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Inverted velocity profile in the cylindrical Couette flow of a rarefied gas

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The cylindrical Couette flow of a rarefied gas is investigated, under the diffuse-specular reflection condition of Maxwell’s type on the cylinders, in the case where the inner cylinder is rotating whereas the outer cylinder is at rest. The inverted velocity profile for small accommodation coefficients, pointed out by Tibbs, Baras, and Garcia [Phys. Rev. E 56, 2282 (1997)] on the basis of a Monte Carlo simulation, is investigated extensively by means of a systematic asymptotic analysis for small Knudsen numbers as well as the direct numerical analysis of the Boltzmann equation, and the parameter range in which the phenomenon appears is clarified.

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I. INTRODUCTION

Cylindrical Couette flow, which is a textbook example in classical fluid dynamics, is also one of the most fundamental problems in the kinetic theory of gases [1] and has been investigated from various points of view. The recent interest extends to topics such as the validity of the principle of the frame indifference and to clarify the parameter range where it appears.

One of such examples is given in a paper by Tibbs, Baras, and Garcia [6], where the cylindrical Couette flow of a rarefied gas is analyzed numerically by means of the direct simulation Monte Carlo (DSMC) method under the diffuse-specular reflection condition of Maxwell’s type on the cylinders in the case where the inner cylinder is rotating whereas the outer one is at rest. Their result shows that when the accommodation coefficients of the cylinders are small (i.e., when the major part of the molecules undergo specular reflection), the flow speed of the gas increases with the distance from the inner cylinder, which is contrary to the ordinary velocity profile of the Couette flow when only the inner cylinder is rotating.

The aim of the present paper is to investigate the phenomenon of the inverted velocity profile more comprehensively and to clarify the parameter range where it appears.

II. FORMULATION OF THE PROBLEM

A. Problem

Let us consider a rarefied gas between two coaxial circular cylinders with common temperature $T_0$: the inner cylinder with radius $r_I$ is rotating at a constant surface speed $V_I$, whereas the outer cylinder with radius $r_O$ is at rest. Assuming the diffuse-specular reflection condition of Maxwell’s type on the cylinders and restricting ourselves to the axially as well as circumferentially uniform case, we investigate the steady behavior of the gas on the basis of the Boltzmann equation. For actual numerical computation, we will assume hard-sphere molecules or the Bhatnagar-Gross-Krook (BGK) model [7,8].

B. Basic equation

We first introduce some notations: $(r, \theta, z)$ is the cylindrical coordinate system with the $z$ axis being the axes of the cylinders; $\xi$ is the molecular velocity and $\xi_r$, $\xi_\theta$, and $\xi_z$ are its $r$, $\theta$, and $z$ components; $f(r, \xi)$ is the velocity distribution function of the gas molecules; $\rho$ is the density, $u_r$, $u_\theta$, and $u_z$ ($v_r=v_\theta=0$) are the $r$, $\theta$, and $z$ components of the flow velocity, $T$ is the temperature, and $p$ is the pressure of the gas; and $R$ is the gas constant per unit mass. The dimensionless quantities corresponding to $V_I$, $r$, $\xi= (\xi_r, \xi_\theta, \xi_z)$, $f$, $\rho$, $u_\theta$, $T$, and $p$, which are denoted by $\hat{V}_I$, $\hat{r}$, $\hat{\xi}= (\hat{\xi}_r, \hat{\xi}_\theta, \hat{\xi}_z)$, $\hat{f}$, $\hat{\rho}$, $\hat{u}_\theta$, $\hat{T}$, and $\hat{p}$, respectively, are defined by

\[
\hat{V}_I = \frac{V_I}{(2RT_0)^{3/2}}, \quad \hat{r} = \frac{r}{r_I}, \quad \hat{\xi} = \frac{\xi}{(2RT_0)^{1/2}},
\]

\[
\hat{f} = \frac{(2RT_0)^{3/2}}{\rho_{av}} f, \quad \hat{\rho} = \frac{\rho}{\rho_{av}}, \quad \hat{u}_\theta = \frac{u_\theta}{(2RT_0)^{1/2}},
\]

\[
\hat{T} = \frac{T}{T_0}, \quad \hat{p} = \frac{p}{R\rho_{av}T_0},
\]

where $\rho_{av}$ is the average density of the gas between the cylinders. Therefore, $\hat{\rho}$ is normalized as

\[
\frac{2}{(r_O/r_I)^2} \int_{r_O/r_I}^{r_O/r_I} r \rho dr = 1.
\]

Then, the Boltzmann equation in the dimensionless form reads

\[
\begin{align*}
\frac{\partial \hat{f}}{\partial \hat{r}} + \frac{\hat{\xi}_r}{\hat{r}} \frac{\partial \hat{f}}{\partial \hat{\xi}_r} + \frac{\hat{\xi}_\theta}{\hat{r}} \frac{\partial \hat{f}}{\partial \hat{\xi}_\theta} - \frac{\hat{\xi}_r \hat{\xi}_\theta}{\hat{r}} & = \frac{2}{\sqrt{\pi Kn}} J(\hat{f}, \hat{\rho}),
\end{align*}
\]

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where \( \hat{J}(\hat{f},\hat{f}) \) is the dimensionless collision term, the explicit form of which is given in Appendix A. Kn is the Knudsen number, and \( l_0 \) is the mean free path of the gas molecules in the equilibrium state at rest at density \( \rho_{av} \) and temperature \( T_0 \). For hard-sphere molecules, \( l_0 \) is given by \( l_0 = (\sqrt{2\pi d^2 \rho_{av}/m})^{-1} \), \( m \) and \( d \) being the mass and diameter of a molecule.

The boundary conditions on the cylinders are written as follows: on the inner cylinder (at \( \hat{r} = 1 \)), for \( \xi_r > 0 \),

\[
\hat{f}(1,\xi) = (1 - \alpha_I) \hat{f}(1, -\xi_r, \xi_\theta, \xi_z) + \alpha_I \frac{\sigma_I}{\pi^{3/2}} \exp[-\xi_r^2 - (\xi_\theta - \hat{v}_\theta)^2 - \xi_z^2],
\]

and on the outer cylinder (at \( \hat{r} = r_O/\hat{r}_I \)), for \( \xi_r < 0 \),

\[
\hat{f}(r_O/\hat{r}_I, \xi) = (1 - \alpha_O) \hat{f}(r_O/\hat{r}_I, -\xi_r, \xi_\theta, \xi_z) + \alpha_O \frac{\sigma_O}{\pi^{3/2}} \exp[-\xi_r^2 - (\xi_\theta - \hat{v}_\theta)^2 - \xi_z^2],
\]

where \( \alpha_I \) and \( \alpha_O \) are the accommodation coefficient of the inner cylinder and that of the outer cylinder, respectively, and \( d\xi = d\xi_r d\xi_\theta d\xi_z \).

