}}/v_{\text{shock}} \) equal to the ion-magnetic field diffusion time ahead of the front \( t_{\text{diff}} = L_{\text{pre}}^{2}/D_{\text{ad}} \) (see Equation (49d)). Because \( V_{\text{shock}} \) equals \( v_{\text{A}} \) in our models, it follows from Equation (31) that \( L_{\text{pre}} \propto n_{i}^{-1} \). Hence, the model with the greater ion density \( n_{i} \) will have the smaller values of \( L_{\text{pre}} \). This explains why the precursors extend further in the NMT model than in the MT model.

Figure 8(d) reveals a consequence of the dependence of precursor width on ion density: the distribution of magnetic flux is noticeably different in the MT and NMT models. Compression of the magnetic field by a shock is made less abrupt by having a precursor of greater width. This accounts for the smaller magnetic field increase at the left shock in the NMT model than in the MT model. (Note, however, that magnetic flux is conserved in both models: the total magnetic flux contained in the region from \( 1.5 \times 10^{16} \) cm to \( 4.5 \times 10^{16} \) cm [\( \approx \) area under the curve] is the same for both.) A by-product of the mass transfer test, then, is that we have demonstrated how loading of mass onto magnetic field lines affects the structure of a magnetic precursor and the magnetic field profile elsewhere in a shock. These effects feed back into the dynamics, because the magnetic pressure gradient is one of the main driving forces in the ion–electron momentum transport equation (5).

The fidelity of the mass-transfer test can be quantified by noting that from the left-hand side of the ion mass equation (4) we can define an ion advection time

\[
\tau_{\text{adv}} \equiv \frac{L}{v_{i}} = 317 \left( \frac{L}{10^{10} \text{ cm}} \right) \left( \frac{10 \text{ km s}^{-1}}{v_{i}} \right) \text{ yr} \tag{61a}
\]

where \( L \) is a characteristic length scale. From the rhs of the same equation we can also define a recombination timescale

\[
\tau_{\text{rec}} \equiv \frac{n_{i}}{m_{i} \alpha_{\text{rec}} n_{i}^{2}} = 66.0 \left( \frac{2 \times 10^{-3} \text{ cm}^{-3}}{n_{i}} \right) \left( \frac{T_{i}}{300 \text{ K}} \right)^{0.69} \text{ yr.} \tag{61b}
\]

and a cosmic-ray ionization time

\[
\tau_{\text{CR}} \equiv \frac{\rho_{i}}{m_{i} \zeta_{\text{CR}} n_{i}} = 63.4 \left( \frac{n_{i}}{2 \times 10^{-3} \text{ cm}^{-3}} \right) \left( \frac{2 \times 10^{4} \text{ cm}^{-3}}{n_{n}} \right) \left( \frac{5 \times 10^{-17} \text{ s}^{-1}}{\zeta_{\text{CR}}} \right) \text{ yr} \tag{61c}
\]

(see Equation (55)). For the physical conditions in the MT model, \( \tau_{\text{adv}} \gtrsim \tau_{\text{rec}}, \tau_{\text{CR}} \) on length scales \( L_{\text{eq}} \gtrsim 5 \times 10^{15} \) cm. For model ages greater than \( \tau_{\text{ref}} \) and \( \tau_{\text{CR}} \) and length scales greater than \( L_{\text{eq}} \), ion advection can be ignored, and there should be approximate equality between the rates of mass creation and destruction. That is, quasi-mass-equilibrium with \( S_{\text{n}} \approx 0 \) should occur with \( n_{i} \) given by Equation (60). It is therefore expected that the ion density in the MT run for the time displayed should be set by the condition of quasi-mass-equilibrium in regions having length scales \( \gtrsim L_{\text{eq}} \).

To test this hypothesis, the values of \( n_{i} \) predicted by the quasi-equilibrium equation (60) are also plotted in Figure 8(b) (dash-dotted curve). It is seen that the predicted values do match the actual MT model results well in most places, except at the shock fronts where the assumption of large length scales becomes invalid. Away from the shock fronts the agreement between the MT model and Equation (60) is generally quite good. For instance in the region between the shocks, including the region about the neutral contact discontinuity, the relative differences of the quasi-equilibrium and model values for \( n_{i} \) range from \( 3 \times 10^{-3} \) to 0.01. The very good agreement between the values predicted by Equation (60) for the ion density with the actual results of the fully dynamical MT model (in the regions where the mass-quasi-equilibrium approximation is valid) illustrates the accuracy of the split-operator scheme for multifluid shocks with mass transfer by ionization and recombination.

4. SUMMARY

Because many protostellar outflow shocks are much younger than the timescale \( t_{\text{ref}} \) to accelerate the neutral gas, it is likely they are not steady flows. A time-dependent treatment is therefore generally required to model outflow shocks and their emission. In this paper we have presented a method for modeling perpendicular, time-dependent, multifluid MHD shocks using operator splitting. This scheme splits the time integration of the evolutionary equations into separate steps, one dealing with the solution of independent homogeneous Riemann problems for the neutral and charged fluids using Godunov’s method, plus other steps where only the equations describing interactions between the fluids are integrated. Our method exploits the fact that the thermal pressure of the charged fluid is usually negligible compared to its magnetic pressure. Neglecting the thermal pressure reduces the number of characteristic waves in the ion–electron fluid to three, the same number as in one-dimensional gas dynamics (Toro 2009). We have shown that under these circumstances the MHD Riemann problem can be solved exactly. Using this exact solution, we have constructed an approximate MHD solver which is used in the RG integration of the charged fluid in our split-operator scheme. The similarity in structure of the Riemann problems for both fluids in our split-operator scheme makes it straightforward to adapt well-established gas-dynamic numerical techniques (TVD slope limiters, other data reconstruction methods such as WENO, etc.) for use in the RG integration of the charged fluid. The symmetry in the RG problems for both the neutral and charged fluids also greatly simplifies the numerical coding. (In fact, following our scheme it should not be too difficult to modify an already existing non-magnetic single-fluid RG hydrodynamics code to one for use in multifluid MHD.)

Since the inertia of a fluid and the wave modes it supports are fundamental to the solution of the Riemann problem, the multifluid split-operator method outlined here has no difficulty dealing with flows involving MHD waves or transients in the charged fluid. Several tests spanning a wide variety of time- and length scales were performed to demonstrate the versatility and accuracy of our algorithm. The numerical results are in very good agreement with analytic solutions in all of our benchmark tests. The latter include a model for the formation of MHD shocks resulting from the collision of two identical clouds, a problem solved analytically by RC07; the split-operator code successfully captures the ion density and magnetic field transients that propagate away from the collision.
surface at early times. Our code also correctly reproduces the multifluid shocks with magnetic precursors which develop at much later times, when the charged fluid is force free and evolves diffusively.

Another benchmark test involved MHD shocks with transfer of mass between the neutral and charged fluids. A numerical model with mass transfer has significantly greater ion densities behind the shocks and in the magnetic precursor compared to a model with no mass transfer. Analysis of the characteristic timescales for cosmic-ray ionization and ion–electron dissociative recombination in the mass transfer model suggested that for sufficiently large length scales the ion density should be close to the value that would be predicted when ionization and recombination balance one another. The ion density profile in the mass transfer model is indeed well fit by the quasi-equilibrium relation, except in the vicinity of the shock fronts where the criterion of large length scales (or, equivalently, negligible ion mass advection) is violated.

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APPENDIX A
THE EXACT MHD RIEMANN SOLUTION

The three-dimensional MHD Riemann problem for an arbitrary flow geometry and a plasma with nonzero thermal pressure is probably too complex to be solved exactly (Brio & Wu 1988; Torrilhon 2003). Because of this, several approximate MHD Riemann solvers have been developed (e.g., Dai & Woodward 1994, 1997; Ryu & Jones 1995; Balsara 1998; Li 2005; Mignone 2007). However, for the special case studied here—one-dimensional flow, perpendicular geometry, and no thermal pressure—the MHD Riemann problem for the charged fluid has a simple, exact solution. We give it here.

A.1. Characteristics

The Riemann problem for the charged fluid is to solve Equation (28),

$$\frac{\partial U_i}{\partial t} + \frac{\partial \mathbf{F}_i(U_i)}{\partial x} = 0,$$

subject to Riemann initial conditions,

$$U_i(x, 0) = \begin{cases} U^L_i \equiv \begin{bmatrix} \rho^L_i, \rho^L_i v^L_i, B^L \\ \rho^L_i v^L_i, B^L \\
\end{bmatrix}^T & \text{if } x < 0 \\ U^R_i \equiv \begin{bmatrix} \rho^R_i, \rho^R_i v^R_i, B^R \\ \rho^R_i v^R_i, B^R \\
\end{bmatrix}^T & \text{if } x > 0 \end{cases},$$

where the constant vectors $U^R_i$ and $U^L_i$ are the initial states on the right and left half planes. The first step is to rewrite Equation (A1) in the equivalent form

$$\frac{\partial U_i}{\partial t} + \mathbf{A} \cdot \frac{\partial U_i}{\partial x} = 0$$

and examine the eigenvalues and eigenvectors of the Jacobian matrix,

$$\mathbf{A}_{k\ell} \equiv \frac{\partial F_{i\ell}}{\partial U_{ik}} = \begin{bmatrix} 0 & 1 & 0 \\ -v_i^2 & 2v_i & B/4\pi \\ -Bv_i/\rho_i & B/\rho_i & v_i \end{bmatrix}.$$

The eigenvalues are

$$\{\lambda_- , \lambda_0 , \lambda_+ \} = \{v_i - V_{IA}, v_i, v_i + V_{IA}\},$$

where $V_{IA}$ is the ion Alfvén speed (Equation (31)). The corresponding right (column) eigenvectors are

$$R_- = [\rho_i, \rho_i (v_i - V_{IA}), B]^T,$$

$$R_0 = [\rho_i, \rho_i v_i, 0]^T,$$

and

$$R_+ = [\rho_i, \rho_i (v_i + V_{IA}), B]^T.$$

The solution also depends on the left (row) eigenvectors, which are

$$L_- = \frac{1}{2v_{IA}} \begin{bmatrix} v_i \\ -\frac{1}{\rho_i} \rho_i V_{IA} \\ B \end{bmatrix}.$$
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![Figure 9. Three characteristics emanating from the initial discontinuity.](image)

Three characteristics emanating from the initial discontinuity.

\[
L_0 = \left[ \begin{array}{c} \frac{1}{\rho_i}, 0, -\frac{1}{B} \end{array} \right].
\]  
(A7b)

and

\[
L_+ = \frac{1}{2v_{iA}} \left[ -\frac{v_i}{\rho_i}, \frac{1}{\rho_i}, \frac{V_{iA}}{B} \right].
\]  
(A7c)

The left and right eigenvectors are orthonormal.

