

## A Numerical Study of Shock Formation in Cylindrical and Two-Dimensional Shock Tubes

*By*

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*Summary:* Shock formation process and shock tube performance are investigated numerically. Numerical method applied is FLIC method proposed by Gentry, Martin and Daly for solving compressible fluid flows. A model of diaphragm opening process which does not base on the usual one-dimensional assumption is presented. Four cases for cylindrical shock tubes and one for two-dimensional were calculated for diaphragm pressure ratios of 100 and 1000. Diaphragm opening times were varied from about 100 to 500  $\mu$ sec. Shock formation process predicted from present calculations shows a qualitative agreement with experimental results. Shock tube performance was also calculated and compared with those predicted by ideal shock tube theory and other models such as formation-from-compression model proposed by White and multi stage model proposed by Ikui and Matsuo. Presented result shows a close agreement with that predicted by multi stage model.

### 1. INTRODUCTION

Most theoretical studies on shock tube flows have been based upon the usual ideal model which assumes; (1) one-dimensional flow, (2) ideal nonviscous, non-heat conducting driver and driven gases and (3) instantaneous diaphragm removal [1],[2]. According to these assumptions, wave speeds and the flow properties of the uniform states are readily determined from the initial conditions in terms of the specific heat ratios of driver and driven gases, diaphragm pressure ratio and initial temperature ratio. Schematic diagrams of this model are illustrated in Fig. 1.

In practice, however, it is not possible to remove instantaneously the diaphragm which separates high and low pressure chambers. Photographic investigations have revealed that the process by which a diaphragm actually ruptures involves formation of a small opening near the center, followed by the outward tearing of the material and the folding back of the leaves so formed. The effective cross-sectional area of the shock tube at the diaphragm location therefore increases gradually with time instead of attaining its maximum value instantaneously as assumed in the usual ideal analysis. Associated with this fact, a constant strength shock wave is not formed immediately but a train of compression waves generated by the

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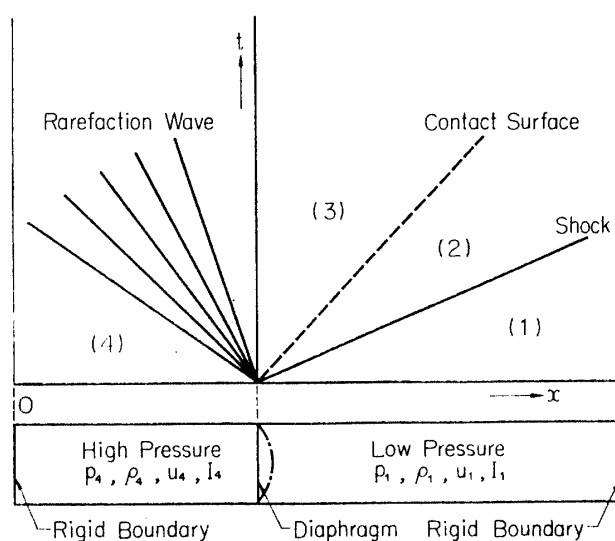


FIG. 1. Schematic diagrams of flow in a shock tube.

gradual opening of the diaphragm are sent out and overtake each other to form a weak shock wave. Then this shock wave is continuously accelerated by succeeding compression waves and finally attains its stationary speed.

White [3] has suggested a model which has taken into account this effect of finite opening time of the diaphragm. His model has assumed, instead of above-mentioned assumption (3), that diaphragm opening time is finite and continuously sent out unsteady compression waves all overtake the first wave at the same point to form a shock wave. When the same driver and driven gases such as air/air are selected, for a given diaphragm pressure ratio his theory always predicts a shock Mach number greater than that given by ideal shock tube theory. His model, however, does not include the effects of shock acceleration process.

Ikui and Matsuo [4] have extended White's model to take into account of this process of shock acceleration. Their model (multi stage model) has based upon the assumptions that continuous compression waves sent out through the diaphragm opening process can be divided into a finite number of groups, each of which is represented by a single compression wave, and a weak shock wave formed by the coalescence of such compression waves belonging to the first group is accelerated stepwise by succeeding compression waves. According to this model, still higher shock Mach numbers than White's model are predicted for the case of the larger diaphragm pressure ratios.

These two models still base on the one-dimensional assumption. As mentioned above, actual diaphragm opening process is not one-dimensional and analysis taking into account of this not-one-dimensional effects has not yet been performed.

In the present work, the process of shock formation in cylindrical and two-dimensional shock tubes is investigated numerically using the Fluid-in-Cell (FLIC) method [5]. A model of the diaphragm opening process not assuming one-dimensionality and computational cycle of FLIC method is described in some detail and results for shock formation distances and shock tube performance

are presented and discussed. FORTRAN programme used in this calculation is listed in appendix. Present calculations were performed on HITAC 5020 Fs at Institute of Space and Aeronautical Science, University of Tokyo and at National Aerospace Laboratory.

## NOTATION

$c$ ;	speed of sound
CPUT;	computational time
$D$ ;	hydraulic diameter of shock tube
$E$ ;	fluid energy per unit volume
$I$ ;	specific internal energy
$IN$ ;	maximum number of cells in $x$ -direction
$IN1$ ;	maximum number of cells in $x$ -direction in high pressure chamber
$IN2$ ;	maximum number of cells in $x$ -direction in low pressure chamber
$JN$ ;	maximum number of cells in $r$ (or $y$ )-direction
$M_s$ ;	shock Mach number
MEMSTP;	parameter controlling the diaphragm opening time
$N$ ;	maximum time cycle number
$P, p$ ;	static pressure
$r$ ;	coordinate perpendicular to shock tube axis of symmetry
$t$ ;	time
$u$ ;	velocity in $x$ -direction
$v$ ;	velocity in $r$ (or $y$ )-direction
$x$ ;	coordinate parallel to shock tube axis of symmetry
$x_f$ ;	shock formation distance
$y$ ;	coordinate perpendicular to shock tube axis of symmetry in two-dimensional case
$\gamma$ ;	specific heat ratio
$\Delta M$ ;	mass flowing across cell boundary during the time increment $\Delta t$
$\Delta r$ ;	cell dimension in $r$ -direction
$\Delta t$ ;	time increment
$\Delta x$ ;	cell dimension in $x$ -direction
$\Delta y$ ;	cell dimension in $y$ -direction
$\rho$ ;	fluid density
$\tau_{op}$ ;	diaphragm opening time

## SUBSCRIPTS

$i$ ;	cell number in $x$ -direction
$j$ ;	cell number in $r$ (or $y$ )-direction
$n$ ;	time cycle number
1;	initial condition of low pressure chamber
2;	state compressed by the shock wave
3;	state formed from the isentropic expansion of the gas in high pressure chamber
4;	initial condition of high pressure chamber

## 2. DIAPHRAGM OPENING PROCESS

As mentioned in the previous section, actual diaphragm opening process is three-dimensional, that is, diaphragm first ruptures from the center then follows continuous opening to full aperture. As the result a shock wave so formed by the coalescence of continuous compression waves is curved and contact surface becomes conical. Formation of compression waves in diaphragm opening process is to be considered to correspond to the rate of diaphragm opening. Therefore, to simulate this process, we must know whether it opens at an even rate or whether it opens more rapidly at some stage than others.

