A KINETIC APPROACH OF THE BI-TEMPERATURE EULER MODEL

STÉPHANE BRULL

Univ. Bordeaux, CNRS, Bordeaux INP, IMB, UMR 5251, F-33400 Talence, France.

BRUNO DUBROCA

Univ. Bordeaux, Laboratoire des Composites ThermoStructuraux (LCTS), UMR 5801: CNRS-Herakles(Safran)-CEA-UBx, 3, Allée de La Boétie, 33600 Pessac, France

CORENTIN PRIGENT

Univ. Bordeaux, CNRS, Bordeaux INP, IMB, UMR 5251, F-33400 Talence, France.

Abstract. We are interested in the numerical approximation of the bi-temperature Euler equations, which is a non conservative hyperbolic system introduced in [4]. We consider a conservative underlying kinetic model, the Vlasov-BGK-Poisson system. We perform a scaling on this system in order to obtain its hydrodynamic limit. We present a deterministic numerical method to approximate this kinetic system. The method is shown to be Asymptotic-Preserving in the hydrodynamic limit, which means that any stability condition of the method is independent of any parameter $\varepsilon$, with $\varepsilon \to 0$. We prove that the method is, under appropriate choices, consistent with the solution for bi-temperature Euler. Finally, our method is compared to methods for the fluid model (HLL, Suliciu).

1. Introduction

Plasma physics is renowned for being a multiscale problem ([13]). Phenomena range from plasma oscillations and charge separation at very small scales to fluid mechanics at macroscopic levels. Different physical models exist in order to describe a plasma according to the scale of interest. The kinetic model, constituted of the Vlasov-Boltzmann equations, coupled with Maxwell equations, are well suited for mesoscopic descriptions. In this model, each population of particles $\alpha$ (in the case of a plasma, these are electrons, ions and neutral atoms and molecules) is represented via the use of a distribution function $f^\alpha$. This function, defined over the phase space, denotes the number of particles that, at a given time $t$ at a given point $x$, possess a velocity $v$. Via an hydrodynamic limit, equations on the macroscopic conserved quantities (mass, momentum and energy) can be derived. These equations form the so-called bi-temperature Euler system (see [4] for the electrostatic case, and [9] for the transverse magnetic case, and [3] for the viscous development). From a numerical point of view, these equations are less accurate than the underlying kinetic model, but they are cheaper to solve.

The bi-temperature Euler system is comprised of an equation over mass, an equation over momentum, and one equation over each species energy. These energy equations are non-conservative,
accounting for the fact that, in a plasma, electrons and ions are exchanging energy through collisions and via electromagnetic interactions. Contrarily to the usual Euler system, for which the exact solution of any Riemann problem is known, these non-conservative terms prevent the knowledge of such solutions, especially when the solution contains a shock. This is a consequence that these non-conservative terms are written as products of distributions, which does not possess a clear definition. [17] presents a way of defining these products using a theory of paths. It is made clear that a certain amount of arbitrariness is needed in the choice of such paths. A major drawback of this property is that, when attempting numerical resolution of this system, every numerical scheme chooses its own path, and the Rankine-Hugoniot relation holding across shocks will be different for every scheme. This phenomenon is highlighted in [4]. More precisely, different schemes have different numerical viscosity, and the shock relations depend on such viscosities.

Several solutions have been proposed in order to derive numerical methods that use the notion of paths to define jump relations ([28], [11]). Other approaches have been proposed for different kinds of non-conservative systems (see [12], [21], [26], [29], [1]). Finally, Bouchut’s book [7] presents an overview of numerical methods for non-conservative Shallow-water equations.

In this paper, the aim is to propose a reference numerical method for such solutions using the underlying kinetic model. This kinetic model is conservative, and hence does not exhibit the drawback of the macroscopic model. Hence, the idea dwells in solving the Vlasov-BGK-Ampère system in the hydrodynamic limit, which will be presented in section 2.3. In order to be able to compare the results with the scheme applied to the bi-temperature Euler system, it is necessary to describe identical scales. By performing a scaling on the kinetic system, dimensionless parameters (scaled Debye length, scaled plasma frequency) are introduced in the system. Taking the hydrodynamic limit then amounts to taking the limit when these parameters tend to zero. In the general case, if a naive numerical approach is used, extremely restrictive stability conditions will appear, rendering the scheme unusable in a decent amount of computation time. Hence, an Asymptotic-Preserving (AP) scheme needs to be derived. Such a scheme possesses stability conditions independent of these small parameters and is then able to compute solutions for both cases when parameters are of the order of one and when these parameters are arbitrarily small.

The adjective Asymptotic Preserving has been introduced by Shi Jin in [23]. In the study of plasma simulation, numerous several Asymptotic Preserving methods have been proposed for various models, such as Euler-Maxwell, Euler-Poisson, Euler-Lorentz, Vlasov-Poisson ([20], [8], [19], [16], [15], [18], [5]).

The paper is organized as follows. Firstly, the macroscopic and kinetic models relative to our problem are introduced. A reduced 1D3V kinetic model is derived, and the relative properties of moment conservation and entropy dissipation are specified. Subsequently, a scaling of the kinetic model is proposed in order to exhibit the limiting behaviour of the system. Then, an Asymptotic-Preserving numerical method is presented. This method is able to compute solutions of the kinetic model in the asymptotic regime. Finally, test cases are presented in order to illustrate the derived method. In particular, the numerical results are compared with an HLL-inspired method and a Suliciu relaxation approach, both derived in [4].

2. Description of the models

2.1. Bi-temperature Euler model. The one-dimensional bi-temperature Euler model describes the behaviour of a two species fluid (constituted of electrons and ions) on a macroscopic scale
(see [14]):
\begin{align}
(1a) \quad & \partial_t \rho + \partial_x (\rho u) = 0, \\
(1b) \quad & \partial_t (\rho u) + \partial_x (\rho u^2 + p^* + p^i) = 0, \\
(1c) \quad & \partial_t (\rho^e \epsilon^e + \frac{1}{2} \rho^i u^2) + \partial_x (u(\rho^e \epsilon^e + \frac{1}{2} \rho^i u^2 + p^*) + u(c^i \partial_x p^e - c^e \partial_x p^i)) \\
& \hspace{1cm} = -\nu_{ei}(T^i - T^e), \\
(1d) \quad & \partial_t (\rho^i \epsilon^i + \frac{1}{2} \rho^i u^2) + \partial_x (u(\rho^i \epsilon^i + \frac{1}{2} \rho^i u^2 + p^i)) - u(c^i \partial_x p^i - c^e \partial_x p^e) \\
& \hspace{1cm} = \nu_{ei}(T^e - T^i).
\end{align}

Superscripts \( e \) and \( i \) denote quantities related to electrons and ions, respectively. \( \rho = \rho^e + \rho^i = m^e n^e + m^i n^i \) is the total density, \( m^\alpha \) being the mass of a particle of species \( \alpha \), and \( n^\alpha \) the corresponding concentration. \( u \) is the macroscopic velocity of the fluid. \( p^\alpha, T^\alpha \) and \( e^\alpha \) denote the partial pressures, temperatures and specific energies. \( c^\alpha = \frac{m^\alpha}{m^\alpha + m^eta} \) denotes the mass fractions. \( \nu_{ei} \) is the temperature exchange rate between the two species, that can depend on space and time.

### 2.2. Macroscopic quantities

Define \( f^\alpha(t, x, v) : \mathbb{R}_+ \times \mathbb{R}^3 \times \mathbb{R}^3 \to \mathbb{R}_+ \), the particle distribution function of species \( \alpha \), where \( t, x, v \) respectively denote the time, space and microscopic velocity variables. Integration with respect to \( v \in \mathbb{R}^3 \) will be denoted as follows, for any function \( g \) that depends on \( v \) such that \( g \in L^1(\mathbb{R}^3) \):

\[
\langle g \rangle = \int_{\mathbb{R}^3} g dv.
\]

Macroscopic quantities \( n^\alpha(t, x), u^\alpha(t, x) \) and \( T^\alpha(t, x) \) are defined as moments of \( f^\alpha \) such that \((1 + v^2)f^\alpha \in L^1(\mathbb{R}^3)\) as follows:

\[
\langle f^\alpha \rangle = n^\alpha, \\
\langle vf^\alpha \rangle = n^\alpha u^\alpha, \\
\langle \frac{m^\alpha v^2 f^\alpha}{2} \rangle = \frac{1}{2} m^\alpha n^\alpha (u^\alpha)^2 + \frac{3}{2} n^\alpha k_B T^\alpha.
\]

Similarly, mixture macroscopic quantities \( \rho, u \) and \( T \) are defined as:

\[
(2a) \quad \langle m^e f^e + m^i f^i \rangle = \rho, \\
(2b) \quad \langle v(m^e f^e + m^i f^i) \rangle = \rho u, \\
(2c) \quad \langle \frac{v^2}{2}(m^e f^e + m^i f^i) \rangle = \frac{1}{2} \rho u^2 + \frac{3}{2} (n^e + n^i) k_B T.
\]

Finally, we define electromagnetic quantities: \( E \in \mathbb{R} \) is the electric field along \( x \), \( q^\alpha \) is the electric charge of species \( \alpha \). Note that \( q^e = -Z q^i \), with \( Z \) the ionization rate. \( j \) and \( \mathbf{p} \) are respectively the current and total electric charge, defined as:

\[
\langle q^e f^e + q^i f^i \rangle = \mathbf{p}, \\
\langle v(q^e f^e + q^i f^i) \rangle = j.
\]

### 2.3. Vlasov-BGK-Ampère system

Consider the Vlasov-BGK-Ampère model, for \( t \in \mathbb{R}, x \in \mathbb{R}, v = (v_1, v_2, v_3) \in \mathbb{R}^3 \) and \( \alpha \in \{e, i\} \):

\[
(4a) \quad \partial_t f^\alpha + v_1 \partial_{v_1} f^\alpha + \frac{q^\alpha E}{m^\alpha} \partial_{v_1} f^\alpha = \frac{1}{\tau^\alpha} (M(f^\alpha) - f^\alpha) + \frac{1}{\tau^{\alpha\alpha}} (\mathcal{M}(f^\alpha) - f^\alpha),
\]

\[
(4b) \quad \partial_t E = -\frac{j}{\varepsilon_0},
\]

\[
(4c) \quad \partial_t \mathbf{p} = \frac{j}{\varepsilon_0},
\]

where \( \tau^\alpha \) and \( \tau^{\alpha\beta} \) are, respectively, the relaxation rate towards intra-species and inter-species equilibria (\( \tau^{\alpha\beta} = \tau^{\beta\alpha} \)). \( \varepsilon_0 \) is the dielectric permittivity of vacuum. The first equation is called
the Vlasov-BGK equation, while the two equations on the electric field are respectively called the Maxwell-Ampère equation and the Maxwell-Gauss equation.

