Coupling lattice Boltzmann model for simulation of thermal flows on standard lattices

Q. Li, 1,2 K. H. Luo, 2,4 Y. L. He, 1,* Y. J. Gao, 3 and W. Q. Tao 1

¹Key Laboratory of Thermal Fluid Science and Engineering of MOE, National Key Laboratory of Multiphase Flow in Power Engineering, School of Energy and Power Engineering, Xi'an Jiaotong University, Xi'an, Shaanxi 710049, China

²Energy Technology Research Group, School of Engineering Sciences, University of Southampton, Southampton SO17 1BJ, United Kingdom

³Department of Zoology, University of Cambridge, Cambridge CB2 3EJ, United Kingdom

⁴Center for Combustion Energy, Key Laboratory for Thermal Science and Power Engineering of Ministry of Education,
Department of Thermal Engineering, Tsinghua University, Beijing 100084, China
(Received 18 September 2011; revised manuscript received 14 November 2011; published 20 January 2012)

In this paper, a coupling lattice Boltzmann (LB) model for simulating thermal flows on the standard two-dimensional nine-velocity (D2Q9) lattice is developed in the framework of the double-distribution-function (DDF) approach in which the viscous heat dissipation and compression work are considered. In the model, a density distribution function is used to simulate the flow field, while a total energy distribution function is employed to simulate the temperature field. The discrete equilibrium density and total energy distribution functions are obtained from the Hermite expansions of the corresponding continuous equilibrium distribution functions. The pressure given by the equation of state of perfect gases is recovered in the macroscopic momentum and energy equations. The coupling between the momentum and energy transports makes the model applicable for general thermal flows such as non-Boussinesq flows, while the existing DDF LB models on standard lattices are usually limited to Boussinesq flows in which the temperature variation is small. Meanwhile, the simple structure and general features of the DDF LB approach are retained. The model is tested by numerical simulations of thermal Couette flow, attenuation-driven acoustic streaming, and natural convection in a square cavity with small and large temperature differences. The numerical results are found to be in good agreement with the analytical solutions and/or other numerical results reported in the literature.

DOI: 10.1103/PhysRevE.85.016710 PACS number(s): 47.11.-j

I. INTRODUCTION

Because of its inherent parallelizability on multiple processors and the avoidance of nonlinear convective terms, the lattice Boltzmann (LB) method, which originates from the lattice-gas automata (LGA) method [1], has been developed to be a very attractive numerical method in the past two decades [2-5]. Unlike conventional numerical methods, which are based on the discretization of macroscopic governing equations, the LB method is based on the mesoscopic kinetic equation. The main advantages of the LB method can be summarized as follows [6]: (1) simple form of the governing equation and the easiness of programming: In conventional numerical methods the convection terms of governing equations are nonlinear, while in the LB method the convection terms are linear and the viscous effect is modeled through the linearized collision operator, such as the Bhatnagar-Gross-Krook (BGK) collision operator [7–10], multiple-relaxation-time (MRT) collision operator [11–16], and the two-relaxation-time (TRT) collision operator [17,18]; (2) easy implementation of complex boundary conditions; (3) the LB method is much easier to parallelize and is far less costly in terms of data exchange due to its explicit scheme and the local interaction. It has been demonstrated that the LB method is far better than conventional numerical methods in a parallel implementation [19].

In the LB community, the first LB model for simulating thermal flows was devised by Alexander *et al.* [20], who

extended the second-order equilibrium distribution of Qian et al. [9] to a third-order model for thermal problems. In the meantime, Qian [21] also proposed a series of higher-order LB models. Since then, several approaches for constructing thermal LB models have been developed, the multispeed approach [20-26], the double-distribution-function (DDF) approach [27-34], and the hybrid approach [35,36]. The multispeed approach is a straightforward extension of the isothermal LB method. In this approach higher-order velocity moments of the density distribution function are used to describe the energy equation. The DDF approach utilizes two different distribution functions, one for the flow field and the other for the temperature field. In the hybrid approach, the flow simulation is accomplished by the LB method, but the temperature field is solved by conventional numerical methods, such as the finite-difference method.

Despite the fact that many efforts have been made from various viewpoints, several limitations still remain in the LB method for thermal flows. The multispeed approach usually suffers from severe numerical instability and the temperature variation simulated is limited to a narrow range [4]. Meanwhile, the boundary treatments become complex when the multispeed lattices are employed: After the streaming process, not only the boundary nodes but also the internal nodes near to the boundaries will possess unknown distribution functions. Lallemand and Luo [35] have found out that the numerical instability of multispeed LB models results from the *spurious* mode coupling in multispeed LB models and cannot be removed by increasing the number of discrete velocities. They suggested that the best approach to remove the spurious mode coupling is to treat the energy-conservation equation

^{*}yalinghe@mail.xjtu.edu.cn

separately from the mass- and momentum-conservation equations. Obviously, such a treatment can be realized by the hybrid and the DDF approaches. Nevertheless, as pointed out by Lallemand and Luo [35], the hybrid approach significantly deviates from the orthodox LB method based on the kinetic theory, and it only provides a compromised solution. Meanwhile, the terms of viscous heat dissipation and compressible work are often ignored in the hybrid approach. Fortunately, these terms can be easily considered in the DDF approach. However, most of the existing DDF LB models are "decoupling" models, i.e., an equation of state with a constant temperature $[p(\rho, T) = \rho c_s^2 = \rho R T_0$, where $c_s = \sqrt{R T_0}$ is the isothermal sound speed, R is the specific gas constant, and T_0 is the reference temperature] is recovered. Such a decoupling causes these models to be limited to Boussinesq flows, in which the temperature variation is small. When these models are applied to the thermal problems in which the temperature field has significant influences on the flow field, the decoupling between momentum and energy transports will result in considerable

Recently, Prasianakis and Karlin [37,38] devised a coupling LB model for simulating thermal flows on standard lattices. Different from previous coupling thermal LB models, Prasianakis and Karlin's model is constructed on the standard two-dimensional nine-velocity (D2Q9) lattice. Hence the basic advantages of the standard LB method are retained. Moreover, the Prandtl number is adjustable in the model. However, the model employs a semiexplicit LB scheme. Quasi-iterations are therefore introduced in the numerical algorithm of the model. In addition, many gradient terms need to be evaluated in numerical computations. Furthermore, the specific-heat ratio cannot be chosen freely. For the two-dimensional model, the specific-heat ratio is fixed at an unphysical value of 2.

In the present study, we aim to propose a coupling LB model for simulating thermal flows on standard lattices in the framework of the DDF approach on the basis of the following considerations. First, it is well known that in the DDF approach the flow and temperature fields are treated separately, which satisfies the previously mentioned criterion suggested by Lallemand and Luo. Second, the viscous heat dissipation and the compression work can be easily included. Moreover, both the Prandtl number and the specific-heat ratio can be made adjustable in the DDF approach. In the proposed model, the discrete equilibrium distribution functions will be obtained from the Hermite expansions of the continuous equilibrium distribution functions. Particularly, the pressure given by the equation of state of perfect gases will be recovered in the macroscopic momentum and energy equations. The rest of this paper is organized as follows: Section II briefly introduces the DDF LB approach. Section III presents the coupling DDF LB model in detail. Both the BGK and MRT versions of the model are shown. In Sec. IV, numerical simulations are carried out for several test problems to validate the proposed model. Finally, a brief conclusion is made in Sec. V.

II. THE DDF LB APPROACH

In this section, we will briefly introduce the DDF LB approach of Guo *et al.* [31]. Historically, the first DDF LB model including the viscous heat dissipation and the

compression work was developed by He *et al.* [28]. Two different distribution functions are employed in He and coworkers' model. The density distribution function is used for the flow field, while an internal energy distribution function is introduced to solve the temperature flied. He and co-workers' DDF LB model has attracted much attention for its excellent numerical stability over the multispeed LB models and the adjustability of the Prandtl number. Following He and co-workers' approach, Guo *et al.* [31] developed a total-energy-distribution-function-based DDF LB approach, which enables DDF LB models to be simpler and makes the inclusion of viscous heat dissipation and compression work easier as compared with He and co-workers' DDF LB approach. The kinetic equations of Guo and co-workers' DDF LB approach are given as follows [31]:

$$\partial_t f + \boldsymbol{\xi} \cdot \boldsymbol{\nabla} f = -\frac{1}{\tau_f} (f - f^{\text{eq}}), \tag{1a}$$

$$\partial_t h + \boldsymbol{\xi} \cdot \boldsymbol{\nabla} h = -\frac{1}{\tau_h} (h - h^{\text{eq}}) + \frac{Z}{\tau_{hf}} (f - f^{\text{eq}}),$$
 (1b)

where f and h are the density distribution function and the total energy distribution function, respectively; τ_f and τ_h are the momentum and total energy relaxation times, respectively; τ_{hf} is a relaxation time related to τ_f and τ_h ; and $Z = \xi \cdot u - u^2/2$, where ξ and u are the particle velocity and the macroscopic velocity, respectively. The equilibrium distribution functions $f^{\rm eq}$ and $h^{\rm eq}$ are given by

$$f^{\text{eq}} = \frac{\rho}{(2\pi RT)^{D/2}} \exp\left[-\frac{(\xi - u)^2}{2RT}\right],$$

$$h^{\text{eq}} = \frac{1}{2}\xi^2 f^{\text{eq}} = \frac{\rho \xi^2}{2(2\pi RT)^{D/2}} \exp\left[-\frac{(\xi - u)^2}{2RT}\right],$$
(2)

where $\xi^2 = \xi \cdot \xi$. By projecting f^{eq} and h^{eq} onto the tensor Hermite polynomial basis in terms of the particle velocity ξ , Guo *et al.* obtained the following Hermite expansions of f^{eq} and h^{eq} at the Navier-Stokes level, respectively:

$$f^{\text{eq,3}}(T) = \rho \omega(\xi, T) \left\{ 1 + \frac{\xi \cdot u}{RT} + \frac{1}{2} \left(\frac{\xi \cdot u}{RT} \right)^2 - \frac{u^2}{2RT} + \frac{\xi \cdot u}{6RT} \left[\left(\frac{\xi \cdot u}{RT} \right)^2 - \frac{3u^2}{RT} \right] \right\},$$
(3a)
$$h^{\text{eq,2}}(T) = \omega(\xi, T) \left\{ \rho E + (p + \rho E) \frac{\xi \cdot u}{RT} + \frac{p}{2} \left(\frac{\xi^2}{RT} - D \right) + \left(p + \frac{\rho E}{2} \right) \left[\left(\frac{\xi \cdot u}{RT} \right)^2 - \frac{u^2}{RT} \right] \right\},$$
(3b)

where $f^{\text{eq},3}(T)$ is the third-order expansion of f^{eq} , $h^{\text{eq},2}(T)$ is the second-order expansion of h^{eq} , E is the macroscopic total energy, $p = \rho RT$ is the pressure, and $\omega(\xi, T)$ is given by

$$\omega(\xi, T) = \frac{1}{(2\pi RT)^{D/2}} \exp\left(-\frac{\xi^2}{2RT}\right).$$
 (4)

The discrete velocity set e_{α} can be obtained by choosing the abscissae of a suitable Gauss-Hermite quadrature with the weight function $\omega(\xi, T)$. However, as pointed out by Guo *et al.*, the temperature in $\omega(\xi, T)$ is a locally changed variable, which means the abscissae of the Gauss-Hermite quadrature are not fixed. Then the discrete velocity e_{α} will change with the local temperature. To overcome this difficulty, Guo *et al.* replace the local temperature T in $f^{\text{eq},3}(T)$ and $h^{\text{eq},2}(T)$ with a reference temperature T_0 [31]:

$$f^{\text{eq},3}(T_0) = \rho \omega(\boldsymbol{\xi}, T_0) \left\{ 1 + \frac{\boldsymbol{\xi} \cdot \boldsymbol{u}}{RT_0} + \frac{1}{2} \left(\frac{\boldsymbol{\xi} \cdot \boldsymbol{u}}{RT_0} \right)^2 - \frac{\boldsymbol{u}^2}{2RT_0} + \frac{\boldsymbol{\xi} \cdot \boldsymbol{u}}{6RT_0} \left[\left(\frac{\boldsymbol{\xi} \cdot \boldsymbol{u}}{RT_0} \right)^2 - \frac{3\boldsymbol{u}^2}{RT_0} \right] \right\}, \quad (5a)$$

$$h^{\text{eq},2}(T_0)$$

$$= \omega(\boldsymbol{\xi}, T_0) \left\{ \rho E + (p_0 + \rho E) \frac{\boldsymbol{\xi} \cdot \boldsymbol{u}}{RT_0} + \frac{p_0}{2} \left(\frac{\boldsymbol{\xi}^2}{RT_0} - D \right) + \left(p_0 + \frac{\rho E}{2} \right) \left[\left(\frac{\boldsymbol{\xi} \cdot \boldsymbol{u}}{RT_0} \right)^2 - \frac{\boldsymbol{u}^2}{RT_0} \right] \right\}, \tag{5b}$$

where $p_0 = \rho RT_0$. Note that the third-order velocity terms in Eq. (5a) can be neglected in the low Mach number limit.

