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NASA Technical Memorandum 105693

Numerical Solution of a Three-Dimensional Cubic Cavity Flow by Using the Boltzmann Equation

Danny P. Hwang
Lewis Research Center
Cleveland, Ohio

Prepared for the
18th International Symposium on Rarefield Gas Dynamics
sponsored by the University of British Columbia
Vancouver, Canada, July 26–31, 1992

NASA

NUMERICAL SOLUTION OF A THREE-DIMENSIONAL CUBIC CAVITY FLOW BY USING THE BOLTZMANN EQUATION

Danny P. Hwang *

National Aeronautics and Space Administration
Lewis Research Center
Cleveland, Ohio 44135

Summary

A three-dimensional cubic cavity flow has been analyzed for diatomic gases by using the Boltzmann equation with the Bhatnagar-Gross-Krook (B-G-K) model. The method of discrete ordinate was applied, and the diffuse reflection boundary condition was assumed. The results, which show a consistent trend toward the Navier-Stokes solution as the Knudson number is reduced, give us confidence to apply the method to a three-dimensional geometry for practical predictions of rarefied-flow characteristics. The CPU time and the main memory required for a three-dimensional geometry using this method seem reasonable.

Symbols

a	defined in equations (5) and (6)
E_{σ}	energy level (equal to the roots of the Laguerre polynomial of degree n)
F	Maxwellian distribution function
f	distribution function
imax	maximum i index
jmax	maximum j index
Kn	Knudson number
k	Boltzmann constant
kmax	maximum k index

* Aerospace Engineer.

$k_\alpha, k_\beta, k_\gamma$	weighting coefficient of the odd equally spaced quadrature
L	length
m	mass of a molecule
n	number particle density
R	gas constant
R_σ	Gauss-Laguerre weighting coefficient
T	temperature
u	macroscopic velocity (flow velocity)
v	molecular velocity
w	velocity of moving surface
Y	vertical distance
Z_R	rotational relaxational parameter
λ	mean free path
μ	viscosity
ν	collision frequency
\wedge	nondimensional quantity
Subscripts	
el	elastic
eq	equilibrium
i, j, k	index of physical space
inel	inelastic
T	total
t	translational
x	x-component
y	y-component
z	z-component
α, β, γ	index of molecular velocity space
σ	discrete energy level
1	reference condition

Introduction

With space vehicles orbiting the world at hypersonic speed, numerical methods to cover all flow regimes become more important. Advances in computer hardware in recent years allow methods based on the kinetic theory to be used for practical applications. Many two-dimensional problems have been solved by the kinetic approach.¹⁻⁷ A limited number of three-dimensional problems have been solved recently by the direct-simulation Monte Carlo (DSMC) method.⁸⁻¹⁰ However, there are no three-dimensional numerical solutions based on the discrete ordinate method¹¹ applied to a Boltzmann equation.

A cubic cavity flow problem has been solved for diatomic gases using the Boltzmann equation with the Bhatnagar-Gross-Krook (B-G-K) model.^{12,13} Because neither theoretical nor experimental data are available for cubic cavity flow in free molecule flow, slip flow, and transition flow regimes, efforts have been made to obtain a continuum flow solution as

closed as possible so that the results from this study can be compared with the results of the Navier-Stokes solution.¹⁴

Governing Equation

The geometry of the problem considered is shown in figure 1. A cubic cavity with the same dimension on all sides is filled with diatomic gas (such as air), and the wall temperature is assumed to be the same on all walls. The top surface ($j = j_{\max}$ surface) is moving at a constant speed, and steady-state conditions are reached and analyzed numerically.

The distribution function at the energy level E_σ , f_σ , is governed by the Boltzmann equation with the Bhatnagar-Gross-Krook-Morse model for the collision terms:

$$v_x \frac{\partial f_\sigma}{\partial x} + v_y \frac{\partial f_\sigma}{\partial y} + v_z \frac{\partial f_\sigma}{\partial z} = \nu_{el}(F_{t\sigma} - f_\sigma) + \nu_{inel}(F_{i\sigma} - f_\sigma) \quad (1)$$

where

$$F_{t\sigma} = n_\sigma \left(\frac{m}{2\pi k T_t} \right)^{3/2} e^{-\frac{m}{2k T_t} ((v_x - u_x)^2 + (v_y - u_y)^2 + (v_z - u_z)^2)} \quad (2)$$

$$F_{i\sigma} = n_{\sigma eq} \left(\frac{m}{2\pi k T_T} \right)^{3/2} e^{-\frac{m}{2k T_T} ((v_x - u_x)^2 + (v_y - u_y)^2 + (v_z - u_z)^2)} \quad (3)$$

The elastic collision frequency $\nu_{el}^{15,16}$ is given by

$$\nu_{el} = \frac{nkT_t}{(1 + a)\mu} \quad (4)$$

whereas the inelastic collision frequency ν_{inel} is related to ν_{el} in the formula of

$$\nu_{inel} = a \nu_{el} \quad (5)$$

$$Z_R = \frac{5(1 + a)}{3a} \quad (6)$$

where the rotational relaxational parameter Z_R must be obtained by experiment. The viscosity-temperature relation of the form

$$\frac{\mu}{\mu_1} = \left(\frac{T_t}{T_1} \right)^s \quad (7)$$

is assumed and the value of s can be found in reference 17; for the current study, a and s are equal to 0.4 and 0.756, respectively, for air. The subscript 1 indicates the reference condition. The reference condition used in this study is the equilibrium flow in the cubic cavity when the top surface is stationary.