The entropy-minimizing distribution function is the local Maxwellian distribution (see [27] for more details), denoted \( M(f^\alpha) \), defined by:

\[
M(f^\alpha)(t, x, v) = \frac{n^\alpha(t, x)}{(2\pi k_B T^\alpha(t, x))^{\frac{3}{2}}} \exp\left(-\frac{|v - u^\alpha(t, x)|^2}{2k_B T^\alpha(t, x)}\right),
\]

\( \overline{M}(f^\alpha) \), the exchange Maxwellian distribution, presented for example in [6], is defined as:

\[
\overline{M}(f^\alpha)(t, x, v) = \frac{n^\alpha(t, x)}{(2\pi k_B T(t, x))^{\frac{3}{2}}} \exp\left(-\frac{|v - u(t, x)|^2}{2k_B T(t, x)}\right).
\]

2.4. Reduced model.

2.4.1. Definition of the reduced distribution functions. The problem is assumed to be one dimensional in space and 3 dimensional in velocity (1D3V). Hence, the distribution functions \( f^\alpha \) only depend on \( x = x_1 \) and \( v = (v_1, v_2, v_3) \). To further simplify the model, we also assume that the function \( f^\alpha \) is even in \( v_2 \) and \( v_3 \), which can be expressed as:

\[
\begin{align*}
\int_{\mathbb{R}^2} f^\alpha(t, x, v)v_2 dv_2 dv_3 &= 0, \\
\int_{\mathbb{R}^2} f^\alpha(t, x, v)v_3 dv_2 dv_3 &= 0.
\end{align*}
\]

(5a)
(5b)

Denote \( u^\alpha = (u^\alpha_1, u^\alpha_2, u^\alpha_3) \), this assumption gives \( u^\alpha = (u^\alpha_1, 0, 0) \). Hence, the notation \( u^\alpha = u^\alpha_1 \) is used. The components \( u_2 \) and \( u_3 \) of the macroscopic velocity being equal to zero, the number of variables can be reduced, as in [2] or [25], in order to reduce the numerical cost of the method, while retaining the actual 3 dimensional total energy. Define the marginals, for \( \alpha = e, i \), by:

\[
\begin{align*}
g^\alpha_0(t, x, v_1) &= \int_{\mathbb{R}^2} f^\alpha(t, x, v)v_2 dv_2 dv_3, \\
g^\alpha_2(t, x, v_1) &= \int_{\mathbb{R}^2} f^\alpha(t, x, v)\frac{m^\alpha(v_2^2 + v_3^2)}{2} dv_2 dv_3,
\end{align*}
\]

and denote integration with respect to \( v_1 \), for any function \( g \) that depends on \( v_1 \), as:

\[
(g) = \int_{\mathbb{R}} g dv_1.
\]

(7)

2.4.2. Reduced system of equations and its properties. The idea is to describe the system of equations and the macroscopic quantities using these reduced functions. The system obtained after reduction reads as follows:

**Proposition 1.** Under the assumptions (5) and (6), the system (4) can be reduced to the following one, with \( \alpha = e, i \) and \( p = 0, 2 \):

\[
\begin{align*}
\partial_t g^\alpha_0 + v_1 \partial_x g^\alpha_0 + \frac{q_p^2 E}{m^\alpha} \partial_{v_1} g^\alpha_p &= \frac{1}{\tau^\alpha} (M^\alpha_p - g^\alpha_p) + \frac{1}{\tau^\alpha} (\overline{M}^\alpha_p - g^\alpha_p), \\
\partial_x E|_{t=t_0} &= \frac{1}{\varepsilon_0} \mathbf{p}|_{t=t_0}, \\
\partial_t E &= -\frac{1}{\varepsilon_0} j,
\end{align*}
\]

(8a)
(8b)
(8c)
where the different Maxwellian distributions are defined as:

\[
\begin{align*}
(9a) \quad M_0^\alpha(t, x, v) &= \frac{n^\alpha(t, x)}{\sqrt{2\pi k_B}} \exp\left( -\frac{(v_1 - u^\alpha)^2}{2k_B T^\alpha(t, x)} \right), \\
(9b) \quad M_2^\alpha(t, x, v) &= k_B T^\alpha(t, x) M_0^\alpha(t, x, v), \\
(9c) \quad \mathcal{M}_0^\alpha(t, x, v) &= \frac{\alpha}{\sqrt{2\pi k_B}} \exp\left( -\frac{(v_1 - u)^2}{2k_B T(t, x)} \right), \\
(9d) \quad \mathcal{M}_2^\alpha(t, x, v) &= k_B T(t, x) \mathcal{M}_0^\alpha(t, x, v).
\end{align*}
\]

Moreover, \( g_0^\alpha \) and \( g_2^\alpha \) satisfy the following relations:

\[
\begin{align*}
(10a) \quad n^\alpha = \langle g_0^\alpha \rangle, \\
(10b) \quad n^\alpha u^\alpha = \langle g_0^\alpha v_1 \rangle, \\
(10c) \quad \frac{3}{2} n^\alpha k_B T^\alpha + \frac{1}{2} m^\alpha n^\alpha (u^\alpha)^2 = \langle g_0^\alpha \frac{m^\alpha v_1^2}{2} \rangle + \langle g_2^\alpha \rangle.
\end{align*}
\]

**Proof.** Firstly, by integrating equation (4a) with respect to \( v_2 \) and \( v_3 \), it comes:

\[
\begin{align*}
\partial_t \int_{\mathbb{R}^2} f^\alpha dv_2 dv_3 + v_1 \partial_x \int_{\mathbb{R}^2} f^\alpha dv_2 dv_3 &= \frac{1}{n^\alpha} \left( \int_{\mathbb{R}^2} M(f^\alpha) dv_2 dv_3 - \int_{\mathbb{R}^2} f^\alpha dv_2 dv_3 \right) \\
&\quad + \frac{1}{n^\alpha} \left( \int_{\mathbb{R}^2} \mathcal{M}(f^\alpha) dv_2 dv_3 - \int_{\mathbb{R}^2} f^\alpha dv_2 dv_3 \right).
\end{align*}
\]

Define \( M_0^\alpha, \mathcal{M}_0^\alpha : \mathbb{R}^+ \times \mathbb{R} \times \mathbb{R} \to \mathbb{R} \) using (9a-9c), which correspond to the formulae obtained from:

\[
\begin{align*}
M_0^\alpha &= \int_{\mathbb{R}^2} M(f^\alpha) dv_2 dv_3, \\
\mathcal{M}_0^\alpha &= \int_{\mathbb{R}^2} \mathcal{M}(f^\alpha) dv_2 dv_3.
\end{align*}
\]

The definition of \( g_0^\alpha \) given in (6a) allows us to rewrite (11) as (8a) with \( p = 0 \). We have obtained an equation on the function \( g_0^\alpha \). Let us express the moments using \( g_0^\alpha \). The 0\textsuperscript{th}-order moment can be expressed as:

\[
\langle f^\alpha \rangle = \langle g_0^\alpha \rangle = n^\alpha.
\]

Similarly, the first order moment writes:

\[
n^\alpha u^\alpha = \langle f^\alpha v \rangle = \left\langle \int_{\mathbb{R}^2} f^\alpha \begin{pmatrix} v_1 \\ v_2 \\ v_3 \end{pmatrix} dv_2 dv_3 \right\rangle.
\]

According to (5a) and (5b), the second and third components of the first order moment are equal to zero. Then, we get, for the first component:

\[
\left\langle v_1 \int_{\mathbb{R}^2} f^\alpha dv_2 dv_3 \right\rangle = (v_1 g_0^\alpha) = n^\alpha u^\alpha.
\]

The third moments reads as:

\[
\langle f^\alpha m^\alpha |v|^2 \rangle = \left\langle g_0^\alpha \frac{m^\alpha v_1^2}{2} \right\rangle + \langle f^\alpha m^\alpha (v_2^2 + v_3^2) \rangle.
\]

Hence, another function \( g_2^\alpha \), defined by (6b), is introduced. Rewrite (12) as:

\[
\frac{3}{2} n^\alpha k_B T^\alpha + \frac{1}{2} m^\alpha n^\alpha (u^\alpha)^2 = \left\langle g_0^\alpha \frac{m^\alpha v_1^2}{2} \right\rangle + \langle g_2^\alpha \rangle.
\]
Finally, in order to close the system, an equation over \( g_3^2 \) is needed. To obtain it, multiply (4a) by \( \frac{m^\alpha(v_2^3 + v_3^2)}{2} \), and integrate with respect to \( v_2 \) and \( v_3 \):

\[
\int_{\mathbb{R}^2} \partial_t f^\alpha \frac{m^\alpha(v_2^3 + v_3^2)}{2} dv_2 dv_3 + \int_{\mathbb{R}^2} v_1 \partial_z f^\alpha \frac{m^\alpha(v_2^3 + v_3^2)}{2} dv_2 dv_3 \\
+ \frac{q^\alpha E}{m^\alpha} \int_{\mathbb{R}^2} \partial_{v_1} f^\alpha \frac{m^\alpha(v_2^3 + v_3^2)}{2} dv_2 dv_3 \\
= \frac{1}{\tau^\alpha} \left( \int_{\mathbb{R}^2} M(f^\alpha) \frac{m^\alpha(v_2^3 + v_3^2)}{2} dv_2 dv_3 - \int_{\mathbb{R}^2} f^\alpha \frac{m^\alpha(v_2^3 + v_3^2)}{2} dv_2 dv_3 \right).
\]

Then we introduce the function \( M_2^2 \) of (9b) corresponding to:

\[
M_2^2 = \int_{\mathbb{R}^2} M(f^\alpha) \frac{m^\alpha(v_2^3 + v_3^2)}{2} dv_2 dv_3.
\]

We get equation (8a) with \( p = 2 \). □

**Remark 1.** Equation (8b) means that, provided that Maxwell-Gauss equation is satisfied at \( t = 0 \) and if the Maxwell-Ampère equation is satisfied for all \( t \), so will the Maxwell-Gauss equation.