Campbell et al [6] has investigated experimentally the time dependence of the aperture size of rupture of aluminium diaphragms and the piercer-assisted bursting of cruciformed copper diaphragms. For both the natural rupture of an aluminium diaphragm and the piercer-assisted rupture of a copper diaphragm, a slow initiation of opening followed by a much more rapid opening to full aperture was observed. Their result shows that the time taken to 10% of full aperture was 100  $\mu\text{sec}$  while that from 10% to full aperture was 80  $\mu\text{sec}$  in the case of aluminium diaphragm and 750  $\mu\text{sec}$  (out of 1050  $\mu\text{sec}$  total) for copper. But analogous experiments for other materials have not yet been performed.

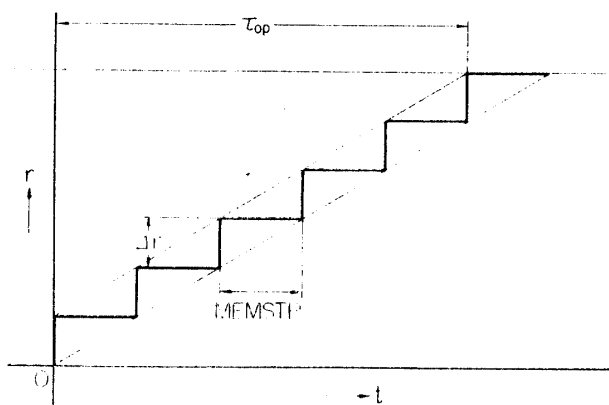


FIG. 2. Suggested model of diaphragm opening process.

In the present study, we simulate the diaphragm opening process as follows; at  $t=0$ , a plane diaphragm begins to rupture from the center, then stepwise increase of aperture is repeated at discrete intervals of time, MEMSTP and finally reaches to full aperture. Increment of the cross sectional area at diaphragm location is  $\Delta r$  or  $\Delta y$ , the cell dimension of  $r$  or  $y$  direction. This model of diaphragm opening process is illustrated in Fig. 2. Choosing an appropriate value of MEMSTP, we can control the rate of diaphragm opening. Throughout this calculation the values of MEMSTP are taken to be constant with time for the convenience of the calculation, therefore, rate of change of cross sectional area opened to high pressure gas is proportional to time in two-dimensional cases and to square of time in cylindrical cases.

### 3. CELL ARRANGEMENT AND BOUNDARY CONDITIONS

ELIC method requires a fixed (Eulerian) grid of rectangular cells of sides  $\Delta x$  and  $\Delta r$  (or  $\Delta y$  for two-dimensional problems) to cover the region under consideration. Fig. 3 indicates how the shock tube configuration fits into the cell system. In both cylindrical and two-dimensional problems the pair of integers  $(i, j)$  denotes the location of cell center with  $i$  and  $j$  increasing in the  $x$  and  $r$  (or  $y$ ) directions, respectively and half integer indices denote the cell boundaries.  $IN$  and  $JN$  are the total number of cells in  $x$  and  $r$  (or  $y$ ) directions.  $IN1$  and  $IN2$  indicate the number of columns of cells located in high pressure and low pressure chambers, respectively. In all cases except one, calculations were performed in cylindrical coordinates. Typical grid sizes were 400 cells in the  $x$ -direction and 10 cells in the  $r$ -direction. The plane  $r=0$  was chosen as a plane of symmetry, thus there being effectively 20 cells in the  $r$ -direction. Cell dimensions were taken to  $\Delta x = 2\Delta r$ .

Boundary conditions on the left, the top and the bottom boundaries were set to be reflective but the right boundary was set to be continuative. Reflective boundary conditions were also set on both sides of the diaphragm not yet opened to the fluid.

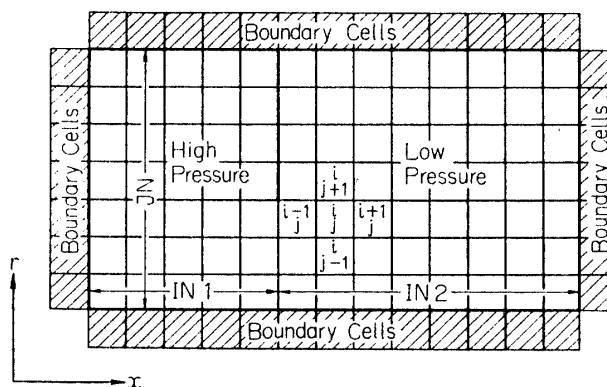


FIG. 3. The computing mesh.

### 4. METHOD OF SOLUTION

A number of numerical methods for time-dependent compressible flows containing shocks have been proposed. Representatives of such schemes are Lax's scheme [7], Godunov's [8] Rusanov's [9], Lax-Wendroff's [10] and Particle-in-Cell (PIC) [11]–[13] or Fluid-in-Cell (FLIC) method. Each of these schemes has characteristic properties which make it suitable for a certain class of problems. Little investigation has been performed as to which of these methods is optimum for a particular problem [14].

In the present calculations, FLIC method proposed by Gentry, Martin and Daly is chosen. The FLIC method is an outgrowth of the work of Rich [15], which is an extension of the PIC method to the case of a continuous fluid. Details of the

method can be found in references [6] and [15]. Here only an outline of the calculational cycles will be presented. The calculation of the time  $(n+1)$  configuration of the fluid from that at time  $n$  proceed in cycles in which all the variables of the cell quantities are advanced to new values slightly removed from the values of the preceding cycle representing a slightly later time.

The governing partial differential equations for compressible fluid flows are as follows;

Conservation of mass,

$$\left( \frac{\partial}{\partial t} + \vec{u} \nabla \right) \rho = -\rho \nabla \vec{u} \quad (1)$$

Conservation of momentum,

$$\rho \left( \frac{\partial}{\partial t} + \vec{u} \nabla \right) \vec{u} = -\nabla p \quad (2)$$

Conservation of energy,

$$\rho \left( \frac{\partial}{\partial t} + \vec{u} \nabla \right) \left\{ I + \frac{1}{2} (u^2 + v^2) \right\} = -\nabla (p \vec{u}) \quad (3)$$

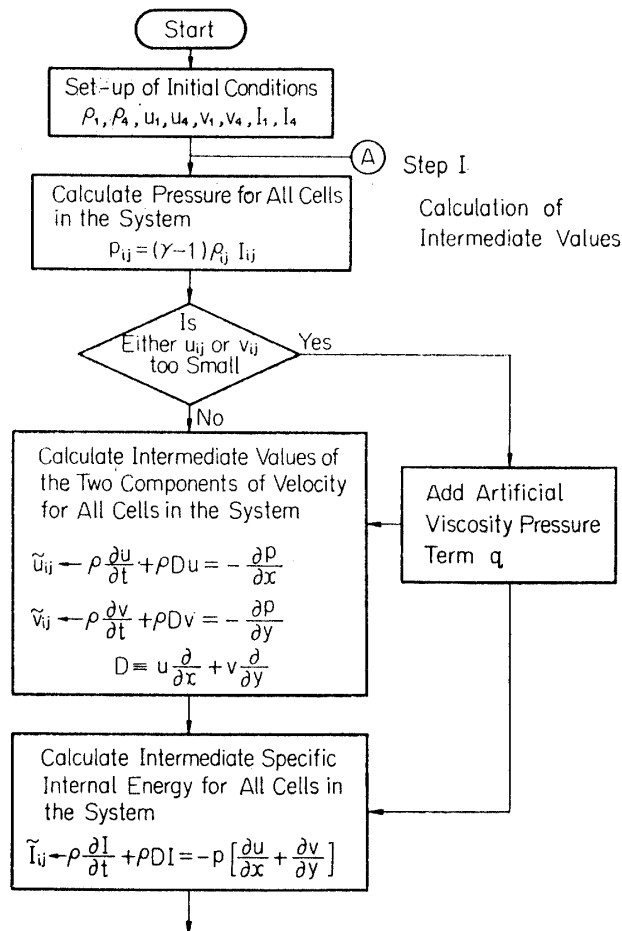


FIG. 4(a).