The reference viscosity is related to the reference mean free path λ_1 by the relation

$$\mu_1 = \frac{5}{16} mn_1 \lambda_1 (2\pi RT_1)^{1/2} \quad (8)$$

Combining equations (4) to (8) gives

$$\nu_{el} = \frac{16}{5(1+a)} \frac{nkT_t}{\lambda_1 mn_1 (2\pi RT_1)^{1/2}} \left(\frac{T_t}{T_1} \right)^{-s} \quad (9)$$

All macroscopic quantities, which will be listed in the nondimensional form, can be calculated in terms of f_σ .

Nondimensionalization

A characteristic velocity used for nondimensionalization is defined as

$$V_1 = (2RT_1)^{1/2} \quad (10)$$

The definitions of nondimensional variables are defined as follows:

$$\hat{n} = \frac{n}{n_1} ; \quad \hat{T} = \frac{T}{T_1} \quad (11)$$

$$\hat{u}_x = \frac{u_x}{V_1} ; \quad \hat{u}_y = \frac{u_y}{V_1} ; \quad \hat{u}_z = \frac{u_z}{V_1}$$

$$\hat{v}_x = \frac{v_x}{V_1} ; \quad \hat{v}_y = \frac{v_y}{V_1} ; \quad \hat{v}_z = \frac{v_z}{V_1}$$

$$\hat{E}_\sigma = \frac{E_\sigma}{kT_1}$$

$$\hat{x} = \frac{x}{L_1} ; \hat{y} = \frac{y}{L_1} ; \hat{z} = \frac{z}{L_1}$$

$$\hat{\nu}_{el} = \frac{\nu_{el} L_1}{V_1}$$

$$\hat{f}_\sigma = \frac{f_\sigma V_1^3}{n_1} ; \hat{F}_{t\sigma} = \frac{F_{t\sigma} V_1^3}{n_1} ; \hat{F}_{i\sigma} = \frac{F_{i\sigma} V_1^3}{n_1}$$

Equation (1) becomes

$$\hat{v}_x \frac{\partial \hat{f}_\sigma}{\partial \hat{x}} + \hat{v}_y \frac{\partial \hat{f}_\sigma}{\partial \hat{y}} + \hat{v}_z \frac{\partial \hat{f}_\sigma}{\partial \hat{z}} = \hat{\nu}_{el} ((\hat{F}_{t\sigma} - \hat{f}_\sigma) + a(\hat{F}_{i\sigma} - \hat{f}_\sigma)) \quad (12)$$

where

$$\hat{F}_{t\sigma} = \hat{n}_\sigma \left(\frac{1}{\pi \hat{T}_t} \right)^{3/2} e^{-\frac{(\hat{v}_x - \hat{u}_x)^2 + (\hat{v}_y - \hat{u}_y)^2 + (\hat{v}_z - \hat{u}_z)^2}{\hat{T}_t}}$$

$$\hat{F}_{i\sigma} = \hat{n}_{\sigma eq} \left(\frac{1}{\pi \hat{T}_T} \right)^{3/2} e^{-\frac{(\hat{v}_x - \hat{u}_x)^2 + (\hat{v}_y - \hat{u}_y)^2 + (\hat{v}_z - \hat{u}_z)^2}{\hat{T}_T}}$$

$$\hat{\nu}_{el} = \frac{8\hat{n}}{5(1+a)\sqrt{\pi Kn}} \hat{T}_t^{1-s}$$

$$Kn = \frac{\lambda_1}{L}$$

$$\hat{n}_\sigma = \int_{-\infty}^{+\infty} \int_{-\infty}^{+\infty} \int_{-\infty}^{+\infty} \hat{f}_\sigma d\hat{v}_x d\hat{v}_y d\hat{v}_z$$

$$\hat{n} = \sum_\sigma \hat{n}_\sigma$$

$$\hat{n}_{\sigma eq} = \hat{n} \left\{ \frac{\exp\left(-\frac{\hat{E}_\sigma}{\hat{T}_T}\right)}{\sum_s \exp\left(-\frac{\hat{E}_s}{\hat{T}_T}\right)} \right\} = \left(\frac{\hat{n}}{\hat{T}_T} \right) R_\sigma \exp\left[\hat{E}_\sigma \left(1 - \frac{1}{\hat{T}_T}\right)\right] \quad (\text{from ref. 1})$$