Regarding the conservation properties of the BGK operators in equations (8a), the following proposition holds:

**Proposition 2.** The intra-species BGK operators defined by the Maxwellians (9a) and (9b) conserve the partial masses, momenta and energies according to the following identities, for \( \alpha = e, i \):

\[
\langle (M_0^\alpha - g_0^\alpha) \left( \frac{1}{v_1} \right) \rangle = 0, \quad \langle (M_0^\alpha - g_0^\alpha) \frac{m^\alpha(v_1^2)}{2} \rangle + \langle (M_2^2 - g_2^2) \rangle = 0.
\]

The exchange Maxwellians (9c) and (9d) define inter-species BGK operators that enjoy the following conservation properties:

\[
\langle M_0^\alpha - g_0^\alpha \rangle = 0, \quad \sum_{\alpha} \langle (M_0^\alpha - g_0^\alpha) m^\alpha v_1 \rangle = 0,
\]

\[
\sum_{\alpha} \langle (M_0^\alpha - g_0^\alpha) \frac{m^\alpha(v_1^2)}{2} \rangle + \sum_{\alpha} \langle M_2^2 - g_2^2 \rangle = 0.
\]

**Proof.** The proof of this proposition is similar to [24]. Using the results from Proposition 1, we first recall equations (10a-10b)

\[
n^\alpha = \langle g_0^\alpha \rangle, \quad n^\alpha u^\alpha = \langle g_0^\alpha v_1 \rangle.
\]

Then, according to the expressions of \( M_0^\alpha \) from (9a), a straightforward computation gives:

\[
n^\alpha = \langle M_0^\alpha \rangle, \quad n^\alpha u^\alpha = \langle M_0^\alpha v_1 \rangle,
\]

which gives the first equation of (13). For the second equation, the following computation holds, according to the expression of \( M_2^2 \) in (9b):

\[
\frac{3}{2} n^\alpha k_B T^\alpha + \frac{1}{2} m^\alpha n^\alpha (u^\alpha)^2 = \langle M_0^\alpha \frac{m^\alpha v_1^2}{2} \rangle + \langle M_2^2 \rangle,
\]

which gives the expected result by substracting equation (10c).
For the results concerning the exchange Maxwellians, the expressions of \( \overline{M}_0^\alpha \) and \( \overline{M}_2^\alpha \) in (9c) and (9d) give the following computations:

\[
\begin{align*}
\langle \overline{M}_0^\alpha \rangle &= n^\alpha, \\
\langle m^\alpha \overline{M}_0^\alpha v_1 \rangle &= \rho^\alpha u, \\
\langle m^\alpha \overline{M}_0^\alpha v_1^2 \rangle &= \frac{1}{2} \rho^\alpha u^2 + \frac{3}{2} n^\alpha k_B T.
\end{align*}
\]

The first equation gives the first equation of (14). By summing the second equation over \( \alpha \), and in favor of the definition of \( \rho \) and \( u \) from (2), one gets the second equation of (14). Similarly, by summing the third equation over \( \alpha \), and in favor of the definition of \( \rho, u \) and \( T \) from (2), equation (15) holds.

\( \square \)

Define now the entropy of the non-reduced system:

\[
H^\alpha(f^\alpha) = \langle f^\alpha \log f^\alpha \rangle.
\]

In order to obtain the entropy of the reduced system, consider the following definition of the Maxwellian distribution of the non-reduced system:

\[
H^\alpha(M(f^\alpha)) = \min_{f^\alpha \geq 0} (\hat{H}^\alpha(f^\alpha)) \text{ s.t. } g_0^\alpha(v_1) = \int f^\alpha(v) dv_2 dv_3,
\]

\[
g_2^\alpha(v_1) = \int f^\alpha(v) \frac{m^\alpha(v_2^2 + v_3^2)}{2} dv_2 dv_3,
\]

with

\[
\hat{H}^\alpha(f^\alpha) = \int f^\alpha \log f^\alpha dv_2 dv_3.
\]

**Proposition 3.** The minimization problem (17) admits the following solution:

\[
M(f^\alpha)(g_0^\alpha, g_2^\alpha) = \frac{(g_0^\alpha)^2 m^\alpha}{2\pi g_2^\alpha} \exp \left( -\frac{g_0^\alpha m^\alpha (v_2^2 + v_3^2)}{2g_2^\alpha} \right).
\]

**Proof.** Define the Lagrangian associated with this minimization problem:

\[
\mathcal{L}(f^\alpha, \lambda_1, \lambda_2) = \int f^\alpha \log f^\alpha dv_2 dv_3 + \lambda_1(g_0^\alpha(v_1)) - \int f^\alpha(v) dv_2 dv_3 \\
+ \lambda_2(g_2^\alpha(v_1)) - \int f^\alpha(v) \frac{m^\alpha(v_2^2 + v_3^2)}{2} dv_2 dv_3,
\]

where \( \lambda_1 \) and \( \lambda_2 \) denote the Lagrange multipliers associated to the constraints

\[
g_0^\alpha(v_1) = \int f^\alpha(v) dv_2 dv_3, \quad g_2^\alpha(v_1) = \int f^\alpha(v) \frac{m^\alpha(v_2^2 + v_3^2)}{2} dv_2 dv_3.
\]

A straightforward computation of the differential of the Lagrangian w.r.t \( f^\alpha \) gives the following equation for the minimum \( M(f^\alpha)(g_0^\alpha, g_2^\alpha) \):

\[
\log(M(f^\alpha)(g_0^\alpha, g_2^\alpha)) + 1 - \lambda_1 - \lambda_2 \frac{m^\alpha (v_2^2 + v_3^2)}{2} = 0.
\]

Then, by using the constraints, the following expressions of \( \lambda_1 \) and \( \lambda_2 \) are found:

\[
\lambda_1 = 1 + \log \frac{(g_0^\alpha)^2 m^\alpha}{2\pi g_2^\alpha}, \quad \lambda_2 = -\frac{g_0^\alpha}{g_2^\alpha},
\]
which give the following expression for the Maxwellian:

\[
M(f^\alpha)(g_0^\alpha, g_2^\alpha) = \frac{(g_0^\alpha)^2 m^\alpha}{2\pi g_2^\alpha} \exp\left(-\frac{g_0^\alpha m^\alpha (v_1^2 + v_3^2)}{2g_2^\alpha}\right).
\]

Hence, by injecting (19) in (16), it comes

\[
H^\alpha(g_0^\alpha, g_2^\alpha) = \left\langle g_0^\alpha \log \frac{(g_0^\alpha)^2}{g_2^\alpha} - 2g_0^\alpha \right\rangle.
\]

Hence, define the following notations for the partial entropies of the reduced model:

\[
H^\alpha(g_0^\alpha, g_2^\alpha) = \left\langle \eta^\alpha(g_0^\alpha, g_2^\alpha) \right\rangle,
\]

with

\[
\eta^\alpha(g_0^\alpha, g_2^\alpha) = g_0^\alpha \log \frac{(g_0^\alpha)^2}{g_2^\alpha} - 2g_0^\alpha,
\]

for \(\alpha = e, i\), and the total entropy

\[
H(g_0^e, g_2^e, g_0^i, g_2^i) = \sum_\alpha H^\alpha(g_0^\alpha, g_2^\alpha).
\]

Firstly, this entropy is shown to be convex:

**Proposition 4.** For \(\alpha = e, i\), under the assumption that \(g_0^\alpha, g_2^\alpha\) are positive, \(H^\alpha\) is a strictly convex function. Consequently, \(H\) is strictly convex as well.

**Proof.** Compute the gradient of \(\eta^\alpha\):

\[
\nabla \eta^\alpha = \left(\frac{\partial \eta^\alpha}{\partial g_0^\alpha}, \frac{\partial \eta^\alpha}{\partial g_2^\alpha}\right) = \left(\log \frac{(g_0^\alpha)^2}{g_2^\alpha}, -\frac{1}{g_0^\alpha}\right).
\]

Compute the Hessian:

\[
\nabla^2 \eta^\alpha = \begin{pmatrix}
\frac{2}{g_0^\alpha} & \frac{1}{g_0^\alpha} \\
-\frac{1}{g_0^\alpha} & \frac{2}{g_0^\alpha}
\end{pmatrix}
\]

By positivity of the distribution functions, \(\text{Tr}(\nabla^2 \eta^\alpha) > 0\) and \(\det(\nabla^2 \eta^\alpha) = \frac{1}{(g_0^\alpha)^2} > 0\). Hence, the Hessian is positive definite and \(\eta^\alpha\) is strictly convex. Consequently, \(H^\alpha\) is strictly convex.