The macroscopic quantities \( \hat{\rho}, \hat{v}_\theta, \hat{T}, \) and \( \hat{p} \) are given by

\[
\hat{\rho} = \int \hat{f} d\xi, \quad \hat{v}_\theta = \frac{1}{\hat{\rho}} \int \xi_\theta \hat{f} d\xi, \quad \hat{T} = \frac{2}{3\hat{\rho}} \int [\xi_r^2 + (\xi_\theta - \hat{v}_\theta)^2 + \xi_z^2] \hat{f} d\xi, \quad \hat{p} = \hat{\rho} \hat{T}.
\]

Here and in what follows, the domain of integration with respect to \( \xi \) is its whole space unless the contrary is stated.

It should be noted that the local Maxwellian distribution corresponding to the solid-body rotation [9] with \( \hat{v}_\theta = \hat{v}_p \hat{r} \), i.e.,

\[
\hat{f} = \frac{\hat{\rho}_0}{\pi^{3/2}} \exp(\hat{V}_p^2 \hat{r}^2) \exp[-\xi_r^2 - (\xi_\theta - \hat{v}_\theta)^2 - \xi_z^2],
\]

where \( \hat{\rho}_0 \) is a constant determined by Eq. (2), is the exact solution of the present problem when \( \alpha_I = 0 \) and \( \alpha_O = 0 \). In view of this fact, one can naturally think that the inverted velocity profile (i.e., \( \hat{v}_\theta \) increasing with \( \hat{r} \)) takes place when \( \alpha_O \) becomes small. Our interest is to see the transition from the normal profile (i.e., \( \hat{v}_\theta \) decreasing with \( \hat{r} \)) to the inverted one at small accommodation coefficients.

### III. ASYMPTOTIC ANALYSIS FOR SMALL KNUDSEN NUMBERS

In this section we consider the case where the accommodation coefficients \( \alpha_I \) and \( \alpha_O \), as well as the Knudsen number Kn, are small and carry out a systematic asymptotic analysis of the boundary-value problem (3) and (5a)–(5b), following Refs. [10–13] as a guideline. To begin with, we assume that \( \alpha_I \) and \( \alpha_O \) are of the order of \( \text{Kn} \); that is, we put

\[
\alpha_I = \beta_I \varepsilon, \quad \alpha_O = \beta_O \varepsilon, \quad \varepsilon = (\sqrt{\pi/2}) \text{Kn} \ll 1,
\]

where \( \beta_I \) and \( \beta_O \) are given constants, and \( \varepsilon \) is a small parameter (of the order of \( \text{Kn} \)) that is mainly used in this section. Our aim is to derive a set of fluid-dynamic equations and its appropriate boundary conditions in the leading order, i.e., the order of \( \text{Kn}^0 \) (or \( \varepsilon^0 \)). Since the method of the asymptotic analysis is described in detail in the recent monograph by Sone [13], we give only the brief outline of the analysis and show the main result.

#### A. Fluid-dynamic equation

First, putting aside the boundary conditions (5a)–(5b), we look for a moderately varying solution \( \hat{f}_H \), satisfying \( \partial \hat{f}_H / \partial \hat{r} = \mathcal{O}(\hat{f}_H) \), in the form of a power series of \( \varepsilon \):

\[
\hat{f}_H = \hat{f}_0 + \hat{f}_1 \varepsilon + \hat{f}_2 \varepsilon^2 + \cdots.
\]

This \( \hat{f}_H \) is called the Hilbert solution or expansion. Let \( \hat{\rho}_H, \hat{v}_\theta, \hat{T}, \) be the macroscopic quantities \( \hat{\rho}, \hat{v}_\theta, \hat{T}, \) corresponding to the Hilbert solution. Then, they are also expanded as

\[
\hat{\rho}_H = h_H + h_{H1} \varepsilon + h_{H2} \varepsilon^2 + \cdots,
\]

where \( h \) represents \( \hat{\rho}, \hat{v}_\theta, \hat{T}, \) or \( \hat{p} \). The explicit expressions of \( h_{Hm} \) in terms of \( \hat{f}_H \) are obtained by substituting \( \hat{f} = \hat{f}_H \) and \( h = h_H \) in Eq. (7) and by equating the coefficients of the same power of \( \varepsilon \).

If we substitute Eq. (10) into Eq. (3) and arrange the power of \( \varepsilon \), we obtain a sequence of integral equations for the coefficients \( \hat{f}_{Hm} \), which can be solved from the lowest order. For the leading-order term \( \hat{f}_{H0} \), we have \( \hat{J}(\hat{f}_{H0}, \hat{f}_{H0}) = 0 \), which means that \( \hat{f}_{H0} \) is a local Maxwellian distribution. As for the higher-order term \( \hat{f}_{Hm} \), we obtain a linear integral equation, containing the collision term \( \hat{J} \) linearized around \( \hat{f}_{H0} \), with inhomogeneous terms consisting of the earlier terms \( \hat{f}_{Hn} (n < m) \) of the Hilbert expansion. Since the corresponding homogeneous equation has the summation invariants (multiplied by \( \hat{f}_{H0} \)) as nontrivial solutions, the inhomogeneous terms should satisfy solvability conditions, which are reduced to the form

\[ \text{Kn} = l_0/\hat{r}_I, \]
The application of this condition to \( f_{\hat{H}} \) and \( f_{\hat{H}} \) equations are derived in the following subsection. The appropriate boundary conditions for this set of equations from the lowest order leads to the fluid-dynamic equations for \( h_{\hat{Hm}} \) in Eq. (11).

To summarize, \( \hat{f}_{\hat{H}} \) is obtained as

\[
\hat{f}_{\hat{H}0} = \frac{\hat{\rho}_{\hat{H}0}}{(\pi \hat{T}_{\hat{H}0}^{3/2})} \exp \left( -\frac{\xi_{\tau}^{2} + (\xi_{\theta} - \nu_{\hat{H}0})^{2} + \xi_{\zeta}^{2}}{\hat{T}_{\hat{H}0}} \right),
\]

and \( \hat{f}_{\hat{H1}} \) as the form given in Appendix B [Eq. (B1)]. The fluid-dynamic equations in the leading order, i.e., those for \( \hat{\rho}_{\hat{H}0}, \hat{\nu}_{\hat{H}0}, \hat{T}_{\hat{H}0}, \) and \( \hat{\rho}_{\hat{H}0} \), are obtained in the following form:

\[
\frac{d\hat{\rho}_{\hat{H}0}}{dr} = \frac{2\hat{\rho}_{\hat{H}0}\nu_{\hat{H}0}^{2}}{r} = 0,
\]

\[\frac{d}{dr} \left[ \gamma_{1}\hat{T}_{\hat{H}0}^{3/2} \left( \frac{d\hat{\nu}_{\hat{H}0}}{dr} - \frac{\hat{\nu}_{\hat{H}0}}{r} \right) \right] = 0,\]

\[\frac{5}{4} \frac{d}{dr} \left( \gamma_{2}\hat{T}_{\hat{H}0}^{1/2} \left( \frac{d\hat{T}_{\hat{H}0}}{dr} + \hat{T}_{\hat{H}0} \right)^{2} \right) = 0,
\]

\[\hat{\rho}_{\hat{H}0} = \hat{\rho}_{\hat{H}0} \hat{\rho}_{\hat{H}0}^{*},\]

where \( \gamma_{1} \) and \( \gamma_{2} \) are functions of \( \hat{T}_{\hat{H}0} \) and their functional form depends on the molecular model (see Appendix B); for hard-sphere molecules, they are constants given by \( \gamma_{1} = 1.270042, \gamma_{2} = 1.922284; \) for the BGK model, \( \gamma_{1} = \gamma_{2} = \hat{T}_{\hat{H0}}^{1/2}. \gamma_{1} = \gamma_{2} = \hat{T}_{\hat{H0}}^{1/2}. \gamma_{1} = \gamma_{2} = \hat{T}_{\hat{H0}}^{1/2}. \)

The appropriate boundary conditions for this set of equations are derived in the following subsection.