$$\hat{u}_x = \frac{1}{\hat{n}} \sum_\sigma \int_{-\infty}^{+\infty} \int_{-\infty}^{+\infty} \int_{-\infty}^{+\infty} \hat{f}_\sigma \hat{v}_x d\hat{v}_x d\hat{v}_y d\hat{v}_z$$

$$\hat{u}_y = \frac{1}{\hat{n}} \sum_\sigma \int_{-\infty}^{+\infty} \int_{-\infty}^{+\infty} \int_{-\infty}^{+\infty} \hat{f}_\sigma \hat{v}_y d\hat{v}_x d\hat{v}_y d\hat{v}_z$$

$$\hat{u}_z = \frac{1}{\hat{n}} \sum_\sigma \int_{-\infty}^{+\infty} \int_{-\infty}^{+\infty} \int_{-\infty}^{+\infty} \hat{f}_\sigma \hat{v}_z d\hat{v}_x d\hat{v}_y d\hat{v}_z$$

$$\hat{T}_t = \frac{2}{3\hat{n}} \left[\sum_\sigma \int_{-\infty}^{+\infty} \int_{-\infty}^{+\infty} \int_{-\infty}^{+\infty} [(\hat{v}_x - \hat{u}_x)^2 + (\hat{v}_y - \hat{u}_y)^2 + (\hat{v}_z - \hat{u}_z)^2] \hat{f}_\sigma d\hat{v}_x d\hat{v}_y d\hat{v}_z \right]$$

$$\hat{T}_i = \frac{1}{\hat{n}} \sum_\sigma \hat{n}_\sigma \hat{E}_\sigma$$

$$\hat{T}_T = \frac{3}{5} \hat{T}_t + \frac{2}{5} \hat{T}_i$$

Discretization of Distribution Function and Numerical Scheme

The distribution function at energy level σ , \hat{f}_σ , was discretized with equal spacing in the physical space, the arguments of the odd equally spaced quadrature (see appendix and ref. 18) was the spacing for the velocity space, and the roots of the Laguerre polynomial of degree 4 was used for the spacing for the internal energy space. The value of the distribution function at these discretized points was calculated by solving the Boltzmann equation using the finite difference method in the physical space.

The difference scheme is the first-order upwind difference based on the molecular velocity; for example,

$$\frac{\partial \hat{f}_\sigma}{\partial \hat{x}} = \frac{\hat{f}_{\sigma,i,j,k,\alpha,\beta,\gamma} - \hat{f}_{\sigma,i-1,j,k,\alpha,\beta,\gamma}}{\Delta \hat{x}} \quad \text{when } \hat{v}_x > 0$$

and

$$\frac{\partial \hat{f}}{\partial \hat{x}} = \frac{\hat{f}_{\sigma,i+1,j,k,\alpha,\beta,\gamma} - \hat{f}_{\sigma,i,j,k,\alpha,\beta,\gamma}}{\Delta \hat{x}} \quad \text{when } \hat{v}_x < 0$$

The distribution functions were then integrated for all macroscopic properties by using the odd equally spaced quadrature, e.g., the particle density at energy level σ and at the physical point (i, j, k) is

$$\int_{-\infty}^{+\infty} \int_{-\infty}^{+\infty} \int_{-\infty}^{+\infty} \hat{f}_\sigma d\hat{v}_x d\hat{v}_y d\hat{v}_z \approx \sum_{\alpha=-n'}^{n'} \sum_{\beta=-n''}^{n''} \sum_{\gamma=-n'''}^{n'''} k_\alpha k_\beta k_\gamma \hat{f}_\sigma(i,j,k,\alpha,\beta,\gamma)$$

where k_α , k_β , and k_γ are the weighting coefficients of the odd equally spaced quadrature for the velocity components \hat{v}_α , \hat{v}_β , and \hat{v}_γ , respectively. These properties were summed over the energy space, and then saved and used for the next iteration.

Boundary Conditions

For a constantly moving surface, perfectly diffuse reflection is assumed to specify the interaction of the molecules with the surface of the moving plate. Molecules which strike the moving surface ($j = j_{\max}$ surface) are emitted with a Maxwellian velocity distribution characterized by the plate temperature \hat{T}_w and the plate velocity \hat{w} ; that is,

$$\hat{f}_\sigma = \hat{n}_w \left(\frac{1}{\pi \hat{T}_w} \right)^{3/2} e^{-\frac{(\hat{v}_x - \hat{w})^2 + \hat{v}_y^2 + \hat{v}_z^2}{\hat{T}_w}}$$

where the density of molecules diffusing from the plate \hat{n}_w is not known a priori and must be found by applying the condition of zero mass flux normal to the plate at the surface, i.e.,

$$\sum_{\sigma} \left\{ \int_{-\infty}^{\infty} \int_0^{\infty} \int_{-\infty}^{\infty} \hat{v}_y \hat{f}_{\sigma}^{+} d\hat{v}_x d\hat{v}_y d\hat{v}_z + \int_{-\infty}^{\infty} \int_{-\infty}^0 \int_{-\infty}^{\infty} \hat{v}_y \hat{f}_{\sigma}^{-} d\hat{v}_x d\hat{v}_y d\hat{v}_z \right\} = 0$$

