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Burnett description for plane Poiseuille flow

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Two recent works have shown that at small Knudsen number ($K$) the pressure and temperature profiles in plane Poiseuille flow exhibit a different qualitative behavior from the profiles obtained by the Navier-Stokes equations. Tij and Santos [J. Stat. Phys. 76, 1399 (1994)] used the Bhatnagar-Gross-Kook model to show that the temperature profile is bimodal and the pressure profile is nonconstant. Malek-Mansour, Baras, and Garcia [Physica A 240, 255 (1997)] qualitatively confirmed these predictions in computer experiments using the direct simulation Monte Carlo method (DSMC). In this paper we compare the DSMC measurements of hydrodynamic variables and non-equilibrium fluxes with numerical solutions of the Burnett equations. Given that they are in better agreement with molecular-dynamics simulations [E. Salomons and M. Mareschal, Phys. Rev. Lett. 69, 269 (1992)] of strong shock waves than Navier-Stokes [F. J. Uribe, R. M. Velasco, and L. S. García-Colín, Phys. Rev. Lett. 81, 2044 (1998)], and that they are in better second order in Knudsen number suggests that the Burnett equations may provide a better description for large $K$. We find that for plane Poiseuille flow the Burnett equations do not predict the bimodal temperature profile but do recover many of the other anomalous features (e.g., nonconstant pressure and nonzero parallel heat flux). [S1063-651X(99)02410-1]

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I. INTRODUCTION

The Navier-Stokes equations provide a sound and robust theoretical method to study flows at the macroscopic level which have innumerable applications. The complexity of the equations is due to the fact that they are nonlinear except for simple cases where, due mainly to the symmetries involved, they reduce to linear equations that can be solved in a closed form. The Navier-Stokes equations can be derived from strictly macroscopic arguments or from a kinetic model, such as the Boltzmann equation for dilute gases; however, their derivation assumes both local equilibrium hypotheses and the small gradient approximation, which allows one to formulate the constitutive equations needed to close the conservation equations. In situations where large gradients exist, as in a shock wave or in a boundary layer, these equations are expected to fail, as has been pointed out in the past. Furthermore, being a macroscopic theory their validity is also restricted to situations in which the Knudsen number ($K$) is small and a crucial point is to know, for a given problem, the range of Knudsen numbers where they remain valid. Another important parameter is the Reynolds number ($Re$), which in some cases determines when the solutions cease to be laminar and the flow becomes turbulent.

The Navier-Stokes equations can, in principle, be solved as an initial value problem. Usually stick boundary conditions are used, but when the Knudsen number is not small ($K>10^{-2}$) the slip at the boundaries must be taken into account [1]. There exist simple examples in which the solutions to the stationary Navier-Stokes equations are infinite (countable) even for a fixed Reynolds number [2], so even when only one solution exists, as in plane Poiseuille flow with stick boundary conditions, one needs to perform a stability analysis of the solutions [3].

Recently Tij and Santos [4] showed that a perturbative solution to the Bhatnagar-Gross-Kook (BGK) model predicted qualitatively different temperature and pressure profiles from those predicted by the Navier-Stokes equations. Malek et al. [5] confirmed this result using the direct simulation Monte Carlo (DSMC) method to model plane Poiseuille flow of a dilute hard sphere gas for $Re=10$ and $K=1/10$. Malek et al. also solved the Navier-Stokes equations numerically and showed that slip corrections could not explain the discrepancy. The results obtained by Tij and Santos are in better quantitative agreement with the DSMC data if the Chapman-Cowling expressions for the transport coefficients are used instead of the those derived from the BGK model, which give an incorrect Euken factor. While the Euken factor is in contradiction with the results from both the Chapman-Enskog method and experiment, it should be pointed out that the BGK model is able to reproduce qualitative features in situations far from equilibrium where the Chapman-Enskog method is not applicable, and so the model is useful as a guide in such situations.

To the authors’ knowledge, Poiseuille flow is the first scenario in which the Navier-Stokes equations have been shown to be susceptible to significant improvement for a flow with relatively small gradients. Since Uribe et al. [6] showed that the Burnett equations can provide an accurate description for strong shock waves, it is then a natural step to compare the Burnett equations with DSMC simulations of Poiseuille flow, and that is what we do here. The next four sections present the theoretical development of the conservation equations, the Navier-Stokes equations, the Burnett...
equations, and the Burnett fluxes. The DSMC algorithm is outlined briefly in Sec. VI; simulation results are presented and discussed in Secs. VII and VIII. Section IX contains final remarks and a discussion of future work.

II. CONSERVATION EQUATIONS

We consider the flow between two infinite and stationary plane parallel walls located at \( y = \pm L/2 \). The hydrodynamic velocity \( \mathbf{e}_0 = \mathbf{u} = u(x) \mathbf{i} \) in the laminar, stationary case (unless stated otherwise, we adopt the notation used by Chapman and Cowling [7]). Following Landau and Lifshitz [8] we assume that all the relevant quantities can only depend on the \( y \) coordinate, in particular \( \mathbf{e}_0 = u(y) \mathbf{i} \). Noting that, for any function, \( f(y) \), \( (\mathbf{e}_0 \cdot \nabla) f(y) = 0 \), it follows that the conservation equations reduce to

\[
\begin{align*}
0 &= \nabla \cdot (\rho(y) u(y) \mathbf{i}), \\
0 &= -\frac{1}{\rho(y)} \frac{dP_{xx}}{dy} + a, \\
0 &= -\frac{1}{\rho(y)} \frac{dP_{yy}}{dy}, \\
0 &= \frac{dP_{yz}}{dy}, \\
0 &= P_{yx} \frac{du(y)}{dy} + \frac{dq_y(y)}{dy},
\end{align*}
\]

where \( \rho(y) \) is the mass density, \( P(y) \) the pressure tensor, \( q(y) \) the heat flux, and \( a \) the acceleration from a constant, external body force (e.g., gravity) acting in the \( x \) direction. Equation (1) is the mass conservation equation, which is automatically satisfied. Equations (2), (3), and (4) are the components of Newton’s second law (momentum conservation), and Eq. (5) is the energy conservation equation. Notice that we can conclude that the components \( P_{yy} \) and \( P_{yz} \) of the pressure tensor are constant, a result that does not depend on the constitutive relations.