**B. Knudsen layers and fluid-dynamic boundary conditions**

Now we take into account the boundary conditions that were put aside in the preceding subsection. The boundary conditions (5a)–(6b) with Eq. (9) can be recast as

\[
\hat{f}(\hat{r}_w, \xi) = (1 - \epsilon\beta_{w})S[\hat{f}(\hat{r}_w, \xi)] + \epsilon\beta_{w}D[\hat{f}(\hat{r}_w, \xi)] \quad \text{for} \quad \delta_{w}\xi > 0,
\]

where \( \hat{r}_w \) represents 1 and \( r_O/r_I \), and \( \beta_{w} \) and \( \delta_{w} \) are as follows:

\[
\beta_{w} = \beta_{1}, \quad \delta_{w} = 1, \quad \text{for} \quad \hat{r}_w = 1,
\]

\[
\beta_{w} = \beta_{O}, \quad \delta_{w} = -1, \quad \text{for} \quad \hat{r}_w = r_O/r_I.
\]

In addition, the operators \( S \) and \( D \) are defined by

\[
S[\hat{f}(\hat{r}_w, \xi)] = \hat{f}(\hat{r}_w, -\xi_{\theta}, -\xi_{\zeta}),
\]

\[
D[\hat{f}(\hat{r}_w, \xi)] = \frac{2}{\pi} \int_{\delta_{w}}^{1} \xi_{\theta} \hat{f}(\hat{r}_w, \xi) d\xi,
\]

where

\[
\hat{V}_{w} = \begin{cases} \hat{V}_{I} & \text{for} \quad \hat{r}_w = 1 \\ 0 & \text{for} \quad \hat{r}_w = r_O/r_I. \end{cases}
\]

Since \( \hat{f}_{\hat{H}0} \) is a local Maxwellian given by Eq. (14), it satisfies the specular-reflection condition on the cylinders, i.e.,

\[
\hat{f}_{\hat{H}0} = S[\hat{f}_{\hat{H}0}].
\]

In other words, it satisfies Eq. (16) in the leading order. However, as in the usual situation [13], the terms of \( O(\epsilon) \) of condition (16) cannot be satisfied by the next-order term \( \hat{f}_{\hat{H}1} \) of the Hilbert expansion. To obtain the solution satisfying the boundary condition, therefore, we need to introduce the so-called Knudsen layers.

Let us seek the solution in the form

\[
\hat{f} = \hat{f}_{H} + \hat{f}_{K},
\]

with

\[
\hat{f}_{K} = \hat{f}_{K1} \epsilon + \hat{f}_{K2} \epsilon^{2} + \cdots.
\]

Here, \( \hat{f}_{K} \) is the correction to the Hilbert solution \( \hat{f}_{H} \) appreciable only in the thin layers of thickness of the order of \( \epsilon \) (or of the mean free path in the dimensional \( r \) variable) adjacent to the cylinders (Knudsen layers). Expansion (22) is started from \( \epsilon \) order because \( \hat{f}_{\hat{H}0} \) could satisfy the boundary condition in the leading order.

To handle the two Knudsen layers on the inner and outer cylinders in a unified way, we introduce the following variables:

\[
y = \hat{r} - 1, \quad \eta = y/\epsilon, \quad \xi_{n} = \xi_{r},
\]

near the inner cylinder (\( \hat{r} = 1 \)); and

\[
y = r_O/r_I - \hat{r}, \quad \eta = y/\epsilon, \quad \xi_{n} = -\xi_{r},
\]

near the outer cylinder (\( \hat{r} = r_O/r_I \)). Here, \( y \) is the normal coordinate measured from each cylinder toward the gas, and
\( \eta \) is the stretched normal coordinate. We suppose that the length scale of variation of \( \hat{f}_K \) is \( \epsilon \), i.e.,

\[
\hat{f}_K = \hat{f}_K(\eta, \xi, \phi, \xi_\phi)
\]

or \( \partial \hat{f}_K / \partial \eta = O(\hat{f}_K) \), and that \( \hat{f}_K \) vanishes rapidly as \( \eta \to \infty \).

If we substitute Eq. (21) [with Eqs. (10) and (22)] into Eqs. (3) and (16) and take into account the explicit forms of \( \hat{f}_{H0} \) and \( \hat{f}_{H1} \) as well as the properties of \( \hat{f}_K \), we obtain the equation and boundary conditions for \( \hat{f}_{K1} \), which are summarized as follows.

\[
\xi_n \frac{\partial \hat{f}_{K1}}{\partial \eta} = 2J(\hat{f}_{H0}) \hat{f}_{K1}, \tag{26a}
\]

\[
\hat{f}_{K1} = S[\hat{f}_{K1}] + S(\hat{f}_{H1}) - (\hat{f}_{H1})_\eta - \beta_{nu} S[\hat{f}_{H0}] + \beta_{nu} D[\hat{f}_{H0}] \quad \text{for} \quad \xi_n > 0, \quad \eta = 0,
\]

\[
\hat{f}_{K1} \to 0 \quad \text{as} \quad \eta \to \infty, \tag{26b}
\]

where \( (\cdot)_\eta \) indicates the value of the Hilbert solution at \( \eta = 0 \), i.e., at \( \hat{r} = \hat{r}_\eta \), and \( J(\hat{f}, \hat{g}) \) is defined by Eq. (A1) in Appendix A. Equations (26a)–(26c) are essentially the half-space boundary-value problem of the linearized Boltzmann equation.