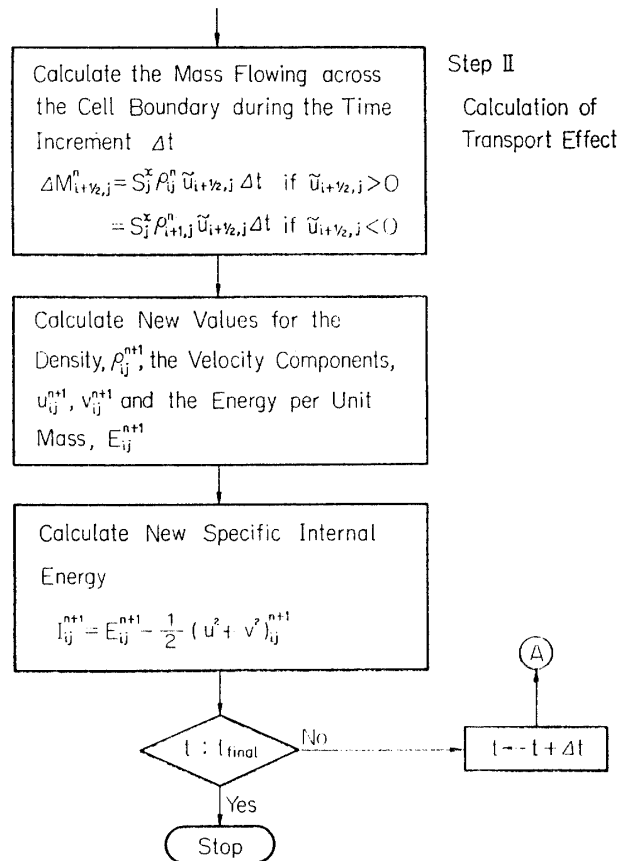


FIG. 4(b). Flow diagram of FLIC method in two-dimensional case  
(a) Phase I; (b) Phase II.

Equations (1), (2) and (3) are solved in two separate phases. Equation (1), that of mass conservation, may be eliminated from further consideration, as it is used only implicitly in the calculation of mass transport. Phase I treats the terms due to pressure forces only, and Phase II accounts for the remaining transport terms. Thus the second terms on the left-hand side of equations (2) and (3) are temporarily dropped in Phase I.

Briefly, Phase I begins with the calculation of new pressures for all cells from cell values of density and specific internal energy at time  $n$  or from those of initial conditions. From the gradients of these pressures are calculated tentative new values of the two components of velocity for each cell. From these new velocity components and those from the preceding cycle or from the initial condition, tentative new values of specific internal energy are computed for all cells. In both instances, the values are tentative because of omission of transport terms from the equations given above. As was seen in Phase I of the calculational cycle, only cell quantities were updated, while the fluid was assumed to be momentarily completely at rest.

In the second stage (Phase II) the term that were omitted in Phase I are accounted for, so that the cell density, momenta and energy are updated to time  $(n+1)$  for the effects due to transport. Mass fluxes across cell boundaries are first computed for each cell, then mass at time  $(n+1)$  is calculated for each cell

from the mass at time  $n$  and mass fluxes across four boundaries. New velocity components and specific energy are calculated from those obtained in Phase I and new density just calculated. To complete the time cycle, the pressure is calculated from the density and specific internal energy and this present the starting condition for the next cycle. To help clear understanding of this calculational cycle, flow diagrams for two-dimensional case are shown in Fig. 4(a) and (b). FORTRAN program for the present calculation is listed in appendix.

## 5. RESULTS AND DISCUSSIONS

Using the numerical technique described in the prior section, four cases for cylindrical shock tube and one for two-dimensional were calculated for various combinations of diaphragm opening times and diaphragm pressure ratios. These cases with some representative results are summarized in Table 1. Both driver and driven gases were assumed to be ideal gases and the value of specific heat ratio,  $\gamma$  was taken to be 1.4. The hydraulic diameter of the shock tube,  $D$  was assumed to be 10 cm. All numerical calculations were performed on HITAC 5020 Fs at Institute of Space and Aeronautical Science, University of Tokyo and National Aerospace Laboratory.

TABLE 1. Summary of calculational conditions and results.

$P_{41}$	MEMSTP	$N$	CPUT	$\tau_{op}$ ( $\mu$ sec)	$X_f/D$	$M_s$
100	6	480	1801	99.3	7.8	2.40
	16	520	1801	264.7	8.2	2.40
	16	"	"	"	8.6	2.39
	30	600	1963	496.3	>10.7	>2.36
1000	16	1200	5119	264.7	12.0	3.30

### Shock formation process

Henshall [16] has explained the process of shock formation as follows; when the bowed diaphragm is ruptured, a number of curved compression waves are propagated into the channel. These compression waves very quickly coalesce to form a curved shock wave. Its profile will become less curved as it progresses down the channel. Regular reflection of this curved shock will take place at the walls of the tube until the angle of incidence of the shock is such that regular reflection of the shock is impossible and the shock then undergoes Mach reflection. The triple points of the Mach reflection move toward the center of the tube, thus producing a plane shock wave. Subsequently, the triple points continue to move back and forth across the tube until the secondary branches of the Mach configuration are very weak and disappear. Finally, an optically flat primary shock is propagated down the channel.



Experimental investigations on the process of shock formation after diaphragm rupture have been performed by Geiger and Mautz [17] for the Cellophane diaphragms in a 2"×7" shock tube using the Shadowgraph technique. Ikui and Matsuo [18] have also investigated the process in a 38×38 mm shock tube by the Schlieren method. Photographs obtained by Geiger and Mautz support Henshall's explanation.

To see the flow patterns near the diaphragm at the initial stage of the opening process, calculated velocity distributions near the diaphragm are shown in Fig. 5 (a)–(c) for the case of  $\tau_{op}=496.3 \mu\text{sec}$  and locations of shock fronts at various radial positions are also displayed in Fig. 6(a)–(c). Fig. 6(a) shows the result for diaphragm pressure ratio,  $P_{41}=100$  and  $\tau_{op}=264.7 \mu\text{sec}$ , 6(b) is for the same diaphragm pressure ratio as 6(a) but  $\tau_{op}=496.3 \mu\text{sec}$ . Result for  $P_{41}=1000$  and  $\tau_{op}=264.7 \mu\text{sec}$  is plotted in Fig. 6(c).