where \hat{f}_{σ}^{+} is the distribution function for $\hat{v}_y > 0$ and \hat{f}_{σ}^{-} is the distribution function for $\hat{v}_y < 0$. In the discrete ordinate form, this gives the relation at (i, j, k)

$$\hat{n}_w = \frac{2(\pi)^{1/2}}{\hat{T}_w^{1/2}} \sum_{\sigma=1}^n \sum_{\alpha=-n'}^{n'} \sum_{\beta=1}^{n''} \sum_{\gamma=-n'''}^{n'''} k_{\alpha} k_{\beta} k_{\gamma} \hat{v}_{\beta} \hat{f}_{\sigma, \alpha, \beta, \gamma}$$

Similarly, a zero-mass-flux boundary condition was applied to all other surfaces, and a Maxwellian velocity distribution function characterized by the plate temperature was used for the molecules emitted from the surface.

Because of the symmetry of the flow inside the cubic cavity, half of the cubic cavity was used for the computation. At the symmetry plane (i.e., $k = k_{\max}$ plane), the symmetry boundary condition was applied; that is, the distribution function of the outgoing molecule is equal to the distribution function of the incoming molecule at the symmetry plane.

Computational Procedures

For the present study, the initial condition assumed was that the flow everywhere was characterized by $\hat{n} = 1$ and $\hat{T} = 1$. As mentioned previously, the $j = j_{\max}$ surface was selected to be the moving surface. The computation began from the $j = 2$ surface with a known value for \hat{f}_{σ} at the $j = 1$ surface. On each $j = \text{constant}$ surface, the computation was performed along the i direction first and then marching toward the symmetry plane ($k = k_{\max}$) starting with \hat{f}_{σ}^{+++} . The symbol \hat{f}_{σ}^{+++} is defined as the subdistribution function at energy level σ for $\hat{v}_x > 0$, $\hat{v}_y > 0$, and $\hat{v}_z > 0$. At the symmetry plane, the symmetry boundary condition was applied (i.e., $\hat{f}_{\sigma}^{++-} = \hat{f}_{\sigma}^{+++}$). Knowing \hat{f}_{σ}^{++-} at the symmetry plane, the computation was marched backward from $k = k_{\max} - 1$ toward $k = 1$. Similarly, the computation continued for \hat{f}_{σ}^{-++} and \hat{f}_{σ}^{+-+} .

After the computation reached the $j = j_{\max}$ surface, the particle density was found at this surface by applying the no-flux boundary condition and then the computation marched downward from $j = j_{\max} - 1$ to $j = 1$ surface. The order of computation was \hat{f}_{σ}^{+-+} , \hat{f}_{σ}^{+--} , \hat{f}_{σ}^{--+} , and \hat{f}_{σ}^{---} . The particle density on the surfaces could then be found by using the no-flux boundary condition and the iteration

continued. The fourth-order Gauss-Laguerre quadrature was used for the energy space, and four sets of the third-order odd equally spaced quadrature were used for the molecular velocity space with the equal spacing of 0.15. (See the appendix.) The convergence was assumed when the maximum particle density increment was less than 10^{-4} .

Results

As mentioned previously, efforts have been made to calculate the flow with as small a Knudson number as possible so that the results could be compared with the solution of the Navier-Stokes equation. The results of this study are all based on $\hat{w} = 0.1$. Comparisons between the Boltzmann solution and the Navier-Stokes solution¹⁴ are given in figures 2 to 4.

Figure 2 shows the velocity profiles along the centerline for free molecule flow (Knudson number (Kn) = 100), slip flow ($Kn = 0.1$), and nearly continuum flow ($Kn = 0.03$). It clearly shows that the flow slips on the moving surface for $Kn = 100$ and $Kn = 0.1$; that is, the flow velocity on the surface is less than the speed of the moving wall. It also shows that the no-slip boundary condition of the Navier-Stokes equation was almost recovered for $Kn = 0.03$.

The comparison of the velocity vector plot on the symmetry plane between the Navier-Stokes solution and the Boltzmann solution is given in figure 3. The general shapes of the primary vortex are similar except that the center of the vortex for the Navier-Stokes solution is slightly farther downstream than that for the Boltzmann solution.

Figure 4 shows the surface static pressure looking from the center of the cubic cavity toward the upper corner for both the Navier-Stokes solution and the Boltzmann solution. The three-dimensional effect of the static pressure distribution is clearly shown qualitatively. The Boltzmann solution, even for the case of $Kn = 0.03$, is still not close to the continuum solution obtained by the Navier-Stokes equation; however, the consistent trend toward the Navier-Stokes solution is encouraging.

The symmetry plane velocity vector plots for two different Knudson numbers are shown in figure 5. The shape of the primary vortex is clearly shown even for $Kn = 100$. The center of the vortex moves upward toward the center obtained by the Navier-Stokes solution (fig. 3) when the Knudson number is reduced. It also clearly shows that the magnitude of the velocity vectors increases as expected when the Knudson number is reduced.