\(\square\)

The following property on entropy dissipation holds:

**Proposition 5.** The reduced system (8) satisfies the H-theorem for the entropy defined by (22):

\[
\frac{\partial H}{\partial t} + \frac{\partial G}{\partial x} \leq 0,
\]

with \(G\) the entropy flux associated to \(H\):

\[
G(g_0^e, g_2^e, g_0^i, g_2^i) = \sum_\alpha G^\alpha(g_0^\alpha, g_2^\alpha) = \sum_\alpha \left\langle v_1 \eta^\alpha(g_0^\alpha, g_2^\alpha) \right\rangle.
\]

**Proof.** Multiply equation (8a) by \(\partial_p \eta^\alpha(g_0^\alpha, g_2^\alpha)\), sum over \(p\) and integrate over \(v_1\):

\[
\frac{\partial H^\alpha}{\partial t} + \frac{\partial G^\alpha}{\partial x} = \frac{1}{\tau^\alpha} \left( (M_0^\alpha - g_0^\alpha) \partial_{g_0^\alpha} \eta^\alpha + (M_2^\alpha - g_2^\alpha) \partial_{g_2^\alpha} \eta^\alpha \right)
+ \frac{1}{\tau^\alpha \beta} \left( (\bar{M}^\alpha - g_0^\alpha) \partial_{g_0^\alpha} \eta^\alpha + (\bar{M}^\alpha - g_2^\alpha) \partial_{g_2^\alpha} \eta^\alpha \right).
\]
Then, summing over $\alpha$ gives:

$$
\frac{\partial H}{\partial t} + \frac{\partial G}{\partial x} = \sum_{\alpha} \frac{1}{\tau_\alpha} \left( \langle (M_0^\alpha - g_0^\alpha) \partial_{g_0^\alpha} \eta^\alpha + (M_2^\alpha - g_2^\alpha) \partial_{g_2^\alpha} \eta^\alpha \rangle \right)
+ \sum_{\alpha} \frac{1}{\tau_\alpha^\beta} \left( \langle (M_0^\alpha - g_0^\alpha) \partial_{g_0^\alpha} \eta^\alpha + (M_2^\alpha - g_2^\alpha) \partial_{g_2^\alpha} \eta^\alpha \rangle \right).
$$

(26)

Consider every term of the right-hand side consecutively.

$$
\langle (M_0^\alpha - g_0^\alpha) \partial_{g_0^\alpha} \eta^\alpha + (M_2^\alpha - g_2^\alpha) \partial_{g_2^\alpha} \eta^\alpha \rangle = \left( \langle (M_0^\alpha - g_0^\alpha) (\partial_{g_0^\alpha} \eta^\alpha (g_0^\alpha, g_2^\alpha) - \partial_{g_0^\alpha} \eta^\alpha (M_0^\alpha, M_2^\alpha)) \rangle \right)
+ \left( \langle (M_2^\alpha - g_2^\alpha) (\partial_{g_2^\alpha} \eta^\alpha (g_0^\alpha, g_2^\alpha) - \partial_{g_2^\alpha} \eta^\alpha (M_0^\alpha, M_2^\alpha)) \rangle \right)
+ \left( \langle (M_0^\alpha - g_0^\alpha) \partial_{g_0^\alpha} \eta^\alpha (M_0^\alpha, M_2^\alpha) \rangle \right)
+ \left( \langle (M_2^\alpha - g_2^\alpha) \partial_{g_2^\alpha} \eta^\alpha (M_0^\alpha, M_2^\alpha) \rangle \right).
$$

(27)

By convexity of $\eta^\alpha$, the sum of the first two terms of the right-hand side of (27) is negative. The remaining terms are

$$
\left( \langle (M_0^\alpha - g_0^\alpha) \partial_{g_0^\alpha} \eta^\alpha (M_0^\alpha, M_2^\alpha) \rangle \right) + \left( \langle (M_2^\alpha - g_2^\alpha) \partial_{g_2^\alpha} \eta^\alpha (M_0^\alpha, M_2^\alpha) \rangle \right)
= \left( \langle (M_0^\alpha - g_0^\alpha) \log \left( \frac{(M_0^\alpha)^2}{M_2^\alpha} \right) \rangle \right) - \left( \langle (M_2^\alpha - g_2^\alpha) \frac{M_0^\alpha}{M_2^\alpha} \rangle \right)
= \frac{1}{k_B T_\alpha} \left( \langle (M_0^\alpha - g_0^\alpha) \log (M_0^\alpha) \rangle - \frac{1}{k_B T_\alpha} \langle (M_2^\alpha - g_2^\alpha) \rangle \right)
= \frac{1}{k_B T_\alpha} \log \left( \frac{n^\alpha}{2\pi k_B T_\alpha \frac{m^\alpha}{2} \right)} \langle (M_0^\alpha - g_0^\alpha) \rangle + \left( \langle (M_0^\alpha - g_0^\alpha) \rangle \left( \frac{n^\alpha}{k_B T_\alpha \frac{m^\alpha}{2} \right)} - \frac{(\alpha T_\alpha)^2}{2k_B T_\alpha \frac{m^\alpha}{2}} \right)
- \frac{1}{k_B T_\alpha} \left( \langle (M_0^\alpha - g_0^\alpha) \rangle \frac{(\alpha T_\alpha)^2}{2k_B T_\alpha \frac{m^\alpha}{2}} \right)
= 0,
$$

according to (13). Then, consider the remaining last term in (26). Firstly, rewrite the term as:

$$
\left( \langle (M_0^\alpha - g_0^\alpha) \partial_{g_0^\alpha} \eta^\alpha + (M_2^\alpha - g_2^\alpha) \partial_{g_2^\alpha} \eta^\alpha \rangle \right) = \left( \langle (M_0^\alpha - g_0^\alpha) (\partial_{g_0^\alpha} \eta^\alpha (g_0^\alpha, g_2^\alpha) - \partial_{g_0^\alpha} \eta^\alpha (M_0^\alpha, M_2^\alpha)) \rangle \right)
+ \left( \langle (M_2^\alpha - g_2^\alpha) (\partial_{g_2^\alpha} \eta^\alpha (g_0^\alpha, g_2^\alpha) - \partial_{g_2^\alpha} \eta^\alpha (M_0^\alpha, M_2^\alpha)) \rangle \right)
+ \left( \langle (M_0^\alpha - g_0^\alpha) \partial_{g_0^\alpha} \eta^\alpha (M_0^\alpha, M_2^\alpha) \rangle \right)
+ \left( \langle (M_2^\alpha - g_2^\alpha) \partial_{g_2^\alpha} \eta^\alpha (M_0^\alpha, M_2^\alpha) \rangle \right).
$$

(28)

By convexity of $\eta^\alpha$, the sum of the first two terms is negative. Then, in favor of Proposition 2, by treating the remaining terms as previously, the dissipation of entropy (25) is obtained.

\[ \square \]

2.5. Scaling of the equations. The regime of interest is the hydrodynamic limit of this model, which corresponds to the scale of the phenomena described by the bi-temperature Euler equations (see [4]). To do so, it is necessary to perform a scaling on this model, in order to exhibit the behaviour of the equations when the macroscopic characteristic scale is much larger than the ones involved by the kinetic equations (collisional scale, Debye scale). Denote $t_0, L, v_{th}, q, m$ respectively the reference time, length, velocity, electric charge and particle mass. Denoting with a tilda the dimensionless quantities, it comes:

$$
\bar{t} = \frac{t}{t_0}, \quad \bar{x} = \frac{x}{L}, \quad \bar{v}_1 = \frac{v_1}{v_{th}}, \quad \bar{q} = \frac{q^\alpha}{q}, \quad \bar{m} = \frac{m^\alpha}{m}.
$$
From this scaling, define also the following dimensionless quantities:
\[
\bar{g}_p^\alpha = \frac{v_{th}}{n_0} g_p^\alpha, \quad \bar{E} = \frac{q_{t0}}{v_{th} m} E, \quad \bar{\rho} = \frac{\bar{\rho}}{q_{n0}}, \quad \bar{T} = \frac{k_B}{mv_{th}^2} T,
\]
where \( n_0 \) is the reference density.

Applying this scaling to our model, and removing the tildas for the sake of readability, it comes, for \( p = 0, 2 \) and \( \alpha = e, i \):
\[
\begin{align*}
\partial_t g_p^\alpha + v_1 \partial_x g_p^\alpha + \frac{q_0}{m^\alpha} E \partial_v v_i g_p^\alpha &= \frac{1}{\varepsilon^\alpha} (M_p^\alpha - g_p^\alpha) + \frac{1}{\varepsilon^\alpha\beta} (\bar{M}_p^\alpha - g_p^\alpha), \\
\partial_t E &= -\frac{j}{\beta^2}, \\
\partial_x E &= \frac{\bar{\rho}}{\beta^2},
\end{align*}
\]
where \( \varepsilon^\alpha = \frac{\varepsilon}{\varepsilon_0} \), \( \varepsilon^\alpha\beta = \frac{\varepsilon_0}{\varepsilon_0} \) and \( \beta = \frac{\lambda_{De}}{L} \), \( \lambda_{De} = (\varepsilon_0 k_B T_0) \frac{1}{T} \), with \( T_0 = \frac{m v_{th}^2}{k_B} \) the reference temperature. \( \lambda_{De} \) is the so-called Debye length, which is the characteristic length of charge separation in a plasma. In the regime of interest, this quantity is extremely small compared to the reference length of the domain \( L \). Similarly, the same-species mean free path related quantities \( \varepsilon^\alpha \) are small compared to the reference length \( L \). Hence, reaching the hydrodynamic limit consists in taking the limit of the model when \( \varepsilon^\alpha \) and \( \beta \) tend to 0. Notice that, however, \( \varepsilon^\alpha\beta \) is finite. In the limit, the Maxwell equations (29b-29c) become \( j = 0 \) and \( \bar{\rho} = 0 \), which can be rewritten as:
\[
\begin{align*}
u^\varepsilon &= \bar{u}^i, \\
n^\varepsilon &= Z \bar{n}^i.
\end{align*}
\]

The equations (30) are called the quasi-neutrality constraints. This form of the system of equations is the singular quasi-neutral limit of our initial model, where two evolution equations degenerate into algebraic relations. The purpose of this work is to derive an Asymptotic-preserving (AP) numerical scheme, that is to say with a stability condition that is independent of every non-dimensional small parameters \( \varepsilon^\alpha, \varepsilon^\alpha\beta \) and \( \beta \).