With regard to the discretization of velocity space, Guo *et al.* adopted a nine-point fifth-degree Gauss-Hermite quadrature, which leads to the following discrete velocity (the D2Q9 lattice):

$$e_{\alpha} = \begin{cases} (0,0) & \alpha = 0, \\ c\left(\cos[(\alpha - 1)\pi/2], \sin[(\alpha - 1)\pi/2]\right) & \alpha = 1,2,3,4, \\ \sqrt{2}c\left(\cos[(2\alpha - 1)\pi/4], \sin[(2\alpha - 1)\pi/4]\right) & \alpha = 5,6,7,8, \end{cases}$$
(6)

where $c = \sqrt{3RT_0}$. According to Eq. (1), the kinetic equations for the discrete density distribution function f_{α} and the discrete total energy distribution function h_{α} are given by [31]

$$\partial_t f_\alpha + \boldsymbol{e}_\alpha \cdot \nabla f_\alpha = -\frac{1}{\tau_f} (f_\alpha - f_\alpha^{\text{eq}}),$$
 (7a)

$$\partial_t h_\alpha + \boldsymbol{e}_\alpha \cdot \nabla h_\alpha = -\frac{1}{\tau_h} (h_\alpha - h_\alpha^{\text{eq}}) + \frac{Z_\alpha}{\tau_{hf}} (f_\alpha - f_\alpha^{\text{eq}}), \quad (7b)$$

where $Z_{\alpha} = \boldsymbol{e}_{\alpha} \cdot \boldsymbol{u} - u^2/2$. At the Navier-Stokes level, Z_{α} can be simplified as $Z_{\alpha} = \boldsymbol{e}_{\alpha} \cdot \boldsymbol{u}$. The discrete equilibrium distribution functions $f_{\alpha}^{\rm eq}$ and $h_{\alpha}^{\rm eq}$ are given by

$$f_{\alpha}^{\text{eq}} = \rho w_{\alpha} \left[1 + \frac{\boldsymbol{e}_{\alpha} \cdot \boldsymbol{u}}{RT_{0}} + \frac{1}{2} \left(\frac{\boldsymbol{e}_{\alpha} \cdot \boldsymbol{u}}{RT_{0}} \right)^{2} - \frac{\boldsymbol{u}^{2}}{2RT_{0}} \right], \quad (8a)$$

$$h_{\alpha}^{\text{eq}} = E f_{\alpha}^{\text{eq}} + w_{\alpha} p_{0} \left[\frac{(\boldsymbol{e}_{\alpha} \cdot \boldsymbol{u})}{RT_{0}} + \left(\frac{\boldsymbol{e}_{\alpha} \cdot \boldsymbol{u}}{RT_{0}} \right)^{2} - \frac{\boldsymbol{u}^{2}}{RT_{0}} + \frac{1}{2} \left(\frac{\boldsymbol{e}_{\alpha}^{2}}{RT_{0}} - D \right) \right]. \quad (8b)$$

The coefficients are $w_0 = 4/9$, $w_{1, 2, 3, 4} = 1/9$, and $w_{5, 6, 7, 8} = 1/36$. Through the Chapman-Enskog analysis, Guo *et al.* demonstrated that Eq. (7) together with Eq. (8) can recover the following equations in the low Mach number limit:

$$\partial_{t}\rho + \nabla \cdot (\rho \mathbf{u}) = 0, \qquad (9a)$$

$$\partial_{t}(\rho \mathbf{u}) + \nabla \cdot (\rho \mathbf{u} \mathbf{u}) = -\nabla p_{0} + \nabla \cdot \mathbf{\Pi}, \qquad (9b)$$

$$\partial_{t}(\rho E) + \nabla \cdot [(\rho E + p_{0})\mathbf{u}] = \nabla \cdot (\lambda \nabla T) + \nabla \cdot (\mathbf{u} \cdot \mathbf{\Pi}),$$

where $\Pi = \mu [\nabla u + (\nabla u)^T]$ is the viscous stress tensor. It can be seen that Eq. (9) is a decoupling Navier-Stokes-Fourier (NSF) equation because the equation of state is given by

 $p(\rho, T) = p_0 = \rho RT_0$, which makes the momentum transport decoupled from the energy transport. The model is therefore a decoupling model and is limited to Boussinesq flows, in which the temperature variation is small.

III. THE COUPLING DDF LB MODEL

A. Hermite expansions of continuous equilibrium distribution functions

From Sec. II it can be seen that the decoupling between the momentum and energy transports in Guo and co-workers' DDF LB model results from replacing the local temperature T in $f^{\text{eq},3}(T)$ and $h^{\text{eq},2}(T)$ with the reference temperature T_0 . It is noted that $f^{\text{eq},3}(T)$ given by Eq. (3a) satisfies the zeroth-through third-order velocity moments of the Maxwell distribution function f^{eq} :

$$\int f^{\text{eq},3}(T)d\boldsymbol{\xi} \equiv \int f^{\text{eq}}d\boldsymbol{\xi} = \rho, \tag{10a}$$

$$\int f^{\text{eq},3}(T)\xi_i d\boldsymbol{\xi} \equiv \int f^{\text{eq}}\xi_i d\boldsymbol{\xi} = \rho u_i, \tag{10b}$$

$$\int f^{\text{eq},3}(T)\xi_i\xi_j d\boldsymbol{\xi} \equiv \int f^{\text{eq}}\xi_i\xi_j d\boldsymbol{\xi} = \rho u_i u_j + \rho RT\delta_{ij}, (10c)$$

$$\int f^{\text{eq},3}(T)\xi_i\xi_j\xi_k d\boldsymbol{\xi}
\equiv \int f^{\text{eq}}\xi_i\xi_j\xi_k d\boldsymbol{\xi}
= \rho u_i u_j u_k + \rho RT(u_k \delta_{ij} + u_i \delta_{jk} + u_j \delta_{ik}), \quad (10d)$$

where δ is the Kronecker delta with two indices. However, when the local temperature in $f^{\text{eq},3}(T)$ is replaced with the reference temperature T_0 , the following relationships will be

(9c)

obtained:

$$\int f^{\text{eq},3}(T_0)\xi_i\xi_j d\boldsymbol{\xi} = \rho u_i u_j + \rho R T_0 \delta_{ij}, \qquad (11a)$$

$$\int f^{\text{eq},3}(T_0)\xi_i\xi_j\xi_k d\boldsymbol{\xi}$$

$$= \rho u_i u_j u_k + \rho R T_0 (u_k \delta_{ij} + u_i \delta_{jk} + u_j \delta_{ik}). \qquad (11b)$$

Similar results will be obtained for $h^{\text{eq},2}(T_0)$. As a consequence, the pressure in the macroscopic momentum and energy equations is given by $p(\rho, T) = p_0 = \rho R T_0$ [see Eq. (9)].

In order to overcome the above problem, in the present study we employ another type of Hermite expansion of the continuous equilibrium distribution function. As summarized in Refs. [39,40], in the LB method there exist three different types of Hermite expansion of f^{eq} . Here we adopt Shan and co-workers' third-order Hermite expansion of f^{eq} , which is given as follows [23]:

$$f^{\text{eq*,3}}(T) = \rho \omega(\xi, T_0) \left\{ 1 + \frac{\xi \cdot u}{RT_0} + \frac{1}{2} \left[\left(\frac{\xi \cdot u}{RT_0} \right)^2 - \frac{u^2}{RT_0} + (\theta - 1) \left(\frac{\xi^2}{RT_0} - D \right) \right] + \frac{\xi \cdot u}{6RT_0} \left[\left(\frac{\xi \cdot u}{RT_0} \right)^2 - \frac{3u^2}{RT_0} + 3(\theta - 1) \left(\frac{\xi^2}{RT_0} - D - 2 \right) \right] \right\}, \quad (12)$$

where $\theta = T/T_0$. It can be proved that $f^{\text{eq*,3}}(T)$ satisfies the zeroth-through third-order velocity moments of f^{eq} . In the low Mach number limit, the third-order velocity terms in $f^{\text{eq*,3}}(T)$ can be neglected. Then we have

$$f^{\text{eq*,3}}(T) = \rho \omega(\boldsymbol{\xi}, T_0) \left\{ 1 + \frac{\boldsymbol{\xi} \cdot \boldsymbol{u}}{RT_0} + \frac{1}{2} \left[\left(\frac{\boldsymbol{\xi} \cdot \boldsymbol{u}}{RT_0} \right)^2 - \frac{\boldsymbol{u}^2}{RT_0} + (\theta - 1) \left(\frac{\boldsymbol{\xi}^2}{RT_0} - D \right) \right] + \frac{\boldsymbol{\xi} \cdot \boldsymbol{u}}{2RT_0} (\theta - 1) \left(\frac{\boldsymbol{\xi}^2}{RT_0} - D - 2 \right) \right\}. \quad (13)$$

Inspired by the approach of Shan *et al.* for the Hermite expansion of f^{eq} [23], after some algebra we obtain the following second-order Hermite expansion of h^{eq} (see Appendix A for details):

$$h^{\text{eq*,2}}(T) = \omega(\boldsymbol{\xi}, T_0) \left\{ \rho E + (p + \rho E) \frac{\boldsymbol{\xi} \cdot \boldsymbol{u}}{RT_0} + \left(p + \frac{\rho E}{2} \right) \right\} \times \left[\left(\frac{\boldsymbol{\xi} \cdot \boldsymbol{u}}{RT_0} \right)^2 - \frac{\boldsymbol{u}^2}{RT_0} \right] + \frac{p\theta}{2} \left(\frac{\boldsymbol{\xi}^2}{RT_0} - D \right) + \frac{\rho E}{2} (\theta - 1) \left(\frac{\boldsymbol{\xi}^2}{RT_0} - D \right) \right\}.$$
(14)

It can be verified that $h^{\text{eq*},2}(T)$ satisfies the zeroth-through second-order velocity moments of h^{eq} :

$$\int h^{\text{eq}*,2}(T)d\boldsymbol{\xi} \equiv \int h^{\text{eq}}d\boldsymbol{\xi} = \rho E,$$

$$\int h^{\text{eq}*,2}(T)\xi_i d\boldsymbol{\xi} \equiv \int h^{\text{eq}}\xi_i d\boldsymbol{\xi} = (\rho E + p)u_i,$$
(15a)

$$\int h^{\text{eq*,2}}(T)\xi_i\xi_j d\xi \equiv \int h^{\text{eq}}\xi_i\xi_j d\xi$$
$$= (\rho E + 2p)u_i u_j + p(E + RT)\delta_{ij},$$
(15c)

where $p = \rho RT$. From Eqs. (5), (12), and (14) we can see that $f^{\text{eq*,3}}(T)$ and $h^{\text{eq*,2}}(T)$ share a similar feature with $f^{\text{eq,3}}(T_0)$ and $h^{\text{eq,2}}(T_0)$. Namely, the weight function $\omega(\xi, T_0)$ is independent of the local temperature. Hence the discrete velocity e_{α} will not change with the local temperature. However, as previously mentioned, $f^{\text{eq,3}}(T_0)$ and $h^{\text{eq,2}}(T_0)$ cannot satisfy the related statistical relationships, which makes the pressure in the macroscopic momentum and energy equations to be given by $p(\rho, T) = \rho RT_0$. Note that $h^{\text{eq*,2}}(T)$ can also be written as follows:

$$h^{\text{eq*,2}}(T) = \rho E \omega(\boldsymbol{\xi}, T_0) \left\{ 1 + \frac{\boldsymbol{\xi} \cdot \boldsymbol{u}}{RT_0} + \frac{1}{2} \left[\left(\frac{\boldsymbol{\xi} \cdot \boldsymbol{u}}{RT_0} \right)^2 - \frac{\boldsymbol{u}^2}{RT_0} + (\theta - 1) \left(\frac{\boldsymbol{\xi}^2}{RT_0} - D \right) \right] \right\} + \omega(\boldsymbol{\xi}, T_0) p$$

$$\times \left[\frac{\boldsymbol{\xi} \cdot \boldsymbol{u}}{RT_0} + \left(\frac{\boldsymbol{\xi} \cdot \boldsymbol{u}}{RT_0} \right)^2 - \frac{\boldsymbol{u}^2}{RT_0} + \frac{\theta}{2} \left(\frac{\boldsymbol{\xi}^2}{RT_0} - D \right) \right]. \tag{16}$$

Therefore, for convenience, we can use the following form of $h^{\text{eq}*,2}(T)$:

$$h^{\text{eq*,2}}(T) = Ef^{\text{eq*,3}}(T) + \omega(\boldsymbol{\xi}, T_0)p \left[\frac{\boldsymbol{\xi} \cdot \boldsymbol{u}}{RT_0} + \left(\frac{\boldsymbol{\xi} \cdot \boldsymbol{u}}{RT_0} \right)^2 - \frac{\boldsymbol{u}^2}{RT_0} + \frac{\theta}{2} \left(\frac{\boldsymbol{\xi}^2}{RT_0} - D \right) \right], \tag{17}$$

where $f^{\text{eq}*,3}(T)$ is given by Eq. (13).