Figure 6 is a plot of the number particle density on the surfaces for three Knudson numbers. The distribution patterns for $Kn = 100$ and $Kn = 0.03$ are completely different. For a free molecular flow ($Kn = 100$), there are not enough molecular collisions to ensure the high number particle density at the upper downstream corner as shown for the case of $Kn = 0.03$. The high number particle density on the upper corner of the downstream vertical surface (i.e., near $i = i_{max}$ and $j = j_{max}$) and the low number

particle density on the upper corner of the upstream vertical surface (i.e., near $i = 1$ and $j = j_{\max}$) is a reasonable solution.

The convergence history is shown in figure 7 for $Kn = 0.03$. Each iteration took about 40 min on the Cray 2 supercomputer. Fortunately, it required less than 50 iterations to reach a convergent solution. In other words, it took less than 30 hr of CPU time for $Kn = 0.03$. The CPU time for $Kn = 100$ was less than 10 hr. To reduce the main memory requirement, the Maxwellian distribution functions and collision frequency were calculated repeatedly; as a consequence, extra CPU time was needed for each iteration. The grid size used in the study was 31 by 31 by 16 in the physical space, 12 by 12 by 12 in the molecular velocity space and 4 levels of internal energy in the energy space. Considerable CPU time can be saved by improving the integration scheme in the molecular velocity space. In any case, the discrete ordinate method used in this study can be applied to simple three-dimensional geometries without using the parallel programming technique.

Concluding Remarks

A three-dimensional cubic cavity flow was solved for diatomic gases by using the Boltzmann equation with the Bhatnagar-Gross-Krook (B-G-K) model. A comparison was made between the Boltzmann solution and the Navier-Stokes solution for the velocity profiles along the centerline, the primary vortex on the symmetry plane, and the surface static pressure. The general trend toward the Navier-Stokes solution as the Knudson number is reduced indicates that the solutions are very reasonable and that the discrete ordinate method can be used with confidence to a three-dimensional geometry for practical predictions of rarefied-flow characteristics. Because of the robustness of this numerical scheme, it requires less than 50 iterations to obtain a converged solution. The present method has a potential to be a practical flow simulation method to cover all flow regimes.

Appendix—Laguerre and Odd Equally Spaced Quadratures

Abscissa and Weight Factors for Laguerre Integration

$$\int_0^{\infty} e^{-E} f(E) dE = \sum_{i=1}^n R_i f(E_i)$$

where E_i are the abscissas and are the roots of Laguerre polynomials.

For $n = 4$,

$$E_1 = 0.322547689619$$

$$\begin{aligned}
E_2 &= 1.745761101158 \\
E_3 &= 4.536620296921 \\
E_4 &= 9.395070912301 \\
R_1 &= 6.03154104342 \times 10^{-1} \\
R_2 &= 3.57418692438 \times 10^{-1} \\
R_3 &= 3.88879085150 \times 10^{-2} \\
R_4 &= 5.39294705561 \times 10^{-4}
\end{aligned}$$

Odd Equally Spaced Quadrature

$$\int_0^1 f(x) dx \simeq \sum_{i=1}^n k_i f(\alpha_i)$$

where the arguments α_i are taken to be $1/(2n)$, $3/(2n)$, ..., $1 - [1/(2n)]$ and the weighting coefficients k_i for $n = 3$ are

$$\begin{aligned}
k_1 &= k_3 = 0.375 \\
k_2 &= 0.25
\end{aligned}$$

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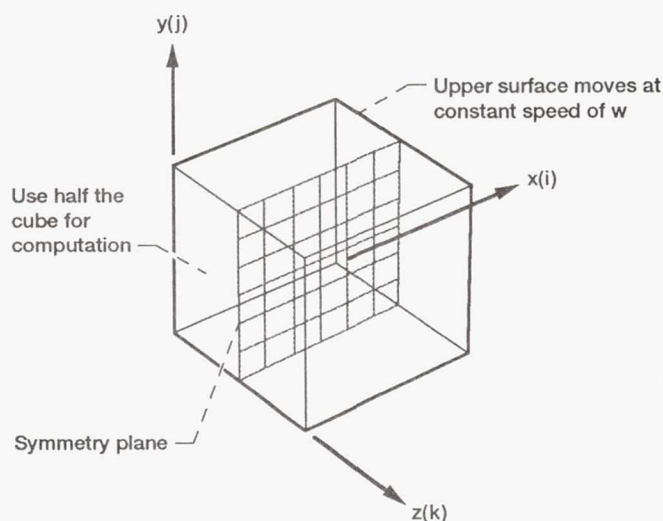


Figure 1.—Geometry.