Introduce the characteristics curves, \(x_m(t)\), defined by

\[
\frac{dx_m}{dt} = \lambda_m, \quad m = -, 0, +.
\]  
(A8)

For Riemann initial conditions, the characteristics emanating from the initial discontinuity at \(x = 0\) are straight lines (Figure 9).

Each represents a discontinuity in the solution which may be a shock wave, rarefaction, or contact discontinuity. Since discontinuities propagate along characteristics, and the initial data are uniform to the left and right of the origin, Figure 9 shows immediately that \(U_i\) to the left of \(\lambda_-\) and \(U_i\) to the right of \(\lambda_+\). The problem therefore reduces to finding \(U_i\) in the “star region” between \(\lambda_-\) and \(\lambda_+\). Let \(\rho_i^L, v_i^L, B_i^L\) denote the three unknowns in the star region between \(\lambda_-\) and \(\lambda_0\) and \(\rho_i^R, v_i^R, B_i^R\) be the three unknowns between \(\lambda_0\) and \(\lambda_+\). These are found by (1) classifying each characteristic as a shock, rarefaction, or contact discontinuity, and (2) joining the solutions in such a way that the Rankine–Hugoniot (RH) jump conditions are satisfied across shocks and the generalized Riemann invariants are conserved across rarefactions and contact discontinuities. For a detailed discussion of the underlying theory, see the monograph by Toro (2009).

A.2. The \(\lambda_0\) Characteristic

The classification of \(\lambda_0\) depends on the “eigenvalue gradient,”

\[
\nabla \lambda_0 \equiv \begin{bmatrix} \frac{\partial \lambda_0}{\partial U_{i1}}, \frac{\partial \lambda_0}{\partial U_{i2}}, \frac{\partial \lambda_0}{\partial U_{i3}} \end{bmatrix}.
\]  
(A9)

It is easy to show that

\[
\nabla \lambda_0 = \frac{1}{\rho_i} \begin{bmatrix} -v_i, 1, 0 \end{bmatrix}
\]  
(A10)

and hence that

\[
\nabla \lambda_0 \cdot R_0 = 0.
\]  
(A11)

According to the theory of hyperbolic PDEs, Equation (A11) establishes that \(\lambda_0\) is always a contact discontinuity. Conservation of the Riemann invariants across the \(m\)th characteristic implies that

\[
\frac{dU_1}{R_{m1}} = \frac{dU_2}{R_{m2}} = \frac{dU_3}{R_{m3}}
\]  
(A12)

(see Toro 2009, Section 2.4.4) and writing this out for \(m = 0\) gives

\[
\frac{d\rho_i}{\rho_i} = \frac{d(\rho_i v_i)}{\rho_i v_i} = \frac{dB}{0}.
\]  
(A13)

Integrating Equation (A13) across the contact discontinuity yields

\[
v_i^L = v_i^R \equiv v_i^*
\]  
(A14)

and

\[
B_i^L = B_i^R \equiv B^*.
\]  
(A15)

Thus the ion density may undergo a jump across the contact discontinuity but the other fluid variables are continuous, in precise analogy with contact discontinuities in gas dynamics. Note that the number of unknowns in the star region has now been reduced to four: \(v_i^*, B^*, \rho_i^{L4}, \) and \(\rho_i^{R4}.

---

1 If \(B \neq 0\) the eigenvalues of \(A\) are real and nondegenerate. This guarantees that Equation (A3) is strictly hyperbolic (e.g., Jeffrey 1976).
A.3. The $\lambda_+ \text{ and } \lambda_- \text{ Characteristics}$

The eigenvalue gradients for the other characteristics are

$$\nabla \lambda_{\pm} = \frac{1}{\rho_i} \left[ -\frac{1}{\rho_i} (v_i \mp V_{iA}/2), \frac{1}{\rho_i} \pm \frac{1}{\sqrt{8\pi \rho_i}} \right],$$  \hspace{1cm} (A16)

from which it follows that

$$\nabla \lambda_{\pm} \cdot R_{\pm} = \mp \frac{3}{2} V_{iA}. \hspace{1cm} (A17)$$

It seems reasonable to assume on physical grounds that $B \neq 0$ everywhere for $t > 0$ (no vacuum) so that the rhs of expression (A17) is always nonzero. Then the theory of hyperbolic PDEs establishes that $\lambda_- \text{ and } \lambda_+$ are never contact discontinuities. The $\lambda_-$ wave is a shock wave if the pressure in the star region exceeds the pressure in the left region ($B^* > B^L$) and a rarefaction wave otherwise. Similarly, $\lambda_+$ is a shock if and only if $B^* > B^R$. These conclusions follow from the fact that shocks are compressive and rarefactions are not, plus our assumption that the pressure of the plasma is entirely magnetic. For a given set of initial conditions, only one combination of shocks and/or rarefactions at $\lambda_- \text{ and } \lambda_+$ gives a solution that satisfies the matching conditions across all three discontinuities. In Appendix A.4, we give a simple criterion for identifying the correct combination.

The matching conditions for shock waves are the RH jump conditions, which require $\mathcal{F}_i$ to be conserved across the shock front in a frame comoving with the shock. We omit details of the derivation and simply give the results. For a shock at $\lambda_-$ (a “left shock”), we find that the ion velocities on opposite sides of the shock front are related by

$$v_i^* = v_i^L + f_{LS} (B^*), \hspace{1cm} (A18)$$

where

$$f_{LS} (B^*) \equiv - \left[ \frac{(P^* - P^L)}{\rho_i} \left( 1 - \frac{B^L}{B^*} \right) \right]^{1/2}. \hspace{1cm} (A19)$$

and $P(B) \equiv B^2/8\pi$ is the magnetic pressure. Similarly, for a right shock we find

$$v_i^* = v_i^R + f_{RS} (B^*), \hspace{1cm} (A20)$$

where

$$f_{RS} (B^*) \equiv + \left[ \frac{(P^* - P^R)}{\rho_i} \left( 1 - \frac{B^R}{B^*} \right) \right]^{1/2}. \hspace{1cm} (A21)$$

The matching conditions for rarefactions are governed by the Riemann invariants. Carrying out steps analogous to the derivation of Equation (A13), we find that

$$\frac{d\rho_i}{\rho_i} = \frac{d (\rho_i v_i)}{\rho_i (v_i \mp V_{iA})} = \frac{dB}{B}, \hspace{1cm} (A22)$$

where the upper sign corresponds to $\lambda_+$. Equating the first and third terms simply gives flux freezing in the charged fluid. This implies that the magnetic field and hence $V_{iA}$ may be viewed as functions of the ion density alone. Knowing this, we equate the first two terms in Equation (A22) to obtain a differential equation for $v_i$:

$$dv_i = \pm V_{iA} (\rho_i) \frac{d\rho_i}{\rho_i}. \hspace{1cm} (A23)$$

Integrating Equation (A23) across $\lambda_-$ gives the matching condition for a left rarefaction:

$$v_i^* = v_i^L + f_{LR} (B^*), \hspace{1cm} (A24)$$

where

$$f_{LR} (B^*) \equiv +2 V_{iA}^L \left[ 1 - \left( \frac{B^*}{B^L} \right)^{1/2} \right]. \hspace{1cm} (A25)$$

and $V_{iA}^L$ is the ion Alfvén speed in the left region. For a right rarefaction one finds similarly that

$$v_i^* = v_i^R + f_{RR} (B^*), \hspace{1cm} (A26)$$

where

$$f_{RR} (B^*) \equiv -2 V_{iA}^R \left[ 1 - \left( \frac{B^*}{B^R} \right)^{1/2} \right]. \hspace{1cm} (A27)$$
the plasma, and this determines the density in both halves of the star region:

\( f \)

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configuration depends only on three dimensionless parameters, \( \lambda \)

shock plus a right rarefaction, and so on. If the configuration was known a priori, one could write the matching conditions across the

all three characteristics, an exercise in the process of elimination. For example, suppose that the flow is of type RR, which requires

by choosing \( f_L \) and \( f_R \) appropriately from the functions \( f_{LS}, f_{LR}, f_{RR} \), etc. Subtracting Equation (A29) from (A28) gives

\[ v_L^1 - v_R^1 = f_L(B^*) - f_R(B^*) = 0, \]

which is a nonlinear equation for \( B^* \). Once \( B^* \) has been determined by solving Equation (A30), the velocity in the star region follows from Equation (A28) or Equation (A29). Finally, the matching conditions for shocks and rarefactions both require flux freezing in the plasma, and this determines the density in both halves of the star region:

\[ \rho^*_L = \rho^*_L \left( \frac{B^*}{B^*_L} \right) \]  

and

\[ \rho^*_R = \rho^*_R \left( \frac{B^*}{B^*_R} \right). \]