Let us denote any of \( 1, \xi, \) and \( \xi^2 \) by \( \Psi \). If we multiply Eq. (26a) by \( \Psi \) and integrate it over the whole space of \( \xi \), we have \( (\partial / \partial \eta) \int \Psi \xi_n \hat{f}_{K1} d\xi = 0 \) because of the property of the collision integral \( J \). Condition (26c) then gives

\[
\int \Psi \xi_n \hat{f}_{K1} d\xi = 0 \quad \text{for} \quad \eta > 0. \tag{27}
\]

We now consider Eq. (27) on the boundary \( \eta = 0 \), using Eq. (26b) with the explicit forms of \( \hat{f}_{H0} \) and \( \hat{f}_{H1} \) for \( \xi_n > 0 \) in the integral. Then we find that Eq. (27) with \( \Psi = \xi_n \) and \( \xi^2 \) gives two compatibility conditions for the boundary values of the leading-order terms \( h_{H0} \) [Eq. (11)] and of their derivatives. [Equation (27) with \( \Psi = 1 \) and \( \xi \) is automatically satisfied, whereas that with \( \Psi = \xi_n \) does not give any additional condition, consisting only of the boundary values of \( h_{H0} \) and \( dh_{H0}/d\eta \).] The two compatibility conditions, arranged in the appropriate form for each cylinder, are given as follows: on the inner cylinder (at \( \hat{r} = 1 \)),

\[
\sqrt{\pi} \gamma_1 \left( \frac{d^2 \hat{v}_{H0}}{d\hat{r}^2} - \hat{v}_{H0} \right) + \beta_1 \hat{\rho}_{H0}(\hat{V}_1 - \hat{v}_{H0}) = 0, \tag{28a}
\]

\[
\frac{5}{4} \sqrt{\pi} \gamma_2 \frac{d\hat{T}_{H0}}{d\hat{r}} - \beta_1 \hat{\rho}_{H0}(\hat{T}_{H0} - 1) + \frac{1}{2} \beta_1 \hat{\rho}_{H0}(\hat{V}_1 - \hat{v}_{H0})^2 = 0, \tag{28b}
\]

and on the outer cylinder (at \( \hat{r} = r_o / r_1 \)),

\[
\sqrt{\pi} \gamma_1 \left( \frac{d^2 \hat{v}_{H0}}{d\hat{r}^2} - \hat{v}_{H0} \right) + \beta_0 \hat{\rho}_{H0}(\hat{v}_{H0}) - \beta_0 \hat{\rho}_{H0} \hat{v}_{H0} = 0, \tag{29a}
\]

\[
\frac{5}{4} \sqrt{\pi} \gamma_2 \frac{d\hat{T}_{H0}}{d\hat{r}} - \beta_0 \hat{\rho}_{H0}(\hat{T}_{H0} - 1) + \frac{1}{2} \beta_0 \hat{\rho}_{H0}(\hat{v}_{H0})^2 = 0. \tag{29b}
\]

In deriving Eqs. (28b) and (29b) from Eq. (27) with \( \Psi = \xi^2 \), Eqs. (28a) and (29a) have been used. Equations (28a)–(29b) give the boundary conditions for the fluid-dynamic equations (15a)–(15d).

In this way, the boundary conditions for the leading-order fluid-dynamic equations (15a)–(15d) are obtained from the first-order Knudsen-layer problem (26a)–(26c). At the same time, it should be noted that the former boundary conditions can be derived without solving the latter problem. This procedure to determine fluid-dynamic boundary conditions in the case of a small accommodation coefficient or specularly reflecting boundary was first devised in Ref. [14], where the thermal creep along a specularly reflecting wall was clarified. Then, it was applied in Refs. [15] and [16], where the fluid-dynamic system for small Knudsen numbers was derived for the general geometry with a specularly [15] or almost specularly [16] reflecting boundary when the system was close to a uniform equilibrium state at rest. The existence and uniqueness of a solution of a half-space problem that contains the problem (26a)–(26c) as a special case have been proved by Golse et al. [17].

As is easily seen in the course of analysis, if the accommodation coefficients are of the order of unity [i.e., \( \alpha_i = O(1) \) and \( \alpha_o = O(1) \)], the boundary conditions for Eqs. (15a)–(15d) become the so-called nonslip conditions:

\[
\hat{v}_{H0} = \hat{V}_1, \quad \hat{T}_{H0} = 1, \quad \text{at} \quad \hat{r} = 1, \tag{30}
\]

\[
\hat{v}_{H0} = 0, \quad \hat{T}_{H0} = 1, \quad \text{at} \quad \hat{r} = r_o / r_1.
\]

These can be obtained formally by letting \( \beta_i \to \infty \) in Eq. (28) and \( \beta_o \to \infty \) in Eq. (29).

C. Flow properties at small Knudsen numbers

We have derived the fluid-dynamic equations (15a)–(15d) and their boundary conditions (28a)–(29b) for the leading order in Kn (or \( \epsilon \)) under assumption (10). Some numerical results of this system for hard-sphere molecules are shown in Fig. 1. That is, the velocity profile \( \hat{v}_{H0} \) is shown in the case where \( \beta_i = \beta_o = \beta, \ r_o / r_1 = 2, \) and \( \hat{V}_1 = V_1 / (2RT_o)^{1/2} = 0.1 \) [\( \alpha \)], 0.3 [\( \beta \)], and 0.5 [\( \alpha \)]. The dashed line in Fig. 1(c) indicates the solution given by Eq. (34) below [see the sentences following Eq. (34)]. In the figures, the inverted velocity profile is observed for \( \beta \leq 0.5 \).

Since the above system cannot be solved analytically, we consider the case where the rotation speed is small (but much larger than the Knudsen number). More specifically, we assume that \( \hat{V}_1 \ll 1 \) and thus \( \hat{v}_{H0} \ll 1 \) (\( \hat{V}_1 \) and \( \hat{v}_{H0} \) are supposed to be positive). Correspondingly, we put \( \hat{\rho}_{H0} = 1 + \omega \),
neglect the higher-order terms, Eqs. \(~\) and on the outer cylinder \(~\) and boundary conditions \(~\). 5

\[ r \frac{\partial}{\partial r} \left( \frac{1}{r} \frac{\partial}{\partial r} \right) \psi = \frac{1}{r} \frac{\partial}{\partial r} \left( \frac{1}{r} \frac{\partial}{\partial r} \right) \psi = 0. \]

\[ \sqrt{\pi} \gamma_1^* \left( \frac{d}{dr} \frac{1}{r} \frac{\partial}{\partial r} - \beta_1 \right) \psi = 0, \] 33a

\[ \frac{5}{4} \sqrt{\pi} \gamma_2^* \frac{d}{dr} \tau - \beta_1 \tau = 0, \] 33b

where \( \gamma_1^* \) and \( \gamma_2^* \) are \( \gamma_1 \) and \( \gamma_2 \) at \( \hat{T}_{H_0} = 1 \), respectively. This system can be solved readily and gives

\[ u = \Lambda \hat{r}^{-1} + B \hat{r}, \quad \omega = \tau = P = 0, \] 34

where

\[ A = \frac{1}{D} \frac{r_0}{r_1} \beta_1 \beta_0 \hat{V}_1, \] 35a

\[ B = \frac{1}{D} \frac{2}{2} \sqrt{\pi} \gamma_1^* \left( \frac{r_1}{r_0} \right)^2 - \frac{r_1}{r_0} \beta_0 \right) \beta_1 \hat{V}_1, \] 35b

\[ D = \frac{r_o}{r_1} - \frac{r_1}{r_o} \beta_1 \beta_0 + 2 \sqrt{\pi} \gamma_1^* \left( \frac{r_1}{r_0} \right)^2 \beta_1 \] 35c

Equation (34) corresponds to the solution of the incompressible Navier-Stokes equations. It should be noted that the pressure gradient due to the centrifugal force appears in the higher order, so that there is no pressure gradient in the leading order [Eq. (31a)]. Solution (34) is also plotted in Fig. 1(a) (dashed line). Even for \( \hat{V}_1 = 0.5 \), the difference between the solution of the compressible system and Eq. (34) is very small. If the latter is plotted in Fig. 1(a), the difference is invisible.