For the discussion that follows it is important to keep in mind that the results of this section depend on the three conditions: (1) \( \mathbf{u}_x = 0 \). (2) \( \mathbf{u}_x = 0 \). (3) All the physical quantities depend only on \( y \). The last condition will be referred to as Landau’s symmetry argument.

In a dilute gas the pressure \( p(y) \) and temperature \( T(y) \) are related by the ideal gas equation of state \( p = \rho kT/m \), where \( m \) is the particle mass and \( k \) is Boltzmann’s constant. To have a closed system in the variables \( u(y) \), \( p(y) \), and \( T(y) \), one needs to introduce constitutive relations. Since the contributions up to second order in the Knudsen number for the pressure tensor and heat flux include the Navier-Stokes terms (first order in the Knudsen number), these terms are developed in Sec. III.

III. NAVIER-STOKES REGIME

The Navier-Stokes expressions for the pressure tensor \( P^{NS} \) and the heat flux \( q^{NS} \) are given by

\[
\begin{align*}
P^{NS} &= \rho I - 2\mu \nabla \mathbf{e}_0, \\
q^{NS} &= -\lambda \nabla T,
\end{align*}
\]

where \( \mu \) is the shear viscosity, \( \lambda \) the thermal conductivity, and \( I \) the unit tensor. The double overbar denotes the symmetric tensor, and the circle means the corresponding traceless tensor. For stationary, plane Poiseuille flow, these constitutive relations simplify to

\[
\begin{align*}
P^{NS}_{yx} &= -\frac{d u(y)}{dy}, \\
P^{NS}_{yy} &= p(y), \\
q^{NS}_{yy} &= -\frac{d T(y)}{dy},
\end{align*}
\]

For a dilute gas of rigid spheres, the transport coefficients are [7]

\[
\mu = \frac{5\epsilon_\mu}{16\sigma^2} \left( \frac{mkT(y)}{\pi} \right)^{1/2}, \quad \lambda = \frac{75\epsilon_\lambda}{64\sigma^2} \left( \frac{k^2 T(y)}{\pi m} \right)^{1/2},
\]

where \( \sigma \) is the particle diameter. While often the transport coefficients are approximated by the first Sonine expansion (\( \epsilon_\mu = \epsilon_\lambda = 1 \)), for rigid spheres more precise values (\( \epsilon_\mu = 1.016034 \) and \( \epsilon_\lambda = 1.02513 \)) are known [7].

Substitution of Eqs. (8) and (9) into Eqs. (1)–(5) leads us to

\[
\begin{align*}
10\epsilon_\mu k^2 T(y) \frac{d^2 u}{dy^2} &= -5\epsilon_\mu k^2 T(y) \frac{dT}{dy} \frac{du}{dy} \\
&\quad -32\rho \sigma^2 \epsilon_\lambda (\pi mkT(y))^{1/2}, \\
30\epsilon_\lambda k^3 T(y) \frac{dT}{dy^2} &= -8\epsilon_\mu k^2 T(y) \left( \frac{du}{dy} \right)^2 -15\epsilon_\lambda k^4 \left( \frac{dT}{dy} \right)^2.
\end{align*}
\]

Note that in the Navier-Stokes regime the pressure is constant, so we have two second order nonlinear equations for \( u \) and \( T \). The boundary conditions for these equations are discussed in Sec. IV.

IV. BURNETT REGIME

The second order Knudsen number corrections for the stress tensor \( P^{(2)} \) and the heat flux \( q^{(2)} \) are given by [7]
\[ \mathbf{P}^{(2)} = \omega_1 \mu_1^2 \Delta \mathbf{e} + \omega_2 \frac{\mu_2^2}{p} D_0 \frac{D_t}{D_t} \mathbf{e} - 2 \omega_3 \frac{\mu_3^2}{\rho T} \mathbf{e} + \omega_4 \frac{\mu_4^2}{p \rho T} \mathbf{e} + \omega_5 \frac{\mu_5^2}{\rho T} \mathbf{e} + \omega_6 \frac{\mu_6^2}{\rho T} \mathbf{e} \]

\[ q^{(2)} = \theta \frac{\mu_1^2}{\rho T} \Delta \nabla T + \theta_2 \frac{\mu_2^2}{\rho T} D_0 \frac{D_t}{D_t} \mathbf{e} - \theta_3 \frac{\mu_3^2}{\rho T} \mathbf{e} + \theta_4 \frac{\mu_4^2}{\rho T} \mathbf{e} + \theta_5 \frac{\mu_5^2}{\rho T} \mathbf{e} T \mathbf{e}. \]

where the numerical values for the \( \omega \)'s and \( \theta \)'s are given below and [7]

\[ e = -c_0, \Delta = -c_0, \]

\[ D_0 \frac{D_t}{D_t} (T) = -\frac{2}{3} (T \cdot c_0) - (c_0 \cdot T), \]

\[ D_0 \frac{D_t}{D_t} (\mathbf{e}) = \left( a - \frac{1}{\rho} \right) - c_0. \]