To implement the steps above it remains only to identify the unique flow configuration that satisfies the matching conditions across all three characteristics, an exercise in the process of elimination. For example, suppose that the flow is of type RR, which requires that \( B^* < B_L^* \) and \( B^* < B_R^* \), for the case \( v_L^1 - v_R^1 > 0 \) (the flows collide). Equation (A30) has a solution only if \( f_L - f_R < 0 \); but examination of Equations (A25) and (A27) shows that \( f_{LR} - f_{RR} \) is strictly positive if \( B^* < B_L^* \) and \( B^* < B_R^* \). We conclude then that the RR configuration never occurs when \( v_L^1 - v_R^1 > 0 \). Similarly, consider a flow of type SS, which has \( B^* > B_L^* \) and \( B^* > B_R^* \), for the situation \( v_L^1 - v_R^1 < 0 \) (the flows diverge). For that situation, according to Equation (A30), there can only be a solution if \( f_L - f_R > 0 \). Equations (A19) and (A21) reveal that, for \( B^* > B_L^* \) and \( B^* > B_R^* \), \( f_L - f_R \) is always <0. We conclude then that the SS configuration cannot occur when \( v_L^1 - v_R^1 < 0 \). Straightforward but lengthy analysis of the other possibilities shows that the configuration depends only on three dimensionless parameters,

\[ \xi \equiv \frac{v_L^1 - v_R^1}{V_{1A}}, \]  

\[ \theta \equiv B_R^*/B_L^*, \]  

and

\[ \delta \equiv \sqrt{\rho^*_L/\rho^*_R}. \]

The flow classification is given in Table 2, where

\[ \Gamma \equiv \frac{1}{\delta \sqrt{2}} \left[ (1 - \theta^2)(1 - \xi^2) \right]^{1/2} - \xi, \]

Note. * The RR-state can exist only if \(-2(1 + \theta/\delta)V_{1A} < v_L^1 - v_R^1 < 0 \) (see Equation (A35)).

### A.4. Flow Configuration and Solution

Let RR, SR, RS, and SS denote the four possible flow configurations, where “RR” is a flow with two rarefactions, “SR” a left shock plus a right rarefaction, and so on. If the configuration was known a priori, one could write the matching conditions across the \( \lambda_- \) wave,

\[ v_i^* = v_{iL}^1 + f_L(B^*), \]

and the \( \lambda_+ \) wave,

\[ v_i^* = v_{iR}^1 + f_R(B^*), \]

Table 2

| Initial Conditions | Flow Type | Note |
|--------------------|-----------|------|
| \( v_L^1 - v_R^1 \) > 0: | \( B_L^* < B_R^*, \Psi^* < 0 \) | SS | left shock + right shock |
| \( B_L^* < B_R^*, \Psi^* > 0 \) | SR | left shock + right rarefaction |
| \( B_L^* > B_R^*, \Gamma > 0 \) | SS | left shock + right shock |
| \( B_L^* > B_R^*, \Gamma > 0 \) | RS | left rarefaction + right shock |
| \( v_L^1 - v_R^1 \) = 0: | \( B_L^* < B_R^*, \Psi^* > 0 \) | SR | left shock + right rarefaction |
| \( B_L^* > B_R^*, \Gamma > 0 \) | RS | left rarefaction + right shock |
| \( v_L^1 - v_R^1 \) < 0: | \( B_L^* < B_R^*, \Phi^* < 0 \) | RR | left rarefaction + right rarefaction |
| \( B_L^* < B_R^*, \Phi^* > 0 \) | SR | left shock + right rarefaction |
| \( B_L^* > B_R^*, \Phi^* < 0 \) | RR | left rarefaction + right rarefaction |
| \( B_L^* > B_R^*, \Phi^* > 0 \) | RS | left rarefaction + right shock |

Initial Conditions Flow Type Note
Figure 10. First MHD Riemann demonstration problem. Dashed lines indicate the initial state of the charged fluid and magnetic field. The solid line is the exact MHD Riemann solution at $t = 0.1$ yr. Circles are the solution calculated using the approximate MHD Riemann solver (some data points have been omitted for clarity). Top panel: magnetic field. Middle: ion number density. Bottom: ion velocity.

\[
\Psi \equiv \frac{1}{\sqrt{2}} [(\theta^2 - 1)(1 - \theta^{-1})]^{1/2} - \xi, \quad (A34b)
\]

\[
\Phi \equiv \frac{2}{\delta} (\theta - \sqrt{\theta}) + \xi, \quad (A34c)
\]

\[
\Upsilon \equiv 2(1 - \sqrt{\theta}) + \xi. \quad (A34d)
\]

Once the flow type is known, which of the expressions (A19), (A21), (A25), and (A27) should be used as the functions $f_L(B^*)$ and $f_R(B^*)$ in Equation (A30) is then also known. The value of $B^*$ can then be determined to any desired level of accuracy from Equation (A30) using an iterative numerical technique such as Newton’s method (Atkinson 1989; Press et al. 1996). We note that for flows with both left and right rarefaction waves (type RR), the ion mass densities in the star region $\rho^L_i$ and $\rho^R_i$ must be $>0$; it follows from Equations (A31) and (A32) that the magnetic field in the star region $B^*$ must also be $>0$. This non-vacuum (or, positivity) condition imposes a lower limit on the value of $\xi$, the dimensionless velocity difference between the left and right states of the initial discontinuity (A33a). Using Equations (A25), (A27), and (A30), one finds that this condition is satisfied so long as $\xi > \xi_{\text{vac}}$, where

\[
\xi_{\text{vac}} = -2 \left(1 + \frac{\theta}{\delta}\right). \quad (A35)
\]

As demonstrative examples, we consider two representative MHD Riemann test problems. In the first problem, the initial state of the plasma and magnetic field has $B^L = 50 \mu G$, $B^R = 25 \mu G$, $v^L_i = 200 \text{ km s}^{-1}$, $v^R_i = 0$, and $\rho^L_i = \rho^R_i = 2.51 \times 10^{-28} \text{ g cm}^{-3}$ (for a cloud with $m_i = 25 m_p$, this corresponds to a number density $n_i = 6 \times 10^{-4} \text{ cm}^{-3}$). For these parameters $B^L > B^R$, and $\Gamma = 0.321 > 0$. Examination of Table 2 indicates that the flow configuration for this first test will be of type RS, a left rarefaction with a right shock. That this is so can be seen in Figure 10 which shows the results for this problem. The initial state of the plasma and magnetic field are displayed as dashed lines, and the solid curves are the MHD Riemann solution at the time $t = 0.1$ yr. The right shock is located at $x_{\text{shk}} = +2.07 \times 10^{14} \text{ cm}$, and the head of the left rarefaction wave is at $x_{\text{rw}} = -2.49 \times 10^{14} \text{ cm}$. A contact discontinuity is
located at $x_c = +8.23 \times 10^{13}$ cm; that is the position the initial discontinuity in the charged fluid’s magnetic flux-to-mass ratio $B/\rho_i$ has moved to by the time shown.

In the second MHD Riemann test problem we have an initial state with $B^L = 45 \mu$G, $B^R = 50 \mu$G, $v^L_i = -200 \text{ km s}^{-1}$, $v^R_i = +200 \text{ km s}^{-1}$, and the same constant ion density as in the first test. This state has $B^L < B^R$, $v^L_i < v^R_i$, and $\Phi = -0.385 < 0$. According to Table 2 this should be an RR flow. This is indeed the case, as can be seen in Figure 11, which displays the exact MHD Riemann solution for this model at $t = 0.1$ yr. At the time shown, the head of the left rarefaction wave is at $x_{L, rw} = -3.14 \times 10^{14}$ cm, and the head of the right rarefaction wave is at $x_{R, rw} = +3.42 \times 10^{14}$ cm. There is also a charged fluid contact discontinuity at $x_c = -1.70 \times 10^{13}$ cm.

APPENDIX B
AN APPROXIMATE MHD RIEMANN SOLVER

B.1. The Star Region

The Riemann solution found in Appendix A can also be used to numerically solve the MHD flow problem of Equation (28) using Godunov’s method (Godunov 1959; Toro 2009). However, doing so can be computationally expensive, because it uses an iterative method to solve for $B^*$ from Equation (A30).

It is desirable then to instead use a more efficient approximate solver to calculate the variables in the star region ($B^*$, $v^*_i$, $\rho^*_i$, $\rho^*_R$, and $\rho^*_L$) between the characteristics $\lambda_-$ and $\lambda_+$ (see Figure 9). This can be done by assuming that the flow motion is approximately linear, that is, a wave, and then using the characteristic relations derived from the Riemann invariants (Equation (A22)) to connect the variables in the star region to those in the adjacent L or R region of the Riemann problem (see Figure 9). But this is just what was done in deriving Equations (A25) and (A27). Solving those two equations for $B^*$ and $v^*_i$ in terms of the L and R region variables yields

$$B^* = \left[ \frac{V^L_{iA} + V^R_{iA} + \frac{1}{2} (v^L_i - v^R_i)}{C^L_i + C^R_i} \right]^2,$$

$$v^*_i = \frac{C^L_i v^R_i + C^R_i v^L_i + 2C^L_i C^R_i (\sqrt{B^L} - \sqrt{B^R})}{C^L_i + C^R_i},$$

24
where
\[ C^L_i = \sqrt{\frac{V_{iA}^L}{B^L}}, \quad C^R_i = \sqrt{\frac{V_{iA}^R}{B^R}}. \] (B3)

With these quantities now known, \( \rho^L_i \) and \( \rho^R_i \) can be calculated directly from Equations (A31) and (A32), respectively.