B. Discrete equilibrium distribution functions

According to Eqs. (13) and (17), the following discrete equilibrium distribution functions can be obtained when discretizing ξ to e_{α} :

$$f_{\alpha}^{\text{eq}} = \rho w_{\alpha} \left\{ 1 + \frac{\boldsymbol{e}_{\alpha} \cdot \boldsymbol{u}}{RT_{0}} + \frac{1}{2} \left[\left(\frac{\boldsymbol{e}_{\alpha} \cdot \boldsymbol{u}}{RT_{0}} \right)^{2} - \frac{\boldsymbol{u}^{2}}{RT_{0}} + (\theta - 1) \left(\frac{\boldsymbol{e}_{\alpha}^{2}}{RT_{0}} - D \right) \right] + \frac{\boldsymbol{e}_{\alpha} \cdot \boldsymbol{u}}{2RT_{0}} (\theta - 1) \left(\frac{\boldsymbol{e}_{\alpha}^{2}}{RT_{0}} - D - 2 \right) \right\}, \quad (18a)$$

$$h_{\alpha}^{\text{eq}} = E f_{\alpha}^{\text{eq}} + w_{\alpha} p \left[\frac{\boldsymbol{e}_{\alpha} \cdot \boldsymbol{u}}{RT_{0}} + \left(\frac{\boldsymbol{e}_{\alpha} \cdot \boldsymbol{u}}{RT_{0}} \right)^{2} - \frac{\boldsymbol{u}^{2}}{RT_{0}} + \frac{\theta}{2} \left(\frac{\boldsymbol{e}_{\alpha}^{2}}{RT_{0}} - D \right) \right]. \quad (18b)$$

Obviously, when $\theta = T/T_0 = 1$ and $p = p_0$, the above discrete equilibrium distribution functions will reduce to those in the decoupling DDF model [see Eq. (8)]. When the D2Q9 lattice is adopted, it can be found that f_{α}^{eq} and h_{α}^{eq} rigorously

satisfy the zeroth- through second-order velocity moments of f^{eq} and h^{eq} , respectively, i.e.,

$$\sum_{\alpha} f_{\alpha}^{\text{eq}} = \rho, \quad \sum_{\alpha} f_{\alpha}^{\text{eq}} e_{\alpha i} = \rho u_{i},$$

$$\sum_{\alpha} f_{\alpha}^{\text{eq}} e_{\alpha i} e_{\alpha j} = \rho u_{i} u_{j} + \rho RT \delta_{ij}, \tag{19a}$$

$$\sum_{\alpha} h_{\alpha}^{\text{eq}} = \rho E, \ \sum_{\alpha} h_{\alpha}^{\text{eq}} e_{\alpha i} = (\rho E + p) u_i,$$

$$\sum_{\alpha} h_{\alpha}^{\text{eq}} e_{\alpha i} e_{\alpha j} = (\rho E + 2p) u_i u_j + p(E + RT) \delta_{ij}. \tag{19b}$$

However, taking the third-order velocity moment of f_{α}^{eq} with the D2Q9 lattice, we will obtain

$$\sum_{\alpha} f_{\alpha}^{eq} e_{\alpha i} e_{\alpha j} e_{\alpha k} = \rho R T (u_k \delta_{ij} + u_j \delta_{ik} + u_i \delta_{jk})$$

$$+ 3\rho R T_0 u_l (1 - \theta) \delta_{ijkl}, \qquad (20)$$

where $\delta_{ijkl} = 1$ when i = j = k = l, otherwise $\delta_{ijkl} = 0$. Equation (20) can be rewritten as follows:

$$\sum_{\alpha} f_{\alpha}^{\text{eq}} e_{\alpha i} e_{\alpha j} e_{\alpha k}$$

$$= \begin{cases} \rho R T_0(u_k \delta_{ij} + u_j \delta_{ik} + u_i \delta_{jk}), & \text{if } i = j = k, \\ \rho R T(u_k \delta_{ij} + u_j \delta_{ik} + u_i \delta_{jk}), & \text{others.} \end{cases}$$
(21)

This means that, when the D2Q9 lattice is used, the off-diagonal elements of the third-order velocity moment of f_{α}^{eq} satisfy the related statistical relationship [see Eq. (10d) and note that the third-order velocity term can be neglected as done in the standard LB method with the low Mach number limit], but the diagonal elements deviate from the relationship. This deviation results from the low symmetry of the D2Q9 lattice and cannot be removed by choosing f_{α}^{eq} different from Eq. (18a).

For standard lattices, as pointed out by Prasianakis and Karlin [37], the deviation of f_{α}^{eq} can be removed only by introducing a correction term into the Boltzmann-BGK equation (7a):

$$\partial_t f_{\alpha} + \boldsymbol{e}_{\alpha} \cdot \nabla f_{\alpha} = -\frac{1}{\tau_f} (f_{\alpha} - f_{\alpha}^{\text{eq}}) + C_{\alpha},$$
 (22)

where C_{α} is a correction term. The constraints on C_{α} can be established with the Chapman-Enskog analysis. By introducing the following multiscale expansions,

$$\begin{aligned}
\partial_t &= \varepsilon \partial_{t_1} + \varepsilon^2 \partial_{t_2}, \quad \nabla = \varepsilon \nabla_1, \\
f_\alpha &= f_\alpha^{\text{eq}} + \varepsilon f_\alpha^{(1)} + \varepsilon f_\alpha^{(2)}, \quad C_\alpha = \varepsilon C_\alpha^{(0)},
\end{aligned} \tag{23}$$

where ε is the expansion parameter, we can rewrite Eq. (22) in the consecutive orders of ε :

$$\varepsilon^{1}: \left(\partial_{t_{1}} + \boldsymbol{e}_{\alpha} \cdot \nabla_{1}\right) f_{\alpha}^{\text{eq}} + \frac{1}{\tau_{f}} f_{\alpha}^{(1)} = C_{\alpha}^{(0)}, \tag{24}$$

$$\varepsilon^{2}: \partial_{t_{2}} f_{\alpha}^{(1)} + (\partial_{t_{1}} + \boldsymbol{e}_{\alpha} \cdot \nabla_{1}) f_{\alpha}^{(1)} + \frac{1}{\tau_{f}} f_{\alpha}^{(2)} = 0. \quad (25)$$

To recover the correct macroscopic equations, the first two velocity moments of C_{α} should satisfy

$$\sum_{\alpha} C_{\alpha} = 0, \quad \sum_{\alpha} e_{\alpha} C_{\alpha} = 0.$$
 (26)

Furthermore, taking the second-order velocity moment of Eq. (24) leads to

$$\Pi_{ij} = -\sum_{\alpha} e_{\alpha i} e_{\alpha j} f_{\alpha}^{(1)} = \tau_{f} \left[\partial_{t_{1}} \left(\sum_{\alpha} e_{\alpha i} e_{\alpha j} f_{\alpha}^{\text{eq}} \right) + \partial_{1k} \left(\sum_{\alpha} e_{\alpha i} e_{\alpha j} e_{\alpha k} f_{\alpha}^{\text{eq}} \right) - \sum_{\alpha} e_{\alpha i} e_{\alpha j} C_{\alpha}^{(0)} \right]. \quad (27)$$

According to Eqs. (20) and (27), we can find that, in order to remove the deviation, the second-order velocity moment of C_{α} should be defined as

$$\sum_{\alpha} e_{\alpha x} e_{\alpha y} C_{\alpha} = 0, \quad \sum_{\alpha} e_{\alpha x}^{2} C_{\alpha} = \partial_{x} [\rho c^{2} (1 - \theta) u_{x}],$$

$$\sum_{\alpha} e_{\alpha y}^{2} C_{\alpha} = \partial_{y} [\rho c^{2} (1 - \theta) u_{y}].$$
(28)

Here $c = \sqrt{3RT_0} = 1$ is the lattice speed. Equations (26) and (28) are the statistical constraints on the correction term C_{α} . With these constraints, the deviation results from the low symmetry of the D2Q9 lattice can be removed. The detailed form of C_{α} will be given in the next section.

The full Chapman-Enskog analysis of Eqs. (22) and (7b) with the equilibrium distribution functions given by Eq. (18) is presented in Appendix B. The following macroscopic equations are obtained at the Navier-Stokes level:

$$\partial_t \rho + \nabla \cdot (\rho \mathbf{u}) = 0, \tag{29a}$$

$$\partial_t(\rho \boldsymbol{u}) + \nabla \cdot (\rho \boldsymbol{u} \boldsymbol{u}) = -\nabla p + \nabla \cdot \boldsymbol{\Pi}, \tag{29b}$$

$$\partial_t(\rho E) + \nabla \cdot [(\rho E + p)\mathbf{u}] = \nabla \cdot (\lambda \nabla T) + \nabla \cdot (\mathbf{u} \cdot \mathbf{\Pi}),$$
(29c)

where $\lambda = \tau_h[(b+2)/2]pR$ is the thermal conductivity, in which *b* is the number of degrees of freedom of the gas. The viscous stress tensor in Eq. (29) is given by

$$\mathbf{\Pi} = \mu \left[\nabla \mathbf{u} + (\nabla \mathbf{u})^T - \frac{2}{3} (\nabla \cdot \mathbf{u}) \mathbf{I} \right] + \mu_{\mathbf{B}} (\nabla \cdot \mathbf{u}) \mathbf{I}, \quad (30)$$

where $\mu = \tau_f p$ is the dynamic viscosity, $\mu_B = 2(1/3 - 1/b)\mu$ is the bulk viscosity, and **I** is the unit tensor. The Prandtl number of the system is given by $\Pr = \tau_f/\tau_h$. An important difference between Eqs. (29) and (9) can be observed although they take the same form: The pressure in Eqs. (29b) and (29c) is given by $p = \rho RT$, while in Eqs. (9b) and (9c) the pressure is $p_0 = \rho RT_0$.

C. Time-discrete evolution equations

1. BGK scheme

The time discretization of the Boltzmann-BGK equation (22) can be made by integrating the equation along the characteristic line e_{α} over a time interval of length

 δ_t , which leads to [28,31]

$$f_{\alpha}(\mathbf{x} + \mathbf{e}_{\alpha}\delta_{t}, t + \delta_{t}) - f_{\alpha}(\mathbf{x}, t)$$

$$= \int_{0}^{\delta_{t}} [\Omega_{f}(\mathbf{x} + \mathbf{e}_{\alpha}t', t + t') + C_{\alpha}(\mathbf{x} + \mathbf{e}_{\alpha}t', t + t')] dt,$$
(31)

where $\Omega_f = (f_{\alpha}^{eq} - f_{\alpha})/\tau_f$. Applying the trapezoidal rule to the integral on the right-hand side yields

$$f_{\alpha}(\mathbf{x} + \mathbf{e}_{\alpha}\delta_{t}, t + \delta_{t}) - f_{\alpha}(\mathbf{x}, t)$$

$$= \frac{\delta_{t}}{2} [\Omega_{f}(\mathbf{x} + \mathbf{e}_{\alpha}\delta_{t}, t + \delta_{t}) + \Omega_{f}(\mathbf{x}, t)]$$

$$+ \frac{\delta_{t}}{2} [C_{\alpha}(\mathbf{x} + \mathbf{e}_{\alpha}\delta_{t}, t + \delta_{t}) + C_{\alpha}(\mathbf{x}, t)]. \quad (32)$$

Following He *et al.* [28] and Guo *et al.* [31], the implicitness of Eq. (32) can be eliminated by

$$\bar{f}_{\alpha} = f_{\alpha} - \frac{\delta_t}{2} (\Omega_f + C_{\alpha}). \tag{33}$$

Then Eq. (32) can be rewritten as

$$\bar{f}_{\alpha}(\mathbf{x} + \mathbf{e}_{\alpha}\delta_{t}, t + \delta_{t}) - \bar{f}_{\alpha}(\mathbf{x}, t)
= -\omega_{f} \left[\bar{f}_{\alpha}(\mathbf{x}, t) - f_{\alpha}^{eq}(\mathbf{x}, t) \right] - \delta_{t} (1 - 0.5\omega_{f}) C_{\alpha}(\mathbf{x}, t),$$
(34)

where $\omega_f = 2\delta_t/(2\tau_f + \delta_t)$. According to Eq. (33), the macroscopic density and velocity can be explicitly computed from \bar{f}_{α} as follows (note that $\sum_{\alpha} C_{\alpha} = \sum_{\alpha} e_{\alpha} C_{\alpha} = 0$):

$$\rho = \sum_{\alpha} \bar{f}_{\alpha}, \quad \rho u = \sum_{\alpha} e_{\alpha} \bar{f}_{\alpha}. \tag{35}$$

Similarly, from Eq. (8b), the following time-discrete evolution equation can be obtained:

$$\bar{h}_{\alpha}(\mathbf{x} + \mathbf{e}_{\alpha}\delta_{t}, t + \delta_{t}) - \bar{h}_{\alpha}(\mathbf{x}, t)
= -\omega_{h} \left[\bar{h}_{\alpha}(\mathbf{x}, t) - h_{\alpha}^{\text{eq}}(\mathbf{x}, t) \right]
+ (\omega_{h} - \omega_{f}) Z_{\alpha} \left[\bar{f}_{\alpha}(\mathbf{x}, t) - f_{\alpha}^{\text{eq}}(\mathbf{x}, t) + \frac{\delta_{t}}{2} C_{\alpha}(\mathbf{x}, t) \right],$$
(36)

where $\omega_h = 2\delta_t/(2\tau_h + \delta_t)$ and $\bar{h}_\alpha = h_\alpha - 0.5\delta_t(\Omega_h - \tau_f Z_\alpha \Omega_f/\tau_{hf})$, in which $\Omega_h = (h_\alpha^{\rm eq} - h_\alpha)/\tau_h$ and $\tau_{hf} = \tau_h \tau_f/(\tau_f - \tau_h)$. The macroscopic total energy can be obtained via $\rho E = \sum_\alpha \bar{h}_\alpha$. According to the definition of the total energy $E = (bRT + u^2)/2$, the macroscopic temperature is given by

$$T = \frac{2}{bR} \left(\sum_{\alpha} \bar{h}_{\alpha} / \rho - \frac{1}{2} \boldsymbol{u}^2 \right). \tag{37}$$

The specific-heat ratio γ is related to b via $\gamma = c_p/c_v = (b+2)/b$, where $c_p = (b+2)R/2$ and $c_v = bR/2$ are the specific-heat coefficients at constant pressure and volume, respectively.