2.6. Hydrodynamic limit. In this section, we prove that the hydrodynamic limit of the scaled kinetic system (29) is the bi-temperature Euler system (1). Firstly, let us rewrite system (29) under the following form, where all parameters are \( \beta, \varepsilon^\alpha \) and \( \varepsilon^\varepsilon \) are equal to a common value \( \tau \):
\[
\begin{align*}
\partial_t g_p^\alpha + v_1 \partial_x g_p^\alpha + \frac{q_0}{m^\alpha} E \partial_v v_i g_p^\alpha &= \frac{1}{\varepsilon^\alpha} (M_p^\alpha - g_p^\alpha) + \frac{1}{\varepsilon^\alpha\beta} (\bar{M}_p^\alpha - g_p^\alpha), \\
\partial_t E &= -\frac{j}{\tau^2}, \\
\partial_x E &= \frac{\bar{\rho}}{\tau^2}.
\end{align*}
\]
As said in the previous section, the Maxwell equations give, in the limit \( \tau \rightarrow 0 \), the following equation:
\[
u^\varepsilon = \bar{u}^i.
\]
Hence, in the limit, electronic and ionic velocities are both equal to the mixture velocity \( u \). Besides, the BGK operators (31a) imply that the distribution functions \( g_p^\alpha \) are equal to their partial maxwellian. Then, for \( \alpha = e, i \) and \( p = 0, 2 \):
\[
\bar{g}_p^\alpha = M_p^\alpha.
\]
By integrating equation (31a) multiplied by \( m^\alpha \) for \( p = 0 \) w.r.t. \( v \) for each \( \alpha \), the following equation is obtained, for \( \alpha = e, i \):
\[
\partial_t \rho^\alpha + \partial_x (\rho^\alpha u^\alpha) = 0,
\]
and, by summing over \( \alpha \), the total mass equation is obtained:
\[
\partial_t \rho + \partial_x (p u) = 0.
\]
To obtain the equation on total momentum, multiply equation (31a) for \( p = 0 \) by \( m^αv^2 \), and integrate w.r.t. \( v \). It comes:

\[
(33) \quad \partial_t (\rho^α u^α) + \partial_x (m^α v^2 g^α_0) + q^α E (v \partial_v g^α_0) = \frac{\rho^α}{\varepsilon_α}(u - u^α).
\]

The third term on the left hand side can be rewritten as follows, by use of a integration by parts:

\[
q^α E (v \partial_v g^α_0) = -q^α E (g^α_0) = -E \bar{p}^α,
\]

where \( \bar{p}^α = q^α n^α \). The second term on the left hand side, in favor of (32), is equal to:

\[
\partial_x (m^α v^2 g^α_0) = \partial_x (m^α v^2 M^α_0) = \partial_x (\rho^α (u^α)^2 + p^α).
\]

By summing over \( α \), and in favor of conservation properties 2, one gets:

\[
\partial (\rho u) + \partial_x (\rho u^2 + p^e + p^i) - E \bar{p} = 0.
\]

Finally, compute the evolution equations on partial energies. For \( α = e, i \), multiply equation (31a) for \( p = 0 \) by \( m^α v^2 / 2 \), and integrate w.r.t. \( v \). It comes:

\[
\partial_t (m^α v^2 g^α_0) + \partial_x (m^α v^3 g^α_0) + q^α E (u^2 \partial_v g^α_0) = \frac{1}{\varepsilon_α}(M^α - g^α_0).
\]

To obtain an equation on total energy for species \( α \), integrate equation (31a) for \( p = 2 \) and sum with the previous equation. The following equation is obtained:

\[
\partial_t (m^α v^2 g^α_0 + g^α_0) + \partial_x (m^α v^3 g^α_0 + v g^α_0) + q^α E (v^2 \partial_v g^α_0) = \frac{1}{\varepsilon_α}(E - E^α).
\]

where the total plasma energy \( E \) is defined by:

\[
E = \sum_α E^α, \quad E^α = \frac{1}{2} \rho^α u^α + \rho^α \varepsilon^α.
\]

According to this definition, the right hand side can be rewritten as:

\[
\frac{1}{\varepsilon_α}(E - E^α) = \frac{3}{\varepsilon_α} m^α k_B (T^α - T^e).
\]

Then, in favor of quasi-neutrality properties and the definition of mixture temperature, we get

\[
T = \frac{T^α + T^e}{2}.
\]

The following expression \( \nu^{αβ} \) is then obtained:

\[
\frac{1}{\varepsilon_α}(E - E^α) = \frac{3}{\varepsilon_α} m^α k_B (T^i - T^e) = \nu^{αβ}(T^β - T^e).
\]

For the left hand side, the third term can be computed:

\[
q^α E (v^2 / 2 \partial_v g^α_0) = -q^α E (v g^α_0) = -E j^α,
\]

where \( j^α = q^α n^α u^α \) is the partial current of species \( α \). Finally, the second term of the left hand side, by use of (32), can be computed as well:

\[
\partial_x (m^α v^3 g^α_0 + v g^α_0) = \partial_x (m^α v^3 (M^α_0 + v M^α_2)) = u (\rho^α u^2 / 2 + \frac{3}{2} n^α k_B T^α + p^α).
\]

Hence, the following equation on total energy of species \( α \) is obtained:

\[
\partial_t E^α + \partial_x (u (E^α + p^α)) - j^α E = \nu^{αβ} (T^β - T^α).
\]

Finally, to compute the third term on the left hand side, rewrite equation (33) as follows:

\[
\partial_t u + u \partial_x u + \frac{\partial_x p^α}{\rho^α} - \bar{F} E = 0.
\]
Substracting this equation for \( \alpha = i \) to the equation for \( \alpha = e \), the generalized Ohm law is obtained, which writes:

\[
\frac{\partial_x p^e}{\rho^e} - \frac{\partial_x p^i}{\rho^i} = \frac{q^e n^e \rho}{\rho^e \rho^i} E.
\]

By denoting the following relation:

\[
j^e E = -j^i E,
\]

and by using the expression of \( E \) given by Ohm’s law, it comes:

\[
j^e E = q^e n^e u^e \left( \frac{\partial_x p^e}{\rho^e} - \frac{\partial_x p^i}{\rho^i} \right) = u^e \left( c^e \partial_x p^e - c^i \partial_x p^i \right).
\]

By injecting this expression into (34), the bi-temperature Euler system (1) is recovered.

2.7. Entropy inequality for hydrodynamic limits. In [4], the existence of an entropy-flux pair is shown for the bi-temperature Euler system (1). In this section, an entropy dissipation inequality in the limit is derived for the reduced model presented in Proposition 1.

Firstly, denote \( \mathcal{U} = (\rho, \rho u, e^e, e^i) \) the unknowns of the bi-temperature model and \((H_{\text{macro}}(\mathcal{U}), G_{\text{macro}}(\mathcal{U}))\) the corresponding macroscopic entropy-flux pair. Note the following properties regarding the link between kinetic and macroscopic entropies:

\[
H(M_0^e, M_2^e, M_0^i, M_2^i) = H_{\text{macro}}(\mathcal{U}), \quad G(M_0^e, M_2^e, M_0^i, M_2^i) = G_{\text{macro}}(\mathcal{U}).
\]

See [4] for further details on these properties.

**Proposition 6.** Let \( \mathcal{U} \) be a limit solution of the kinetic model from Proposition 1, i.e moments of solution distribution functions \((g_0^e, g_2^e)\) in the hydrodynamic limit. Then, the following inequality on entropy holds:

\[
\frac{\partial H_{\text{macro}}}{\partial t} + \frac{\partial G_{\text{macro}}}{\partial x} \leq -\frac{\nu_{\alpha}}{k_B T^e} (T^i - T^e)^2,
\]

with

\[
\nu_{\alpha} = \frac{3k_B n^\alpha n^\beta}{2\tau_{\alpha\beta} (n^\alpha + n^\beta)}.
\]

**Proof.** Multiply equation (8a) by \( \partial_x \eta^\alpha (g_0^e, g_2^e) \), sum over \( p \) and integrate over \( v_1 \). It comes, for \( \alpha = e, i \):

\[
\frac{\partial H^\alpha}{\partial t} + \frac{\partial G^\alpha}{\partial x} = \frac{1}{\varepsilon} \left( (M_0^\alpha - g_0^\alpha) \partial_{g_0^\alpha} \eta^\alpha (g_0^e, g_2^e) + (M_2^\alpha - g_2^\alpha) \partial_{g_2^\alpha} \eta^\alpha (g_0^e, g_2^e) \right)
+ \frac{1}{\varepsilon} \left( (\overline{M}_0^\alpha - \overline{g}_0^\alpha) \partial_{\overline{g}_0^\alpha} \eta^\alpha (g_0^e, g_2^e) + (\overline{M}_2^\alpha - \overline{g}_2^\alpha) \partial_{\overline{g}_2^\alpha} \eta^\alpha (g_0^e, g_2^e) \right).
\]

Using the entropy dissipation properties of the BGK operator shown in Section 2.4.2:

\[
\frac{\partial H^\alpha}{\partial t} + \frac{\partial G^\alpha}{\partial x} \leq \frac{1}{\varepsilon} \left( (M_0^\alpha - g_0^\alpha) \partial_{g_0^\alpha} \eta^\alpha (g_0^e, g_2^e) + (\overline{M}_0^\alpha - \overline{g}_0^\alpha) \partial_{\overline{g}_0^\alpha} \eta^\alpha (g_0^e, g_2^e) \right).
\]

By formally passing to limit \( \varepsilon \rightarrow 0 \), it comes \( g_0^\alpha = M_0^\alpha \), \( n^\alpha = n^i \) and \( u = u^e = u^i \). Hence, summing over \( \alpha \) and using (36), the inequality becomes

\[
\frac{\partial H_{\text{macro}}(\mathcal{U})}{\partial t} + \frac{\partial G_{\text{macro}}(\mathcal{U})}{\partial x} \leq \frac{1}{\varepsilon} \sum_{\alpha} \left( (M_0^\alpha - M_0^\alpha) \partial_{g_0^\alpha} \eta^\alpha (M_0^0, M_2^0) + (\overline{M}_0^\alpha - \overline{M}_0^\alpha) \partial_{\overline{g}_0^\alpha} \eta^\alpha (M_0^0, M_2^0) \right).
\]
The right hand side reads
\[ -\sum_{\alpha} \left( (\overline{m}_0^\alpha - m_0^\alpha) \frac{(v-u)^2}{2k_B T_2} \right) = -\sum_{\alpha} \frac{1}{k_B T^{\alpha}} \left( m^{\alpha} (\overline{m}_0^\alpha - m_0^\alpha) \frac{v^2}{2} \right). \]

The total right hand side is then equal to
\[ -\sum_{\alpha} \left( (\overline{m}_0^\alpha - m_0^\alpha) \partial_g \eta (m_0^\alpha, m_2^\alpha) + (\overline{m}_2^\alpha - m_2^\alpha) \partial_g \eta (m_0^\alpha, m_2^\alpha) \right) = -\sum_{\alpha} \frac{1}{k_B T^{\alpha}} \left( m^{\alpha} (\overline{m}_0^\alpha - m_0^\alpha) \frac{v^2}{2} \right). \]

According to the definition of $T$, it comes:
\[ T - T^{\alpha} = \frac{n^{\alpha}}{n^{\alpha} + n^\beta} (T^{\beta} - T^{\alpha}). \]

Hence,
\[ -\frac{3}{2}\varepsilon \sum_{\alpha} n^{\alpha} \frac{T - T^{\alpha}}{T^{\alpha}} = -\frac{3}{2\varepsilon} \frac{n^{\epsilon} n^{i} (T^{\epsilon} - T^i)^2}{n^{\epsilon} T^{i} - T^{i}}, \]
and we get (37).