If we assume the specular-reflection condition on both cylinders from the beginning [i.e., if \( \alpha_I = \alpha_O = 0 \) in Eqs. (5a) and (6a)], the solution is indeterminate because the solid-body rotation with an arbitrary angular speed [Eq. (8) with \( \hat{V}_1 \) replaced by an arbitrary constant] is a solution. On the contrary, if we take the limit \( \beta_1 \to 0 \) and \( \beta_0 \to 0 \), keeping the ratio \( \chi = \beta_0 / \beta_1 \) fixed, in Eq. (34), then we have \( u \to \hat{V}_1 \hat{r}/[1 + (r_0/r_1)^3 \chi] \), that is, the limiting flow is the solid-body rotation with an angular speed determined uniquely by ratio \( \chi \).

**D. Continuum limit**

In this subsection, we comment briefly on the continuum limit. The continuum limit is the limit where the Knudsen number Kn (or \( \epsilon \)) vanishes. In this limit, because of Eq. (9), the boundary conditions on both cylinders approach the specular reflection. On the other hand, we have \( \hat{\rho} \to \hat{\rho}_{H_0} \), \( \hat{\psi} \to \hat{\psi}_{H_0} \), \( \hat{T} \to \hat{T}_{H_0} \), and \( \hat{\rho} \to \hat{\rho}_{H_0} \) in this limit because the Knudsen-layer correction appears in the order of \( \epsilon \). This means that the continuum limit of the Couette flow for the specular-reflection condition is uniquely determined if the limit is taken appropriately, i.e., in such a way that \( \alpha_I \) and \( \alpha_O \) vanish in proportion to \( \epsilon \) [Eq. (9)], and the limiting solution depends on the proportionality constants (\( \beta_I \) and \( \beta_O \)).
In contrast, if \( \alpha_I \) and \( \alpha_O \) are set to be zero from the beginning, the solution is indeterminate for any \( \text{Kn} \) including the continuum limit. Finally, if the continuum limit \( \text{Kn} \to 0 \) is taken for any fixed \( \alpha_I \) and \( \alpha_O \), then the limiting solution is given by the solution of Eqs. (15a)–(15d) with the nonslip condition (30) even when \( \alpha_I \) and \( \alpha_O \) are small.

### IV. CASE OF FREE-MOLECULAR FLOW

Next, we consider the other extreme case where the Knudsen number is infinitely large, i.e., the free-molecular flow. In this case, one can obtain the exact solution to the original system, Eq. (3) (with the right-hand side equal to zero) and boundary conditions (5a)–(6b), in the following form:

\[
\hat{f}(r, \xi) = \frac{C}{\pi^{3/2}} \exp(-r^2 - \xi^2) \quad (\varphi < |\hat{\theta}| < \pi - \varphi), \quad (36a)
\]

\[
\hat{f}(r, \xi) = \frac{\alpha_f(1 - \alpha_O)}{\alpha_I + \alpha_O - \alpha_I \alpha_O} \frac{C}{\pi^{3/2}} \exp(-\xi^2 + 2\hat{V}_I \xi \sin \theta - \xi^2) - \hat{V}_I \left( \frac{\alpha_O}{\alpha_I + \alpha_O - \alpha_I \alpha_O} \frac{C}{\pi^{3/2}} \exp(-\xi^2 - \xi^2) \right) \quad (0 < |\hat{\theta}| < \varphi), \quad (36b)
\]

where

\[
\xi = (\xi_r + \xi_\theta) \frac{1}{2}, \quad \theta = \arctan(\xi_\theta / \xi_r), \quad (37)
\]

\[
\varphi = \arcsin(1/r), \quad (37)
\]

and \( C \) is a constant. Then, the macroscopic quantities are obtained from Eqs. (7) and (36) as follows:

\[
\hat{\rho}(r) = \frac{C}{\pi} \left[ \pi - 2\varphi + \left( \frac{2\alpha_O - \alpha_I \alpha_O}{\alpha_I + \alpha_O - \alpha_I \alpha_O} \right) \varphi \right] + \frac{2\alpha_I - \alpha_I \alpha_O}{\alpha_I + \alpha_O - \alpha_I \alpha_O} \frac{C}{\pi} \exp(-\hat{V}_I^2) \times \left[ \varphi + \sqrt{\pi} \int_0^\varphi \Theta \exp(\Theta^2) \text{erf} \Theta d\theta \right], \quad (38)
\]

\[
\hat{\rho}(r) = \frac{2\alpha_I - \alpha_I \alpha_O}{\alpha_I + \alpha_O - \alpha_I \alpha_O} \frac{C}{\pi} \exp(-\hat{V}_I^2) \times \left[ \varphi + \sqrt{\pi} \int_0^\varphi \Theta \exp(\Theta^2) \text{erf} \Theta d\theta \right], \quad (39)
\]

\[
\hat{\rho}(r) = \frac{2\alpha_I - \alpha_I \alpha_O}{\alpha_I + \alpha_O - \alpha_I \alpha_O} \frac{C}{\pi} \exp(-\hat{V}_I^2) \times \left[ \varphi + \sqrt{\pi} \int_0^\varphi \Theta \exp(\Theta^2) \text{erf} \Theta d\theta \right], \quad (40)
\]

where \( \Theta = \hat{V}_I \sin \hat{\theta}, \) and \( \text{erf} z = (2/\sqrt{\pi}) \int_0^z \exp(-t^2) dt \) is the error function. Constant \( C \) in Eqs. (36) and (38)–(40) is determined by Eq. (2). In the derivation of the above solution, we have assumed that \( \alpha_O \neq 0 \). It should be noted that when \( \alpha_I = \alpha_O \), all the macroscopic quantities are independent of the accommodation coefficient, though the velocity distribution function still depends on it. Results (38)–(40) will be shown in some figures in Sec. V.

### V. NUMERICAL ANALYSIS FOR WHOLE RANGE OF KNUDSEN NUMBER

Finally, we consider the case of arbitrary Knudsen numbers. Here, we use two different numerical approaches: one is the stochastic method known as the direct simulation Monte Carlo (DSMC) method [18,19] and the other is a deterministic finite-difference method based on the BGK model. We hereafter restrict ourselves to the case where the accommodation coefficients are common to both cylinders and put \( \alpha_I = \alpha_O = \alpha \).