The followings can be seen from Figs. 5 and 6; immediately after the diaphragm rupture, radial flow of the high pressure gas is initiated and a weak curved shock first formed by coalescence of compression waves becomes gradually plane as it propagates down the tube, which shows that the one-dimensional assumption does

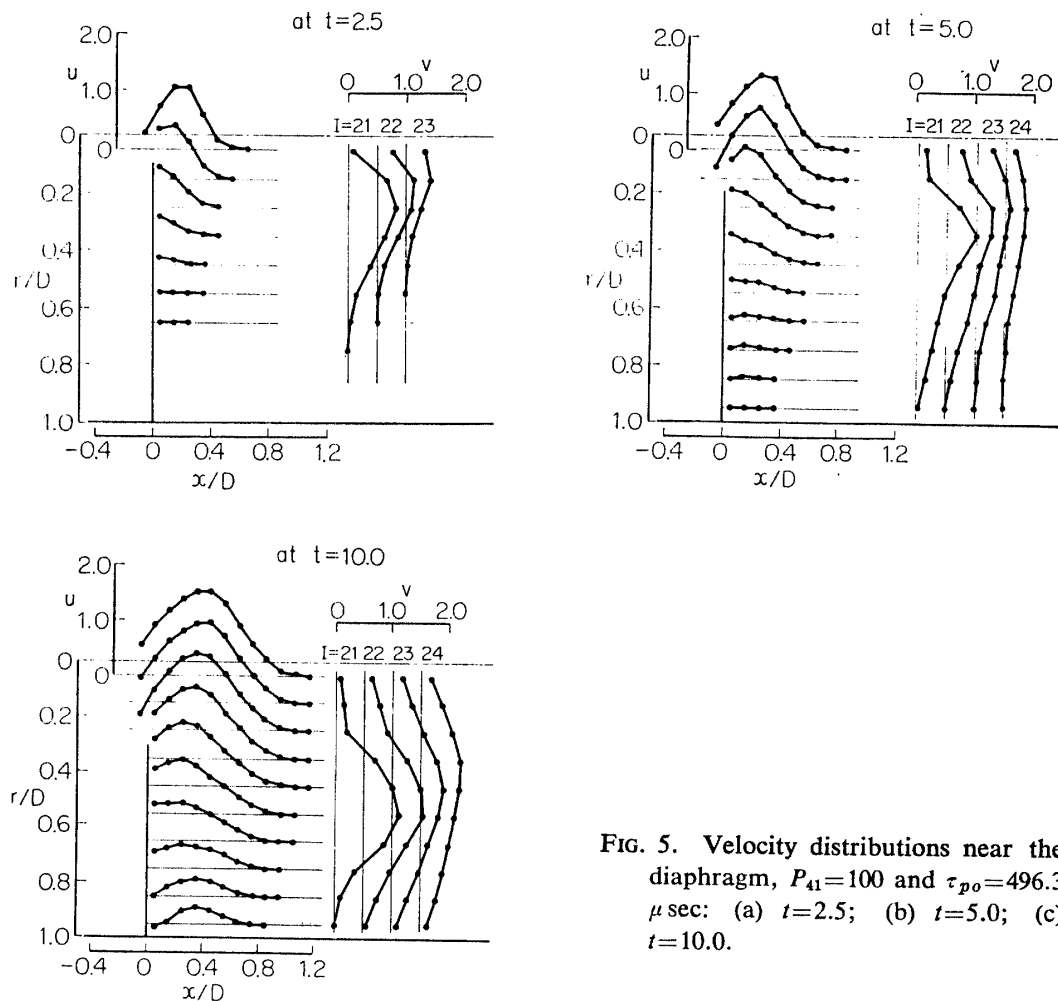


FIG. 5. Velocity distributions near the diaphragm,  $P_{41}=100$  and  $\tau_{po}=496.3 \mu\text{sec}$ : (a)  $t=2.5$ ; (b)  $t=5.0$ ; (c)  $t=10.0$ .

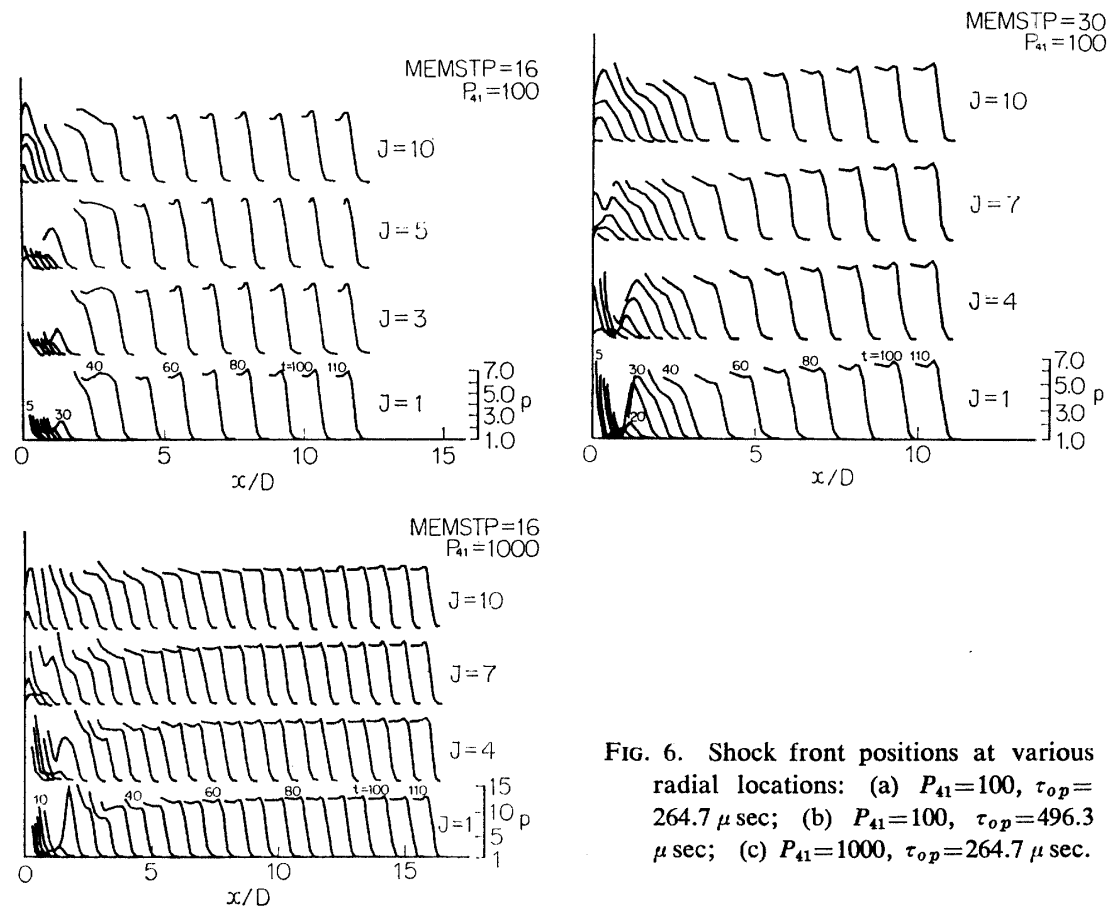
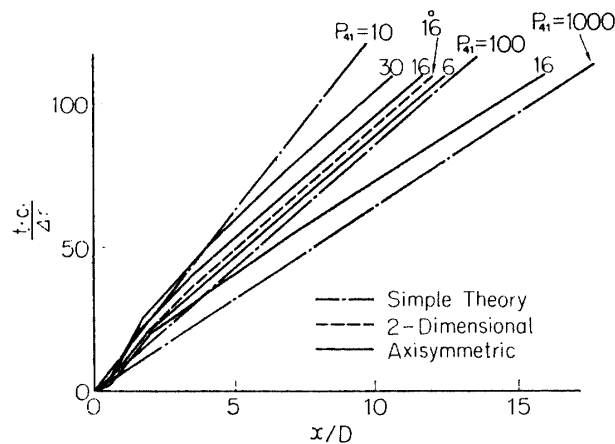