### B.2. Calculation of Fluxes at the Discontinuity

To solve the MHD Riemann problem for a state having the discontinuous initial conditions (A2) using the method of Godunov (1959), we need to calculate the flux vector \( \mathbf{F}_i(0) \) at the interface separating the two states, \( \chi = 0 \) (or, equivalently, along the ray with the similarity variable \( x/t = 0 \), which corresponds to the \( t \)-axis in Figure 9). To calculate \( \mathbf{F}_i(0) \) from Equation (20), we need the array of variables \( U_i(0) \) (Equation (18)).

A very thorough discussion of how to obtain the state variables at the interface \( \chi = 0 \) for the one-dimensional gas dynamic Riemann problem using Godunov’s method is presented in chapter 6 of Toro (2009); the fundamental ideas described there—in which the RG flux is determined by examining the characteristic wave structure at the discontinuity—remain valid and carry over directly to the MHD Riemann problem that we are concerned with here. The first step in the assignment of \( U_i(0) \) is made by finding the values of the velocity \( v^*_i \) (\( = \lambda \delta \)) of the contact discontinuity and the magnetic field of the star region \( B^* \) from Equations (B1) and (B2).

For \( v^*_i > 0 \), the contact discontinuity is moving to the right. If \( B^* > B^R \) the left wave is a shock wave with velocity
\[ s^L_i = v^*_i - Q^L_i / \rho^L_i \] (B4)
in the stationary frame. This relation was obtained by applying the RH conditions in a frame moving with the shock front. The quantity
\[ Q^L_i = \frac{(B^L)^2/8\pi - (B^*)^2/8\pi}{v^*_i - v^L_i} \] (B5)
is the ion mass flux across the front. If \( s^L_i > 0 \), the shock is moving to the right and \( U_i(0) = U^L_i \). If \( s^L_i < 0 \), the shock is instead heading to the left, and \( U_i(0) = U^{\lambda L}_i \equiv [\rho^L_i, \rho^{\lambda L}_i v^*_i, B^\lambda]^T \).

If \( B^* > B^L \), the left wave will instead be a rarefaction. As is widely known (e.g., Courant & Friedrichs 1948; Landau & Lifshitz 1959; Zeldovich & Raizer 1966; Toro 2009), a rarefaction emanating from a discontinuity has a fan-like phase-space structure (a centered simple wave) with a head and a tail. The velocities of the head \( (h^L_i) \) and tail \( (t^L_i) \) of the left rarefaction fan are
\[ h^L_i = v^*_i - v^L_i, \quad t^L_i = v^*_i - v^{\lambda L}_i, \] (B6)
respectively, where
\[ v^{\lambda L}_i \equiv B^* \sqrt{\pi / \rho^L_i}. \] (B7)

For \( h^L_i > 0, t^L_i > 0 \), the entire rarefaction fan is heading to the right and \( U_i(0) = U^L_i \), while for \( h^L_i < 0, t^L_i < 0 \) the entire fan is traveling left and \( U_i(0) = U^{\lambda L}_i \). If \( h^L_i < 0 \) and \( t^L_i > 0 \) values inside the rarefaction fan are needed (e.g., see chapter 4 of Toro 2009). In that situation, \( U_i(0) = U^{\lambda \text{fan}}_i \equiv [\rho^{\lambda \text{fan}}_i, \rho^{\lambda \text{fan}}_i v^{\lambda \text{fan}}_i, B^{\lambda \text{fan}}]^T \). Using the Riemann invariant relation (A22) to integrate across the \( \lambda \)-characteristic to connect quantities in the left rarefaction fan to those in the left state, along with the relations (A5), (A8), and (A31), and setting \( x/t = 0 \) in the resulting expressions yields
\[ v^{\lambda \text{fan}}_i = \frac{2}{3} \left( V^L_i + \frac{1}{2} v^L_i \right), \quad B^{\lambda \text{fan}} = B^L \left( \frac{2}{3} + \frac{v^L_i}{3 V^L_i} \right)^2, \quad \rho^{\lambda \text{fan}}_i = \frac{\rho^L_i B^{\lambda \text{fan}}}{B^L}. \] (B8)

If \( v^*_i < 0 \), the contact discontinuity is traveling to the left. For \( B^* > B^R \) the right wave is a shock wave. The velocity of the shock in the stationary frame is
\[ s^R_i = v^R_i + Q^R_i / \rho^R_i, \] (B9)
where
\[ Q^R_i = \frac{(B^*)^2/8\pi - (B^R)^2/8\pi}{v^*_i - v^R_i} \] (B10)
is the ion mass flux across the shock front. If \( s^R_i > 0 \) the right shock is traveling to the right and \( U_i(0) = U^{*R}_i \equiv [\rho^{*R}_i, \rho^{*R}_i v^*_i, B^*]^T \); if \( s^R_i < 0 \) the shock is heading left, and \( U_i(0) = U^{\text{R}}_i \) instead.

For \( B^* < B^R \) the right wave will be a rarefaction. The right rarefaction fan has head \( (h^R_i) \) and tail \( (t^R_i) \) velocities
\[ h^R_i = v^R_i + V^R_{iA}, \quad t^R_i = v^*_i + V^R_{iA}, \] (B11)
where
\[ V^R_{iA} = \frac{B^*}{\sqrt{4\pi \rho^R_i}}. \] (B12)
For $h_i^R > 0$, $\gamma_i^R > 0$, the entire fan is rightward traveling and $U_i(0) = U_i^R$, while, for $h_i^R < 0$, $\gamma_i^R < 0$ the complete fan is heading to the left, and therefore $U_i(0) = U_i^L$. If $h_i^R < 0$ and $\gamma_i^R > 0$ values from inside the right rarefaction fan are required and $U_i(0) = U^{\text{Rfan}}_i \equiv \begin{bmatrix} \rho_i^{\text{Rfan}}, \rho_i^{\text{Rfan}}/\gamma_i^{\text{Rfan}}, B^{\text{Rfan}}_i \end{bmatrix}^T$. Integrating the Riemann invariant (A22) across $\lambda_+$ to link quantities in the tail of the right rarefaction fan to quantities in the right state, also using Equations (A5), (A8), and (A32), and taking $x/t = 0$ in the resulting relations, gives

$$v_i^{\text{Rfan}} = \frac{2}{3} \left( -\frac{V_i^{\Lambda}}{3} + \frac{1}{3} v_i^R \right), \quad B^{\text{Rfan}} = B^R \left( \frac{2}{3} - \frac{v_i^R}{3V_i^{\Lambda}} \right)^2, \quad \rho_i^{\text{Rfan}} = \frac{\rho_i^R B^{\text{Rfan}}}{B^R}. \quad (B13)$$

As we have stated earlier, approximate Riemann solvers for the one-dimensional neutral gas dynamic RG problem have a very similar structure to that of the MHD solver we present here. In several instances the algorithm for the MHD and neutral gas Riemann solvers are basically the same when a suitable substitution is made. For example, the relations we derived for quantities in the left and right rarefaction fans (Equations (B8) and (B13)) are identical to the left and right fan relations at $x/t = 0$ of the neutral gas Riemann solver. The zero-order value which simply expresses freezing of magnetic flux in the charged fluid. Inserting this result in Equation (C1) and using $n_i = \rho_i/m_i$, Equations (48b), (49b), and (50b) follow directly.

**APPENDIX C**

**ANALYTIC BENCHMARK SOLUTIONS**

We present here a derivation of the analytical small-amplitude wave solutions used to test the split-operator models discussed in Sections 3.1.1–3.1.3. For the governing fluid equations we linearize the system in the standard way, assuming for any physical quantity $f$ that

$$f(x, t) = f_0 + \delta f(x, t); \quad (C1)$$

the zero-order value $f_0$ is taken to be constant and uniform, and the first-order perturbation $|\delta f(x, t)| \ll f_0$. Inserting these perturbations into Equations (1)–(6) and retaining terms up to first order give the linearized system of equations:

$$\frac{\partial}{\partial t} (\delta \rho_n) = -\rho_0 \frac{\partial}{\partial x} (\delta v_n), \quad (C2)$$

$$\frac{\rho_0}{\tau_m} \frac{\partial}{\partial t} (\delta v_n) = -\frac{\partial}{\partial x} (\delta P_n) - \frac{\rho_0}{\tau_m} \frac{\partial}{\partial t} \delta v_n + \frac{\rho_0}{\tau_m} \frac{\partial}{\partial t} \delta v_i, \quad (C3)$$

$$\frac{1}{T_n} \frac{\partial}{\partial t} (\delta T_n) = (\gamma - 1) \frac{1}{\tau_m} \frac{\partial}{\partial t} (\delta \rho_n), \quad (C4)$$

$$\frac{\partial}{\partial t} (\delta \rho_n) = -\rho_0 \frac{\partial}{\partial x} (\delta v_i), \quad (C5)$$

$$\frac{\rho_0}{\tau_m} \frac{\partial}{\partial t} (\delta v_i) = -\frac{B_0}{4\pi} \frac{\partial}{\partial x} (\delta B) + \frac{\rho_0}{\tau_m} \frac{\partial}{\partial t} \delta v_n - \frac{\rho_0}{\tau_m} \frac{\partial}{\partial t} \delta v_i, \quad (C6)$$

$$\frac{\partial}{\partial t} (\delta B) = -\frac{B_0}{\tau_m} \frac{\partial}{\partial x} (\delta v_i), \quad (C7)$$

where we have used the fact that the charged and neutral fluids are at rest in the zero-order reference state ($v_{n0} = v_{i0} = 0$), and also that $F_{n,\text{inel}} = S_n = S_i = G_n = \Lambda_n = 0$ for all of the model tests in Section 3.1. Equation (13) was also used to substitute for $F_n$ in the perturbed force equations (C3) and (C6).