Note that, if the viscous heat dissipation and the compression work are negligible in the flow, we can directly adopt an internal energy distribution function g_{α} to solve the

temperature field. Its evolution equation can be given by

$$g_{\alpha}(\mathbf{x} + \mathbf{e}_{\alpha}\delta_{t}, t + \delta_{t}) - g_{\alpha}(\mathbf{x}, t)$$

$$= -\omega_{g} [g_{\alpha}(\mathbf{x}, t) - g_{\alpha}^{eq}(\mathbf{x}, t)], \tag{38}$$

where $g_{\alpha}^{\rm eq}=bRTf_{\alpha}^{\rm eq}/2$ is the equilibrium internal energy distribution and $\omega_g=2\delta_t/(2\tau_g+\delta_t)$, in which τ_g is the relaxation time for g_{α} . The macroscopic temperature and the thermal diffusivity are calculated by $T=2\sum_{\alpha}g_{\alpha}/(\rho bR)$ and $\chi=\tau_g p$, respectively.

2. MRT scheme

In the LB community, it has been accepted that the MRT collision operator can improve the numerical stability of the LB equation as compared with the BGK collision operator. The relaxation times for hydrodynamic and nonhydrodynamic moments can be separated when employing the MRT collision operator [11–16]. By replacing the BGK collision operator with the MRT collision operator, Eq. (22) can be rewritten as

$$\partial_t f_{\alpha} + \boldsymbol{e}_{\alpha} \cdot \nabla f_{\alpha} = -\Lambda_{\alpha\beta} (f_{\beta} - f_{\beta}^{\text{eq}}) + C_{\alpha},$$
 (39)

where $\mathbf{\Lambda} = \mathbf{M}^{-1}\mathbf{S}\mathbf{M}$ is the collision matrix, \mathbf{M} is an orthogonal transformation matrix (see Appendix B), and S is a diagonal matrix given by

$$\mathbf{S} = \operatorname{diag}(\tau_{\rho}^{-1}, \, \tau_{e}^{-1}, \, \tau_{\varsigma}^{-1}, \, \tau_{j}^{-1}, \, \tau_{q}^{-1}, \, \tau_{j}^{-1}, \, \tau_{q}^{-1}, \, \tau_{f}^{-1}, \, \tau_{f}^{-1}). \tag{40}$$

Through the transformation matrix, the density distribution function f_{α} can be projected onto the moment space with $\mathbf{m} = \mathbf{Mf}$, where $\mathbf{f} = (f_0, f_1, \dots, f_8)^{\mathrm{T}}$. For the D2Q9 lattice, \mathbf{m} is defined as $\mathbf{m} = (\rho, e, \varsigma, j_x, q_x, j_y, q_y, p_{xx}, p_{xy})^T$, where e is the energy mode, ς is related to energy square, (j_x, j_y) are the momentum components, (q_x, q_y) correspond to the energy flux; and (p_{xx}, p_{xy}) are related to the diagonal and off-diagonal components of the stress tensors [12]. Similarly, f_{α}^{eq} can be projected onto the moment space with $\mathbf{m}^{\mathrm{eq}} = \mathbf{Mf}^{\mathrm{eq}}$, where $\mathbf{f}^{\mathrm{eq}} = (f_0^{\mathrm{eq}}, \dots, f_8^{\mathrm{eq}})^{\mathrm{T}}$. According to Eq. (18a), after some algebra we can obtain the equilibria \mathbf{m}^{eq} of the present model as follows:

$$\mathbf{m}^{\text{eq}} = (\rho, e^{\text{eq}}, \varsigma^{\text{eq}}, j_x, q_x^{\text{eq}}, j_y, q_y^{\text{eq}}, p_{xx}^{\text{eq}}, p_{xy}^{\text{eq}})^T$$

$$= \rho [1, (-4 + 3u^2 + 2\theta), (3 - 3u^2 - 2\theta), u_x,$$

$$\times (-2u_x + u_x\theta), u_y, (-2u_y + u_y\theta), u_x^2 - u_y^2, u_x u_y]^T.$$
(41)

When $\theta = T/T_0 = 1$, the equilibria will reduce to those in the isothermal D2Q9 MRT model [41,42]. It should be noted that the moment corresponding to the energy, $m_1 = e$, is not a conserved moment. The energy conservation is separately described by the total energy distribution function.

The time discretization of Eq. (39) can also be made by integrating the equation along the characteristic line. Nevertheless, the MRT collision term needs to be integrated explicitly. Then the following equation will be obtained:

$$f_{\alpha}(\mathbf{x} + \mathbf{e}_{\alpha}\delta_{t}, t + \delta_{t}) - f_{\alpha}(\mathbf{x}, t)$$

$$= -\delta_{t}\Lambda_{\alpha\beta}.\left(f_{\beta} - f_{\beta}^{\text{eq}}\right)\Big|_{(x, t)} + \frac{\delta_{t}}{2}(C_{\alpha}|_{(\mathbf{x} + \mathbf{e}_{\alpha}\delta_{t}, t + \delta_{t})} + C_{\alpha}|_{(\mathbf{x}, t)}).$$
(42)

By using $\hat{f}_{\alpha} = f_{\alpha} - 0.5\delta_t C_{\alpha}$, we can rewrite Eq. (42) as follows:

$$\widehat{f}_{\alpha}(\mathbf{x} + \mathbf{e}_{\alpha}\delta_{t}, t + \delta_{t})
= \widehat{f}_{\alpha}(\mathbf{x}, t) - \delta_{t}\Lambda_{\alpha\beta}(\widehat{f}_{\beta} - f_{\beta}^{\text{eq}})\big|_{(\mathbf{x}, t)}
+ \delta_{t}(C_{\alpha} - 0.5\Lambda_{\alpha\beta}C_{\beta})\big|_{(\mathbf{x}, t)}.$$
(43)

The right-hand side of Eq. (43) can be carried out in the moment space:

$$\hat{\mathbf{m}}^* = \hat{\mathbf{m}} - \delta_t \mathbf{S}(\hat{\mathbf{m}} - \mathbf{m}^{\text{eq}}) + \delta_t \left(\mathbf{I} - \frac{\mathbf{S}}{2} \right) \hat{\mathbf{C}}, \tag{44}$$

where $\widehat{\mathbf{m}} = \mathbf{M}\widehat{\mathbf{f}}$ and $\widehat{\mathbf{C}} = \mathbf{M}\mathbf{C}$, in which $\mathbf{C} = (C_0, C_1, \dots, C_8)^{\mathbf{T}}$. The streaming process is given by

$$\widehat{f}_{\alpha}(\mathbf{x} + \mathbf{e}_{\alpha}\delta_{t}, t + \delta_{t}) = \widehat{f}_{\alpha}^{*}(\mathbf{x}, t), \tag{45}$$

where $\hat{\boldsymbol{f}}^* = \mathbf{M}^{-1} \hat{\mathbf{m}}^*$. The macroscopic density and velocity are calculated by

$$\rho = \sum_{\alpha} \hat{f}_{\alpha}, \quad \rho u = \sum_{\alpha} e_{\alpha} \hat{f}_{\alpha}. \tag{46}$$

With $\hat{f}_{\alpha} = f_{\alpha} - 0.5\delta_t C_{\alpha}$, the evolution equation for the total energy distribution function is given by

$$\bar{h}_{\alpha}(\mathbf{x} + \mathbf{e}_{\alpha}\delta_{t}, t + \delta_{t}) - \bar{h}_{\alpha}(\mathbf{x}, t)
= -\omega_{h} \left[\bar{h}_{\alpha}(\mathbf{x}, t) - h_{\alpha}^{\text{eq}}(\mathbf{x}, t) \right] + (1 - 0.5\omega_{h}) \frac{\delta_{t} Z_{\alpha}}{\tau_{hf}}
\times \left[\hat{f}_{\alpha}(\mathbf{x}, t) - f_{\alpha}^{\text{eq}}(\mathbf{x}, t) + \frac{\delta_{t}}{2} C_{\alpha}(\mathbf{x}, t) \right].$$
(47)

According to the Chapman-Enskog analysis, the shear viscosity in the energy equation (contained in Π) will be given by $\mu = (\tau_f \tau_h / \tau_{hf} + \tau_h) p$, while the shear viscosity in the macroscopic momentum equation from Eq. (43) is $\mu = (\tau_f - 0.5\delta_t)p$. To ensure these two viscosities are the same, the relaxation time τ_{hf} should be defined as $\tau_{hf} = \tau_f / (\Pr - 1)$. The Prandtl number is now given by $\Pr = \mu c_p / \lambda = (\tau_f - 0.5\delta_t) / \tau_h$.

Furthermore, theoretically speaking, the relaxation time τ_e in Eq. (40) should be equal to τ_f to ensure that the bulk viscosity μ_B in the energy equation is consistent with that in the momentum equation. Actually, if $\tau_e = \tau_\varsigma = \tau_f$, the MRT collision operator will become a TRT collision operator, which is also better than the BGK collision operator in terms of numerical stability [14]. In practical applications, different values of τ_e and τ_f can be adopted since the influence of the bulk viscosity on the energy equation can be neglected in most low Mach number flows.

3. Correction term

Now we give the detailed form of the correction term. Actually, in the moment space the correction term $\widehat{\mathbf{C}}$ can be directly determined via the Chapman-Enskog analysis. Details about the Chapman-Enskog analysis of isothermal MRT models can be found in Refs. [40,43]. The Chapman-Enskog

analysis of the present MRT model can be conducted in a similar way, and we can obtain

$$\hat{\mathbf{C}} = (0, \, \varphi_x, \, 0, \, 0, \, 0, \, 0, \, \varphi_y, \, 0)^T,
\varphi_x = 3(\partial_x \, Q_x + \partial_y \, Q_y),
\varphi_y = (\partial_x \, Q_x - \partial_y \, Q_y),$$
(48)

where $Q_x = \rho u_x(1-\theta)$ and $Q_y = \rho u_y(1-\theta)$. Q_x and Q_y are the order of Ma ΔT , where Ma is the local Mach number and $\Delta T = (T-T_0)/T_0$. The correction term in the velocity space, C_α , can be obtained via the relation $\mathbf{C} = \mathbf{M}^{-1} \widehat{\mathbf{C}}$. The transformation matrix \mathbf{M} of the D2Q9 lattice and its inverse matrix \mathbf{M}^{-1} are given in Appendix \mathbf{C} . According to Eqs. (48) and (C2), C_α is given by

$$C_0 = -\frac{1}{9}\varphi_x$$
, $C_1 = -\frac{1}{36}\varphi_x + \frac{1}{4}\varphi_y$, $C_2 = -\frac{1}{36}\varphi_x - \frac{1}{4}\varphi_y$, $C_3 = C_1$, $C_4 = C_2$, $C_5 = C_6 = C_7 = C_8 = \frac{1}{19}\varphi_x$. (49)

It can be easily verified that C_{α} given by Eq. (49) satisfies the constraints in Eqs. (26) and (28). In numerical simulations, the gradient terms $\partial_x Q_x$ and $\partial_y Q_y$ in Eq. (48) can be evaluated as follows:

$$\partial_x Q_x|_{(I,J)} = \frac{1}{2\delta_x} [Q_x|_{(I+1,J)} - Q_x|_{(I-1,J)}], \ \partial_y Q_y|_{(I,J)}$$

$$= \frac{1}{2\delta_y} [Q_y|_{(I,J+1)} - Q_y|_{(I,J-1)}], \tag{50}$$

where $\delta_x = \delta_y$ are the spatial steps in the *x* and *y* directions, respectively. If needed, one can also adopt some other schemes [41,42]. In fact, for low Mach number flows with small or moderate temperature variations, the correction term is usually negligible.

When a force \boldsymbol{a} is exerted on the flow, a forcing term F_{α} should be included in Eqs. (22) and (39). F_{α} can be defined as $F_{\alpha} = w_{\alpha}[(\boldsymbol{e}_{\alpha} \cdot \boldsymbol{a})/(RT_0) + \boldsymbol{u}\boldsymbol{a} : (\boldsymbol{e}_{\alpha}\boldsymbol{e}_{\alpha} - RT_0\mathbf{I})/(RT_0)^2]$ [31]. The forcing term in the moment space can be obtained via $\widehat{\mathbf{F}} = \mathbf{MF}$, i.e.,

$$\mathbf{F} = [0, 6u \cdot a, -6u \cdot a, a_x, -a_x, a_y, -a_y, 2(u_x a_x - u_x a_y), (u_x a_y + u_y a_x)]^T.$$
 (51)

In the evolution equations, F_{α} and $\widehat{\mathbf{F}}$ can be absorbed into C_{α} and $\widehat{\mathbf{C}}$, respectively. When the forcing term is included, the macroscopic velocity will be given by $\rho \mathbf{u} = \sum_{\alpha} \mathbf{e}_{\alpha} \widehat{f}_{\alpha} + 0.5 \delta_t \mathbf{a}$.

Finally, a brief comparison between the present model and Prasianakis and Karlin's coupling BGK-LB model [37,38] is made. First, it is noted that these two models share some similar features: Both models are based on the standard D2Q9 lattice, the Prandtl number is adjustable, and a correction term is introduced into the kinetic equation due to the low symmetry of the D2Q9 lattice. On the other hand, the differences between the two models are also apparent. The present model is constructed in the framework of the DDF approach. Hence only the deviation of the third-order velocity moment of f_{α}^{eq} needs to be considered, while Prasianakis and Karlin's model is based on the single density distribution function and should consider the deviations of the third-order and fourth-order

velocity moments of f_{α}^{eq} . As a result, more than ten gradient terms need to be computed in Prasianakis and Karlin's model.