For plane Poiseuille flow, we are led to

\[ \mathbf{P}_{y}^{(2)} = 0, \quad \mathbf{P}_{z}^{(2)} = 0, \]

\[ \mathbf{q}_{y}^{(2)} = \frac{2}{3} \frac{\mu_1^2}{\rho(y) p(y)} \frac{dp}{dy} \left( \frac{dy}{dy} \frac{dy}{dy} + \frac{2}{3} \frac{\mu_1^2}{\rho(y) p(y)} \frac{d^2 p}{dy^2} \right) - \frac{2}{3} \frac{\mu_2^2}{\rho(y) p(y)}, \]

\[ \times \left( \frac{d u}{dy} \right)^2 + \frac{2}{3} \frac{\mu_1^2}{\rho(y) p(y)} \frac{d^2 T}{dy^2} + \frac{2}{3} \frac{\mu_2^2}{\rho(y) p(y)} \frac{d^2 T}{dy^2}, \]

\[ \times \frac{dp}{dy} \frac{d T}{dy} + \frac{2}{3} \frac{\mu_3^2}{\rho(y) T(y)} \frac{d T}{dy} + \frac{1}{12} \frac{\mu_4^2}{\rho(y) T(y)} \frac{d u}{dy}, \]

\[ q_{z}^{(2)} = \frac{1}{2} \frac{\mu_1^2}{\rho(y) p(y)} \frac{du}{dy} + \frac{1}{2} \frac{\mu_4^2}{\rho(y) T(y)} \frac{du}{dy} + \frac{3}{2} \frac{\mu_5^2}{\rho(y) T(y)} \frac{d T}{dy}, \quad q_{x}^{(2)} = 0, \quad q_{y}^{(2)} = 0. \]

Note that there is a heat flow along the \( x \) direction, but it does not make a contribution to the conservation equations, the reason being that in the energy conservation equation one has to evaluate the divergence of the heat flow, and only the \( y \) component of the heat flux can contribute.

The \( x \) component in the momentum conservation is the same as the Navier-Stokes result because \( \mathbf{P}_{x}^{(2)} = 0 \). Similarly the \( z \) component of momentum conservation is not altered, so only the \( y \) component of momentum conservation is altered in the Burnett regime. Since this component must be a constant, we see from previous equations that the result is a second order ordinary differential equation for \( p \). In the previous equation for \( \mathbf{P}_{y} \), there appears a second derivative for the temperature, which can be obtained from Eq. (10). Using this result, we are led to solve the following reduced second order system of ordinary differential equations:

\[ \frac{d^2 u}{ds^2} = -\frac{1}{2T^*(s)} \frac{dt}{ds} \frac{du}{ds} + \frac{b_0 p^*}{T^* T^*(s)}, \]

\[ \frac{d^2 T^*}{ds^2} = -8 \frac{c_\mu}{150 \lambda} \left( \frac{du}{ds} \right)^2 - \frac{1}{2T^*(s)} \left( \frac{d t}{ds} \right)^2, \]

\[ \frac{d^2 p^*}{ds^2} = \frac{b_1 p^*}{T^* T^*(s)} + \frac{b_2 p^*}{T^* T^*(s)} - \frac{b_1}{T^*(s)} \frac{p}{p^*} \left( \frac{dp}{ds} \right)^2 \]

\[ - \frac{1}{T^*(s)} \frac{dp}{ds} \frac{d^2 T^*}{ds^2} - \frac{2}{T^*(s)} \frac{dp}{ds} \left( \frac{du}{ds} \right)^2 \]

\[ - \frac{8}{150} \frac{p^*}{s^*} \frac{c_\mu}{T^* T^*(s)} \left( \frac{du}{ds} \right)^2 - \frac{p^*}{s^*} \frac{\omega_3}{2} \left( \frac{dT^*}{ds} \right)^2 \]

\[ + \frac{\omega_4}{\omega_2 T^*(s)} \frac{dp}{ds} \frac{d^2 T^*}{ds^2} + \frac{\omega_5 p^*}{\omega_2 T^* T^*(s)} \left( \frac{dT^*}{ds} \right)^2 \]

\[ + \frac{\omega_6 p^*}{4 \omega_2 T^* T^*(s)} \left( \frac{du}{ds} \right)^2, \]

where \( s = y/L, p^* = p(y)/p(0), T^* = T(y)/T_R \), with \( T_R \) the temperature of the walls and \( u^*(y) = u(y)/(2k T_R/m)^{1/2} \). The coefficients are

\[ b_0 = -\frac{8}{5} \frac{L^2 a \sqrt{2 \pi p(0) m a^2}}{c_\mu k^2 T_R^2}, \quad b_1 = -\frac{384}{25} \frac{L^2 p(0)^2 \pi \sigma^4}{\omega_2 c_\mu T^2 T_R^2}, \]

\[ b_2 = -\frac{L^2 d^2 p}{p(0) T_R^2}. \]

for symmetric profiles. Though the physical system is geometrically symmetry about the \( x \) axis, this does not guarantee a symmetric profile since spontaneous symmetry breaking may occur [2]. However, the profiles measured in our DSMC computer experiments are in good agreement with this assumed symmetry.

In terms of the variables: \( Y_1(s) = u^*(s), Y_2(s) = T^*(s), Y_3(s) = p^*(s), Y_4(s) = du/ds, Y_5(s) = (dT*/ds)(s), \) and \( Y_6(s) = (dp*/ds)(s) \), the previous system of three second order differential equations can be expressed as a system of six coupled first order differential equations, \( \dot{Y} = F(Y) \), where the dot denotes the derivative with respect to \( s \). The vector field \( F \) is given by
\[
F_1(Y) = Y_4, \quad F_2(Y) = Y_5, \quad F_3(Y) = Y_6, \\
F_4(Y) = -\frac{1}{2} Y_4 Y_5 + b_0, \quad F_5(Y) = -\frac{1}{2} Y_5 Y_6 + b_0, \\
F_6(Y) = b_1 Y_4^2 + b_2 Y_5^2 - b_3 Y_4^3 + b_4 Y_5^3 - b_5 Y_4 Y_5 - \frac{2}{3} \frac{Y_4 Y_5}{Y_2}.
\]

where for simplicity an explicit \(s\) dependence is omitted.

To have a well posed problem requires the specification of initial or boundary conditions. For the Burnett equations, several issues regarding boundary conditions are still unresolved. Rather than looking at this point in a general setting, it is important to know first if the Burnett equations can indeed reproduce the features observed in problems for which the Navier-Stokes equations are susceptible to improvement, and this is what we would like to test here for plane Poiseuille flow.