The internal energy for an ideal gas (11) having a ratio of specific heats $\gamma$ was used in the energy equation of the neutral gas (3) to derive (C4); integrating that latter equation and inserting the result into the linearized ideal gas law (9),

$$\frac{\delta P_n}{P_{n0}} = \frac{\rho_0}{\rho_{n0}} \frac{\delta \rho_n}{\rho_{n0}} + \frac{\delta T_n}{T_{n0}}, \quad (C8)$$

yields the adiabatic relation between density and pressure perturbations in the neutral fluid,

$$\delta P_n = \frac{\gamma P_{n0}}{\rho_{n0}} \delta \rho_n = C_n^2 \delta \rho_n, \quad (C9)$$

where $C_n$ is the adiabatic speed of sound (44).

Combining Equations (C5) and (C7) and integrating we immediately obtain the result

$$\frac{\delta \rho_i}{\delta B} = \frac{\rho_{n0}}{B_0}, \quad (C10)$$

which simply expresses freezing of magnetic flux in the charged fluid. Inserting this result in Equation (C1) and using $n_i = \rho_i/m_i$, Equations (48b), (49b), and (50b) follow directly.
C.1. Solution for $L_G \ll L_{\text{ims}}$

In this limit, the neutral fluid is unable to respond to the very short time- and length scales of the perturbations; therefore, for this situation the neutrals are effectively a fixed stationary background with $\delta v_n = \delta \rho_n = 0$, and the only equations relevant to these particular modes are (C6) and (C7).

The Fourier transform $\hat{f}$ of any function $f$ is given by

$$\hat{f}(k, t) = \frac{1}{\sqrt{2\pi}} \int_{-\infty}^{\infty} f(x, t)e^{-ikx}dx,$$

where $k$ is the wavenumber. The inverse transform is

$$f(x, t) = \frac{1}{\sqrt{2\pi}} \int_{-\infty}^{\infty} \hat{f}(k, t)e^{ikx}dk$$

Fourier transforming equations (C6) and (C7) gives a system of linear ordinary differential equations with constant coefficients:

$$\frac{\partial \hat{y}_i}{\partial t} = M_i \hat{y}_i, \quad \text{(C13)}$$

where the vector of transformed variables $\hat{y}_i(k, t) ≡ [\delta \hat{v}_i(k, t), \delta \hat{B}(k, t)]^T$ and

$$M_i = \begin{bmatrix} -1/\tau_{in} & -i B_0k/4\pi \rho_0 \\ -i B_0k & 0 \end{bmatrix}. \quad \text{(C14)}$$

The ODE system (C13) has the solution

$$\hat{y}_i = a_{i+} \mathcal{E}_{i+} e^{\xi_{i+} t} + a_{i-} \mathcal{E}_{i-} e^{\xi_{i-} t} \quad \text{(C15a)}$$

(e.g., see chap. 3 of Braun 1983), where the eigenvalues and eigenvectors of $M_i$ are, respectively,

$$\xi_{i\pm} = -\frac{1}{\tau_{in}} \pm i V_{\text{ims}}k \left[ 1 - \left( \frac{k_{\text{ims}}}{k} \right)^2 \right]^{1/2} \quad \text{(C15b)}$$

and

$$\mathcal{E}_{i\pm} ≡ \begin{bmatrix} 1, & -i \frac{B_0k}{g_{i\pm}} \end{bmatrix}^T. \quad \text{(C15c)}$$

In these expressions we have used Equation (31) and the fact that $V_{\text{ims}} = V_A$ for all conditions of interest in our model interstellar clouds. We also introduced the ion magnetosound cutoff wave number

$$k_{\text{ims}} ≡ \frac{1}{2V_{\text{ims}} \tau_{in}} = \frac{2\pi}{L_{\text{ims}}}; \quad \text{(C16)}$$

modes with $k > k_{\text{ims}}$ propagate as ion magnetosound waves.

The constants $a_{i+}$ and $a_{i-}$ in Equation (C15a) are determined by applying the initial conditions $\delta \hat{v}_i(k, 0) = 0$ (ions initially at rest) and the Fourier transform of the initial Gaussian pulse in the magnetic field (see Equation (46)),

$$\delta \hat{B}(k, 0) = \frac{B_0 A_p L_G}{\sqrt{2}} \exp \left( -\frac{L_G^2 k^2}{4} \right). \quad \text{(C17)}$$

Doing that yields

$$\delta \hat{B}(k, t) = \frac{B_0 A_p L_G}{2\sqrt{2}} \exp \left( -\frac{t}{2\tau_{in}} \right) \times \left\{ \begin{array}{l}
1 - \frac{ik_{\text{ims}}/k}{[1 - (k_{\text{ims}}/k)^2]^{1/2}} \exp \left( -\frac{L_G^2 k^2}{4} + i V_{\text{ims}} k t \left[ 1 - \left( \frac{k_{\text{ims}}}{k} \right)^2 \right]^{1/2} \right) \\
+ \frac{ik_{\text{ims}}/k}{[1 - (k_{\text{ims}}/k)^2]^{1/2}} \exp \left( -\frac{L_G^2 k^2}{4} - i V_{\text{ims}} k t \left[ 1 - \left( \frac{k_{\text{ims}}}{k} \right)^2 \right]^{1/2} \right) \end{array} \right\} \quad \text{(C18)}$$

27
\[
\delta \hat{v}_i(k, t) = -\frac{V_{ims} A_p L_G}{2\sqrt{2}} \exp \left( -\frac{t}{2\tau_{in}} \right) \left\{ \exp \left( -\frac{L_G^2 k^2}{4} + i V_{ims} t \left[ 1 - \left( \frac{k}{k_{ims}} \right)^2 \right]^{1/2} \right) 
- \exp \left( -\frac{L_G^2 k^2}{4} - i V_{ims} t \left[ 1 - \left( \frac{k}{k_{ims}} \right)^2 \right]^{1/2} \right) \right\}.
\]

(C19)

\[
\delta B(x, t) \text{ and } \delta \hat{v}_i(x, t) \text{ are obtained by inverse transforming (C18) and (C19). Performing the integration in the inverse transforms is greatly aided by noting that because the initial perturbation in the magnetic field excites wavelengths primarily in the region } \sim L_G, \text{ the condition } L_G \ll L_{ims} \text{ implies that the contributions in the Gaussian packet are principally from waves with } k \gg k_{ims}. \text{ Taking the limit } (k_{ims}/k)^2 \to 0 \text{ will therefore introduce only a negligible error when inverting the transforms. Applying this limit gives the final solutions}
\]

\[
\delta B(x, t) = B_0 \frac{A_p}{2} \exp \left( -\frac{t}{2\tau_{in}} \right) \left[ \exp \left( -\frac{(x - V_{ims} t)^2}{L_G^2} \right) + \exp \left( -\frac{(x + V_{ims} t)^2}{L_G^2} \right) \right] 
+ B_0 \frac{A_p \sqrt{\pi} L_G}{8V_{ims} \tau_{in}} \exp \left( -\frac{t}{2\tau_{in}} \right) \left[ \text{erf} \left( \frac{x + V_{ims} t}{L_G} \right) - \text{erf} \left( \frac{x - V_{ims} t}{L_G} \right) \right].
\]

\]

(C20)

\[
\delta \hat{v}_i(x, t) = V_{ims} \frac{A_p}{2} \exp \left( -\frac{t}{2\tau_{in}} \right) \left[ \exp \left( -\frac{(x - V_{ims} t)^2}{L_G^2} \right) - \exp \left( -\frac{(x + V_{ims} t)^2}{L_G^2} \right) \right].
\]

(C21)

The solution corresponds to left- and right-traveling wave pulses propagating at the ion magneto-sound speed \( V_{ims} \) through a background of stationary neutrals. The decay timescale for the pulses is \( 2\tau_{in} \), a value that is consistent with the decay of individual Fourier wave modes at wavelengths \( \ll L_{ims} \) (e.g., see Section 3.1.1 of Ciolek et al. 2004, or Section 3.2.1 of Mouschovias et al. 2011). Equations (48a) and (48c) follow immediately from inserting (C20) and (C21) into (C1).