Moreover, from Eq. (27) it can be seen that, in the present model, the deviation of the third-order velocity moment of f_{α}^{eq} is removed via the second-order velocity moment of the correction term. Such a treatment is consistent with the philosophy of the Chapman-Enskog analysis. However, in Prasianakis and Karlin's model, the related deviation is removed by the first-order velocity moment of the correction term, which makes the computational scheme of the macroscopic velocity u implicit. Quasi-iterations are therefore introduced in the numerical algorithm of Prasianakis and Karlin's model [37]. In addition, the specific-heat ratio γ , which is fixed at an unphysical value of (D+2)/D in Prasianakis and Karlin's model, can be freely chosen in the present model.

IV. NUMERICAL TESTS

In this section, we conduct a series of numerical simulations to validate the proposed coupling DDF LB model. Without loss of generality, the MRT version of the model is used in the simulations. The testing problems include thermal Couette flow, attenuation-driven acoustic streaming, and natural convection in a square cavity.

A. Thermal Couette flow

Thermal Couette flow is a classical heat transfer problem which can provide a good test of new LB models to describe the viscous heat dissipation. The existing LB simulations of Couette flows are mostly conducted under the assumption of constant viscosity and thermal conductivity coefficients. In the present work, we pay attention to thermal Couette flows with temperature-dependent transport coefficients, in which the shear viscosity and the thermal conductivity vary with the temperature through a power law or the Sutherland law or some other empirical functions. Since the viscosity depends on the temperature, the velocity and temperature fields are coupled together.

Consider the viscous fluid flow between two infinite parallel plates, the lower plate is fixed and the upper one is moving at speed U. The ratio $\lambda/\mu=c_p/\mathrm{Pr}$ is assumed to be constant, and this is reasonable for gases [44]. In fact, there are no closed-form solutions for a general viscosity model. However, under the conditions of $\mu/\mu_0=(T/T_0)^n$ with n=1 and of an adiabatic lower wall, there will be an analytical solution as follows [44]:

$$\frac{T}{T_0} = 1 + \Pr{\frac{\gamma - 1}{2} \operatorname{Ma}^2 \left[1 - \left(\frac{u_x}{U} \right)^2 \right]},$$
 (52a)

$$\left(1 + \operatorname{Pr} \frac{\gamma - 1}{3} \operatorname{Ma}^{2}\right) \frac{y}{H}$$

$$= \frac{u_{x}}{U} + \operatorname{Pr} \frac{\gamma - 1}{2} \operatorname{Ma}^{2} \left[\frac{u_{x}}{U} - \frac{1}{3} \left(\frac{u_{x}}{U}\right)^{3}\right], \quad (52b)$$

where T_0 is the temperature of the upper wall and H is the distance between the two plates. The Mach number is defined as Ma = $U/\sqrt{\gamma RT_0}$.

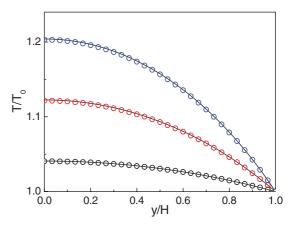


FIG. 1. (Color online) Dimensionless temperature profiles of thermal Couette flow at $\gamma=5/3$ and Ma = 0.35. From top to bottom: Pr = 5, 3, and 1. Solid lines and the circles are the analytical and numerical results, respectively.

Numerical simulations are carried out with different values of Pr and Ma: (1) Pr = 1, 3, and 5 with $\gamma = 5/3$ (b = 3) and Ma = 0.35; (2) Ma = 0.15, 0.25, and 0.35 (in the low Mach number regime) with $\gamma = 5/3$ and Pr = 5. A grid size of $N_x \times N_y = 10 \times 60$ (N_x and N_y exclude the extra layers outside the boundaries) is employed, and the nonequilibrium extrapolation method [31,45] is used to treat the velocity and temperature boundary conditions of the upper and lower walls, while the periodic boundary condition is applied in the x direction.

In simulations, we set $\rho_0 = 1$, $T_0 = 1$, $\mu_0 = 0.25$, $c = \sqrt{3RT_0} = 1$, and $\delta_x = \delta_y = \delta_t = 1$. The relaxation times are chosen as follows: $\tau_e^{-1} = \tau_5^{-1} = \tau_q^{-1} = 1.5$ and $\tau_\rho = \tau_j = 1$. From Eq. (52a) it can be seen that $(\theta - 1) = (T/T_0 - 1)$ is the order of Ma². Therefore the correction term is omitted since it will be the order of Ma³. The temperature profiles are shown in Figs. 1 and 2, in which the analytical results are also presented for comparison. It is clearly seen that the numerical results are in good agreement with the analytical ones. To quantify the results, the numerical (T) and analytical (T_a) temperatures at the lower wall are listed in Table I. The relative error defined by $E_r = |(\theta - 1) - (\theta_a - 1)|/(\theta_a - 1)$

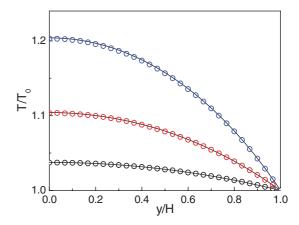


FIG. 2. (Color online) Dimensionless temperature profiles of thermal Couette flow at $\gamma=5/3$ and Pr=5. From top to bottom: Ma = 0.35, 0.25, and 0.15. Solid lines and the circles are the analytical and numerical results, respectively.

TABLE I. Comparison of the numerical (T) and analytical (T_a) temperatures at the lower wall of thermal Couette flow in different cases.

Cases	Pr = 1, $Ma = 0.35$	· · · · · · · · · · · · · · · · · · ·	Pr = 5, $Ma = 0.35$		Pr = 5, $Ma = 0.15$
T/T_0 T_a/T_0 $E_r (\%)$	1.0405	1.1217	1.2032	1.1036	1.0373
	1.0408	1.1225	1.2042	1.1042	1.0375
	0.74	0.65	0.49	0.58	0.53

is computed, where $\theta_a = T_a/T_0$. As shown in Table I, the difference between the present results and the analytical results are within 1.0% for the cases considered.

B. Attenuation-driven acoustic streaming

Now we consider an acoustic problem, the acoustic streaming, which describes a flow field superimposed upon the oscillatory motion of a sound wave propagating in a fluid. It is a nonlinear effect which occurs due to the presence of boundaries or because of damping of waves [46]. Acoustic streaming has important and possibly undesirable effects in the study of acoustic levitation and the operation of thermoacoustic systems. There are two basic types of acoustic streaming. One is the Rayleigh streaming, and the other is the attenuation-driven acoustic streaming, which results from the attenuation of the wave in a fluid.

In this test, the attenuation-driven acoustic streaming is simulated. A plane acoustic wave of wave vector $k = 2\pi/\lambda'$ and angular frequency $\omega = kc_s$ ($c_s = \sqrt{\gamma RT}$ is the speed of sound) propagating in the positive x direction is considered. The wave propagates from a source at x = 0. The problem is initialized with a zero velocity and undisturbed density and temperature everywhere. The sound wave is generated by adding sinusoidal density and temperature variations:

$$\rho = \rho_0 + \delta\rho \sin(\omega t), \quad T = T_0 + (\gamma - 1)\frac{\delta\rho}{\rho_0} T_0 \sin(\omega t), \quad (53)$$

at x = 0. Then the analytical velocity is given by [47]

$$u_x(x, t) = Ue^{-\beta x} \sin(\omega t - kx), \tag{54}$$

where $U = c_s \delta \rho / \rho_0$ and β is the absorption coefficient. The velocity amplitude is given by

$$\Gamma(x) = U e^{-\beta x}. (55)$$

The absorption coefficient β can be obtained via the perturbation solution of the NSF equations [47]:

$$\beta = \frac{2\pi^2 f'^2}{\rho c_s^3} \left[\frac{4}{3} \mu + \mu_B + \left(\frac{1}{c_v} - \frac{1}{c_p} \right) \lambda \right], \tag{56}$$

where f' is the frequency. Since Eq. (56) is based on a perturbation solution, the density perturbation $\delta \rho$ should be small enough to ensure Eq. (56) is valid. In the present study, we set $\delta \rho = 0.01 \rho_0$.

In simulations, we choose the region $x \in [0, 500]$ as the measurement region. The length along the x direction, $L_x = N_x \times \delta_t$, where N_x is the lattice number in the x direction, should be large enough so that the density disturbance will

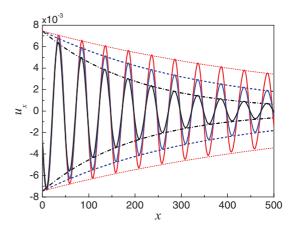


FIG. 3. (Color online) Velocity profiles of attenuation-driven acoustic streaming at $\gamma=5/3$ with different Prandtl numbers. The solid lines are the numerical results. The dotted lines (Pr = 5), dashed lines (Pr = 0.5), and dash-dotted lines (Pr = 0.2) are the analytical velocity amplitude [see Eq. (55) and note that $\pm\Gamma(x)$ are plotted].

not reach $x = L_x$ during the measurement procedure, namely $L_x > c_s t_m > 500$, where t_m is the measurement time. In the computation, we set $\rho_0 = 1$, $T_0 = 1$, $\mu = 0.1$, $\lambda' = 50$, c = 0.11, $\delta_x = \delta_y = \delta_t = 1$, $\tau_e = \tau_\varsigma = \tau_f$, $\tau_q^{-1} = 1.5$, $\tau_\rho = \tau_j = 1$, $t_m = 800$, and $L_x = 600$. A grid size of $N_x \times N_y = 600 \times 4$ is employed. The inlet boundary condition is implemented by constantly resetting the equilibrium distributions with the desired values of density, velocity, and temperature. At the outlet, the zero-gradient condition is imposed by simply extrapolating the information (ρ, \mathbf{u}, T) from the fluid field near the outlet boundary, and then the outlet boundary condition can be implemented by using the nonequilibrium extrapolation method. The periodic boundary condition is applied in the y directions. The velocity profiles at $\gamma = 5/3$ with Pr = 0.2, 0.5, and 5 are presented in Fig. 3, and the velocity profiles at Pr = 2 with $\gamma = 5/3$, 7/5, and 9/7 are shown in Fig. 4. For comparison, the analytical velocity amplitudes are also shown in Figs. 3 and 4. It can be seen that the numerical results agree well with the analytical solutions.

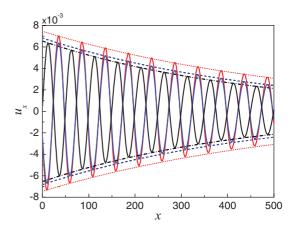


FIG. 4. (Color online) Velocity profiles of attenuation-driven acoustic streaming at Pr=2 with different specific-heat ratios. The solid lines are the numerical results. The dotted lines ($\gamma=5/3$), dashed lines ($\gamma=7/5$), and dash-dotted lines ($\gamma=9/7$) are the analytical velocity amplitude.

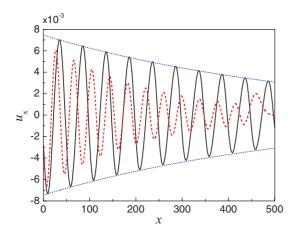


FIG. 5. (Color online) Comparison of the velocity profiles of attenuation-driven acoustic streaming at Pr = 2 and $\gamma = 5/3$ between the present model (the solid line) and the decoupling model (the dashed line). The dotted lines are the analytical velocity amplitude.

A comparison of the numerical results predicted by the present model and the decoupling model ($\theta=1$) is presented in Fig. 5, from which we can observe that the numerical results predicted by the decoupling model significantly deviate from the analytical ones. Actually, in this test the correction term is found to be negligible; hence we have omitted it. Then the only difference between the present model and the decoupling model lies in that the terms related to θ are included in the equilibrium distribution functions of the present model. From Eq. (53) it can be derived that the maximum value of θ is $\theta_{\text{max}} = 1 + (\gamma - 1)\delta\rho/\rho_0$, which is approximately equal to 1 since $\delta\rho/\rho_0 = 0.01$. However, Fig. 5 clearly illustrates that directly setting $\theta=1$ will result in considerable errors.

In fact, when θ is set to be unity, the momentum and energy transports will be decoupled. As a result, the term related to the thermal conductivity λ will vanish in Eq. (56). In addition, the sound speed c_s will be given by $c_s = \sqrt{RT_0}$ rather than $c_s = \sqrt{\gamma RT_0}$. Then the efficient wavelength should be defined as $\lambda'_{\rm eff} = \lambda'/\sqrt{\gamma}$. In other words, the wavelength predicted by the decoupling model will be smaller than the one predicted by the coupling model, which can be clearly seen in Fig. 5. This test demonstrates that in acoustic problems the temperature field will have significant influences on the flow field even though the temperature variation is very small.

C. Natural convection in a square cavity

Natural convection in a square cavity with adiabatic top and bottom walls and with side walls maintained at constant but different temperatures is often taken as one of the standard cases to test new computational schemes. This flow is characterized by two nondimensional parameters: the Prandtl number Pr and the Rayleigh number Ra, which is defined as

$$Ra = \frac{g\beta(T_h - T_l)H^3 Pr}{\mu_r^2},$$
 (57)

where g is the gravity acceleration, β is the thermal expansion coefficient, T_h and T_l are the temperatures of left and right walls $(T_h > T_l)$, H is the distance between the walls, and μ_r is the shear viscosity evaluated at the reference temperature

 $T_r = (T_h + T_l)/2$. For perfect gases, the thermal expansion coefficient β is defined as

$$\beta = -\frac{1}{\rho} \left(\frac{\partial \rho}{\partial T} \right)_{P} = \frac{1}{T}.$$
 (58)

The thermal expansion coefficient in Eq. (57) is also evaluated at T_r .