#### A. Monte Carlo simulation

We start with the DSMC computation. Since the method is widely used and described in many places (see, e.g., Refs. [20–22] in addition to Refs. [18,19]), we give only the result of analysis, omitting the description of the solution process. In this subsection, we assume that the gas molecules are hard spheres. In the actual computation, we use 200 uniform cells in \( r_I < r < r_O \) and 200 simulation particles per cell on the average.

Figure 2 shows some results for \( r_O/r_I = 2 \) and \( \hat{V}_I/(2RT_o)^{1/2}(-\hat{V}_I) = 0.5 \). More specifically, the velocity profile \( \hat{v}_I/(2RT_o)^{1/2}(-\hat{v}_I) \) is plotted for various values of \( \alpha \) (\( \alpha_I = \alpha_O \)) in the case of \( \text{Kn} = 0.02, 0.05, 0.1, 1, \) and 100. Note that the Reynolds number \( \text{Re} = (4/\sqrt{\pi}) \gamma_I (\hat{V}_I/\text{Kn}) \) (cf. Ref. [13]) in the present computation, which is less than 45, is below the critical Reynolds number for the Taylor-Couette instability [23] so the flow is axially...
and circumferentially uniform (see also Ref. [24]). In Figs. 2(a)–2(c) [Kn = 0.02, 0.05, and 0.1], the numerical solution of the fluid-dynamic system for small α and Kn, i.e., Eqs. (15a)–(15d) and (28a)–(29b), is shown by the dashed line for α = 0.01, 0.05, 0.1, and 0.2, and the numerical solution of Eqs. (15a)–(15d) and (30) (nonslip conditions) is shown by the dotted line. In Fig. 2(f) (Kn = 100), the free-molecular-flow result, Eq. (39), is shown by the thick dot-dashed line. For Kn = 1, the inverted velocity profile is observed for α = 0.1, but it is limited to smaller α when Kn is either small or large. For small Kn and α, the fluid-dynamic solution shows good agreement with the DSMC result, so that the dashed lines in Figs. 2(a)–2(c) are not seen clearly. For Kn = 100, the profile is weakly dependent on α, which is consistent with the free-molecular-flow result, and thus the inverted velocity profile is not observed even for α = 0.01. The range of α for the inverted profile will be discussed in the following subsection.

**B. Finite-difference analysis of the BGK model**

In order to obtain more detailed information on the parameter range for the inverted velocity profile, we carry out a deterministic numerical analysis using the BGK model rather than the original Boltzmann equation. We employ the finite-difference method developed by Sone and Sugimoto in their study of strong evaporation from spherical and cylindrical condensed phases [25–27].

A difficulty inherent in the finite-difference analysis is caused by the fact that the molecular velocity distribution function around a convex body generally contains discontinuities [28]. The method mentioned above is capable of describing the behavior of the discontinuity around a spherical or cylindrical body. In the present problem, the situation is slightly more complicated because the discontinuity caused by the inner cylinder, in general, reaches the outer cylinder and is reflected there by the specular-reflection part of the boundary condition (6a). Therefore, we need to adjust the above-mentioned scheme to the present problem. But, since the method is essentially the same as that described in detail in Refs. [25–27], we give only the result of the analysis.

Figure 3 shows the velocity profile for the BGK model, corresponding to Fig. 2. The meaning of the dashed and dotted lines in Figs. 3(a)–3(c) is the same as in Fig. 2, but the relation \( \gamma_1 = \gamma_2 = \gamma = 1/2 \), which corresponds to the BGK model, is used in Eqs. (15a)–(15d) and (28a)–(29b). Figures 4 and 5 show, respectively, the density and temperature profiles for the BGK model for typical Kn in the same case as in Fig. 3. Figure 3 is very similar to Fig. 2, though there is a small quantitative discrepancy.

As is seen from Figs. 2 and 3, the velocity profile is monotonically decreasing when α is close to unity. If α is decreased for a fixed Kn, the monotonicity ceases at a relatively small α, say, \( \alpha_c \), i.e., the profile exhibits a local minimum when \( \alpha < \alpha_c \). Then, with the further decrease of α, the profile becomes monotonically increasing. We show the critical value \( \alpha_c \) versus Kn for the BGK model in Fig. 6, where \( r_O/r_I = 2 \) and \( V_I/(2RT_O)^{1/2} = 0.1 \) and 0.5. The \( \alpha_c \) is almost the same for both values of \( V_I/(2RT_O)^{1/2} \) and becomes largest at \( Kn = 0.7 \sim 0.8 \). From Eq. (34), we have the following expression for small Kn and small \( V_I/(2RT_O)^{1/2}[Kn \ll V_I/(2RT_O)^{1/2} \ll 1] \):

\[
\alpha_c = (\pi/2) (r_I/r_O) \gamma^{1/2} Kn. \tag{41}
\]

This result for the BGK model (\( \gamma^{1/2} = 1 \)) is also shown by the solid line in Fig. 6. The range of parameters, α and Kn, for which the profile exhibits a local minimum is within the reach of laboratory experiments [29].

Here, we comment on the comparison between the result for hard-sphere molecules and that for the BGK model. As is well known, the way of comparison between the results for different molecular models is not unique. One of the standard ways is the following. The viscosity coefficient \( \mu_0 \) corresponding to the reference density \( \rho_{av} \) and temperature \( T_O \) is...
FIG. 3. The numerical result by the finite-difference method for the BGK model: velocity profile. The $v/(2RT_0)^{1/2}$ is shown for various values of $a_i = \alpha_0 = \alpha$ and Kn in the case of $r_O/r_I = 2$ and $V_i/(2RT_0)^{1/2} = 0.5$. (a) $Kn = 0.02$, (b) $Kn = 0.05$, (c) $Kn = 0.1$, (d) $Kn = 1$, (e) $Kn = 10$, (f) $Kn = 100$. In (a)–(c), the dashed line indicates the numerical solution of Eqs. (15a)–(15d) with boundary conditions (28a)–(29b), and the dotted line that of the same equations with the nonslip conditions (30). In (f), the thick dot-dashed line indicates the free-molecular flow solution (39), which is independent of $\alpha$.

FIG. 4. The numerical result by the finite-difference method for the BGK model: density profile. The $\rho/\rho_0$ is shown for various values of $a_i = \alpha_0 = \alpha$ and typical values of Kn in the case of $r_O/r_I = 2$ and $V_i/(2RT_0)^{1/2} = 0.5$. (a) $Kn = 0.02$, (b) $Kn = 0.1$, (c) $Kn = 1$, (d) $Kn = 10$. In (a) and (b), the dashed line indicates the numerical solution of Eqs. (15a)–(15d) with boundary conditions (28a)–(28b), and the dotted line that of the same equations with the nonslip conditions (30). In (d), the thick dot-dashed line indicates the free-molecular flow solution (39), which is independent of $\alpha$.