FIG. 6. Shock front positions at various radial locations: (a)  $P_{41}=100$ ,  $\tau_{op}=264.7 \mu \text{ sec}$ ; (b)  $P_{41}=100$ ,  $\tau_{op}=496.3 \mu \text{ sec}$ ; (c)  $P_{41}=1000$ ,  $\tau_{op}=264.7 \mu \text{ sec}$ .

not hold in this case. Distance travelled by the primary shock before it becomes plane is about 2 to 3 diameters. This result agrees well with that of 2.4 times hydraulic diameters obtained by above-mentioned experiment of Ikui and Matsuo. Results presented in Figs. 5 and 6 seem to support Henshall's explanation. Although, transition from regular reflection to Mach one cannot be fully identified due to insufficient resolution. Photographic investigation of Geiger and Mautz also proves this process as was indicated before. On the other hand, Ikui and Matsuo report that such reflected waves cannot be observed in their experiment but it should be considered that sensitivity of the Schlieren apparatus was not satisfactory.

### Shock formation distance

As is generally known, shock formation distance is one of the most important parameters in the shock tube problems but it is very difficult to determine this theoretically. In this section we shall discuss the effect of the diaphragm opening time on shock formation distance. Fig. 7 shows  $x-t$  diagrams calculated by FLIC method for various combinations of diaphragm pressure ratio,  $P_{41}$  and diaphragm opening time,  $\tau_{op}$ . In Fig. 7  $x-t$  diagrams predicted by the ideal shock tube theory are also shown. Since the slope of tangent of  $x-t$  diagram is reciprocal of shock speed, the smaller incrementation shows the higher shock Mach number. As is


 FIG. 7. Calculated  $x-t$  diagram for  $j=1$  cells.

clearly seen from Fig. 7 the shock front is accelerated by succeeding compression waves gradually to its final speed. This seems to prove that multi stage model proposed by Ikui and Matsuo is reasonable. Although shock formation distances tend to increase with diaphragm opening time, for the same diaphragm pressure ratio shock Mach numbers converge to a value which is greater than that predicted by ideal shock tube theory. As in our model of diaphragm opening process applied to this calculation, rate of change of cross-sectional area opened to high pressure gas is set to be proportional to time in two-dimensional case and to square of time in cylindrical case for the convenience of calculation, for the same diaphragm opening time, shock front is accelerated faster in two-dimensional case than in cylindrical case, which can be clearly seen in Fig. 7.

From  $x-t$  diagrams displayed in Fig. 7 variations of shock Mach numbers along the shock tube axis are obtained and plotted in Fig. 8 for several diaphragm opening time for the case of  $P_{41}=100$ . Fig. 8 also shows the tendency that although acceleration of shock fronts is slower for larger diaphragm opening time, shock Mach numbers for all cases converge to the same value. This shows a qualitative agreement with the result of Shtemenko's experiment [19]. In his experiment,

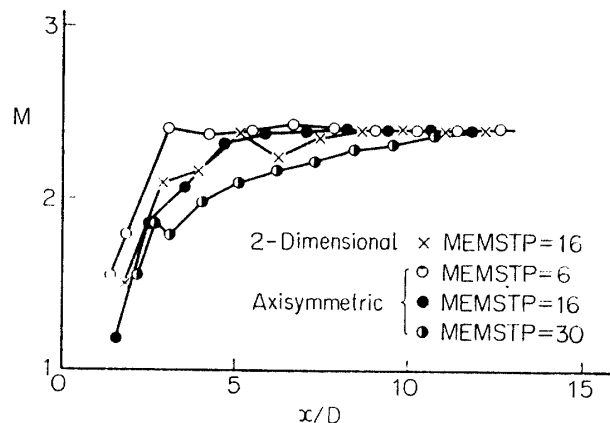


FIG. 8. Variation of shock Mach numbers with distance along shock tube axis.

maximum shock Mach numbers decrease slightly with increasing diaphragm opening time but this may be due to the effect of viscosity and should be a constant value neglecting such effect. As is clearly seen from Figs. 7 and 8, for the case of  $\tau_{op}=496.3 \mu\text{sec}$ , the largest value of diaphragm opening time, shock front is still accelerating. Shock formation distances calculated from Fig. 8 are tabulated in Table 1.

### Shock tube performance

According to the ideal shock tube theory, shock Mach numbers are easily calculated from specific heat ratios of driver and driven gases, diaphragm pressure ratio and initial temperature ratio. In actual shock tubes, however, shocks with higher Mach numbers than predicted by the ideal theory have been observed for diaphragm pressure ratios of the order of  $10^4$ . Ikui and Matsuo [18] have reported that even for the diaphragm pressure ratios less than  $10^2$ , stronger shocks than predicted by ideal theory were observed in a shock tube with large diameter. The formation-from-compression model proposed by White predicts significantly stronger shock waves than the usual ideal model for the larger diaphragm pressure ratios. Multi stage model proposed by Ikui and Matsuo which is an extension of White's model to take into account the effects of shock acceleration gives still stronger shocks than White's model.

In the present calculation by FLIC numerical method we did not base upon the one-dimensional assumption on which all the above-mentioned models were based. In Fig. 9 calculated shock tube performance is plotted and compared with those predicted by ideal, White's and multi stage model. Present result shows a close agreement with multi stage model for the cases of  $P_{41}=10$  and 100 but significantly higher shock Mach number is predicted for  $P_{41}=1000$ . Calculated data are so insufficient to draw any conclusion about this that further investigations should be needed.

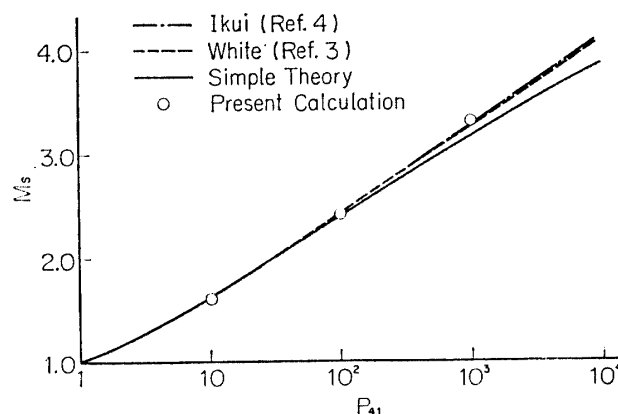


FIG. 9. Predicted shock tube performance for air/air.