Firstly, we simulate the natural convection with a small temperature difference: T_h and T_l are set to be 315 and 285 K, respectively. With such a temperature difference, the Boussinesq assumption is approximately satisfied. However, since the present model is a coupling model, we do not need to use the Boussinesq assumption. We can directly set the force a in Eq. (51) as follows: $a_x = 0$ and $a_y = -\rho g$. The viscous heat dissipation and the compression work are negligible. Hence Eq. (38) is adopted. The gradient terms in the correction term are evaluated by Eq. (50) in this test (including the case with a large temperature difference given below). The relaxation times are chosen as follows: $\tau_e^{-1} = \tau_s^{-1} = \tau_q^{-1} = 1.5$ and $\tau_\rho = \tau_j = 1$. The Prandtl number is set to be Pr = 0.71. The temperature T_0 in the lattice speed can be chosen as $T_0 = T_r$, and then R is determined by $R = c^2/(3T_0)$. The

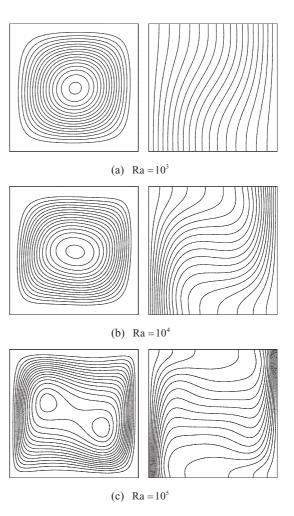


FIG. 6. Streamlines (left) and isotherms (right) of natural convection in a square cavity at Ra = 10^3 , 10^4 , and 10^5 ($T_h = 315$ K and $T_l = 285$ K).

nonequilibrium extrapolation scheme [31,45] is employed to treat different boundary conditions of \hat{f}_{α} and \bar{h}_{α} .

Through the grid-dependence examination of the numerical results, the grid sizes of 100×100 for Ra = 10^3 , 150×150 for Ra = 10^4 , and 200×200 for Ra = 10^5 are adopted. The streamlines and isotherms are shown in Fig. 6. From the figure it can be observed that for low values of Rayleigh number, a vortex appears in the center of the cavity. The vortex gradually becomes elliptic as the Rayleigh number increases, and breaks up into two vortices at $Ra = 10^5$. Meanwhile, the isotherms clearly show that, for the case of $Ra = 10^3$, heat is transferred mainly by conduction between the hot and cold walls. Hence the isotherms are almost vertical. When the Rayleigh number increases, the dominant heat transfer mechanism changes from conduction to convection, and then the isotherms become horizontal in the center of the cavity. These observations are in good agreement with the results reported in the literature [31,48].

To quantify the results, the maximum horizontal velocity component $|u|_{\max}$ in the vertical midplane x=H/2 and its location y_{\max} , the maximum vertical component $|v|_{\max}$ in the horizontal midplane y=H/2 and its location x_{\max} , and the average Nusselt number Nu along the cold wall are computed. The x and y coordinates are normalized with the cavity width y. The velocities are normalized with the diffusion velocity y where y is the thermal diffusivity. The results are listed in Table II together with the data from previous studies. As shown, the present results agree well with the available data.

Furthermore, the case of natural convection with a large temperature difference is also considered: $T_h = 960 \text{ K}$ and $T_l = 240 \text{ K}$. The shear viscosity is defined by Sutherland's law:

$$\frac{\mu(T)}{\mu^*} = \left(\frac{T}{T^*}\right)^{\frac{3}{2}} \frac{(T^* + S)}{(T + S)},\tag{59}$$

where $T^* = 273\,$ K, $S = 110.5\,$ K, and μ^* is the shear viscosity evaluated at T^* . The Prandtl number is assumed to remain constant, i.e., Pr = 0.71, and this is reasonable for gases. For example, for air from $0\,^{\circ}F$ to $1000\,^{\circ}F$, the ratio λ/μ increases only 20% [44].

In simulations, the relaxation time τ_f is determined by $\tau_f = \mu/p + 0.5\delta_t$, where μ is calculated by Eq. (59). The viscosity

TABLE II. Comparisons of the average Nusselt number and the maximum velocity components across the cavity center ($T_h = 315$ K and $T_l = 285$ K).

Ra		Nu	$ u _{\max}$	y_{max}	$ v _{\max}$	x_{max}
10^{3}	Present	1.1207	3.664	0.8100	3.699	0.1800
	Reference [48]	_	3.649	0.8130	3.697	0.1780
	Reference [31]	1.1195	3.643	0.8047	3.692	0.1719
104	Present	2.2528	16.351	0.8200	19.589	0.1200
	Reference [48]	2.2448	16.180	0.8265	19.630	0.1193
	Reference [31]	2.2545	16.125	0.8203	19.558	0.1172
105	Present	4.5350	35.703	0.8550	68.536	0.0650
	Reference [48]	4.5216	34.740	0.8558	68.640	0.0657
	Reference [31]	4.5278	34.603	0.8516	68.082	0.0703

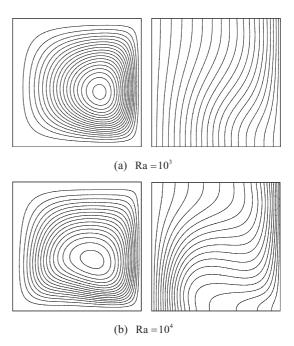


FIG. 7. Streamlines (left) and isotherms (right) of natural convection in a square cavity at Ra = 10^3 and 10^4 ($T_h = 960$ K and $T_l = 240$ K).

 μ_r in Eq. (57) is set to be $\mu_r = 0.14$, 0.1, and 0.06 for Ra = 10^3 , 10^4 , and 10^5 , respectively. The viscosity μ^* in Eq. (59) is determined from $\mu(T_r)$. The grid sizes of 100×100 , 150×150 , and 250×250 are adopted for Ra = 10^3 , 10^4 , and 10^5 , respectively. The streamlines and isotherms are presented in Figs. 7 and 8. The present results agree well with the numerical results reported in Ref. [49]. The asymmetry feature of natural convection with a large temperature difference can be clearly seen in Figs. 7 and 8: There is an apparent shift of the center of the primary vortex both towards the right wall and downwards towards the bottom of the cavity.

The comparisons of the average Nusselt number and the maximum velocity components (the reference velocity $V_{\rm ref}=\mu_r\,{\rm Pr}^{-1}\,H^{-1}$) across the cavity center between the present results and the results reported in Ref. [49] are made in Table III. In the table, "w" and "w/o" denote "with correction term" and "without correction term," respectively. The average Nusselt number of the left wall is defined as ${\rm Nu}=\int_0^H {\rm Nu}(y)dy/H$, where ${\rm Nu}(y)=H\lambda.\partial_x T|_{\rm wall}/[\lambda_r(T_h-T_l)]$ [49]. From the table we can see that in the present case the correction term

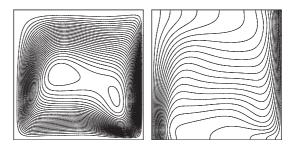


FIG. 8. Streamlines (left) and isotherms (right) of natural convection in a square cavity at Ra = 10^5 ($T_h = 960$ K and $T_l = 240$ K).

TABLE III. Comparisons of the average Nusselt number and the maximum velocity components across the cavity center ($T_h = 960 \text{ K}$ and $T_l = 240 \text{ K}$).

Ra		Nu	$ u _{\max}$	y_{max}	$ v _{\max}$	x_{max}
10^{3}	Present (w/o)	1.105	3.971	0.7800	3.991	0.9000
	Present (w)	1.111	3.726	0.1600	4.429	0.9100
	Reference [49]	1.108	3.702	0.1618	4.324	0.9036
104	Present (w/o)	2.217	18.165	0.7933	20.407	0.9267
	Present (w)	2.217	16.786	0.7800	20.404	0.9267
	Reference [49]	2.218	16.777	0.7821	20.327	0.9270
10 ⁵	Present (w/o)	4.446	47.079	0.8440	71.992	0.0920
	Present (w)	4.454	43.984	0.8360	71.070	0.0960
	Reference [49]	4.480	43.692	0.8364	71.084	0.0948

has an important influence on the velocity field: The relative errors of $|u|_{\text{max}}$ and $|v|_{\text{max}}$ are mostly around 8% when the correction term is not considered, while the results obtained with the correction term are in excellent agreement with the data in Ref. [49]. Furthermore, it is observed that there are only slight differences for the average Nusselt number. This means the correction term has a little influence on the temperature field in the present case. The above points can also be seen in Fig. 9, in which the streamlines and isotherms obtained without the correction term at $Ra = 10^5$ are shown. By comparing Fig. 9(a) with Fig. 8(a), we can find that, without the correction term, the degree of the asymmetry of the velocity field is reduced. This is expected since the local temperature in the diagonal elements of the third-order velocity moment of f_{α}^{eq} is replaced by a constant temperature. Quantitatively, without the correction term, the positions of the left and right vortices will change from (x/H, y/H) = (0.352, 0.509) to (0.362, 0.529)and from (0.786, 0.324) to (0.783, 0.355), respectively.

By comparing the present case and the previous case with a small temperature difference, we can observe that the temperature gradients become higher near the right wall but lower near the left wall. This is because the local Rayleigh number near the right wall is larger than the defined Rayleigh number, while the local Rayleigh number near the left wall is smaller than the defined Rayleigh number. For $Ra = 10^5$, the maximum Rayleigh number near the right wall is around 9.5×10^5 , whereas the minimum Rayleigh number near the left wall is only about 3.5×10^4 . Figure 10 illustrates the temperature profiles along the horizontal line crossing

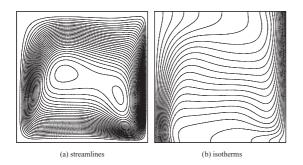
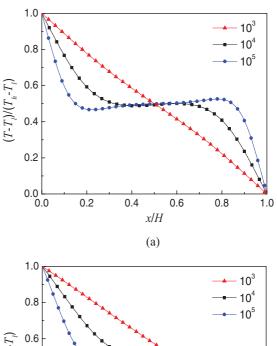


FIG. 9. Streamlines and isotherms of natural convection in a square cavity at Ra = 10^5 without the correction term ($T_h = 960$ K and $T_l = 240$ K).



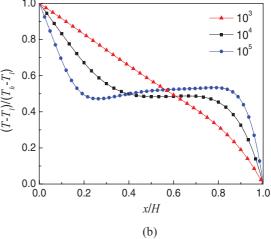


FIG. 10. (Color online) Comparions of the temperature profiles along the horizontal line crossing the center of the cavity: (a) $T_h = 315$ K and $T_l = 285$ K; and (b) $T_h = 960$ K and $T_l = 240$ K.

the center of the cavity. From the figure the symmetry and asymmetry features of natural convection in different cases can be clearly observed.

Finally, the comparison of numerical stability between the BGK and MRT schemes of the proposed model is made at Ra = 10^5 with a higher temperature ratio: $T_h/T_l = 8$ $(T_h \approx 1066.7 \text{ K} \text{ and } T_l \approx 133.3 \text{ K})$. The related local Rayleigh number ranges from 2.7×10^4 to 4.8×10^6 . In simulations, the parameters and the grid system are the same as those used in the case with $T_h/T_l = 4$. The results are presented in Fig. 11. From the figure it can be seen that at $t = 1000\delta_t$ and $t = 3500\delta_t$ the temperature contours obtained from the BGK scheme are still smooth, however, at $t = 7000\delta_t$ obvious oscillations can be observed near the right (cold) wall, which leads to numerical blowup around $t = 7300\delta_t$. The oscillations appear near the right wall only because the local viscosity near the right wall is much lower than the viscosities in other regions and the transient heat transfer near the right wall is very strong due to the high local Rayleigh number. Under the same simulation condition, the MRT scheme is found to work well by separating the relaxation times for hydrodynamic and nonhydrodynamic moments, and the stable solution is shown

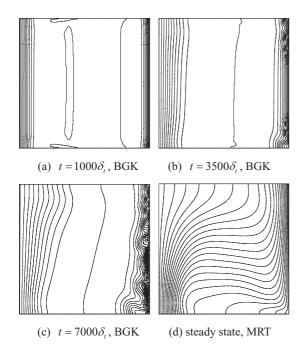


FIG. 11. Isotherms of natural convection in a square cavity at $Ra = 10^5$ with $T_h/T_l = 8$.

in Fig. 11(d), from which we can see that, with the increase of the temperature ratio, the asymmetry feature is obviously enhanced: The differences between the left and right regions are more apparent.

V. CONCLUSION

In this paper, a coupling LB model has been developed for simulating thermal flows on the standard D2Q9 lattice in the framework of the DDF LB approach. In the model, the discrete equilibrium distribution functions are derived from the Hermite expansions of the continuous equilibrium distribution functions. When the D2Q9 lattice is used, these discrete equilibrium distribution functions can rigorously satisfy the zeroth- through second-order velocity moments of the corresponding continuous equilibrium distribution functions. However, the third-order velocity moment of the discrete equilibrium density distribution function deviates from the related statistical relationship. To remove this deviation, a correction term has been introduced into the kinetic equation. The statistical constraints on the correction term and the detailed form of the correction term have been determined with the Chapman-Enskog analysis. The BGK and MRT versions of the present model have been both proposed.

Numerical simulations have been performed for thermal Couette flow with temperature-dependent transport coefficients, attenuation-driven acoustic streaming with different Prandtl numbers and specific-heat ratios, and natural convection in a square cavity with small and large temperature differences. The numerical experiments show that the results predicted by the present model are in good agreement with the analytical solutions and/or the numerical results reported in the literature.