3. Numerical treatment

In this section, numerical treatment of system (29) is addressed. The method to compute the evolution of the particle distribution is presented. For the sake of simplicity, in the Ampère equation and BGK operator, every non-dimensional quantity $\varepsilon^\epsilon, \varepsilon^i$ and $\beta$ are chosen equal to $\tau$. Reaching the hydrodynamic limit is then equivalent to the limit $\tau \rightarrow 0$.

3.1. Vlasov-BGK equations: general method. A time-splitting method is used on the Vlasov-BGK equations, in order to separate the Vlasov equations from the BGK operators. More precisely, the following equations are going to be consecutively solved:

(38) \[ \partial_t \rho_p^\alpha + v_1 \partial_x \rho_p^\alpha + \frac{q^\alpha}{m^\alpha} E \partial_{v_1} \rho_p^\alpha = 0, \]

the Vlasov-Maxwell equations for $\alpha = \epsilon, i$, coupled with the Ampère equation:

(39) \[ \partial_t E = -\frac{j}{\tau^2}. \]
Then, the effects of the inter-species collisions are computed:

\[ \partial_t g_{p}^\alpha = \frac{1}{\varepsilon_{\alpha \beta}} (M_p^\alpha - g_{p}^\alpha), \]

and finally, intra-species collisions:

\[ \partial_t g_{p}^\alpha = \frac{1}{\tau} (M_p^\alpha - g_{p}^\alpha), \]

where \( \tau \rightarrow 0 \).

A uniform discretization of the phase space is considered. Concerning the velocity discretization, in the present case, only static grids are considered, contrarily to [10]. A suitable subset of \( \mathbb{R} \) that contains the most part of the distribution function has to be chosen in order to perform correct numerical evaluations of the moments.

However, in this model, two particle distributions are considered, in order to follow the dynamics of both electrons and ions via a kinetic description. The major difference between these two species is the value of the mass of a particle that can have a ratio of the order of one thousand. Consequently, the distribution functions have very different supports. Thus, if only one velocity grid is employed, then this grid has to be large enough to encompass the largest distribution and has to be fine enough to describe the smallest one, which is numerically very costly. Hence, two velocity grids, one for each species, are considered. The boundaries of the domains are chosen as follows: for a given initial condition, let us denote the initial moments by \( n_0^\alpha, u_0^\alpha \) and \( T_0^\alpha \). For \( \alpha = e, i \), we have:

\[ V^\alpha = \min_x (u_0^\alpha (x)) - l^\alpha \max_x \sqrt{\frac{k_B T_0^\alpha (x)}{m^\alpha}}, \max_x (u_0^\alpha (x)) + l^\alpha \max_x \sqrt{\frac{k_B T_0^\alpha (x)}{m^\alpha}}, \]

where \( l^\alpha \) is chosen to fit the initial maxwellian distribution. Note that this grid is the same for all points in space and is constant in time.

Notations for the discretization are chosen as follows: \( \Delta x \) denotes the discretization step of the domain \([L_{\min}, L_{\max}] \subset \mathbb{R} \) in \( N_x \) cells. For \( \alpha = e, i \), the velocity domains \([V_{\min}^\alpha, V_{\max}^\alpha] \subset \mathbb{R} \) are both discretized using \( N_v \) cells and the discretization steps are denoted \( \Delta v^e \) and \( \Delta v^i \). Then, we use indices \( k \) and \( j \) to denote quantities computed at the \( k \)-th cell in space and \( j \)-th cell in velocity. Moreover, we denote \( \Delta t \) the time step and we use a superscript \( n \) to denote the quantity computed at time \( t^n = t^0 + n \Delta t \).

**Step 1: Computation of \( g \) and \( E \)**

The following discretization is performed, for \( \alpha = e, i \):

\[ g_{0,k,j}^{\alpha,n+1} = g_{0,k,j}^{\alpha,n} - \frac{\Delta t}{\Delta x} (\phi_{0,k+\frac{1}{2},j}^{\alpha,n} - \phi_{0,k-\frac{1}{2},j}^{\alpha,n}) - \frac{\Delta t}{2 \Delta v^\alpha} \frac{q^2 E_{k,n+1}}{m^\alpha} (g_{0,k,j+1}^{\alpha,n} - g_{0,k,j-1}^{\alpha,n}), \]

coupled with

\[ E_{k,n+1} = E_k^n - \frac{\Delta t}{\tau} j_{k,n+1}. \]

Note that the term \( \partial_t g_0 \) has been discretized using a centred scheme. For the space discretization, \( \phi_0 \) represented the flux function of the spatial discretization which will be specified in the next section.

Then, the following procedure is applied to obtain \( g_{2,n+1}^\alpha \):

\[ g_{2,k,j}^{\alpha,n+1} = g_{2,k,j}^{\alpha,n} - \frac{\Delta t}{\Delta x} (\phi_{2,k+\frac{1}{2},j}^{\alpha,n} - \phi_{2,k-\frac{1}{2},j}^{\alpha,n}) - \frac{\Delta t}{\Delta v^\alpha} (\psi_{2,k,j+\frac{1}{2}}^{\alpha,n} - \psi_{2,k,j-\frac{1}{2}}^{\alpha,n}), \]

where

\[ \phi_{2,k+\frac{1}{2},j}^{\alpha,n} = \frac{v^\alpha}{2} (g_{2,\alpha,n+1,j}^{\alpha} + g_{2,\alpha,n,j}^{\alpha}) - \frac{|v^\alpha|}{2} (g_{2,\alpha,n+1,j}^{\alpha} - g_{2,\alpha,n,j}^{\alpha}), \]

and finally, intra-species collisions:

\[ \partial_t g_{p}^{\alpha} = \frac{1}{\tau} (M_p^{\alpha} - g_{p}^{\alpha}), \]
and
\[ q_{2,k,j+1}^{\alpha,n} = \frac{q_0^{\alpha n+1}}{2m_0^{\alpha}} (g_{2,k,j+1}^{\alpha,n} + g_{2,k,j}^{\alpha,n}) - \frac{|q_0^{\alpha n+1}|}{2m_0^{\alpha}} (g_{2,k,j+1}^{\alpha,n} - g_{2,k,j}^{\alpha,n}). \]

Equations (43) are stable under the CFL condition:
\[ \Delta t \leq \min_{\alpha} \left( \frac{1}{\|V^{\alpha}\|_{\infty}} + \frac{1}{m_0^{\alpha} \Delta v_{\infty}} \right). \]

Moments of \( \bar{g}_p^{\alpha,n+1} \) and \( \bar{g}_p^{\alpha,n+1} \) are then computed using the midpoint quadrature formula. \( \forall k \in \{1, \ldots, N_x\} \):
\[
\bar{n}_k^{\alpha,n+1} = \sum_{j=1}^{N_v} \bar{g}_0^{\alpha,n+1} \Delta v^{\alpha},
\]
\[
\bar{u}_k^{\alpha,n+1} = \frac{1}{\bar{n}_k^{\alpha,n+1}} \sum_{j=1}^{N_v} \bar{g}_0^{\alpha,n+1} v_j \Delta v^{\alpha},
\]
\[
\bar{T}_k^{\alpha,n+1} = \frac{1}{3 \bar{n}_k^{\alpha,n+1} k_B} \left[ \sum_{j=1}^{N_v} m_0^{\alpha} \bar{g}_0^{\alpha,n+1} v_j^2 \Delta v^{\alpha} + 2 \sum_{j=1}^{N_v} \bar{g}_2^{\alpha,n+1} \Delta v^{\alpha} - m_0^{\alpha} \bar{n}_k^{\alpha,n+1} \bar{u}_k^{\alpha,n+1} \right].
\]

In order to compute the mixture Maxwellian distribution in the next step, mixture quantities must also be computed:
\[
\bar{n}_k^{\alpha,n+1} = \sum_{\alpha} n_k^{\alpha,n+1},
\]
\[
\bar{u}_k^{\alpha,n+1} = \frac{1}{\bar{n}_k^{\alpha,n+1}} \sum_{\alpha} \bar{u}_k^{\alpha,n+1},
\]
\[
\bar{T}_k^{\alpha,n+1} = \frac{1}{\bar{n}_k^{\alpha,n+1} k_B} \left[ \sum_{\alpha} \frac{1}{3} \rho_k^{\alpha,n+1} ((\bar{u}_k^{\alpha,n+1})^2 - (\bar{u}_k^{\alpha,n+1})^2) + \sum_{\alpha} \bar{n}_k^{\alpha,n+1} k_B \bar{T}_k^{\alpha,n+1} \right].
\]

Step 2: inter-species relaxation

We apply a backward Euler method to the equations (40):
\[
\bar{g}_p^{\alpha,n+1} = \frac{1}{\Delta t + \tau^{\alpha}} \left[ \Delta t \bar{M}_{p,k,j}^{\alpha,n+1} + \tau^{\alpha} \bar{g}_p^{\alpha,n+1} \right]
\]
which is unconditionally stable. \( \bar{M}_{p,k,j}^{\alpha,n+1} \) is the mixture Maxwellian defined by the moments \( \bar{n}_k^{\alpha,n+1}, \bar{u}_k^{\alpha,n+1} \) and \( \bar{T}_k^{\alpha,n+1} \), which are conserved through this step. Hence, the mixture Maxwellian is known beforehand, which allows this implication. Partial velocities and temperatures (denoted in the next section with a hat) are then recomputed using the previous formulae.