C.2. Solution for \( L_{ims} \ll L_G \ll L_{nms} \)

In this limit, the neutrals still remain motionless. Also, the collisional drag of the neutrals on the ions essentially balances the driving magnetic-pressure gradient in the charged fluid force equation. The solution in this wavelength regime can be derived from the Fourier-transformed magnetic field and velocity relations (C18) and (C19), by rewriting them as

\[
\delta \hat{B}(k, t) = B_0 A_p L_G \left( \frac{2}{\sqrt{2}} \right) \exp \left( -\frac{t}{2\tau_{in}} \right)
\times \left\{ \left[ 1 - \frac{1}{1 - (k/k_{ims})^2} \right] \exp \left( -\frac{L_G^2 k^2}{4} - V_{ims} k_{ims} t \left[ 1 - \left( \frac{k}{k_{ims}} \right)^2 \right]^{1/2} \right) 
+ \left[ 1 + \frac{1}{1 - (k/k_{ims})^2} \right] \exp \left( -\frac{L_G^2 k^2}{4} + V_{ims} k_{ims} t \left[ 1 - \left( \frac{k}{k_{ims}} \right)^2 \right]^{1/2} \right) \right\}.
\]

(C22)

and

\[
\delta \hat{v}_i(k, t) = -i \frac{V_{ims} A_p L_G k}{2\sqrt{2} k_{ims} \left[ 1 - (k/k_{ims})^2 \right]^{1/2}} \exp \left( -\frac{t}{2\tau_{in}} \right) \left\{ \exp \left( -\frac{L_G^2 k^2}{4} - V_{ims} k_{ims} t \left[ 1 - \left( \frac{k}{k_{ims}} \right)^2 \right]^{1/2} \right) 
- \exp \left( -\frac{L_G^2 k^2}{4} + V_{ims} k_{ims} t \left[ 1 - \left( \frac{k}{k_{ims}} \right)^2 \right]^{1/2} \right) \right\}.
\]

(C23)

The condition \( L_G \gg L_{ims} \) for the modes considered here is equivalent to taking the limit \( k/k_{ims} \to 0 \). Expanding (C22) and (C23) in the ratio \( k/k_{ims} \) and keeping terms only up to second-order will then suffice to give the principal solution when inverting the transforms. Doing this, and using the relations (C16) and (49d) to replace the cutoff wavenumber \( k_{ims} \), with the diffusion
The diffusion coefficient \( D_{ad} \) gives

\[
\delta B(x, t) = \frac{B_0 A_p}{(1 + 4D_{ad} t / L_G^2)^{3/2}} \exp \left( -\frac{x^2 / L_G^2}{1 + 4D_{ad} t / L_G^2} \right) 
+ \frac{2B_0 A_p D_{ad} \tau_m / L_G^2}{(1 + 4D_{ad} \tau_m / L_G^2)^{3/2}} \left( 1 - \frac{2x^2 / L_G^2}{1 + 4D_{ad} \tau_m / L_G^2} \right) \exp \left( -\frac{x^2 / L_G^2}{1 + 4D_{ad} \tau_m / L_G^2} \right) 
- \frac{2B_0 A_p D_{ad} \tau_m / L_G^2}{(1 - 4D_{ad} \tau_m / L_G^2)^{3/2}} \left( 1 - \frac{2x^2 / L_G^2}{1 - 4D_{ad} \tau_m / L_G^2} \right) \exp \left( \frac{-x^2 / L_G^2}{1 - 4D_{ad} \tau_m / L_G^2} \right),
\]

(C24)

\[
\delta v_i(x, t) = \frac{2D_{ad} A_p x}{L_G^2 (1 + 4D_{ad} t / L_G^2)^{3/2}} \exp \left( -\frac{x^2 / L_G^2}{1 + 4D_{ad} t / L_G^2} \right) 
- \frac{2D_{ad} A_p x}{L_G^2 (1 - 4D_{ad} \tau_m / L_G^2)^{3/2}} \exp \left( \frac{-x^2 / L_G^2}{1 - 4D_{ad} \tau_m / L_G^2} \right).
\]

(C25)

The solution here is one of ambipolar diffusion, in which the charged fluid and magnetic field diffuse through a sea of fixed neutrals, as reflected by the first terms after the equalities in Equations (C24) and (C25). The diffusion coefficient \( D_{ad} \) reflects the balance of the magnetic field pressure (since \( V_{ims}^2 \propto B^2 \)) and collisional forces (represented by \( \tau_m \)). From expressions (C24) and (C25) the characteristic decay timescale \( \tau_{dec} \) for this solution is found to be the ambipolar diffusion time \( \tau_{ad} = L_G^2 / 4D_{ad} \).

**C.3. Solution for \( L_G \gg L_{ims} \)**

In this limit, which has length scales \( \gg L_{ims} \) (since \( L_{ims} \gg L_{ims} \)) for our clouds, the charged fluid continues to move in a force-free manner, for this situation we can make the approximation \( \rho_0 \partial (\delta v_i) / \partial t \simeq 0 \) in the linearized ion momentum equation (C6). Hence, the ions will be traveling at the terminal drift velocity

\[
\delta v_i = -\frac{B_0 \tau_m}{4\pi \rho_0} \frac{\partial}{\partial x}(\delta B).
\]

(C26)

Inserting expression (C26) into the linearized force equation for the neutral fluid (C3) and also the linearized magnetic induction equation (C7), then Fourier transforming those equations along with the perturbed mass continuity equation (C2) yields the approximate system of ODEs for the wave modes in this wavelength region,

\[
\frac{\partial \hat{y}_n}{\partial t} = M_n \hat{y}_n,
\]

(C27)

where \( \hat{y}_n = [\delta \hat{\rho}_n(k, t), \delta \hat{v}_n(k, t), \delta \hat{B}(k, t)]^T \) is the vector of transformed variables, and the \( 3 \times 3 \) array of constant coefficients

\[
M_n \equiv \begin{bmatrix} 0 & -i \rho_0 k & 0 \\ -i k C_n^2 / \rho_0 & 0 & -i k B_0 / 4 \pi \rho_0 \\ 0 & -i B_0 k & -D_{ad} k^2 \end{bmatrix}.
\]

(C28)

In typical clouds with \( (C_n / V_{ims})^2 \ll 1 \) (see Equations (43)–(45)), the system (C27) has the solution

\[
\hat{y}_n = a_{ns} \mathcal{E}_{ns} e^{g_{ns} t} + a_{nt} \mathcal{E}_{nt} e^{g_{nt} t} + a_{ns} \mathcal{E}_{nt} e^{g_{nt} t},
\]

(C29a)

where

\[
g_{ns} = -\frac{1}{2} D_{ad} k^2 \pm i V_{ims} k \left[ 1 - \left( \frac{k}{k_{ims}} \right)^2 \right]^{1/2},
\]

(C29b)

\[
g_{nt} = -\left( \frac{C_n}{V_{ims}} \right)^2 D_{ad} k^2 = -D_{th} k^2,
\]

(C29c)

are the eigenvalues of \( M_n \), and

\[
\mathcal{E}_{ns} = \begin{bmatrix} 1, i \frac{g_{ns}}{\rho_0 k}, \frac{g_{ns} B_0}{\rho_0} \end{bmatrix}^{T} \approx \begin{bmatrix} 1, +V_{ims} \frac{B_0}{\rho_0}, \frac{B_0}{\rho_0} \end{bmatrix}^{T},
\]

(C29d)
are their associated eigenvectors. In the above,
\[ k_{\text{rms}} \equiv \frac{2\pi}{L_{\text{rms}}} = \frac{2V_{\text{rms}}}{V_{nms}^2 \tau_{ni}} = \frac{2V_{\text{rms}}}{D_{\text{ad}}} \]  
(C30)
is the neutral magnetosound cutoff wave number (Equations (49d), (31), (45), and (15) were used to derive the latter equalities). For \( k < k_{\text{rms}} \) the modes with eigenvalues \( g_{n0} \) are neutral magnetosound waves; these waves decay by ambipolar diffusion of the charged fluid and magnetic field with respect to the neutrals. The \( g_{00} \) mode is a neutral pressure-driven diffusion mode (e.g., see Figures 1(b) and (d) of Ciolek et al. 2004; or Section 3.2.1 of Mouschovias et al. 2011), in which the thermal-pressure-gas force of the neutral fluid drives the neutrals through a stationary background of the charged fluid and magnetic field. The balance of thermal-pressure and neutral–ion collisional drag in this mode has the neutrals moving at a terminal drift speed, resulting in neutral diffusion with the coefficient \( D_{\text{th}} \) defined in Equation (53). Simplifications resulting in the latter equalities of (C29d) and (C29e) were made by noting that for \( (C_n/V_{\text{rms}})^2 \ll 1, |g_{00}| \ll |g_{n0}| \). The limit \( k \ll k_{\text{rms}} \) was also used.