## 6. CONCLUDING REMARKS

It may be summarised from the calculated result using the FLIC method that shock formation process is not one-dimensional as assumed in ideal shock tube theory and other models proposed by various authors, and shock formation distances tend to increase with diaphragm opening time and with diaphragm pressure ratios, which shows qualitative agreement with experimental results. For  $P_{41}=100$ , calculated shock formation distances varies from about 8 diameters to above 11 according to the diaphragm opening time. For  $\tau_{op}=496.3 \mu\text{sec}$ , the largest value of diaphragm opening time shock front is still accelerating. Calculated shock tube performance shows a close agreement with that predicted by multi stage model for the cases of  $P_{41}=10$  and 100. For  $P_{41}=1000$ , however, significantly higher shock Mach number is predicted by FLIC method. Calculated data are so insufficient to draw any conclusion that further investigation should be needed.

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## APPENDIX

C ***	***		
C	COMMENT		
C ***	***	C	1
C ***	***		
C	HITACHI CORD		
C	FLIC2C	C	2
C	NUMERICAL CALCULATION FOR AERODYNAMICS	C	3
C	FLUID-IN-CELL (FLIC) CALCULATION	C	6
C	TWO-DIMENSIONAL FORMULATION IN RECTANGULAR AND CYLINDRICAL COORDINATES	C	7
C	TWO DIMENSIONAL SHOCK FORMATION IN A SHOCK TUBE	C	8
C	ITYPE = 1 RECTANGULAR COORDINATES	C	9
C	ITYPE = 2 CYLINDRICAL COORDINATES		
C	ITAPE = 1 R/T EXECUTION		
C	ITAPE = 0 R/T NON EXECUTION		
C	COMPUTATION TIME SAVING PROGRAMME		
C	ISTP(J).....CALCULATION FROM I=1 TO ISTP(J)	C	
C	PRINT PAGE SAVING PROGRAMME		
C	ISTPP(J).....PRINT FROM I=1, TO ISTPP(J)	C	
C ***	***		
C ***	***		
C	INITIAL SET-UP	SUP	1
C ***	***		
C ***	***		
	DIMENSION RGU(440,10),P(440,10),C(440,10),U(440,10),V(440,10),UTD(	SUP	2H
	1440,10),VTD(440,10),EI(440,10),ETD(440,10),E(440,10),ETD(440,10),	SUP	3H
	2QU(440),DMUL(440),ISTP(10),ISTPP(10)	SUP	4H
	EQUIVALENCE (P,E)	SUP	5
	CALL DPON	SUP	6
	CALL OFCOFF(3)	SUP	61
70	WRITE(6,1)	SUP	7
60	READ(5,2)DX,DR,GAM,STEP	SUP	8
	IF(DX.EQ.0.) GO TO 70	SUP	9
	IF(DX.LT.0.) STOP	SUP	10
	READ(5,3)UINF,PINF,CINF,RATIO	SUP	11
	READ(5,4)IN1,IN2,IN3,JN,ITYPE	SUP	12
	READ(5,5)LJ1,LJ2,LJ3,NPR1,NPR2,ILAST,ISTART,IRUN	SUP	13
	READ(5,6)AK,B	SUP	14
	READ(5,7)ITAPE	SUP	14A
	IF(ITAPE.EQ.1) REWIND 3	SUP	14B
1	FORMAT(1H1,30HFLUID IN CELL CALCULATION	SUP	15
2	FORMAT(4F12.0)	SUP	16
3	FORMAT(5F12.0)	SUP	17
4	FORMAT(4I8)	SUP	18
5	FORMAT(8I8)	SUP	19
6	FORMAT(2F12.0)	SUP	20
7	FORMAT( 18)	SUP	201
	READ(5,50)REMSTP,IMT	SUP	21
50	FORMAT(2I8)	SUP	22
	PAI=3.14159	SUP	23
	ROINF=GAM*PINF/CINF**2	SUP	24
	EIINF=PINF/((GAM-1.)*ROINF)	SUP	25
	EINF=EIINF+(UINF**2+VINF**2)/2.	SUP	26
	IN=IN1+IN2	SUP	27
	T=0.	SUP	28
	ITIME=0	SUP	29
	ITOPP=0	SUP	29A
	IPRINT=LJ1	SUP	30