To avoid the difficulty of specifying boundary conditions at \(y = \pm L/2\) \((s = \pm 1/2)\), we use the centerline \((s = 0)\) profile values measured in the DSMC simulations as the initial values for the previous first order system [Eq. (17)]. The explicit initial conditions used are

\[ u^*(0) = u_0, \quad T^*(0) = T_0, \quad p^*(0) = 1, \]

with the derivatives at \(s = 0\) equal to zero and the values \(u_0, \quad p(0), \quad p(0), \quad T_0, \quad T_0\) taken from the DSMC data. The dynamical system [Eq. (17)], is given in terms of quotients of differentiable functions, so that it is differentiable in \(U = R \times R^+ \times R^+ \times R^3\), \(R\) being the real numbers and \(R^+\) the positive real numbers. In particular, it is locally Lipshitz in that set [9], so that the usual local existence and uniqueness theorem holds for the initial values used as long as they are elements of \(U\) [9,10]. We would like to emphasize that the Burnett equations will be solved with initial conditions provided by the DSMC data; in this way the question of slip at the boundaries is bypassed, but the calculated profiles at the boundaries allow us to compare the predicted slip with the DSMC data.

Finally, the values of the coefficients are

\[ \omega_1 = 1.014 \times 4 \left( \frac{7}{3} \frac{1}{2} \right), \quad \omega_2 = 2 \times 1.014, \]

\[ \omega_3 = 3 \times 0.806, \quad \omega_4 = 0.681, \quad \omega_5 = \frac{3}{2} \times 0.806 - 0.99, \]

\[ \omega_6 = 8 \times 0.928, \quad \theta_3 = -3.09, \quad \theta_4 = 2.418, \]

\[ \theta_3 = 8.3855. \]

Notice that \(\theta_3\)'s are not needed for solving the differential equations, but are used to evaluate the heat flux; also \(\theta_1\) and \(\theta_2\) do not appear in the expressions for the fluxes given by Eqs. (13). The Navier-Stokes equations result from taking \(F_3 = 0\) and \(F_6 = 0\) in Eqs. (17).

V. BURNETT DESCRIPTION FOR THE FLUXES

In Sec. IV we saw that, unlike the Navier-Stokes equations, the Burnett description predicts a heat flow along the direction that has no effect on the temperature profile. Since the computer experiments can measure momentum and heat flux, it is useful to evaluate the Burnett expressions explicitly for the various components of these fluxes.

First consider the expression of the \(x\) component of the heat flux; from Eqs. (13), we have

\[ q_x^{(2)} = \frac{1}{2} \theta_3 \mu^2 \frac{dp}{dy} + \frac{1}{2} \theta_4 \mu^2 \frac{d^2 u}{dy^2} + \frac{3}{2} \theta_5 \mu^2 \frac{dT}{dy} - \frac{2}{2} \rho(y) \frac{dy}{dy} \]

\[ \frac{dT}{dy} dT \frac{du}{dy} \frac{dy}{dy}. \]

The second derivative of the velocity can be evaluated from the Navier-Stokes results and the mass density can be eliminated from the equation of state, giving

\[ q_x^{(2)} = 20 \frac{\theta_3 \mu^2 k^2 T(y)^2 \frac{dp}{dy} + 2 \theta_4 \mu^2 k^2 T(y)^2 \frac{d^2 u}{dy^2} + 3 \theta_5 \mu^2}{2} \frac{dT}{dy} \frac{du}{dy} \frac{dy}{dy}. \]

In terms of the notation used for the dynamical system, the reduced form of the \(x\) component of the heat flux reads

\[ q_x^{(2)*} = \frac{d_1 Y_4 Y_5 Y_6}{\sigma^3} + \frac{d_2 Y_5 Y_4 Y_6}{Y_3}, \]

\[ \frac{d_3 Y_2 + d_4 Y_4 Y_5}{Y_3}, \]

where

\[ d_1 = \frac{25}{1024} \frac{k T_R \theta_3 \mu^2}{\rho(0) L^2}, \quad d_2 = \frac{25}{2048} \frac{k T_R \theta_4 \mu^2}{\rho(0) L^2}, \]

\[ d_3 = \frac{5}{128} \frac{k T_R \alpha \mu^2}{\rho(0) L^2}, \quad d_4 = \frac{75}{1024} \frac{k T_R \theta_5 \mu^2}{\rho(0) L^2}. \]

Since according to Eq. (7) there is no \(x\) component of the heat flux at the Navier-Stokes regime, we conclude that the \(x\) component of the total heat flux \((q_x')\), meaning the Navier-
Stokes plus the Burnett contributions, is given by \( \mathbf{q}_t^T = \mathbf{q}_{NS}^T + \mathbf{q}_x^{(2)} = \mathbf{q}_{NS}^{(2)}. \) Similarly we obtain that

\[
\mathbf{q}_t^T = \frac{2}{2(kT/\mu)^{3/2}} = d_5 \bar{Y}_2 Y_5,
\]

\[
d_5 = -\frac{75}{256} \frac{c_\lambda \sigma}{L^2}.
\]

(23)

V. DIRECT SIMULATION MONTE CARLO METHOD

The DSMC method is a well-established algorithm for computing gas dynamics at the level of kinetic theory. For completeness, this section presents a brief summary of the method; it is described in detail in Ref. [11]; and see Ref. [12] for a tutorial, and Refs. [13,14] for reviews. Wagner [15] proved that the DSMC method is equivalent to a stochastic evaluation of the Boltzmann equation.

In the DSMC method, the state of the system is given by the positions and velocities of particles, \( \{ \mathbf{r}_i, \mathbf{v}_i \} \). First, the particles are moved as if they did not interact; that is, their positions are updated to \( \mathbf{r}_i + \mathbf{v}_i t + \frac{1}{2} \mathbf{a} t^2 \), where \( t \) is the time step. A particle that strikes the thermal walls at \( y = \pm L/2 \) has its velocity replaced with a random value generated from the biased Maxwell-Boltzmann distribution; the system is periodic in the \( x \) and \( z \) directions. Second, after all particles have moved, a given number are randomly selected for collisions. This splitting of the evolution between streaming and collisions is accurate if the time step, \( t \ll \tau \), where \( \tau \) is the mean collision time for a particle.