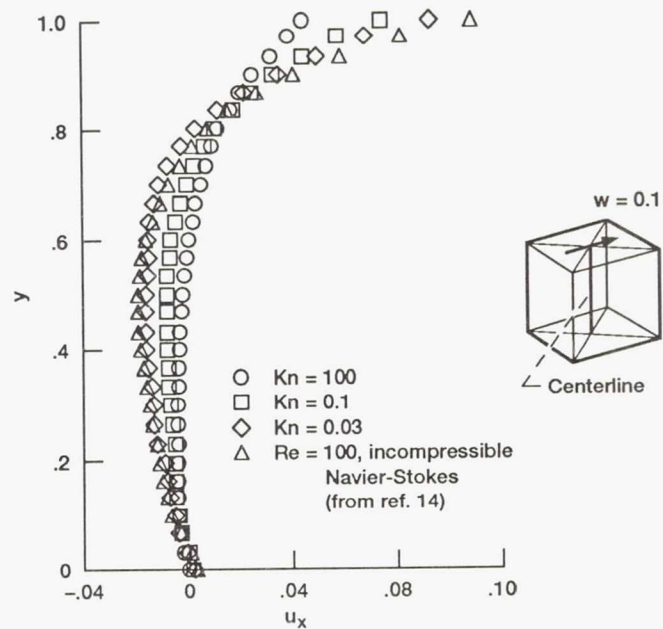
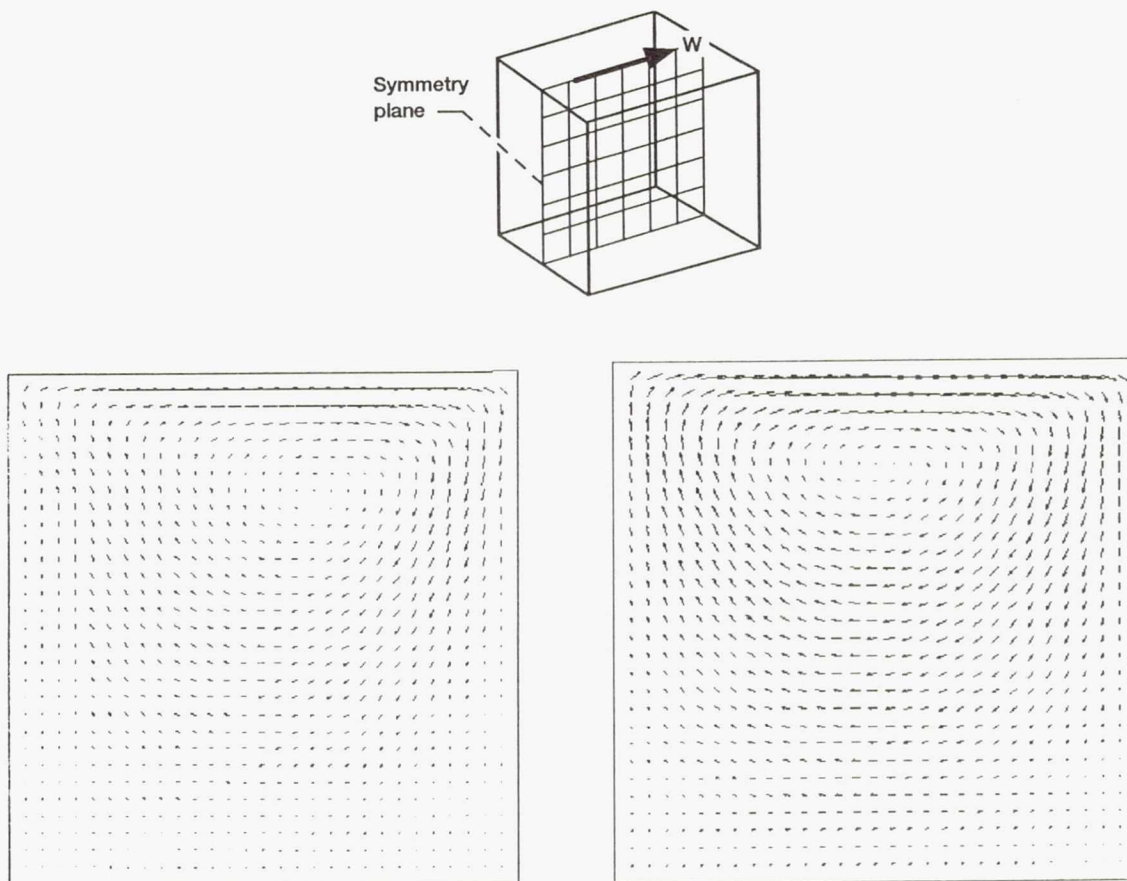


Figure 2.—Velocity profiles along centerline.



(a) Navier-Stokes solution ($Re = 100$).

(b) Boltzmann solution ($Kn = 0.03$).

Figure 3.—Velocity vectors on symmetry plane for Navier-Stokes solution and Boltzmann solution.