MEMR=0	SUP	31
DT=STEP*DX/CINF	SUP	32
IF(DR.LT.DX) DT=STEP*DR/CINF	SUP	33
DO 8 J=1,JN	SUP	34
DO 9 I=1,IN	SUP	35
ROU(I,J)=ROINF	SUP	36
IF(I.LE.IN1) ROU(I,J) =RATIO*ROINF	SUP	37
EI(I,J)=EIINF	SUP	38
U(I,J)= UINF	SUP	39
V(I,J)= VINF	SUP	40
9 CONTINUE	SUP	41
8 CONTINUE	SUP	42
CMTOT=0.	SUP	43
ETOT=0.	SUP	44
VOL=DX*DR	SUP	45
DO 10 J=1,JN	SUP	46
IF(ITYPE.EQ.2) VOL=2. *PAI*(FLOAT(J)-0.5)*DX*DR**2	SUP	47
DO 11 I=1,IN	SUP	48
CMTOT=CMTOT+ROU(I,J)*VOL	SUP	49
E(I,J)=EI(I,J)+(U(I,J)**2+V(I,J)**2)/2.	SUP	50
ETOT=ETOT+E(I,J)*ROU(I,J)*VOL	SUP	51
11 CONTINUE	SUP	52
10 CONTINUE	SUP	53
ETH=ETOT	SUP	54
CMTH=CMTOT	SUP	55
EPSE=10.0*EINF/ETOT	SUP	56
EPSE=1. E-3	SUP	57
IF(ITER.GE.2) GO TO 888	SUP	58
WRITE(6,14)	SUP	59
14 POWER(140,40)INITIAL TIME FLOW FIELD PROPERTY PLOT )	SUP	60
DO 12 J=1,JN	SUP	61M
ISTP(J)=IN1+1	SUP	62M
ISTPP(J)=IN1+1	SUP	62M
12 CONTINUE	SUP	63M
GO TO 605	SUP	64M
C ***		
C ***		
C I-1 PRESSURE AND SOUND SPEED CALCULATION	I -1	1
C ***		
C ***		
888 READ(3) ITIME, IPRINT, CMTH, ETH,DT	I -1	1
READ(3) ((ROU(I,J),U(I,J),V(I,J),EI(I,J),I=1,IN),J=1,JN)	I -1	2
IF(ITIME.EQ.1START) GO TO 102	I -1	3
GO TO 888	I -1	4
102 DO 103 J=1,JN	I -1	6
DO 104 I=1,IN	I -1	7
P(I,J)=(GAM-1. )*ROU(I,J)*EI(I,J)	I -1	8
C(I,J)=SQR(GAM*(GAM-1. )*EI(I,J))	I -1	9
104 CONTINUE	I -1	10
103 CONTINUE	I -1	11
IF(ITIME.EQ.1START) GO TO 118	I -1	11A
IF(ROD(ITIME,IPRINT).EQ.0) GO TO 105	I -1	12
GO TO 118	I -1	13
105 DO 106 J=1,JN	I -1	14
WRITE(6,108) J	I -1	15
WRITE(6,109)	I -1	16
ITER=ITER+1	I -1	161



```

DO 107 I=1,INTPP1
  ABVEL=SQRT(U(I,J)**2+V(I,J)**2)
  AMACH=ABVEL/C(I,J)
  SPEED=ABVEL/SINF
  IF (ABVEL.EQ.0.) GO TO 151
  IF(U(I,J)**2.LT.V(I,J)**2) GO TO 150
  ANGLE=ATAN(V(I,J)/U(I,J))*57.29578
  IF(U(I,J).LT.0. .AND.V(I,J).GE.0.) ANGLE=ANGLE+180
  IF(U(I,J).LT.0. .AND.V(I,J).LT.0.) ANGLE=ANGLE-180
  GO TO 152
150 IF(V(I,J).GE.0.) GO TO 149
  ANGLE=-90.-ATAN(U(I,J)/V(I,J))*57.29578
  GO TO 152
149 ANGLE=90.-ATAN(U(I,J)/V(I,J))*57.29578
  GO TO 152
151 ANGLE=0.
152 NMOD=MOD(I,3)+1
  GO TO (111,109,110),NMOD
109 WRITE(6,113)I, P(I,J),AMACH,SPEED,ANGLE
  GO TO 107
110 WRITE(6,114)I, P(I,J),AMACH,SPEED,ANGLE
  GO TO 107
111 WRITE(6,115)I, P(I,J),AMACH,SPEED,ANGLE
107 CONTINUE
106 CONTINUE
299 >FORMAT(1H0,3(40H I PRESSURE AMACH SPEED ANGLE ))
108 FORMAT(1H0,2HJ=I5)
113 FORMAT(1H , 14,4F9.4 )
114 FORMAT(1H+,40X, 14,4F9.4)
115 FORMAT(1H+,50X, 14,4F9.4)
118 IF(ITIME.EQ.ILAST) GO TO 119
  IF(P(IN,1).GE.1.5*PINF) GO TO 119
  ITIME=ITIME+1
  IF(ITIME.GE.NPR1.AND.ITIME.LT.NPR2) IPRINT=LJ2
  IF(ITIME.GE.NPR2) IPRINT=LJ3
  IF(MOD(ITIME,MEMSTP).EQ.1) MEMR=MEMR+1
  IF(MEMR.GE.JN) MEMR=JN
  IF(ITIME.NE.IDT) GO TO 199
  ITOFF=IDT-1
  DT=2.0*DT
  TOFF=T
  GO TO 199
119 IF(ITAPE.NE.1) GO TO 60
  REWIND 3
  GO TO 60
C ***
C ***
C I-2 THE FILDE VELOCITY CALCULATION
C ***
C ***
199 VOL=DX*DR
  S2=DX
  S4=DX
  J=1
200 IF(ITYPE.NE.2) GO TO 201
  VOL=2.*PAI*(FLOAT(J)-.5)*DX*DR**2
  S2=2.*PAI*FLOAT(J-1)*DR*DX

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I -1 17
I -1 18
I -1 19
I -1 20
I -1 21
I -1 22
I -1 23
I -1 24
I -1 25
I -1 26
I -1 27
I -1 28
I -1 29
I -1 30
I -1 31
I -1 32
I -1 33
I -1 34
I -1 35
I -1 36
I -1 37
I -1 38
I -1 39
I -1 40
I -1 41
I -1 42
I -1 43
I -1 44
I -1 45
I -1 46
I -1 47
I -1 471
I -1 48
I -1 49
I -1 50
I -1 51
I -1 52
I -1 52A
I -1 52B
I -1 52C
I -1 52D
I -1 53
I -1 54
I -1 55
I -1 56
I -2 1
I -2 2
I -2 3
I -2 4
I -2 5
I -2 6
I -2 7
I -2 8

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	S4=2. *PAI*FLOAT(J)*DR*DX	I -2	9
201	I=1	I -2	10
202	IF(I.EQ.1.OR.(I.EQ.IN1+1.AND.J.GT.MEMR)) GO TO 203	I -2	11
	P1=P(I-1,J)	I -2	12
	Q1=QRL	I -2	13
	GO TO 204	I -2	14
203	P1=P(I,J)	I -2	15
	Q1=0.	I -2	16
	IF((AK*(U(I,J)**2+V(I,J)**2).LT.C(I,J)**2).AND.(U(I,J).LT.O. ))	I -2	17
	1 Q1=-2. *B*C(I,J)*U(I,J)*ROU(I,J)	I -2	18
204	IF(J.EQ.1) GO TO 205	I -2	19
	P2=P(I,J-1)	I -2	20
	Q2=QUL(I)	I -2	21
	GO TO 206	I -2	22
205	P2=P(I,J)	I -2	23
	Q2=0.	I -2	24
	IF((TYPE.EQ.1).AND.(AK*(U(I,J)**2+V(I,J)**2).LT.C(I,J)**2).AND.(V(I	I -2	25
	1(I,J).LT.O. )) Q2=-2. *B*C(I,J)*ROU(I,J)*V(I,J)	I -2	26
206	IF(I.EQ.IN.OR.(I.EQ.IN1.AND.J.GT.MEMR)) GO TO 207	I -2	27
	P3=P(I+1,J)	I -2	28