The concept of “collision” implies that the interaction potential between particles is short ranged. In the simulations presented here the particles are taken to be rigid spheres of diameter \( \sigma \), so \( \tau = m/\rho \sigma^2 \langle \mathbf{v}_r \rangle \) where \( \langle \mathbf{v}_r \rangle \) is the average relative speed among the particles. The number of collisions among \( N \) particles during a time step is \( M = \frac{1}{2} N \tau/\tau \). We use Bird’s “no time counter” method [11] for determining \( M \), since it avoids the explicit evaluation of \( \langle \mathbf{v}_r \rangle \).

Particles are randomly selected as collision partners with the restriction that their mean separation be a fraction of a mean free path \( l = \langle \mathbf{v} \rangle \tau [16] \). This restriction is typically enforced by sorting the particles into cells whose dimensions are smaller than a mean free path. For hard spheres, the probability of selecting a given pair is proportional to the relative speed between the particles. DSMC evaluates individual collisions stochastically, conserving momentum and energy and selecting the post-collision angles from their kinetic theory distributions. For hard spheres, the center of mass velocity and relative speed are conserved in the collision with the direction of the relative velocity uniformly distributed in the unit sphere. This Markov approximation of the collision process is statistically accurate so long as the number of particles in a collision cell is sufficiently large, typically over 20 [17,18].

We define the average density of \( f(\mathbf{v}) \) as

\[
\langle f(\mathbf{v}) \rangle = \frac{1}{N_t} \sum_i \frac{1}{V_c} \sum_{i \in c} f(\mathbf{v}_i(t)),
\]

where \( \mathbf{v}_i \) is the velocity of particle \( i \), and \( N_t \) is the number of time steps over which samples are taken. The inner sum is over all particles within cell \( c \), which has volume \( V_c \) and is centered at position \( y \). The hydrodynamic variables are defined as

\[
\rho(y) = \langle m \rangle_y, \quad \langle \mathbf{v} \rangle_y = \langle m \mathbf{v} \rangle_y / \rho,
\]

\[
T(y) = \frac{2m}{3kT} \left( \frac{1}{2} \langle \mathbf{v}^2 \rangle_y - \frac{1}{2} \rho u^2 \right).
\]

and \( p = \rho kT/\mu \). The fluxes are defined as

\[
\mathbf{P}_{\alpha\beta}(y) = \langle m(\mathbf{v}_a - \mathbf{u}_a)(\mathbf{v}_\beta - \mathbf{u}_\beta) \rangle_y = \langle m \mathbf{v}_a \mathbf{v}_\beta \rangle_y - \rho \mathbf{u}_a \mathbf{u}_\beta,
\]

\[
\mathbf{q}_a(y) = \frac{1}{2} \langle m(\mathbf{v}_a - \mathbf{u}_a)(\mathbf{v} - \mathbf{u}) \rangle_y
\]

\[
= \frac{1}{2} \langle m \mathbf{v}_a \mathbf{v} \rangle_y - \langle m \mathbf{v}_a \mathbf{v}_a \rangle_y \mathbf{u}_a + \rho \mathbf{u}_a u^2 - \frac{1}{2} \mathbf{u}_a(m \mathbf{v}^2) \mathbf{u}_a,
\]

(32)
where $\alpha$, $\beta$, and $\gamma$ are $x$, $y$, or $z$ with implied summation over $\gamma$.

VII. SIMULATION RESULTS

To simplify comparison with the DSMC data in Ref. [5], for all the simulations $m = \sigma = T_R = 1$ and $k = \frac{1}{2}$. The sound speed at equilibrium is $\frac{5kT_R}{2m} = 1.29$. The equilibrium density is $\rho_0 = 1.21 \times 10^{-3}$, giving a mean free path of $l_0 = 186$.

A. Ten mean free path case

To simplify comparison with the DSMC data in Ref. [5], the pressure is normalized by $(7.13394 \times 10^{-4})^2 \frac{2kT_R}{\sigma^5}$ to make $p^*(0) = 1$. The simulations were done for an acceleration of $a = 1.6 \times 10^{-4} \frac{2kT_R}{m\sigma}$ and a system size of $L = 1860 \sigma = 10l_0$. Using these values, we obtain

\begin{align*}
b_0 &= -6.235007643, \quad b_1 = -162.3309911, \\
d_1 &= -0.50220358 \times 10^{-5}, \\
d_2 &= -0.39493245 \times 10^{-5}, \\
d_3 &= -0.245073776 \times 10^{-4}, \\
d_4 &= +0.408857099 \times 10^{-4}, \\
d_5 &= -0.128339708 \times 10^{-3}, \\
d_6 &= -0.09546098954, \\
d_9 &= +0.008456678889.
\end{align*}

The coefficient $b_2$ was determined in such a way that the solution to the differential equations reproduced the qualitative behavior for the DSMC pressure profiles. The explicit value used is

\begin{equation}
b_2 = 0.75,
\end{equation}

from which we obtain

\begin{equation}
d_7 = 0.9953798101, \quad d_8 = 0.002310094933.
\end{equation}

Finally the initial conditions used are

\begin{equation}
u^*(0) = 0.90287, \quad T^*(0) = 1.2052, \quad p^*(0) = 1.0,
\end{equation}

with the derivatives at the center of the channel equal to zero. The simulation contained $2.5 \times 10^4$ particles, and evaluated $2.5 \times 10^8$ collisions with statistical samples taken during the latter $1.2 \times 10^8$ collisions. The results for the profiles of the normal variables and some moments are given in Figs. 1–14.