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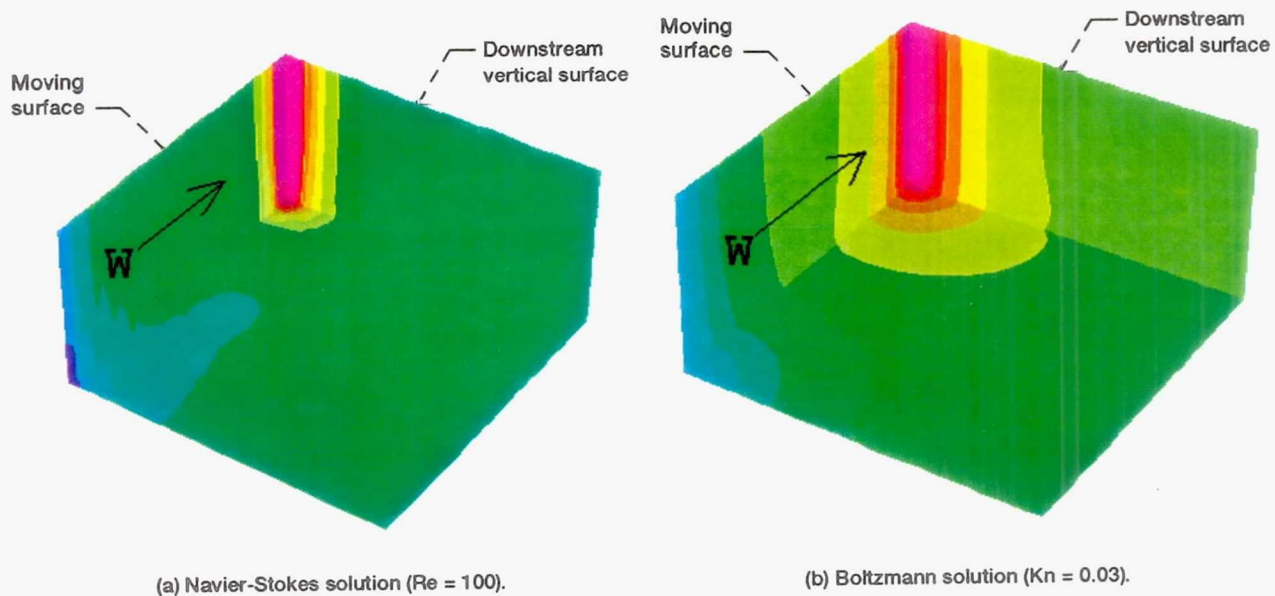
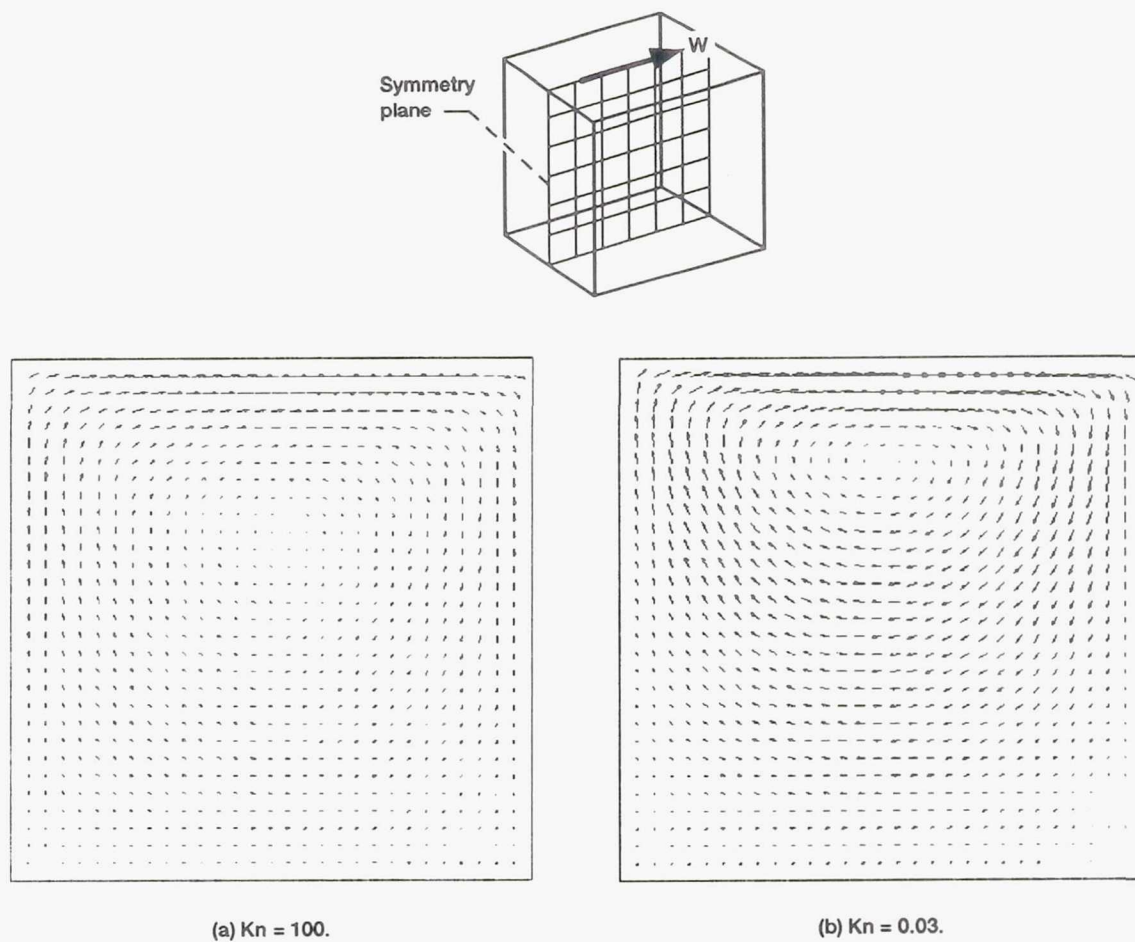