The coefficients \( a_{n0}, a_{n-}, \text{ and } a_{0} \) in (C29a) are found by applying the initial conditions \( \delta \rho_{n}(k, 0) = \delta \upsilon_{n}(k, 0) = 0, \) and \( \delta \mathcal{B}(k, 0) \) given by (C17). Doing this, we have the solutions
\[
\delta \mathcal{B}(k, t) = \frac{B_0 A_p L_G}{2\sqrt{2}[1 + (C_n/V_{\text{rms}})^2]} \times \left\{ \begin{array}{l}
1 + i2 \left( \frac{C_n}{V_{\text{rms}}} \right)^2 \frac{k}{k_{\text{rms}}} \exp \left( -\left[ \frac{L_G^2}{4} + D_{\text{ad}} t \right] \frac{k^2}{2} + i V_{\text{rms}} k t \left[ 1 - \left( \frac{k}{k_{\text{rms}}} \right)^2 \right]^{1/2} \right) \\
1 - i2 \left( \frac{C_n}{V_{\text{rms}}} \right)^2 \frac{k}{k_{\text{rms}}} \exp \left( -\left[ \frac{L_G^2}{4} + D_{\text{ad}} t \right] \frac{k^2}{2} - i V_{\text{rms}} k t \left[ 1 - \left( \frac{k}{k_{\text{rms}}} \right)^2 \right]^{1/2} \right) \\
+ 2 \left( \frac{C_n}{V_{\text{rms}}} \right)^2 \exp \left( -\left[ \frac{L_G^2}{4} + D_{\text{ad}} t \right] \frac{k^2}{2} \right) \end{array} \right
\]  
(C31)
\[
\delta \upsilon_{n}(k, t) = \frac{A_p V_{\text{rms}} L_G}{2\sqrt{2}[1 + (C_n/V_{\text{rms}})^2]} \times \left\{ \begin{array}{l}
1 + i2 \left( \frac{C_n}{V_{\text{rms}}} \right)^2 \frac{k}{k_{\text{rms}}} \exp \left( -\left[ \frac{L_G^2}{4} + D_{\text{ad}} t \right] \frac{k^2}{2} + i V_{\text{rms}} k t \left[ 1 - \left( \frac{k}{k_{\text{rms}}} \right)^2 \right]^{1/2} \right) \\
1 - i2 \left( \frac{C_n}{V_{\text{rms}}} \right)^2 \frac{k}{k_{\text{rms}}} \exp \left( -\left[ \frac{L_G^2}{4} + D_{\text{ad}} t \right] \frac{k^2}{2} - i V_{\text{rms}} k t \left[ 1 - \left( \frac{k}{k_{\text{rms}}} \right)^2 \right]^{1/2} \right) \\
+ i4 \left( \frac{C_n}{V_{\text{rms}}} \right)^2 \frac{k}{k_{\text{rms}}} \exp \left( -\left[ \frac{L_G^2}{4} + D_{\text{ad}} t \right] \frac{k^2}{2} \right) \end{array} \right
\]  
(C32)
\[
\delta \rho_{n}(k, t) = \frac{A_p \rho_{00} L_G}{2\sqrt{2}[1 + (C_n/V_{\text{rms}})^2]} \times \left\{ \begin{array}{l}
1 + i2 \left( \frac{C_n}{V_{\text{rms}}} \right)^2 \frac{k}{k_{\text{rms}}} \exp \left( -\left[ \frac{L_G^2}{4} + D_{\text{ad}} t \right] \frac{k^2}{2} + i V_{\text{rms}} k t \left[ 1 - \left( \frac{k}{k_{\text{rms}}} \right)^2 \right]^{1/2} \right) \\
1 - i2 \left( \frac{C_n}{V_{\text{rms}}} \right)^2 \frac{k}{k_{\text{rms}}} \exp \left( -\left[ \frac{L_G^2}{4} + D_{\text{ad}} t \right] \frac{k^2}{2} - i V_{\text{rms}} k t \left[ 1 - \left( \frac{k}{k_{\text{rms}}} \right)^2 \right]^{1/2} \right) \\
-2 \exp \left( -\left[ \frac{L_G^2}{4} + D_{\text{ad}} t \right] \frac{k^2}{2} \right) \end{array} \right
\]  
(C33)
Inverting the transforms give the solutions for \( \delta \rho_{n}(x, t), \delta \upsilon_{n}(x, t), \text{ and } \delta \mathcal{B}(x, t) \). The inversion is aided by the fact that for the initial perturbation we consider here, with \( L_G \gg L_{\text{rms}} \), the overwhelming bulk of the wave modes contributing to the wave packet have
k \ll k_{\text{num}}. Hence, taking the limit (k/k_{\text{num}})^2 \to 0 will introduce only a minor error when inverting the Fourier transforms (C31), (C32), and (C33). Doing this yields:

\[ \frac{\partial \hat{δ} B(x, t)}{\partial t} = \frac{A_p B_0 G_1}{2} \left[ \left\{ \begin{array}{l} \frac{-2(x + V_{\text{num}t})D_{th}/L_G^2}{V_{\text{num}}[1 + (2D_{\text{ad}t}/L_G^2)]} \exp \left( \frac{-(x + V_{\text{num}t})^2}{L_G^2[1 + (2D_{\text{ad}t}/L_G^2)]} \right) \\
\frac{-2(x - V_{\text{num}t})D_{th}/L_G^2}{V_{\text{num}}[1 + (2D_{\text{ad}t}/L_G^2)]} \exp \left( \frac{-(x - V_{\text{num}t})^2}{L_G^2[1 + (2D_{\text{ad}t}/L_G^2)]} \right)
\end{array} \right\} + \left\{ \begin{array}{l}
\frac{-2(x + V_{\text{num}t})D_{th}/L_G^2}{V_{\text{num}}[1 + (2D_{\text{ad}t}/L_G^2)]} \\
\frac{-2(x - V_{\text{num}t})D_{th}/L_G^2}{V_{\text{num}}[1 + (2D_{\text{ad}t}/L_G^2)]}
\end{array} \right\} \right], \]  

\( \frac{\partial \hat{δ} \rho_n(x, t)}{\partial t} = \frac{A_p \rho_0 G_1}{2} \left[ \left\{ \begin{array}{l} \frac{-2(x + V_{\text{num}t})D_{th}/L_G^2}{V_{\text{num}}[1 + (2D_{\text{ad}t}/L_G^2)]} \exp \left( \frac{-(x + V_{\text{num}t})^2}{L_G^2[1 + (2D_{\text{ad}t}/L_G^2)]} \right) \\
\frac{-2(x - V_{\text{num}t})D_{th}/L_G^2}{V_{\text{num}}[1 + (2D_{\text{ad}t}/L_G^2)]} \exp \left( \frac{-(x - V_{\text{num}t})^2}{L_G^2[1 + (2D_{\text{ad}t}/L_G^2)]} \right)
\end{array} \right\} + \left\{ \begin{array}{l}
\frac{-2(x + V_{\text{num}t})D_{th}/L_G^2}{V_{\text{num}}[1 + (2D_{\text{ad}t}/L_G^2)]} \\
\frac{-2(x - V_{\text{num}t})D_{th}/L_G^2}{V_{\text{num}}[1 + (2D_{\text{ad}t}/L_G^2)]}
\end{array} \right\} \right], \]  

\( \frac{\partial \hat{δ} \rho_i(x, t)}{\partial t} = \frac{A_p \rho_0 G_2}{2} \exp \left( \frac{-x^2}{L_G^2[1 + (4D_{\text{ad}t}/L_G^2)]} \right), \]

where the two functions \( G_1 \) and \( G_2 \) are, respectively, defined by Equations (51) and (52). The solutions clearly show that for initial perturbation with \( L_G \gg L_{\text{num}} \) left- and right-propagating neutral pulses traveling at the magnetoacoustic speed \( V_{\text{num}} \). There is some damping in the pulses because of ambipolar diffusion, occurring on a timescale \( \tau_{\text{dec}} = 2\tau_{\text{ad}} = L_G^2/2D_{\text{ad}} \). There is also a part to the above solutions that indicates an effect due to pressure-driven diffusion of the neutrals, as can be seen by the terms in the above equations involving \( D_{\text{th}} \). Those terms show that pressure and density gradients are created in the neutrals (by collisional drag from the ions) that can also affect the motion of the neutral fluid. However, these effects are small compared to that which results from the magnetically driven collisional drag on the ions on the neutrals.

Finally, Equations (50a)−(50d) are obtained by combining (C1), (C34)−(C36), and (C10), and using the relations \( n_n = \rho_n/m_n \) and \( n_i = \rho_i/m_i \). Equation (50e) follows from taking the partial derivative \( \partial/\partial x \) of (C34) and then inserting the result into (C26).

**APPENDIX D**

**NUMERICAL CONVERGENCE**

Here we provide quantitative data verifying the scaling and convergence of the numerical algorithm described in this paper. We do this by modeling the evolution of small-amplitude disturbances and comparing the numerical model results against exact solutions of the linearized system of equations (C2)–(C7). Inserting (C8) and (C9) into those equations, and then Fourier transforming them in space (Equation (C11)) gives

\[ \frac{\partial \hat{δ} \hat{y}_{\text{tot}}}{\partial t} = \hat{M}_{\text{tot}} \hat{δ} \hat{y}_{\text{tot}}, \]  

where \( \hat{y}_{\text{tot}} \equiv [\hat{δ} \hat{ρ}_n(k, t), \hat{δ} \hat{v}_n(k, t), \hat{δ} \hat{ρ}_i(k, t), \hat{δ} \hat{v}_i(k, t), \hat{δ} \hat{B}(k, t)]^T \) is the vector of transformed variables for the full system, and

\[ \hat{M}_{\text{tot}} \equiv \begin{bmatrix}
0 & -i\rho_0 & 0 & 0 & 0 \\
-1/\tau_{\text{in}} & -iB_0/4\pi & 0 & 0 & 0 \\
0 & 0 & 0 & 0 & -i\rho_0 & k \\
0 & 0 & 0 & 0 & 0 & 0 \\
1/\tau_{\text{in}} & 0 & 0 & -iC_n & 0 & 1/\tau_{\text{in}} & -1/\tau_{\text{in}}
\end{bmatrix}. \]

This particular ODE system has the solution

\[ \hat{y}_{\text{tot}}(k, t) = \sum_{m=1}^{5} a_me^{E_mt}e^{i\omega t}. \]
Figure 12. Rightward-traveling ion magnetosound wave with $\lambda_p = 9.35 \times 10^{11} \text{ cm} < L_{\text{ims}}$ at time $t = 1.37 \times 10^{-4} \text{ yr}$. (a) Ion velocity. Solid line is the exact eigensolution (Equation (D9)) at that time, dashed-line is the initial state. Also shown are numerical code results for three models having different numbers of mesh cells per wavelength of the initial perturbation: 16 cells (plus signs), 64 cells (squares), and 128 cells (circles). (b) Zoom-in of the ion velocity data about the origin. (c) Relative error in the ion velocity for the same three numerical models. (d) Numerically determined error measure of the ion velocity $L_1(v_i)$ as a function of the number of cells per wavelength in a model (curve with triangles). The dash-dotted line is a reference curve with a logarithmic slope of $-2$.