FIG. 5. The numerical result by the finite-difference method for the BGK model: temperature profile. The $T/T_0$ is shown for various values of $a_i = \alpha_0 = \alpha$ and typical values of Kn in the case of $r_O/r_I = 2$ and $V_i/(2RT_0)^{1/2} = 0.5$. (a) $Kn = 0.02$, (b) $Kn = 0.1$, (c) $Kn = 1$, (d) $Kn = 10$. See the caption of Fig. 4.

given by $\mu_0 = (\sqrt{\pi}/4)\gamma_0^s\rho_0/(2RT_0)^{1/2}l_0$ (cf. Sec. 3.9 of Ref. [13]). If we suppose that $\mu_0$ is a fundamental quantity and is common to all the molecular models, then we obtain the relations among the mean free paths or the Knudsen numbers for the different molecular models. In the case of hard-sphere molecules and the BGK model, we have the following relationship:

$$l_0^{(BGK)} = 1.270 042 l_0^{(HS)},$$

$$Kn^{(BGK)} = 1.270 042 Kn^{(HS)},$$

where the suffixes (BGK) and (HS) indicate the quantities for the BGK model and those for hard-sphere molecules, respectively. In Fig. 7, we compare the velocity profile for the BGK model with that for hard-sphere molecules in the case where $r_O/r_I = 2$, $V_i/(2RT_0)^{1/2} = 0.5$, and $Kn^{(HS)} = 0.1$.
FIG. 6. The critical accommodation coefficient $\alpha_c$ versus Kn for the BGK model in the case of $r_o/r_i=2$ and $V_f/(2RT_0)^{1/2}=0.1$ and 0.5. The solid curve indicates Eq. (41).

using the conversion formula (42). More specifically, the result at Kn = 0.127 004 2 for the BGK model is compared with that at Kn = 0.1 for hard-sphere molecules. The figure shows very good agreement.

The data for the computational system in the finite-difference analysis are as follows. Region $1 \leq \tilde{r} \leq r_o/r_i$ is divided into 240 nonuniform intervals (the minimum size is $4.284 \times 10^{-5}$ at the cylinders, and the maximum size is $6.545 \times 10^{-3}$ around $\tilde{r} = 1.5$); region $0 \leq \tilde{\zeta}_p < \infty$ [see Eq. (37)] is replaced by $0 \leq \tilde{\zeta}_p < 6.912$ and then divided into 48 nonuniform intervals (the minimum size is $6.25 \times 10^{-5}$ at $\tilde{\zeta}_p = 0$, and the maximum size 0.4231 at $\tilde{\zeta}_p = 6.912$); region $-\pi \leq \theta_c \leq \pi$ [see Eq. (37)] is divided into 272 uniform intervals (note that $\zeta$ can be eliminated by suitable integration in the case of the BGK model [25,27]).

VI. CONCLUDING REMARKS

In the present study, we investigated the cylindrical Couette flow of a rarefied gas between two coaxial cylindrical in the case where the inner cylinder is rotating whereas the outer one is at rest. The diffuse-specular reflection condition of Maxwell’s type was assumed on the cylinders. Special attention was focused on the inverted velocity profile (the velocity profile increasing with the distance from the inner cylinder) for small accommodation coefficients observed previously by the Monte Carlo simulation [6]. After the formulation of the problem in Sec. II, we first considered the near continuum case in Sec. III, where we derived the fluid-dynamic system for small accommodation coefficients by means of a systematic asymptotic analysis of the Boltzmann equation and showed that the system describes the inverted velocity profile. Next, we considered the other extreme case, the free-molecular-flow limit, in Sec. IV and derived the analytical solution. In this case, the macroscopic quantities do not depend on the accommodation coefficients if they are common to both cylinders. Finally in Sec. V, we carried out numerical analysis of the original Boltzmann system for a wide range of the Knudsen number by using two different approaches: one is the Monte Carlo simulation (DSMC method) for hard-sphere molecules and the other is a deterministic finite-difference analysis of the BGK model. On the basis of these numerical results, the range of the accommodation coefficients that gives the inverted velocity profiles (more precisely, the profile that is not monotonically decreasing) was clarified for certain cases of the driving parameters [ $\alpha_f = \alpha_o = \alpha$, $r_o/r_i = 2$, and $V_f/(2RT_0)^{1/2} = 0.1$ and 0.5].

To conclude the paper, we discuss the inverted velocity profile from a qualitative, physical point of view. The rotating inner cylinder imparts the circumferential component of the momentum to the diffusely reflected molecules. When $\alpha_f$ is small, the rate of the transfer of the circumferential momentum is small, but it is supplied continuously. When Kn is small, the circumferential momentum is transferred to the entire gas by molecular collisions. If the outer cylinder is specularly reflecting ($\alpha_o = 0$), the circumferential momentum is not transferred to the outer cylinder. In other words, there is no braking effect by the outer cylinder. Then, the gas is accelerated until solid-body rotation, in which the transfer of the circumferential momentum vanishes, is established. This situation, in which the velocity profile is inverted since $v_o \propto r$, corresponds to Eq. (8). When $\alpha_o$ is small but non-zero, some of the circumferential momentum is transferred to the outer cylinder by the diffusely reflected molecules, and its reaction brakes the rotation of the gas near the outer cylinder. But, if the braking effect is small, the velocity profile remains inverted, as in solid-body rotation. The numerical result shows that this is the case when $\alpha_f = \alpha_o$ and they are sufficiently small. When Kn is sufficiently large ($Kn > 1$), a competing effect arises: because collisions in the gas are rare, most of the circumferential momentum imparted to the diffusely reflected molecules on the inner cylinder is transferred not to the gas but to the outer cylinder directly. Therefore, in contrast to the case of small Kn, the gas away from the inner cylinder is not accelerated. In this case, the number of the molecules with higher speed in the circumferential direction is larger near the inner cylinder, so that the inversion of the velocity profile does not appear.

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APPENDIX A: COLLISION INTEGRAL

The collision integral \( \mathcal{J}(\hat{f}, \hat{f}) \) is defined by the following bilinear integral operator \( \mathcal{J}(\hat{f}, \hat{g}) \):  
\[
\mathcal{J}(\hat{f}, \hat{g}) = \frac{1}{2} \int (\hat{f}^\ast \hat{g}' + \hat{f}' \hat{g}^\ast - \hat{f} \hat{g}' - \hat{f}' \hat{g}^\ast) \hat{B} d\Omega d\xi \ast ,
\]
(A1a)
\[
\hat{B} = \hat{B}(|V \cdot e|/|V|), \quad V = \xi \ast - \xi ,
\]
(A1b)
\[
\xi = \xi + (V \cdot e) e, \quad \xi \ast = \xi - (V \cdot e) e,
\]
(A1c)
where \( \hat{f} \), \( \hat{f}' \), \( \hat{f}^\ast \), and \( \hat{g} \) stand for \( \hat{f}(\xi) \) with \( \xi = \xi \ast , \xi \ast , \xi \), and \( \xi \ast \), respectively, and \( \xi \ast \) is the variable of integration corresponding to \( \xi \) and \( e \) is a unit vector; \( \hat{B} \) is a non-negative function of \( |V \cdot e|/|V| \) and \( \xi \ast \) depending on the molecular model (for hard-sphere molecules, \( \hat{B} = |V \cdot e|/4\sqrt{2} \pi \); \( d\Omega \) is the solid-angle element around \( e \), and \( \xi \ast = d\xi \ast, d\xi \ast \xi \ast \xi \ast ; \) the domain of integration in Eq. (A1a) is all directions of \( e \) and the whole space of \( \xi \ast \). (See Sec. 2.9 in Ref. [13]).