Figure 4.—Surface static pressure for Navier-Stokes solution and Boltzmann solution.



(a) $Kn = 100$. (b) $Kn = 0.03$.

Figure 5.—Velocity vectors on symmetry plane for two Knudson numbers.

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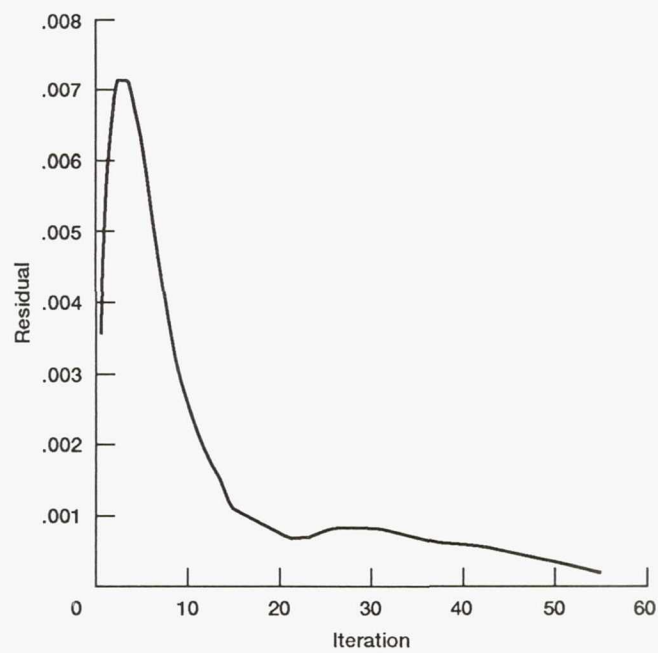


Figure 7.—Convergence history for $Kn = 0.03$.

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1. AGENCY USE ONLY (Leave blank)		2. REPORT DATE July 1992		3. REPORT TYPE AND DATES COVERED Technical Memorandum
4. TITLE AND SUBTITLE Numerical Solution of a Three-Dimensional Cubic Cavity Flow by Using the Boltzmann Equation			5. FUNDING NUMBERS WU-505-62-20	
6. AUTHOR(S) Danny P. Hwang				
7. PERFORMING ORGANIZATION NAME(S) AND ADDRESS(ES) National Aeronautics and Space Administration Lewis Research Center Cleveland, Ohio 44135-3191			8. PERFORMING ORGANIZATION REPORT NUMBER E-7080	
9. SPONSORING/MONITORING AGENCY NAMES(S) AND ADDRESS(ES) National Aeronautics and Space Administration Washington, D.C. 20546-0001			10. SPONSORING/MONITORING AGENCY REPORT NUMBER NASA TM-105693	
11. SUPPLEMENTARY NOTES Prepared for the 18th International Symposium on Rarefield Gas Dynamics, sponsored by the University of British Columbia, Vancouver, Canada, July 26-31, 1992. Responsible person, Danny P. Hwang, (216) 433-2187.				
12a. DISTRIBUTION/AVAILABILITY STATEMENT Unclassified - Unlimited Subject Category 34			12b. DISTRIBUTION CODE	
13. ABSTRACT (Maximum 200 words) A three-dimensional cubic cavity flow has been analyzed for diatomic gases by using the Boltzmann equation with the Bhatnagar-Gross-Krook (B-G-K) model. The method of discrete ordinate was applied, and the diffuse reflection boundary condition was assumed. The results, which show a consistent trend toward the Navier-Stokes solution as the Knudsen number is reduced, give us confidence to apply the method to a three-dimensional geometry for practical predictions of rarefied-flow characteristics. The CPU time and the main memory required for a three-dimensional geometry using this method seem reasonable.				
14. SUBJECT TERMS Boltzmann equation; Cubic cavity flow; Kinetic theory; B-G-K model, Discrete ordinate method			15. NUMBER OF PAGES 20	
			16. PRICE CODE A03	
17. SECURITY CLASSIFICATION OF REPORT Unclassified	18. SECURITY CLASSIFICATION OF THIS PAGE Unclassified	19. SECURITY CLASSIFICATION OF ABSTRACT Unclassified	20. LIMITATION OF ABSTRACT	