(Braun 1983, chap. 3), where $g_m$ is the $m$th eigenvalue of $M_{\text{tot}}$ and $E_m$ is its corresponding eigenvector. The constants $a_m$ are determined by the initial conditions. Making the substitution

$$g_m = -i \omega_m,$$  \hspace{1cm} (D4a)

where

$$\omega_m = \omega_{m,R} + i \omega_{m,I}$$  \hspace{1cm} (D4b)

is a complex frequency (with real and imaginary parts $\omega_{m,R}$ and $\omega_{m,I}$, respectively) and inverting the transform (Equation (C12)), yields

$$y_{\text{tot}}(x, t) = \frac{1}{\sqrt{2\pi}} \sum_{m=1}^{5} \int_{-\infty}^{+\infty} a_m E_m e^{i(kx - \omega_m t)} dk;$$  \hspace{1cm} (D5)

$y_{\text{tot}} = [\delta p_\alpha(x, t), \delta v_\alpha(x, t), \delta p_i(x, t), \delta v_i(x, t), \delta B(x, t)]^T$ is the total vector of perturbed variables.

A single mode with $m = l$ at wavelength $\lambda_p$ and wavenumber $k_p = 2\pi/\lambda_p$ is excited from the initial complex perturbation

$$y_{\text{tot}}(x, 0) = b_p E_l e^{i(k_p x + \chi_p)}.$$  \hspace{1cm} (D6)

The constant $b_p$ is taken to be real, and $\chi_p$ is the phase constant of the initial state. Setting $t = 0$ in (D5) and equating it to (D6) reveals that for the former relation to describe the evolution of the latter, we must have

$$a_m = \sqrt{2\pi} b_p \delta_{lm} \delta(k - k_p) e^{i\chi_p}.$$  \hspace{1cm} (D7)

$E_l$ is also complex, and can be written as

$$E_l = E_{l,R} + i E_{l,I}$$  \hspace{1cm} (D8)

where $E_{l,R}$ and $E_{l,I}$ are both real vectors. Inserting (D4b) and (D8) into (D7), it follows that the real part of the singly excited eigenmode of the system is given by

$$\Re[y_{\text{tot}}(x, t)] = b_p [E_{l,R} \cos(k_p x - \omega_{l,R} t + \chi_p) - E_{l,I} \sin(k_p x - \omega_{l,R} t + \chi_p)] e^{\omega_{l,R} t}.$$  \hspace{1cm} (D9)
To test convergence we excite a single mode of the multifluid system and compare the resulting evolution of our numerical code to the exact solution (D9). A $\lambda_p$ is selected and the eigenvalues and eigenvectors of the array $M_{tot}$ at that wavelength are calculated using widely available numerical routines (e.g., EISPACK, LAPACK) or software packages (e.g., MATLAB, Maple, IDL). To ensure that an excited mode always remains linear, the value of the numerical constant $b_p$ of the initial state (the real part of Equation (D6), or equivalently, Equation (D9) at $t = 0$) is chosen so that its magnetic field perturbation has a relative amplitude that is $10^{-6}$ that of the background magnetic field. For our convergence test runs we assume the same background reference state and conditions for all of the physical variables ($n_{i0}, n_{i0}, B_0, T_n$, etc.) as in the models described in Section 3.1. All of the speeds ($V_{ims}, V_{num}$, and $C_n$), collision times ($\tau_{ni}$ and $\tau_{ni}$), and characteristic length scales ($L_{ims}$ and $L_{num}$) are therefore also the same as for those models. Additionally, we set the eigenmode wavelengths $\lambda_p$ equal to the widths $L_G$ of the Gaussian test packets of Section 3.1; hence, the underlying physics of the waves discussed for those models also applies to the models presented here.

Our first excited eigenmode model has $\lambda_p = 9.35 \times 10^{11}$ cm, which is less than the upper ion magnetosound wave cutoff length $L_{ims}$. The selected mode is an ion magnetosound wave traveling in the +$x$-direction with velocity $+V_{ims}$. At this wavelength the mode has $\omega_{iR} = 5.84 \times 10^{-4}$ s$^{-1}$ and $\omega_{iT} = -1.26 \times 10^{-6}$ s$^{-1}$, the latter value corresponding to a decay rate $\tau_{dec} = 0.0251$ yr = $2\tau_{ni}$. Figure 12(a) shows the exact solution for the ion velocity (solid curve) at $t = 1.37 \times 10^4$ yr, by which time the wave has advanced by more than one-fourth of a wavelength from its initial state (dashed curve). Also shown are the results of three models having different spatial resolution: the first model (displayed as plus signs) has 16 mesh cells within each wavelength of the initial perturbation, the second (squares) 64 cells per wavelength, and the third (circles) 128 cells per wavelength. For this perturbation the numerical time step is determined by the CFL time step (29); the CFL number $\nu = 0.8$ in each of these models. Figure 12(b) displays a zoom-in of the same data about $x = 0$. The local relative error in the ion velocity $\|v_i,\text{exact}(x, t) - v_i,\text{num}(x, t)\|/v_i,\text{exact}(x, t)$ is presented in Figure 12(c), showing how the numerical accuracy increases with finer mesh spacing. Finally, Figure 12(d) shows the error measure in the ion velocity across the entire computational domain of N cells

$$L_1(v_i) \equiv \frac{1}{N} \sum_{j=1}^{N} |v_{i,\text{exact}}(x_j, t) - v_{i,\text{num}}(x_j, t)|$$

(D10)

(curve with triangles) for several different model runs, all with $\nu = 0.8$, as a function of each model’s numerical resolution. The range is from 4 to 512 cells per wavelength of the perturbation. Also shown (dash-dotted curve) is a reference curve with a logarithmic slope = $-2$, the value expected for a code that is second-order accurate in space and time. Our numerical results are seen to be consistent with second-order accuracy.

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Figure 13. Ambipolar diffusion mode with $\lambda_p = 3.74 \times 10^{15}$ cm ($L_{ims} < \lambda_p < L_{num}$) at $t = 3.17$ yr. All symbols and curves have the same meaning as in Figure 12.
The second excited eigenmode is at λ_p = 3.74 × 10^{15} cm, which is in the range \( L_{\text{ims}} < \lambda_p < L_{\text{nms}} \). The ambipolar diffusion mode that exists within this range is chosen for this test: at this wavelength that particular mode has \( \omega_{l,R} = 1.30 \times 10^{-26} \text{ s}^{-1} \) and \( \omega_{l,I} = -8.47 \times 10^{-9} \text{ s}^{-1} \). Its decay time is therefore \( \tau_{\text{dec}} = 3.74 \text{ yr} \). Figure 13(a) displays the initial state and the exact solution for the ion velocity at \( t = 3.17 \text{ yr} \). Also shown are the results of numerical runs having resolutions of 16, 64, and 128 cells per wavelength of the perturbation. For these model runs, the upper limit to the maximum stable numerical time is limited by the timescales occurring in the source terms (here, the ion–neutral collision time \( \tau_{\text{in}} \)) instead of that which would be calculated with a traditional CFL time step with \( v \sim 0.1–1 \) (see Section 2.3). Stability of the time integration for these models thus requires that their time step \( \Delta t = \Delta t_{Sj} \leq \tau_{\text{in}} \).

In analogy to convergence tests in which the CFL number \( v \propto \Delta t/\Delta x \) is kept fixed, the models for this test have \( \Delta t/\Delta x = \text{constant} \), such that

\[
\Delta t = \frac{(4/\lambda_p)}{(N_p/\lambda_p)} \tau_{\text{in}}, \tag{D11}
\]

where \( N_p (\geq 4) \) is the number of cells within one wavelength of the perturbation. A close-up showing the results for this eigenmode about the origin is presented in Figure 13(b), and the local relative error is displayed in Figure 13(c). The \( L_1(v_i) \) error measure of this mode can be seen in Figure 13(d). Examination of that panel indicates that our code tends to scale slightly better than second-order accuracy in this instance.

The final excited eigenmode has \( \lambda_p = 2.24 \times 10^{17} \text{ cm} > L_{\text{nms}} \). For this test we chose the mode that corresponds to a neutral magnetosound wave propagating in the +x-direction. At this wavelength \( \omega_{l,R} = 1.02 \times 10^{-12} \text{ s}^{-1} \) and \( \omega_{l,I} = -7.93 \times 10^{-13} \text{ s}^{-1} \); hence, the decay time by ambipolar diffusion for this mode is \( \tau_{\text{dec}} = 3.99 \times 10^4 \text{ yr} \). Figure 14(a) shows the eigenmode solution for the ion velocity at \( t = 5.03 \times 10^4 \text{ yr} \), by which time the wave has moved about 1/4 of a wavelength from its initial state. Three model runs with different numerical resolution are also shown: one with 64 cells per wavelength (squares), another with 128 cells per wavelength (circles), and the last with 256 cells per wavelength (diamonds). For these eigenmode runs it is again the case that the maximum stable numerical time step cannot exceed \( \tau_{\text{in}} \), so the relation (D11) was used to fix the ratio \( \Delta t/\Delta x \) to the same value in each model to test convergence in the usual way. An expanded view of the ion velocity data near \( x = 0 \) is shown in Figure 14(b), and the local relative error in \( v_i \) as a function of position for the three models is found in Figure 14(c). The error measure \( L_1(v_i) \) as a function of numerical resolution is provided in Figure 14(d); the data there show for the large part that our code again exhibits second-order accuracy.

Figure 14. Right-traveling neutral magnetosound wave with \( \lambda_p = 2.24 \times 10^{17} \text{ cm} > L_{\text{nms}} \) at \( t = 5.03 \times 10^4 \text{ yr} \). (a) Ion velocity: exact eigensolution (solid curve) and initial state (dashed). Also displayed are results for models having different numerical resolutions: 64 cells per initial perturbation wavelength (squares), 128 cells per wavelength (circles), and 256 cells per wavelength (diamonds). (b) Close-up of the ion velocity data about the origin. (c) Relative error in the ion velocity for the three numerical models. (d) Error measure \( L_1(v_i) \) as a function of model resolution. The dash-dotted reference curve has a logarithmic slope = −2.