In the BGK model, the collision term \( \mathcal{J}(\hat{f}, \hat{f}) \) is replaced by the following \( \mathcal{J}_{\text{BGK}}(\hat{f}) \):
\[
\mathcal{J}_{\text{BGK}}(\hat{f}) = \rho \hat{f} \delta(\hat{f} - \hat{f}),
\]
(A2a)
\[
\hat{f} = \rho \left( \frac{\xi}{(\pi T)^{1/2}} \right)^{3/2} \exp \left( -\frac{\xi^2 + (\xi \theta - \hat{v}\theta)^2 + \xi \bar{z}^2}{T} \right),
\]
(A2b)
where \( \rho \), \( \hat{v}\theta \), and \( T \) are given in Eq. (7) [\( \hat{v}\theta = \hat{v}_z = 0 \) is used in Eq. (A2b)].

APPENDIX B: HILBERT SOLUTION \( \hat{f}_{H1} \)

The first-order Hilbert solution \( \hat{f}_{H1} \) is obtained in the following form:
\[
\hat{f}_{H1} = \hat{f}_{H0} \left[ \frac{\rho_{H1}}{\rho_{H0}} + \frac{2 \hat{v}\theta_{H1} \xi \theta}{T_{H1}^{1/2}} \hat{f}_{H1} \xi \theta - \hat{f}_{H0} \xi \frac{3}{2} \right] - \frac{1}{\rho_{H0}} \frac{d\hat{T}_{H0}}{d\tau} \bar{z}, \mathcal{A}(\xi, \hat{T}_{H0}) = \hat{T}_{H0}^{1/2} \frac{\hat{T}_{H1}^{1/2}}{\rho_{H0}} \frac{d\hat{v}\theta_{H0}}{d\tau} - \frac{\hat{v}\theta_{H0}}{\hat{T}_{H1}^{1/2}} \xi \theta \xi \theta \mathcal{B}(\xi, \hat{T}_{H0}),
\]
(B1)
where
\[
\xi \theta = \xi \frac{\xi}{\hat{T}_{H0}^{1/2}}, \quad \xi \frac{\xi}{\hat{T}_{H0}^{1/2}} = \xi \frac{\xi}{\hat{T}_{H0}^{1/2}}, \quad \xi \frac{\xi}{\hat{T}_{H0}^{1/2}}, (B2a)
\]
\[
\xi = (\xi^2 + \xi^2 + \xi^2)^{1/2}.
\]
(B2b)

The functions \( A \) and \( B \) are the solutions of the following integral equations:
\[
\begin{align*}
\mathcal{L}_T(\xi, \mathcal{A}(\xi, \hat{T}_{H0})) &= -\xi (\xi^2 - \frac{5}{2}), \\
\text{subsidiary condition:} & \int_0^\infty z^2 A(\xi, \hat{T}_{H0}) \exp(-\xi^2) d\xi = 0, \\
\mathcal{L}_T(\xi, \mathcal{B}(\xi, \hat{T}_{H0})) &= -2 \xi \xi \frac{\xi}{\hat{T}_{H0}},
\end{align*}
\]
(B3)
where \( \xi = (\xi^2 + \xi^2 + \xi^2)^{1/2} \) [Eq. (13)], and \( \mathcal{L}_T(\phi(\xi)) \) is the linearized collision operator defined by \( \mathcal{L}_A(\phi(\xi)) \) in Ref. [13] [Eq. (A.23) in Sec. A.2 of Appendix A there] with \( a = \hat{T}_{H0} \), i.e.,
\[
\mathcal{L}_T(\phi(\xi)) = \int E(\xi)(\phi + \phi' - \phi) \frac{\hat{B} d\Omega d\xi}{(\xi)},
\]
(B5a)
\[
\hat{B} = \hat{T}_{H0}^{1/2} \hat{B}(|V \cdot e|/|V|, \hat{T}_{H0}^{1/2}|V|),
\]
(B5b)
\[
E(\xi) = \pi^{3/2} \exp(-\xi^2),
\]
(B5c)
\[
\hat{\xi} = (\xi^2 + \xi^2 + \xi^2)^{1/2},
\]
(B5d)
where \( \phi', \phi'' \), \( \phi \), and \( \phi_0 \) stand for \( \phi(\xi) \) with \( \xi = \xi \ast , \xi \ast , \xi \), and \( \xi \ast \) [cf. Eq. (A1c)] respectively. The coefficients \( \gamma_1 \) and \( \gamma_2 \) occurring in the fluid-dynamic equations [Eqs. (15b) and (15c)] are expressed in terms of the functions \( A \) and \( B \) as follows:
\[
\gamma_1(\hat{T}_{H0}) = \frac{8}{15} B \mathcal{L}_T(\xi, \mathcal{B}(\xi, \hat{T}_{H0}) \exp(-\xi^2) d\xi,
\]
(B6a)
\[
\gamma_2(\hat{T}_{H0}) = \frac{16}{15} B \mathcal{L}_T(\xi, \mathcal{B}(\xi, \hat{T}_{H0}) \exp(-\xi^2) d\xi.
\]
(B6b)
Since \( \mathcal{L}_T \) generally depends on \( \hat{T}_{H0} \) and the molecular model, functions \( A \) and \( B \) and thus \( \gamma_1 \) and \( \gamma_2 \) also depend on them. In the case of hard-sphere molecules, \( A, B, \gamma_1 \), and \( \gamma_2 \) are independent of \( \hat{T}_{H0} \) and are given by [13]
\[
A(\xi, \hat{T}_{H0}) = A(\xi), \quad B(\xi, \hat{T}_{H0}) = B(\xi),
\]
(B7a)
\[
\gamma_1 = 1.270 042, \quad \gamma_2 = 1.922 284.
\]
(B7b)
Functions \( A(\xi) \) and \( B(\xi) \) are given in Table 3.1 in Ref. [13]. For the BGK model, these quantities are given as [13]
\[
A(\xi, \hat{T}_{H0}) = \hat{T}_{H0}^{1/2} \left( \xi^2 - \frac{5}{2} \right), \quad B(\xi, \hat{T}_{H0}) = 2 \hat{T}_{H0}^{1/2},
\]
(B8a)
\[
\gamma_1 = \gamma_2 = \hat{T}_{H0}^{1/2}.
\]
(B8b)