B. 20 mean free path case

For comparison with the DSMC data in Ref. [5], the pressure is normalized by $(6.727 \times 10^{-4})^2 \frac{2kT_R}{\sigma^5}$ to make $p^*(0) = 1$. The simulations were done for an acceleration of $a = 4.0 \times 10^{-5} \frac{2kT_R}{m\sigma}$ and a system size of $L = 3720 \times \sigma = 20l_0$. Using these values, we obtain

\begin{align*}
b_0 &= -5.879345273, \quad b_1 = -577.3582581, \\
d_1 &= -0.1331459123 \times 10^{-5}, \\
d_2 &= -0.5209495406 \times 10^{-6}, \\
d_3 &= -0.6125684441 \times 10^{-5}, \\
d_4 &= +0.1083975775 \times 10^{-4}, \\
d_5 &= -0.6441653536 \times 10^{-4}, \\
d_6 &= -0.05061788107, \\
d_9 &= 0.002377693652.
\end{align*}

The coefficient $b_2$ was determined in such a way that the solution to the differential equations reproduced the qualitative behavior for the DSMC pressure profiles. The explicit value used is

\begin{equation}
b_2 = 0.75,
\end{equation}

from which we obtain

\begin{equation}
d_7 = 0.9953798101, \quad d_8 = 0.002310094933.
\end{equation}

Finally the initial conditions used are

\begin{equation}
u^*(0) = 0.90287, \quad T^*(0) = 1.2052, \quad p^*(0) = 1.0,
\end{equation}

with the derivatives at the center of the channel equal to zero. The simulation contained $2.5 \times 10^4$ particles, and evaluated $2.5 \times 10^8$ collisions with statistical samples taken during the latter $1.2 \times 10^8$ collisions. The results for the profiles of the normal variables and some moments are given in Figs. 1–14.
The coefficient $b_2$ was determined in such a way that the solution to the differential equations reproduced the qualitative behavior for the DSMC pressure profiles. The explicit value used is

$$b_2 = 0.2047,$$

from which we obtain,

$$d_7 = 0.9996454541, \quad d_8 = 0.0001772729471.$$  \hspace{1cm} (39)

Finally the initial conditions used are:

$$u^*(0) = 0.7575, \quad T^*(0) = 1.1293, \quad p^*(0) = 1.0,$$

$$d_7 = 0.9996454541, \quad d_8 = 0.0001772729471.$$  \hspace{1cm} (39)

Finally the initial conditions used are:

$$u^*(0) = 0.7575, \quad T^*(0) = 1.1293, \quad p^*(0) = 1.0,$$

with the derivatives at the center of the channel equal to zero. The simulation contained $5.0 \times 10^8$ particles and evaluated $5.0 \times 10^8$ collisions with statistical samples taken during the latter $2.5 \times 10^8$ collisions. The results for the profiles of the normal variables and some moments are given in Figs. 15-25.

C. 40 mean free path case

For comparison with the DSMC data in Ref. [5], the pressure is normalized by $(6.555 \times 10^{-4}) \times 2kT_p\sigma^3$ to make $p^*(0) = 1$. The simulations were done for an acceleration of $a = 1.0 \times 10^{-5} \times 2kT_p/m\sigma$ and a system size of $L = 7440 \times \sigma = 40l_0$. Using these values, we obtain

FIG. 3. Reduced velocity profile for the $z$ component of the velocity vector $u_z^*$ vs $s$ for the $L=10l_0$ system. Circles: DSMC. Solid line: assumption for plane Poiseuille flow.

FIG. 4. Reduced temperature profile $T^*$ vs $s$ for the $L=10l_0$ system. Circles: DSMC; the error bars are of approximately the same size as the symbols. Solid line: Burnett.

FIG. 5. Reduced pressure profile $p^*$ vs $s$ for the $L=10l_0$ system. Circles: DSMC; the error bars are of approximately the same size as the symbols. Solid line: Burnett.

FIG. 6. Reduced $x$ component of the heat flux, $q_x^*$, vs $s$ for the $L=10l_0$ system. Circles: DSMC. Solid line: the error bars are approximately three-fourths of the size of the symbols. (Burnett) Long dashed line: Navier-Stokes.
FIG. 7. Reduced y component of the heat flux, $q_y^*$, vs $s$ for the $L = 10l_0$ system. Circles: DSMC; the error bars are approximately half of the size of the symbols. Solid line: Burnett.

$$b_0 = -5.729018622, \quad b_1 = -2192.8448741,$$
$$d_1 = -0.3415989901 \times 10^{-6},$$
$$d_2 = -0.1336547505 \times 10^{-6},$$
$$d_3 = -0.153142111 \times 10^{-5},$$
$$d_4 = +0.2781046924 \times 10^{-5},$$
$$d_5 = -0.00003220826771,$$
$$d_6 = -0.02597303478,$$
$$d_9 = 0.000626027441.$$

The coefficient $b_2$ was determined in such a way that the solution to the differential equations reproduced the qualitative behavior for the DSMC pressure profiles. The explicit value used is

$$b_2 = +0.051335336,$$

from which we obtain

$$d_7 = 0.9999765896, \quad d_8 = 0.00001170519095.$$

Finally the initial conditions used are

$$u^*(0) = 0.689452, \quad T^*(0) = 1.09814, \quad p^*(0) = 1.0,$$

with the derivatives at the center of the channel equal to zero. The simulation contained $6.0 \times 10^4$ particles and evaluated $9.0 \times 10^8$ collisions with statistical samples taken during the

FIG. 9. Reduced yy component of the pressure tensor, $P_{yy}^*$, vs $s$ for the $L = 10l_0$ system. Circles: DSMC. Solid line: Burnett.

FIG. 8. Reduced z component of the heat flux, $q_z^*$, vs $s$ for the $L = 10l_0$ system. Circles: DSMC. Solid line: Burnett and Navier-Stokes.
latter $3.0 \times 10^8$ collisions. The results for the profiles of the normal variables and some moments are given in Figs. 26–36.