In summary, unlike previous DDF LB models based on standard lattices, the present model can ensure the momentum and energy transports are physically coupled, which makes it applicable for general thermal flows such as non-Boussinesq flows. Furthermore, different from most of the existing coupling LB models which are usually constructed on multispeed lattices, the present model is based on the DDF approach with the standard D2Q9 lattice. The simple structure and general features of the DDF approach are retained in the model. In addition, compared with Prasianakis and Karlin's coupling BGK-LB model which is also based on the standard D2Q9 lattice, the present model has the following advantages: The correction term is much simpler, the computational scheme of the macroscopic velocity is explicit, and the specific-heat ratio is adjustable. With the formulations in the present paper, the model can be readily extended to three-dimensional (3D) space and other systems such as flows of non-Newtonian fluids [50] and axisymmetric flows in cylindrical coordinates [15]. Finally, it should be noted that, with the low Mach number approximation and the standard collision-streaming process, the present model is limited to low Mach number thermal flows.

ACKNOWLEDGMENTS

This work was supported by the National Natural Science Foundation of China (Grants No. 50736005 and No. 51176155) and the Engineering and Physical Sciences Research Council of the United Kingdom under Grants No. EP/I012605/1.

APPENDIX A: HERMITE EXPANSION OF h^{eq}

Following Shan and co-workers' approach for expanding f^{eq} [23,40], we can obtain the Nth-order Hermite expansion of h^{eq} as follows:

$$h^{\text{eq},N} = \omega(\xi, T_0) \sum_{n=0}^{N} \frac{1}{n!} b_{\text{eq}}^{(n)} H^{(n)}(\bar{\xi}) d\xi,$$
 (A1)

where $\bar{\boldsymbol{\xi}} = \boldsymbol{\xi}/\sqrt{RT_0}$, $\boldsymbol{b}_{\rm eq}^{(n)} = \int h^{\rm eq} \, \boldsymbol{H}^{(n)}(\bar{\boldsymbol{\xi}}) d\boldsymbol{\xi}$, and $\boldsymbol{H}^{(n)}(\bar{\boldsymbol{\xi}})$ is the *n*th-order Hermite polynomial:

$$\boldsymbol{H}^{(0)}(\bar{\boldsymbol{\xi}}) = 1, \quad \boldsymbol{H}_{i}^{(1)}(\bar{\boldsymbol{\xi}}) = \bar{\xi}_{i}, \quad \boldsymbol{H}_{ij}^{(2)}(\bar{\boldsymbol{\xi}}) = \bar{\xi}_{i}\bar{\xi}_{j} - \delta_{ij}. \quad (A2)$$

According to $h^{\text{eq}} = \xi^2 f^{\text{eq}}/2$ $(\xi^2 = \xi \cdot \xi)$, $b_{\text{eq}}^{(n)}$ can be written as follows:

$$\boldsymbol{b}_{\text{eq}}^{(n)} = \int \frac{\rho \boldsymbol{\xi}^2}{2(2\pi RT)^{D/2}} \exp\left[-\frac{(\bar{\boldsymbol{\xi}} - \bar{\boldsymbol{u}})^2}{2\theta}\right] \boldsymbol{H}^{(n)}(\bar{\boldsymbol{\xi}}) d\boldsymbol{\xi}, \quad (A3)$$

where $\theta = T/T_0$ and $\bar{u} = u/\sqrt{RT_0}$. By defining $\eta = (\bar{\xi} - \bar{u})/\sqrt{\theta}$, we can rewrite Eq. (A3) as

$$\begin{aligned} \boldsymbol{b}_{\text{eq}}^{(n)} &= \int \frac{\rho \boldsymbol{\xi}^{2} (RT)^{-D/2}}{2(2\pi)^{D/2}} \exp\left(-\frac{\boldsymbol{\eta}^{2}}{2}\right) \boldsymbol{H}^{(n)}(\bar{\boldsymbol{\xi}}) d\boldsymbol{\xi}, \\ &= \frac{\rho}{2} \int RT_{0}(\theta \boldsymbol{\eta}^{2} + 2\sqrt{\theta} \boldsymbol{\eta} \cdot \bar{\boldsymbol{u}} + \bar{\boldsymbol{u}}^{2}) (RT)^{-D/2} \boldsymbol{\varpi}(\boldsymbol{\eta}) \\ &\times \boldsymbol{H}^{(n)}(\sqrt{\theta} \boldsymbol{\eta} + \bar{\boldsymbol{u}}) d\boldsymbol{\xi}, \end{aligned}$$
(A4)

where $\varpi(\eta)$ is given by

$$\varpi(\eta) = \frac{1}{(2\pi)^{D/2}} \exp\left(-\frac{\eta^2}{2}\right). \tag{A5}$$

Note that $d\xi_i = \sqrt{RT_0}d\bar{\xi}_i = \sqrt{RT}d\eta_i$. Hence $d\xi \equiv d\xi_1 d\xi_2 \cdots d\xi_D = (RT)^{D/2}d\eta$. Therefore, $\boldsymbol{b}_{eq}^{(n)}$ is given by

$$\boldsymbol{b}_{\text{eq}}^{(n)} = \frac{\rho R T_0}{2} \int (\theta \boldsymbol{\eta}^2 + 2\sqrt{\theta} \boldsymbol{\eta} \cdot \bar{\boldsymbol{u}} + \bar{\boldsymbol{u}}^2) \varpi(\boldsymbol{\eta}) \times \boldsymbol{H}^{(n)}(\sqrt{\theta} \boldsymbol{\eta} + \bar{\boldsymbol{u}}) d\boldsymbol{\eta}. \tag{A6}$$

According to Eq. (A6), we can obtain

$$\mathbf{b}_{\text{eq}}^{(0)} = \rho E, \quad \mathbf{b}_{eq,i}^{(1)} = (\rho E + p)\bar{u}_i,
\mathbf{b}_{eq,ij}^{(0)} = (\rho E + 2p)\bar{u}_i\bar{u}_j + p\theta\delta_{ij} + (\theta - 1)\rho E\delta_{ij}.$$
(A7)

The following results, which can be found in mathematical handbooks, are used in deriving Eq. (A7):

$$\int_{-\infty}^{+\infty} \exp(-ax^2) dx = \sqrt{\frac{\pi}{a}},$$

$$\int_{-\infty}^{+\infty} x^2 \exp(-ax^2) dx = \frac{1}{2a} \sqrt{\frac{\pi}{a}},$$

$$\int_{-\infty}^{+\infty} x^4 \exp(-ax^2) dx = \frac{3}{4a^2} \sqrt{\frac{\pi}{a}}.$$
(A8)

With Eqs. (A1), (A2), and (A7), we can obtain the following second-order Hermite expansion of h^{eq} :

$$h^{\text{eq*,2}}(T) = \omega(\xi, T_0) \left\{ \rho E + (p + \rho E) \frac{\xi \cdot \mathbf{u}}{RT_0} + \left(p + \frac{\rho E}{2} \right) \right.$$

$$\times \left[\left(\frac{\xi \cdot \mathbf{u}}{RT_0} \right)^2 - \frac{\mathbf{u}^2}{RT_0} \right] + \frac{p\theta}{2} \left(\frac{\xi^2}{RT_0} - D \right) \right.$$

$$\left. + \frac{\rho E}{2} (\theta - 1) \left(\frac{\xi^2}{RT_0} - D \right) \right\}. \tag{A9}$$

It should be noted that, in the above derivations, $h^{\rm eq}$ is defined as $h^{\rm eq} = \xi^2 f^{\rm eq}/2$ and $E = (DRT + u^2)/2$. For polyatomic gases, $h^{\rm eq}$ can be defined as $h^{\rm eq} = [\xi^2 + (b-D)RT] f^{\rm eq}/2$ [31] and the total energy in Eqs. (A8) and (A9) will be given by $E = (bRT + u^2)/2$.

APPENDIX B: CHAPMAN-ENSKOG ANALYSIS

In this Appendix, the Chapman-Enskog analysis of Eqs. (22) and (7b) with the equilibrium distribution functions in Eq. (18) is presented. Substituting the expansions given by Eq. (23) and the expansion $h_{\alpha} = h_{\alpha}^{\text{eq}} + \varepsilon h_{\alpha}^{(1)} + \varepsilon h_{\alpha}^{(2)}$ into Eqs. (22) and (7b), we can obtain a series of equations in terms of the order of ε :

$$\varepsilon^{1}: \left(\partial_{t_{1}} + \boldsymbol{e}_{\alpha} \cdot \nabla_{1}\right) f_{\alpha}^{\text{eq}} + \frac{1}{\tau_{f}} f_{\alpha}^{(1)} = C_{\alpha}^{(0)}, \quad (B1)$$

$$\varepsilon^2 : \partial_{t_2} f_{\alpha}^{(1)} + (\partial_{t_1} + \boldsymbol{e}_{\alpha} \cdot \nabla_1) f_{\alpha}^{(1)} + \frac{1}{\tau_f} f_{\alpha}^{(2)} = 0,$$
 (B2)

and

$$\varepsilon^{1}: \left(\partial_{t_{1}} + \boldsymbol{e}_{\alpha} \cdot \nabla_{1}\right) h_{\alpha}^{\text{eq}} + \frac{h_{\alpha}^{(1)}}{\tau_{h}} = Z_{\alpha} \frac{f_{\alpha}^{(1)}}{\tau_{hf}}, \quad (B3)$$

$$\varepsilon^{2}: \partial_{t_{2}}h_{\alpha}^{(1)} + (\partial_{t_{1}} + \boldsymbol{e}_{\alpha} \cdot \nabla_{1})h_{\alpha}^{(1)} + \frac{1}{\tau_{h}}h_{\alpha}^{(2)} = Z_{\alpha}\frac{f_{\alpha}^{(2)}}{\tau_{hf}}.$$
 (B4)

1. Mass conservation

Summations of Eqs. (B1) and (B2) over α give, respectively,

$$\partial_{t_1} \rho + \partial_{1k}(\rho u_k) = 0, \tag{B5}$$

$$\partial_t, \rho = 0.$$
 (B6)

Combining Eq. (B5) with Eq. (B6) leads to the continuum equation

$$\partial_t \rho + \nabla \cdot (\rho \mathbf{u}) = 0. \tag{B7}$$

2. Momentum conservation

The first-order velocity moments of Eqs. (B1) and (B2) give, respectively,

$$\partial_{t_1}(\rho u_j) + \partial_{1i}\left(\sum_{\alpha} e_{\alpha i} e_{\alpha j} f_{\alpha}^{\text{eq}}\right) = 0,$$
 (B8)

$$\partial_{t_2}(\rho u_j) + \partial_{1i}\left(\sum_{\alpha} e_{\alpha i} e_{\alpha j} f_{\alpha}^{(1)}\right) = 0.$$
 (B9)

According to Eqs. (19a) and (B8) is given by

$$\partial_{t_1}(\rho u_i) + \partial_{1k}(\rho u_k u_i) = -\partial_{1i} p, \tag{B10}$$

where $p = \rho RT$. From Eq. (B1), we can obtain

$$\sum_{\alpha} e_{\alpha i} e_{\alpha j} f_{\alpha}^{(1)} = -\tau_{f} \left[\partial_{t_{1}} \left(\sum_{\alpha} e_{\alpha i} e_{\alpha j} f_{\alpha}^{\text{eq}} \right) + \partial_{1k} \left(\sum_{\alpha} e_{\alpha i} e_{\alpha j} e_{\alpha k} f_{\alpha}^{\text{eq}} \right) - \sum_{\alpha} e_{\alpha i} e_{\alpha j} C_{\alpha}^{(0)} \right].$$
(B11)

With Eqs. (19a) and (20) as well as the constraints on the correction term C_{α} , Eq. (B11) can be rewritten as

$$\sum_{\alpha} e_{\alpha i} e_{\alpha j} f_{\alpha}^{(1)} = -\tau_f \{ \partial_{t_1} (\rho u_i u_j + p \delta_{ij}) + \partial_{1k} [p(u_k \delta_{ij} + u_j \delta_{ik} + u_i \delta_{jk})] \}.$$
(B12)

Combining Eq. (B10) with Eq. (B5), we can obtain

$$\partial_{t_1}(\rho u_i u_j) = -u_i \partial_{1j} p - u_j \partial_{1i} p - \partial_{1k}(\rho u_i u_j u_k). \tag{B13}$$

In the low Mach number limit, the third-order velocity term in Eq. (B13) can be neglected. To get the expression of $\partial p/\partial t_1$, the energy conservation should be considered. The summation of Eq. (B3) gives

$$\partial_{t_1}(\rho E) + \partial_{1k}[(\rho E + p)u_k] = 0. \tag{B14}$$

The total energy E in h_{α}^{eq} is defined by $E = (bRT + u^2)/2$, where b is the number of degrees of freedom of a gas. Then the left-hand side of Eq. (B14) can be rewritten as

$$\partial_{t_1} \left(\frac{1}{2} \rho u^2 + \rho \frac{b}{2} RT \right) + \partial_{1k} \left(\frac{1}{2} \rho u^2 u_k + \rho \frac{b}{2} RT u_k + p u_k \right)$$

$$= u_j \left[\partial_{t_1} (\rho u_j) + \partial_{1k} (\rho u_k u_j) + \partial_{1j} p \right]$$

$$+ \partial_{t_1} \left(\rho \frac{b}{2} RT \right) + \partial_{1k} \left(\rho \frac{b}{2} RT u_k \right) + p \partial_{1k} u_k. \quad (B15)$$

With the aid of Eq. (B10), we can get

$$\partial_{t_1} p = -\partial_{1k}(pu_k) - \frac{2}{h} p \partial_{1k} u_k. \tag{B16}$$

Substituting Eqs. (B13) and (B16) into Eq. (B12) gives

$$\sum_{\alpha} e_{\alpha i} e_{\alpha j} f_{\alpha}^{(1)} = -\tau_f p \left[(\partial_{1i} u_j + \partial_{1j} u_i) - \frac{2}{b} \partial_{1k} u_k \delta_{ij} \right].$$
(B17)

Combining Eqs. (B10) and (B9) with Eq. (B17), the following momentum equation can be obtained:

$$\partial_t(\rho \mathbf{u}) + \nabla \cdot (\rho \mathbf{u} \mathbf{u}) = -\nabla p + \nabla \cdot \mathbf{\Pi},$$
 (B18)

where $\mathbf{\Pi} = \mu [\nabla \mathbf{u} + (\nabla \mathbf{u})^T - (2/3)(\nabla \cdot \mathbf{u})\mathbf{I}] + \mu_B(\nabla \cdot \mathbf{u})\mathbf{I},$ $\mu = \tau_f p$, and $\mu_B = 2(1/3 - 1/b)\mu$.