Step 3: intra-species relaxation

The next step consists in the relaxation towards Maxwellian equilibrium. Since the BGK operator preserves moments, we have the following properties: \( \forall k \in \{1, \ldots, N_x\}, \)
\[
\bar{n}_k^{\alpha,n+1} = n_k^{\alpha,n+1}, \quad \bar{u}_k^{\alpha,n+1} = u_k^{\alpha,n+1}, \quad \bar{T}_k^{\alpha,n+1} = T_k^{\alpha,n+1}.
\]

Similarly to the previous section, this property allows us to use an implicit scheme, the backward Euler method: \( \forall k \in \{1, \ldots, N_x\}, \forall j \in \{1, \ldots, N_v\} \) and \( p = 0, 2, \)
\[
\bar{g}_p^{\alpha,n+1} = \bar{g}_p^{\alpha,n+1} + \Delta t \left( M_{p,k,j}^{\alpha,n+1} - \bar{g}_p^{\alpha,n+1} \right),
\]
where
\[
M_{0,k,j}^{\alpha,n+1} = \frac{n_{k}^{\alpha,n+1}}{\sqrt{2\pi} k_{B}^{\alpha} \frac{T_{m}^{\alpha}}{m^{\alpha}}} \exp\left(-\frac{(v_{j} - u_{k}^{\alpha,n+1})^{2}}{2k_{B}^{\alpha}}\right),
\]
\[
M_{2,k,j}^{\alpha,n+1} = k_{B}^{\alpha} \frac{T_{m}^{\alpha}}{\tau_{k}^{\alpha}} M_{0,k,j}^{\alpha,n+1},
\]
which are expressed using quantities known at the end of step 2, according to the properties (47). Taking the limit \(\tau \to 0\) in (48) leads to:
\[g_{p,k,j}^{\alpha,n+1} = M_{p,k,j}^{\alpha,n+1}.\]

In the next section, enforcement of quasineutrality constraints (30) will be addressed. More precisely, specific spatial fluxes will be given and a reformulation of the Maxwell-Ampère equation will be performed in order to derive an Asymptotic-Preserving method, valid in the limit \(\tau \to 0\). Note that if these constraints are enforced through the hyperbolic step of the scheme (step 1), then they will be trivially preserved by both step 2 and 3. Hence, quasineutrality only has to be enforced through step 1.

3.2. Modified-viscosity upwind scheme. The space divergence is discretized explicitly, via the following modified upwind scheme:

\[
\phi_{0,k+\frac{1}{2},j}^{\alpha,n} = \frac{\Delta t}{\Delta x} \left( \phi_{0,k+\frac{1}{2},j}^{\alpha,n} - \phi_{0,k-\frac{1}{2},j}^{\alpha,n} \right) - \frac{\Delta t}{2\Delta x} q^{\alpha} \frac{E_{k}^{n+1}}{m^{\alpha}} (g_{0,k,j}^{\alpha,n} - g_{0,k,j-1}^{\alpha,n}),
\]

where \(V_{\max} = \max(\|V_{\min}\|, |V_{i}\|, |V_{\max}|, |V_{\min}|)\).

This method is the classical upwind scheme, where the numerical viscosity \(|v_{j}|\), that depends directly of the microscopic velocity, has been replaced by the smallest velocity that ensures stability for all equations. The discrete system writes as follows:

\[
\frac{\Delta t}{\Delta x} (\phi_{0,k+\frac{1}{2},j}^{\alpha,n} - \phi_{0,k-\frac{1}{2},j}^{\alpha,n}) - \frac{\Delta t}{2\Delta x} q^{\alpha} \frac{E_{k}^{n+1}}{m^{\alpha}} (g_{0,k,j}^{\alpha,n} - g_{0,k,j-1}^{\alpha,n}),
\]

and

\[
E_{k}^{n+1} = E_{k}^{n} - \frac{\Delta t}{\tau} j_{k}^{n+1}.
\]

This level of implicitness allows for a reformulation of the equation on the electric field. More precisely, the equation used to compute the electric field in (51) is not valid in the limit \(\tau \to 0\). To solve this problem, we follow the idea developped in [22]. By multiplying equation (50) by \(q^{\alpha} v_{j}^{\alpha} \Delta v^{\alpha}\) and summing over \(j\), it comes:

\[
q^{\alpha} \sum_{j} v_{j}^{\alpha} g_{0,k,j}^{\alpha,n+1} \Delta v^{\alpha} = q^{\alpha} \sum_{j} v_{j}^{\alpha} g_{0,k,j}^{\alpha,n} \Delta v^{\alpha} - q^{\alpha} \frac{\Delta t}{\Delta x} \sum_{j} v_{j}^{\alpha} (\phi_{0,k+\frac{1}{2},j}^{\alpha,n} - \phi_{0,k-\frac{1}{2},j}^{\alpha,n})
\]

\[
- \frac{\Delta t}{2} \sum_{j} v_{j}^{\alpha} (q^{\alpha})^{2} \frac{E_{k}^{n+1}}{m^{\alpha}} (g_{0,k,j+1}^{\alpha,n} - g_{0,k,j-1}^{\alpha,n})).
\]

According to the definition of the current
\[
j_{k}^{n+1} = \sum_{\alpha} q^{\alpha} \sum_{j} v_{j}^{\alpha} g_{0,k,j}^{\alpha,n} \Delta v^{\alpha},
\]

and summing over \(j\), it comes:

\[
j_{k}^{n+1} = \sum_{\alpha} q^{\alpha} \sum_{j} \Delta v^{\alpha} v_{j}^{\alpha} (\phi_{0,k+\frac{1}{2},j}^{\alpha,n} - \phi_{0,k-\frac{1}{2},j}^{\alpha,n})
- \frac{\Delta t}{2} \sum_{\alpha} \frac{(q^{\alpha})^{2}}{m^{\alpha}} \sum_{j} v_{j}^{\alpha} (g_{0,k,j+1}^{\alpha,n} - g_{0,k,j-1}^{\alpha,n}).
\]
By injecting this equation in (51), we get the following expression for the electric field:

\[
E_{k}^{n+1} = \frac{\tau E_{k}^{n} - \Delta t j_{k}^{n} + \frac{\Delta t^{2}}{2} \sum_{\alpha} q_{\alpha} \sum_{j} \Delta v_{\alpha} \psi_{j}^{\alpha} (\phi_{0,k+\frac{1}{2},j}^{\alpha,n} - \phi_{0,k-\frac{1}{2},j}^{\alpha,n})}{\tau - \frac{\Delta t^{2}}{2} \sum_{\alpha} \frac{q_{\alpha}^{2}}{m_{\alpha}} \sum_{j} v_{j}^{\alpha} (g_{0,k,j+1}^{n} - g_{0,k,j-1}^{n})},
\]

which is valid for all values of \( \tau \). In particular, we are interested in the limit \( \tau \to 0 \), which is:

\[
E_{k}^{n+1} = \frac{\frac{1}{\Delta x} \sum_{\alpha} q_{\alpha} \sum_{j} \Delta v_{\alpha} \psi_{j}^{\alpha} (\phi_{0,k+\frac{1}{2},j}^{\alpha,n} - \phi_{0,k-\frac{1}{2},j}^{\alpha,n})}{\sum_{\alpha} \frac{(q_{\alpha}^{n})^{2} \tau_{\alpha,k}^{n+1}}{m_{\alpha} \alpha,n}},
\]

where it has been assumed that \( j_{k}^{n} = 0 \). Such a choice for \( E_{k}^{n+1} \) ensures that \( j_{k}^{n+1} = 0 \), which is one of the two quasi-neutrality constraints (30). The choice of \( \phi_{0} \) (49) that has been made ensures that the other constraint \( \rho_{k}^{n+1} = 0 \) is also verified. More precisely, compute the discrete equation on \( \rho_{k}^{n+1} \). \( \forall k \in \{1, ..., N_{x}\} \), provided that

\[
\sum_{j} \left( \psi_{0,k,j+\frac{1}{2}}^{\alpha,n} - \psi_{0,k,j-\frac{1}{2}}^{\alpha,n} \right) = 0,
\]

the electric charge is

\[
\sum_{j} q_{\alpha} g_{0,k,j}^{n} \Delta v_{\alpha} = \sum_{j} q_{\alpha} g_{0,k,j}^{n} \Delta v_{\alpha} - \frac{\Delta t}{\Delta x} \sum_{j} (\phi_{0,k+\frac{1}{2},j}^{\alpha,n} - \phi_{0,k-\frac{1}{2},j}^{\alpha,n}) \Delta v_{\alpha}.
\]

So, the following equation on the total electric charge is obtained:

\[
\bar{\rho}_{k}^{n+1} = \rho_{k}^{n} - \frac{\Delta t}{\Delta x} \sum_{\alpha} q_{\alpha} \sum_{j} (\phi_{0,k+\frac{1}{2},j}^{\alpha,n} - \phi_{0,k-\frac{1}{2},j}^{\alpha,n}) \Delta v_{\alpha}.
\]

According to (49), we get:

\[
\bar{\rho}_{k}^{n+1} = \rho_{k}^{n} - \frac{\Delta t}{\Delta x} \sum_{\alpha} q_{\alpha} \sum_{j} (\phi_{0,k+\frac{1}{2},j}^{\alpha,n} - \phi_{0,k-\frac{1}{2},j}^{\alpha,n}) \Delta v_{\alpha}.
\]

By rearranging the terms, we obtain:

\[
\bar{\rho}_{k}^{n+1} = \rho_{k}^{n} - \frac{\Delta t}{2 \Delta x} (j_{k+1}^{n} - j_{k-1}^{n}) + \frac{\Delta t}{2 \Delta x} \sum_{\alpha} q_{\alpha} \sum_{j} \left| V_{\max} (g_{0,k+1,j}^{n} - 2g_{0,k,j}^{n} + g_{0,k-1,j}^{n}) \right| \Delta v_{\alpha}.
\]

By assuming that \( j_{k+1}^{n} = j_{k-1}^{n} = 0 \), it comes:

\[
\bar{\rho}_{k}^{n+1} = \frac{\Delta t}{2 \Delta x} \left( V_{\max} (\bar{\rho}_{k+1}^{n} - 2\bar{\rho}_{k}^{n} + \bar{\rho}_{k-1}^{n}) = 0.
\]

Hence, quasineutrality is conserved at all time. However, applying this modification to our scheme amounts to overevaluating the numerical viscosity by its highest value. This will create an important amount of dissipation in the numerical solutions obtained with this scheme. Hence, the discretization of the space domain needs to be very precise to compensate this viscosity, which will result in a longer computation time.
4. Numerical results

In this section, numerical resolution of Riemann problems with our method is presented. Such problems are characterized by initial conditions constituted of two constant states separated by a discontinuity. In the study of such problems, an exact solution can be provided for the conserved quantities (density, momentum and total energy). However, it is not true for non-conserved quantities, and no reference results exist.