VIII. ANALYSIS OF THE RESULTS

This section discusses the general features observed in the above DSMC data and the comparison with the continuum equations. In some cases the effects are subtle, which is the reason we provide numerous graphs where the reader can obtain a more detailed picture. In several cases (e.g., $u_x$, $T$, $P_{yx}$, and $q_y$) the profiles obtained from the Navier-Stokes equations and Burnett equations are practically indistinguishable so only the latter are shown.

From Figs. 1, 15, and 26, we conclude that the profile for the $x$ component of the velocity is in reasonable agreement with the Burnett equations. The largest difference is observed at the highest Knudsen number, $K = l_0/L = 1/10$, where the discrepancy in the slip at the boundary is the most significant, in agreement with previous DSMC and molecular-dynamics (MD) results [19]. The $u_x$ profiles predicted by the Burnett and Navier-Stokes equations are nearly identical, so the latter are omitted in Figs. 1, 15, and 26.

For plane Poiseuille flow we assumed that the other components of the hydrodynamic velocity to be zero, and the measured profiles for $u_y$ and $u_z$ are shown in Figs. 2, 3, 16, 17, 27, and 28. For the velocity component perpendicular to the walls the DSMC data are consistent with the assumption that $u_y = 0$. For the parallel component $u_z$, there is noticeable variation, especially in the $L = 40l_0$ system (Fig. 28), because the long wavelength fluctuations in the periodic directions are very long lived. A similar effect is expected in $u_x$, but is masked by the mean flow, which is nearly four orders of magnitude larger. Though it is possible that the
stationary situation has not been established, this explanation seems unlikely since the DSMC data are assembled from a sequence of simulation runs, discarding any runs exhibiting remnants of the initial transient.

The most striking feature in the DSMC temperature profiles (Figs. 4, 18, and 29) is the appearance of two local maxima, which is quite noticeable for the ten mean free path case. The Burnett and Navier-Stokes equations give nearly identical predictions (the latter is omitted from Figs. 4, 18 and 29), and fail to capture this bimodal behavior. However, the maximum difference between the DSMC data and the Burnett prediction is about 1% for the ten mean free path case and the agreement improves with decreasing Knudsen number. This bimodal behavior is predicted in analyses using the BGK model [4] and Grad’s expansion [26]. The central minimum in the temperature profile is well accounted for in a 19-moment-equations approximation to the Boltzman equation as presented by Hess and Malek Mansour [44].

The pressure profiles (Figs. 5, 19, and 30), all show good agreement between the DSMC data and the Burnett equations. This agreement, however, is constructed because in lieu of determining the initial conditions the coefficient $b_2$ is selected so as to fit the data. This coefficient depends on the second derivative of $p$ at $s=0$ but the DSMC data are not sufficiently accurate to extract its value. It turns out that the pressure profile obtained with the Burnett equations is sensitive to the value of $b_2$ but the velocity and temperature profiles are not affected. The Navier-Stokes equations predict that $p^*(s) = 1$.

Next we examine the components of the heat flux and the
complete pressure tensor. One goal is to infer, if possible, the origin of the bimodal temperature profile. From the general assumptions for plane Poiseuille flow ($u_y = u_z = 0$ and Landau's symmetry argument) the velocity, temperature, and pressure profiles can depend only on $q_y$, $P_{yx}$, and $P_{yy}$. However, to have a better understanding of the problem it is useful to examine all the components of $q$ and $P$ to obtain the leading moments of the distribution function. Furthermore, the three assumptions given in Sec. II need to be tested.

The $x$ component of the heat flux is zero according to the Navier-Stokes equations, but is different from zero according to Burnett. Figures 6, 20, and 31 compare the Burnett prediction with the results from the DSMC method. The qualitative behavior of $q_x$ is very similar, in particular the change in sign of the heat flux, and the agreement between Burnett and the DSMC method is better at a lower Knudsen number.

The discrepancy is greatest at the boundaries where larger gradients are present. Notice that even at the lowest Knudsen number the DSMC data clearly indicate that $q_x \neq 0$. A non-zero parallel heat flux is also predicted in the BGK model [4] and Grad's expansion [26].

Figures 7, 21, and 32 show that $y$ component of the heat flux, as predicted by Burnett and Navier-Stokes, is in good agreement with the DSMC data, the differences being greatest at large Knudsen numbers and near the boundaries where larger gradients are present. However, near the center of the system, where the bimodal behavior of the temperature is observed, the heat flow given by the DSMC method and Burnett are in good agreement, which indicates that the discrepancy in the temperature profiles probably cannot be explained in terms of the differences for $q_y$. The DSMC results confirm that the $z$ component of the heat flux is zero; results from the ten mean free path system are shown in Fig. 8 and
FIG. 23. Reduced $xx$ component of the pressure tensor, $P_{xx}^*$, vs $s$ for the $L=20l_0$ system. Circles: DSMC. Solid line: Burnett.

FIG. 24. Reduced $zz$ component of the pressure tensor, $P_{zz}^*$, vs $s$ for the $L=20l_0$ system. Circles: DSMC. Solid line: Burnett.

FIG. 25. Reduced $yx$ component of the pressure tensor, $P_{yx}^*$, vs $s$ for the $L=20l_0$ system. Circles: DSMC; the error bars are approximately one-fifth of the size of the symbols. Solid line: Burnett.