3. Energy conservation

Taking the summation of Eq. (B4) over α , we can obtain

$$\partial_{t_2}(\rho E) + \partial_{1j}\left(\sum_{\alpha} e_{\alpha j} h_{\alpha}^{(1)}\right) = 0.$$
 (B19)

Multiplying Eq. (B3) by $e_{\alpha j}$ and then taking the summation over α , we have

$$\sum_{\alpha} e_{\alpha j} h_{\alpha}^{(1)} = -\tau_h \left[\partial_{t_1} [(\rho E + p) u_j] + \partial_{1i} \left(\sum_{\alpha} e_{\alpha i} e_{\alpha j} h_{\alpha}^{\text{eq}} \right) \right] + \frac{\tau_h}{\tau_{hf}} u_i \sum_{\alpha} e_{\alpha i} e_{\alpha j} f_{\alpha}^{(1)}.$$
 (B20)

Combining Eq. (B5) with Eq. (B8) leads to

$$\partial_{t_1} u_j = -u_i \partial_{1i} u_j - \frac{1}{\rho} \partial_{1j} p. \tag{B21}$$

With Eqs. (B14), (B16), and (B21), we can obtain

$$\begin{aligned} \partial_{t_1}[(\rho E + p)u_j] &= -(\rho E + p)u_i\partial_{1i}u_j - (E + RT)\partial_{1j}p \\ &- u_j\partial_{1i}[(\rho E + p)u_i] - u_j\partial_{1i}(pu_i) \\ &- \frac{2}{b}pu_i\delta_{ij}\partial_{1k}u_k. \end{aligned} \tag{B22}$$

According to Eq. (19b), the following equation can be obtained after some standard algebra:

$$\partial_{t_{1}}[(\rho E + p)u_{j}] + \partial_{1i} \left(\sum_{\alpha} e_{\alpha i} e_{\alpha j} h_{\alpha}^{\text{eq}} \right) \\
= p u_{i} (\partial_{1j} u_{i} + \partial_{1i} u_{j}) + (c_{v} + R) p \partial_{1j} T \\
- \frac{2}{b} p u_{i} \delta_{ij} \partial_{1k} u_{k}, \tag{B23}$$

where $c_v = bR/2$. With Eqs. (B17) and (B23), we can rewrite Eq. (B19) as

$$\begin{split} \partial_{t_{2}}(\rho E) \\ &= \partial_{1j} \left[\left(\tau_{h} + \frac{\tau_{f} \tau_{h}}{\tau_{hf}} \right) p u_{i} \left(\partial_{1j} u_{i} + \partial_{1i} u_{j} - \frac{2}{b} \delta_{ij} \partial_{1k} u_{k} \right) \right. \\ &+ \tau_{h} (c_{v} + R) p \partial_{1j} T \right]. \end{split} \tag{B24}$$

Note that $(\tau_h + \tau_f \tau_h / \tau_{hf}) = \tau_f$ when $\tau_{hf} = \tau_h \tau_f / (\tau_f - \tau_h)$. Finally, combining Eq. (B14) with Eq. (B24), we can obtain the macroscopic energy equation as follows:

$$\partial_t(\rho E) + \nabla \cdot [(\rho E + p)\mathbf{u}] = \nabla \cdot (\lambda \nabla T) + \nabla \cdot (\mathbf{u} \cdot \mathbf{\Pi}),$$
(B25)

where $\lambda = \tau_h p(c_v + R)$ is the thermal conductivity.

APPENDIX C: THE TRANSFORMATION MATRIX AND ITS INVERSE MATRIX

$$\mathbf{M} = \begin{bmatrix} 1 & 1 & 1 & 1 & 1 & 1 & 1 & 1 & 1 \\ -4 & -1 & -1 & -1 & -1 & 2 & 2 & 2 & 2 & 2 \\ 4 & -2 & -2 & -2 & -2 & 1 & 1 & 1 & 1 \\ 0 & 1 & 0 & -1 & 0 & 1 & -1 & -1 & 1 \\ 0 & -2 & 0 & 2 & 0 & 1 & -1 & -1 & 1 \\ 0 & 0 & 1 & 0 & -1 & 1 & 1 & -1 & -1 \\ 0 & 0 & -2 & 0 & 2 & 1 & 1 & -1 & -1 \\ 0 & 1 & -1 & 1 & -1 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 & 0 & 1 & -1 & 1 & -1 \end{bmatrix}.$$
(C1)

$$\mathbf{M}^{-1} = \begin{bmatrix} \frac{1}{9} & -\frac{1}{9} & \frac{1}{9} & 0 & 0 & 0 & 0 & 0 & 0 \\ \frac{1}{9} & -\frac{1}{36} & -\frac{1}{18} & \frac{1}{6} & -\frac{1}{6} & 0 & 0 & \frac{1}{4} & 0 \\ \frac{1}{9} & -\frac{1}{36} & -\frac{1}{18} & 0 & 0 & \frac{1}{6} & -\frac{1}{6} & -\frac{1}{4} & 0 \\ \frac{1}{9} & -\frac{1}{36} & -\frac{1}{18} & -\frac{1}{6} & \frac{1}{6} & 0 & 0 & \frac{1}{4} & 0 \\ \frac{1}{9} & -\frac{1}{36} & -\frac{1}{18} & 0 & 0 & -\frac{1}{6} & \frac{1}{6} & -\frac{1}{4} & 0 \\ \frac{1}{9} & \frac{1}{18} & \frac{1}{36} & \frac{1}{6} & \frac{1}{12} & \frac{1}{6} & \frac{1}{12} & 0 & \frac{1}{4} \\ \frac{1}{9} & \frac{1}{18} & \frac{1}{36} & -\frac{1}{6} & -\frac{1}{12} & \frac{1}{6} & \frac{1}{12} & 0 & -\frac{1}{4} \\ \frac{1}{9} & \frac{1}{18} & \frac{1}{36} & -\frac{1}{6} & -\frac{1}{12} & -\frac{1}{6} & -\frac{1}{12} & 0 & -\frac{1}{4} \end{bmatrix}$$

$$(C2)$$

U. Frisch, B. Hasslacher, and Y. Pomeau, Phys. Rev. Lett. 56, 1505 (1986).

^[2] R. Benzi, S. Succi, and M. Vergassola, Phys. Rep. 222, 145

^[3] S. Chen and G. D. Doolen, Annu. Rev. Fluid Mech. 30, 329 (1998).

^[4] S. Succi, *The Lattice Boltzmann Equation for Fluid Dynamics and Beyond* (Clarendon Press, Oxford, 2001).

^[5] H. Lai and C. Ma, J. Stat. Mech.: Theory Exp. (2010) P04011; Physica A 388, 1405 (2009).

^[6] S. Succi, Eur. Phys. J. B 64, 471 (2008).

^[7] J. M. V. A. Koelman, Europhys. Lett. 15, 603 (1991).

- [8] S. Chen, H. Chen, D. Martínez, and W. Matthaeus, Phys. Rev. Lett. 67, 3776 (1991).
- [9] Y. H. Qian, D. d'Humières, and P. Lallemand, Europhys. Lett. 17, 479 (1992).
- [10] X. He and L.-S. Luo, Phys. Rev. E 56, 6811 (1997).
- [11] D. d'Humières, in *Rarefied Gas Dynamics: Theory and Simulations*, edited by B. D. Shizgal and D. P. Weaver, Prog. Astronaut. Aeronaut., Vol. 159 (AIAA, Washington, DC, 1992).
- [12] P. Lallemand and L.-S. Luo, Phys. Rev. E 61, 6546 (2000).
- [13] D. d'Humières, I. Ginzburg, M. Krafczyk, P. Lallemand, and L.-S. Luo, Phil. Philos. Trans. R. Soc., A **360**, 437 (2002).
- [14] L.-S. Luo, W. Liao, X. Chen, Y. Peng, and W. Zhang, Phys. Rev. E 83, 056710 (2011).
- [15] Q. Li, Y. L. He, G. H. Tang, and W. Q. Tao, Phys. Rev. E 81, 056707 (2010).
- [16] F. Chen, A. Xu, G. Zhang, Y. Li, and S. Succi, Europhys. Lett. 90, 54003 (2010).
- [17] I. Ginzburg, Adv. Water Resour. 28, 1196 (2005).
- [18] I. Ginzburg, Commun. Comput. Phys. 3, 427 (2008).
- [19] D. O. Martínez, W. H. Matthaeus, S. Chen, and D. C. Montgomery, Phys. Fluids 6, 1285 (1994).
- [20] F. J. Alexander, S. Chen, and J. D. Sterling, Phys. Rev. E 47, R2249 (1993).
- [21] Y. H. Qian, J. Sci. Comput. 8, 231 (1993).
- [22] Y. Chen, H. Ohashi, and M. Akiyama, Phys. Rev. E 50, 2776 (1994).
- [23] X. Shan, X.-F. Yuan, and H. Chen, J. Fluid Mech. **550**, 413 (2006).
- [24] S. S. Chikatamarla and I. V. Karlin, Phys. Rev. Lett. 97, 190601 (2006); Comput. Phys. Commun. 179, 140 (2008).
- [25] L. Zheng, B. Shi, and Z. Guo, Phys. Rev. E 78, 026705 (2008).
- [26] Y. Gan, A. Xu, G. Zhang, X. Yu, and Y. Li, Physica A 387, 1721 (2008); Y. Gan, A. Xu, G. Zhang, and Y. Li, Phys. Rev. E 83, 056704 (2011).
- [27] X. Shan, Phys. Rev. E 55, 2780 (1997).
- [28] X. He, S. Chen, and G. D. Doolen, J. Comput. Phys. 146, 282 (1998).

- [29] Y. Peng, C. Shu, and Y. T. Chew, Phys. Rev. E 68, 026701 (2003).
- [30] Y. Shi, T. S. Zhao, and Z. L. Guo, Phys. Rev. E 70, 066310 (2004).
- [31] Z. Guo, C. Zheng, B. Shi, and T. S. Zhao, Phys. Rev. E 75, 036704 (2007).
- [32] S. Chen, J. Tölke, and M. Krafczyk, Phys. Rev. E 79, 016704 (2009).
- [33] Q. Li, Y. L. He, Y. Wang, and W. Q. Tao, Phys. Rev. E 76, 056705 (2007).
- [34] Q. Li, Y. L. He, G. H. Tang, and W. Q. Tao, Phys. Rev. E 80, 037702 (2009).
- [35] P. Lallemand and L.-S. Luo, Phys. Rev. E 68, 036706 (2003).
- [36] P. Lallemand and L.-S. Luo, Int. J. Mod. Phys. C 17, 41 (2003).
- [37] N. I. Prasianakis and I. V. Karlin, Phys. Rev. E 76, 016702 (2007).
- [38] N. I. Prasianakis and I. V. Karlin, Phys. Rev. E 78, 016704 (2008).
- [39] P. C. Philippi, L. A. Hegele Jr., L. O. E. dos Santos, and R. Surmas, Phys. Rev. E 73, 056702 (2006).
- [40] Z. L. Guo and C. G. Zheng, *Theory and Applications of Lattice Boltzmann Method* (Science Press, Beijing, 2009).
- [41] T. Lee and P. F. Fischer, Phys. Rev. E 74, 046709 (2006).
- [42] C. M. Pooley and K. Furtado, Phys. Rev. E 77, 046702 (2008).
- [43] M. E. McCracken and J. Abraham, Phys. Rev. E 71, 036701 (2005).
- [44] H. W. Liepmann and A. Roshko, *Elements of Gasdynamics* (Wiley, New York, 1957).
- [45] Z. L. Guo, C. Zheng, and B. Shi, Chin. Phys. 11, 0366 (2002).
- [46] D. Haydock and J. M. Yeomans, J. Phys. A: Math. Gen. 36, 5683 (2003).
- [47] G. Emanuel, Int. J. Eng. Sci. 36, 1313 (1998).
- [48] M. Hortmann, M. Perić, and G. Scheuerer, Int. J. Numer. Methods Fluids 11, 189 (1990).
- [49] J. Vierendeels, B. Merci, and E. Dick, Int. J. Numer. Meth. Heat Fluid Flow 13, 1057 (2003).
- [50] G. H. Tang, P. X. Ye, and W. Q. Tao, J. Non-Newtonian Fluid Mech. 165, 435 (2010).