Comparative results obtained by an HLL-type scheme and a Suliciu relaxation method on the bi-temperature Euler system are displayed along the solutions from the kinetic scheme that has been derived in this paper (see [4] for more details on these methods).

The spatial domain is chosen as the interval \([0,1]\), supplemented with homogeneous Neumann boundary conditions. Physical constants \((k_B, Z, q_i)\) are chosen equal to one. All test cases are done in the hydrodynamic limit \(\tau \to 0\). Concerning the choice of \(l^\alpha\) in (42), a convergence test has been performed on the initial conditions of the test cases in order to obtain a appropriate value that would optimize computational cost. Hence, in all test cases, \(l^\alpha = 8\), for \(\alpha \in \{e,i\}\).

A mass ratio of 10 \((m_e = 0.1\) and \(m_i = 1)\) is considered. The greater the mass ratio, the greater the value of \(V_{\text{max}}\) in (49). Hence, a high mass ratio implies an important amount of numerical viscosity. Then, we are restricting ourselves to a mass ratio of 10, which is already demanding in terms of number of discretization points. However, in a proper physical context, mass ratios of the order of 1000 should be considered.

4.1. Shock tube with identical initial temperatures. The Riemann problem of the shock tube with identical ionic and electronic temperature is given by the following initial condition:

\[
\begin{align*}
\rho(t_0, x) &= 1, & u^\alpha(t_0, x) &= 0, & T^\alpha(t_0, x) &= 1 \quad \text{if } x \in [0, 0.5], \\
\rho(t_0, x) &= 0.125, & u^\alpha(t_0, x) &= 0, & T^\alpha(t_0, x) &= 1 \quad \text{if } x \in [0.5, 1].
\end{align*}
\]

Parameters are chosen as \(N_x = 120000\), \(N_v = 40\) and \(l = 8\). The inter-species collision relaxation time is \(\tau^{ei} = 0.1\). Results are computed at time \(t = 0.1\). For fluid methods, 4000 discretization points are used.

The results are provided in figure 1 and 2.

![Figure 1](image1.png)

**Figure 1.** Density and velocity solutions of shock tube test case with a mass ratio of 10 with 120000 space points, 40 velocity points and a domain length of 8

The quantities described by conservative equations (densities and velocities) are consistent with the fluid methods HLL and Suliciu. Moreover, quasi-neutrality is shown to be achieved, since electronic and ionic densities are superposed, as well as electronic and ionic velocities.
For the partial temperatures presented in figure 2, the intermediate state reached between the rarefaction wave and the contact discontinuity is identical for all schemes, which is expected since this part of the solution is smooth. However, the states obtained between the contact discontinuity and the shock are different for every scheme. Figures 3 display a zoom on this part of the solution. The results can be seen to be different for all three methods. The behaviour of the different schemes should be investigated further.

On figure 4, we display the electric field at the final time $t = 0.1$, obtained in two different ways. The first method is the electric field obtained directly by the kinetic scheme, by formula (53). The second one corresponds to the electric field computed by the generalized Ohm’s law (35), computed thanks to the macroscopic quantities obtained by the scheme. As can be seen, the behaviour of the scheme is completely consistent with the hydrodynamic limit.

4.2. Shock tube with different initial temperatures. This test case is similar to the previous one, the only difference being the initial left ionic and electronic temperatures. The following
Figure 4. Electric field obtained by the kinetic scheme and by Ohm’s law for the Sod tube test case with identical initial temperatures.

Initial conditions are:
\[
\begin{align*}
\rho(t_0, x) &= 1, \quad u^\alpha(t_0, x) = 0, \quad T^\alpha(t_0, x) = 1 \quad \text{if } x \in [0, 0.5], \\
\rho(t_0, x) &= 0.125, \quad u^\alpha(t_0, x) = 0, \quad T^e(t_0, x) = 2 \quad T^i(t_0, x) = 3 \quad \text{if } x \in [0.5, 1].
\end{align*}
\]

Parameters are chosen as \(N_x = 120000\), \(N_v = 40\) and \(l = 8\). The inter-species collision relaxation time is \(\tau^{ei} = 0.1\). Results are computed at time \(t = 0.1\). The results are provided in figure 5 and 6. The behaviour of the result is very similar to the previous one. Quasi-neutrality is achieved and temperatures exhibit different jump relations across discontinuities for each method.

Figure 5. Density and velocity solutions of shock tube test case with a mass ratio of 10 with 120000 space points, 40 velocity points and a domain length of 8.

On figure 8, similarly to the previous test case, the behaviour of the electric field is shown. The results are consistent with the hydrodynamic limit.

4.3. Rarefaction wave. This test case is constituted of two rarefaction waves going in opposite directions. It is given by the following data:
\[
\begin{align*}
\rho(t_0, x) &= 1, \quad u^\alpha(t_0, x) = -1, \quad T^\alpha(t_0, x) = 1 \quad \text{if } x \in [0, 0.5], \\
\rho(t_0, x) &= 1, \quad u^\alpha(t_0, x) = 1, \quad T^\alpha(t_0, x) = 1 \quad \text{if } x \in [0.5, 1].
\end{align*}
\]
We have $N_x = 120000$, $N_v = 40$ and $l = 8$. The inter-species collision relaxation time is $\tau^{ei} = 0.1$. Results are computed at time $t = 0.1$.

The result are provided in figure 9 and 10. In this test case, the solution is smooth. There are no shock waves, and then, all methods are expected to give the same results. Figure 11 shows that this is what happens. Figure 12 shows the behaviour of the electric field.

4.4. Shock wave. This test case is constituted of two shock waves. It is given by the following data:

$$\begin{align*}
\rho(t_0, x) = 1, u^\alpha(t_0, x) = 1, T^\alpha(t_0, x) &= 1 & \text{if } x \in [0, 0.5], \\
\rho(t_0, x) = 1, u^\alpha(t_0, x) = -1, T^\alpha(t_0, x) &= 1 & \text{if } x \in [0.5, 1].
\end{align*}$$

We have $N_x = 120000$, $N_v = 40$ and $l = 8$. The inter-species collision relaxation time is $\tau^{ei} = 0.1$. Results are computed at time $t = 0.1$.

The results are provided in Figure 13 and 14. Quasi-neutrality is achieved on the conserved quantities. The solution contains two shock waves and different jump relations can be observed for each numerical method. In Figure 15, a zoom is performed on the constant values reached.
Figure 8. Electric field obtained by the kinetic scheme and by Ohm's law for the Sod tube test case with different initial temperatures.

Figure 9. Density and velocity solution of a rarefaction wave test case with a mass ratio of 10 with 120000 space points, 40 velocity points and a domain length of 8.

Figure 10. Temperature solutions of a rarefaction wave test case with a mass ratio of 10 with 120000 space points, 40 velocity points and a domain length of 8.
by each method between the two shocks. This shows the different behaviour between the three methods. The electric field is displayed on figure 16.

4.5. Sod tube with realistic mass ratio. In this last test case, we consider the same initial data as in the test case from section 4.2, but the electronic mass is chosen equal to $m^e = 0.001$, which is close to a realistic mass ratio. The initial data are:

\[
\begin{align*}
\rho(t_0, x) &= 1, & \alpha(t_0, x) &= 0, & T^\alpha(t_0, x) &= 1 & \text{if } x \in [0, 0.5], \\
\rho(t_0, x) &= 0.125, & \alpha(t_0, x) &= 0, & T^e(t_0, x) &= 2 & T^i(t_0, x) = 3 & \text{if } x \in (0.5, 1].
\end{align*}
\]

Parameters are chosen as $N_x = 100000$, $N_v = 40$ and $l = 8$. The inter-species collision operator is ignored for this test case, which amounts to $\nu^{ij} = 0$ in the bi-temperature Euler equations. Results are computed at time $t = 0.05$. The result are provided in figure 17 and 18. With the absence of a temperature relaxation term, the jump relations for this test case are the same for our method and the HLL method. The electric field is displayed on figure 20 and can be seen to be perfectly consistent with the hydrodynamic limit.
5. Conclusion

In this article, a kinetic numerical method able to provide reference results for a non-conservative hyperbolic system, the bi-temperature Euler system, is proposed. This method is able to enforce all the desired properties that are relevant for comparisons with methods applied directly to the hyperbolic system. Such methods lack an unambiguous definition of the non-conservative products and exhibit different Rankine-Hugoniot relation when the solution contains shock waves. However, the proposed method introduces a important amount of numerical viscosity, which renders the method costly. A second-order method able to circumvent such a drawback is a current on-going work, which should be the subject of a subsequent article.

References


Figure 15. Electronic and ionic temperature of a shock wave test case with a mass ratio of 10 with 120000 space points, 40 velocity points and a domain length of 8

Figure 16. Electric field obtained by the kinetic scheme and by Ohm’s law for the double shock test case

FIGURE 17. Density and velocity solutions of shock tube test case with a mass ratio of 1000 with 100000 space points, 40 velocity points and a domain length of 8

FIGURE 18. Electronic and ionic temperatures of a shock tube test case with a mass ratio of 1000 with 100000 space points, 40 velocity points and a domain length of 8


A KINETIC APPROACH OF THE BI-TEMPERATURE EULER MODEL

Figure 19. Different jump relations of electronic and ionic temperatures for a shock tube test case with a mass ratio of 1000 with 100000 space points, 40 velocity points and a domain length of 8

Figure 20. Electric field obtained by the kinetic scheme and by Ohm’s law for a Sod tube test with different initial temperatures