FIG. 26. Reduced velocity profile for the $x$ component of the velocity vector, $u_x^*$, vs $s$ for the $L=40l_0$ system. Circles: DSMC; the error bars are a small fraction of the size of the symbols. Solid line: Burnett.

the data from the other cases is similar. From Landau’s symmetry argument and given that $u_y = u_z = 0$, it follows from the conservation equations that $P_{yy}$ must be a constant. Figure 9 shows that the DSMC data for the $L=10l_0$ system exhibit a weak nonconstant behavior, and are about 0.2% below the Burnett prediction. This could be an explanation for the discrepancy for the temperature profile, though for $L=20l_0$, the DSMC temperature profile exhibits a bimodal behavior while $P_{yy}$ is, within statistical error, a constant function (see Figs. 22 and 33). A possible explanation for $P_{yy}$ not being exactly constant in the $L=10l_0$ system is that while, on average, $u_y = u_z = 0$, these components of velocity fluctuate and these non-equilibrium fluctuations have long-range spatial correlations [20]. The other diagonal components of the pressure tensor are not zero; the DSMC data and Burnett predictions for $P_{xx}$ (Figs. 10, 23, and 34) and $P_{zz}$ (Figs. 11, 24, and 35) are in qualitative agreement. Finally, we turn our attention to the off-diagonal elements of $P$. The DSMC data confirm the Burnett and Navier-Stokes predictions that the components $P_{xz}$ and $P_{yz}$ are zero, as seen in Figs. 12 and 13 for the $L=10l_0$ system with similar results for the other cases. From Figs. 14, 25, and 36, we see that $P_{yx}$ is basically a linear function of $s$ for both the DSMC data and the Burnett equations. The slopes differ slightly, and while we do not expect that this explains the bimodal temperature profile and a more detailed analysis is needed.

IX. FINAL REMARKS

From our results we conclude that the bimodal temperature profile observed by the DSMC method cannot be explained with the Burnett equations, but that the nonconstant
pressure profile is a second order Knudsen effect. This conclusion is consistent with the fact that as the Knudsen number is lowered the bimodality in the temperature profile becomes unobservable, while a significant nonconstant pressure is still measured by the DSMC method. The bimodal character of the temperature profiles is perhaps a super-Burnett effect; however, it should be pointed out that the higher order Chapman-Cowling gradient expansions are very complicated. While some work has been done with the super-Burnett equations, this work is restricted to linearized corrections to the fluxes. For these reasons it seems more promising to look for alternatives such as the regularization given by Rosenau or the partial summation to the Chapman-Enskog method given by Gorban and Karlin. These methods, however, have some restrictions, such as the use of a linearized collision operator.

Tij et al. investigated the present problem using molecular methods for Maxwellian molecules, and obtained the bimodal behavior for the temperature profiles. Unfortunately simulations of Poiseuille flow using Maxwellian molecules are still lacking, so it is not known whether their prediction is quantitative. For the model used by Tij et al., Ikenberry and Truesdell evaluated a broad set of kinetic quantities that allow one to compute many moments, but a similar analysis has not been performed for other potentials, such as rigid spheres. Tij et al. predicted a heat flux in the $x$ direction, in qualitative agreement with the DSMC data and the Burnett results presented here. As we pointed out, this heat flux has no effect on the conservation equations, which are used to obtain the profiles. Recently Risso and Cordero used MD simulations and Grad’s expansion method to reach some of the same conclusions as in this work. There exist molecular-dynamics calculations for dense systems in which the inadequacy of the Navier-Stokes equations has also been
pointed out [27], in particular for plane Poiseuille flow [28]. Although these works are somewhat inapplicable to the present research for dilute gases, they support the conclusion that there exist simple flows for which the Navier-Stokes method is capable of improvement.

While moment methods recover many of the qualitative features observed in Poiseuille flow, their robustness has been criticized not only for the closure problem but also in evaluating shock waves. As Weiss [29] stated (we have consistently changed the numbering of the references), ‘‘the shock structure in monatomic gases is not satisfactorily described by the Navier-Stokes-Fourier theory, e.g., see Ref. [30]. But Grad’s 13-moment theory is even worse; indeed, Grad [31,32] himself found that no continuous shock structure exists beyond Mach 1.65, which is the maximum speed of propagation of the 13-moment theory.’’ Holway [33] showed that no matter how many moments are taken into account there is no continuous continuous shock structure beyond Mach 1.85, though this result has been challenged by Weiss himself [34]. In contrast, Gilbarg and Paolucci [30] remarked that, ‘‘while the evidence from kinetic theory and the theory of nonlinear viscosity indicate that the Navier-Stokes equations yield values of the shock thickness that are perhaps smaller than the actual one, there is nothing in the evidence to suggest that the classical theory is far wrong.’’

For shock waves there is indeed substantial evidence that the Navier-Stokes theory is susceptible to improvement [11,13,35,36]. However, the evidence provided by Fiscko and Chapman [37] that the Burnett equations give better predictions than the Navier-Stokes equations, which was quoted in Refs. [13,36], was flawed; an explanation of the errors and further extension of this important work can be find in the paper by Zhong et al. [38]. Two classical accounts of the Burnett equations applied to shock waves are the works by...
Wang-Chang and Uhlenbeck [21] and Foch [39]. The DSMC calculations for the shock wave problem were done a long time ago (see Ref. [11] for references), but the MD calculations are more recent [40] and in particular the first MD calculations for a dilute gas of rigid spheres were made in 1992 by Salomons and Mareschal [41]. Salomons and Mareschal [41] provided strong evidence that the Burnett equations can indeed provide a substantial improvement over the Navier-Stokes equations at high Mach numbers, but curiously enough a few months later a communication by these authors and Holian’s group was published [42] leaving aside the previous claim about the Burnett regime. Uribe et al. [6] followed Salomons and Mareschal’s remark, and found indeed that Burnett equations provided an accurate description for strong shock waves; however, in their qualitative analysis of the Burnett dynamical system they found a “Hopf-like bifurcation” at the upstream critical point, suggesting that there may not be a heteroclinic trajectory for Mach numbers greater than about 2.69. Rather than continuing to list all points of view on this topic, let us conclude by stating that a vigorous debate exists and that further experiments and simulations are needed to resolve these arguments.

FIG. 35. Reduced \( \frac{P_{zz}}{\rho U^2} \), vs \( s \) for the \( L=40l_0 \) system. Circles: DSMC. Solid line: Burnett.

FIG. 36. Reduced \( \frac{P_{yx}}{\rho U^2} \), vs \( s \) for the \( L=40l_0 \) system. Circles: DSMC; the error bars are about one fifth the size of the symbols. Solid line: Burnett.

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