	Q3=0.	I -2	29
	IF((AK*(U(I,J)**2+U(I+1,J)**2+V(I,J)**2+V(I+1,J)**2).LT.(C(I,J)**2I	I -2	30
	1+C(I+1,J)**2)).AND.(U(I,J).GT.U(I+1,J))) Q3=B*(C(I,J)+C(I+1,J))*	I -2	31
	2(ROU(I,J)+ROU(I+1,J))*(U(I,J)-U(I+1,J))/4.	I -2	32
	QRL=Q3	I -2	33
	GO TO 208	I -2	34
207	P3=P(I,J)	I -2	35
	Q3=0.	I -2	36
	IF((AK*(U(I,J)**2+V(I,J)**2).LT.C(I,J)**2).AND.(U(I,J).GT.O. ))	I -2	37
	1 Q3=2. *B*C(I,J)*ROU(I,J)*U(I,J)	I -2	38
208	IF(J.EQ.JN) GO TO 209	I -2	39
	P4=P(I,J+1)	I -2	40
	Q4=0.	I -2	41
	IF((AK*(U(I,J)**2+U(I,J+1)**2+V(I,J)**2+V(I,J+1)**2).LT.(C(I,J)**2I	I -2	42
	1+C(I,J+1)**2)).AND.(V(I,J).GT.V(I,J+1))) Q4=B*(C(I,J)+C(I,J+1))*	I -2	43
	2(ROU(I,J)+ROU(I,J+1))*(V(I,J)-V(I,J+1))/4.	I -2	44
	QUL(I)=Q4	I -2	45
	GO TO 210	I -2	46
209	P4=P(I,J)	I -2	47
	Q4=0.	I -2	48
	IF((AK*(U(I,J)**2+V(I,J)**2).LT.C(I,J)**2).AND.(V(I,J).GT.O. ))Q4=1	I -2	49
	1 2. *B*C(I,J)*ROU(I,J)*V(I,J)	I -2	50
210	UTD(I,J)=U(I,J)-DT*(.5*(P3-P1)+Q3-Q1)/(ROU(I,J)*LX)	I -2	51
	VTD(I,J)=V(I,J)-DT*((S4*(P4-P(I,J))-S2*(P2-P(I,J)))/(2.*VOL)+(Q4	I -2	52
	1 -Q2)/DR)/ROU(I,J)	I -2	53
	IF((UTD(I,J).GE.DX/DT).OR.(VTD(I,J).GE.DR/DT)) GO TO 213	I -2	54
	IF(UTD(I,J)**2+VTD(I,J)**2.LE.O.O).AND.(I.J).LE.ROINF)) GO	I -2	55M
	1 TO 250	I -2	56M
	IF(I.GE.IN) GO TO 251	I -2	57M
211	I=I+1	I -2	58M
	GO TO 202	I -2	59M
250	ISTP(J)=I+3	I -2	60M
	GO TO 252	I -2	61M
251	ISTP(J)=I	I -2	62M
252	IF(J.GE.JN) GO TO 300	I -2	63M
	J=J+1	I -2	64M
	GO TO 200	I -2	65M

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213 DT=DT/2
WRITE(6,6666)DT,ITIME
6666 FORMAT(1H0,22H TIME STEP IS TOO LARGE //1H ,18H NEW TIME STEP DT=
1 F12.6,18H AT TIME CYCLE NO.= 18)
GO TO 199
C ***
C ***
C I-3 THE TILDE INTERNAL ENERGY CALCULATION
C ***
C ***
300 VOL=DX*DR
S1=DR
S2=DX
S4=DX
S2H4=DX
S4PH=DX
S2MH=DX
J=1
301 IF(ITYPE.NE.2) GO TO 302
FJ=J
VOL=2. *PAI*(FJ - .5)*DX*DR**2
S1=2. *PAI*(FJ- .5)*DR**2
S2=2. *PAI*(FJ-1. )*DR*DX
S4=2. *PAI*FJ*DR*DX
S2H4=2. *PAI*(FJ- .5)*DR*DX
S4PH=2. *PAI*(FJ+ .5)*DR*DX
S2MH=0.
IF(J.GT.1) S2MH=2. *PAI*(FJ-1.5)*DR*DX
I=1
302 IF(I.EQ.1.OR.(I.EQ.IN1+1.AND.J.GT.MEMR)) GO TO 304
U1=U(I-1,J)+UTD(I-1,J)
Q1=QUL
GO TO 305
304 U1=-(U(I,J)+UTD(I,J))
Q1=0.
IF((AK*(U(I,J)**2+V(I,J)**2).LT.C(I,J)**2).AND.(U(I,J).LT.0. ))
1 Q1=-2. *B*C(I,J)*ROU(I,J)*U(I,J)
305 IF(J.EQ.1) GO TO 306
V2=V(I,J-1)+VTD(I,J-1)
Q2=QUL(I)
GO TO 307
306 V2=-(V(I,J)+VTD(I,J))
Q2=0.
IF((ITYPE.EQ.1).AND.(AK*(U(I,J)**2+V(I,J)**2).LT.C(I,J)**2).AND.(V
1(I,J).LT.0. )) Q2=-2. *B*C(I,J)*ROU(I,J)*V(I,J)
307 IF(I.EQ.IN1.OR.(I.EQ.IN1.AND.J.GT.MEMR)) GO TO 308
U3=U(I+1,J)+UTD(I+1,J)
Q3=0.
IF((AK*(U(I,J)**2+U(I+1,J)**2+V(I,J)**2+V(I+1,J)**2).LT.(C(I,J)**2
1+C(I+1,J)**2)).AND.(U(I,J).GT.U(I+1,J))) Q3=B*(C(I,J)+C(I+1,J))*
2(ROU(I,J)+ROU(I+1,J))*(U(I,J)-U(I+1,J))/4.
Q3L=Q3
GO TO 309
308 IF(I.EQ.IN) GO TO 350
U3=-(U(I,J)+UTD(I,J))
Q3=0.
IF((AK*(U(I,J)**2+V(I,J)**2).LT.C(I,J)**2).AND.(U(I,J).GT.0. ))

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103=2. *B*C(I,J)*ROU(I,J)*U(I,J)
GO TO 309
350 U3= U(I,J)+UTD(I,J)
Q3= 0.
309 IF(J.EQ.JN) GO TO 310
V4=V(I,J+1)+VTD(I,J+1)
Q4=0.
IF((AK*(U(I,J)**2+U(I,J+1)**2+V(I,J)**2+V(I,J+1)**2).LT.(C(I,J)**2
1+C(I,J+1)**2)).AND.(V(I,J).GT.V(I,J+1))) Q4=B*(C(I,J)+C(I,J+1))*
2 (ROU(I,J)+ROU(I,J+1))*(V(I,J)-V(I,J+1))/4.
Q41(I)=Q4
GO TO 311
310 V4=- (V(I,J)+VTD(I,J))
Q4=0.
IF((AK*(U(I,J)**2+V(I,J)**2).LT.C(I,J)**2).AND.(V(I,J).GT.O. ))Q4=I
1 2. *B*C(I,J)*ROU(I,J)*V(I,J)
311 UTEMP=U(I,J)+UTD(I,J)
VTEMP=V(I,J)+VTD(I,J)
ETD(I,J)=E1(I,J)-DT*(P(I,J)*(S4*(V4+VTEMP)-S2* (V2+VTEMP))/4.
1 +(Q4*(SAPH*V4+S2H4*VTEMP)-Q2*(S2H4*VTEMP+S2MH*V2))/4.
2 -(VTEMP*S2H4*(Q4-Q2)+UTEMP*S1*(Q3-Q1))/2.
3 +S1*(UTEMP+U3)*(P(I,J)+Q3)-(UTEMP+U1)*(P(I,J)+Q1))/4. )
4 /(ROU(I,J)*VOL)
ETD(I,J)=ETD(I,J)+(UTD(I,J)**2+VTD(I,J)**2)/2.
IF(I.GE.ISTP(J)) GO TO 312
I=I+1
GO TO 303
312 IF(J.EQ.JN) GO TO 400
J=J+1
GO TO 301
C ***
C ***
C II-1 CALCULATION OF TRANSPORT EFFECTS
C ***
C ***
400 VOL=DX*DR
S1=DR
S2=DX
S4=DX
J=1
401 IF(ITYPE.NE.2) GO TO 402
FJ=J
VOL=2. *PAI*(FJ - .5)*DX*DR**2
S1=2. *PAI*(FJ- .5)*DR**2
S2=2. *PAI*(FJ-1. )*DR*DX
S4=2. *PAI*FJ*DR*DX
402 I=1
403 IF(I.EQ.1.OR.(I.EQ.IN1+1.AND.J.GT.MEMR)) GO TO 404
U1=UTD(I-1,J)
V1=VTD(I-1,J)
E1=ETD(I-1,J)
L1=LDR1
INT1=0
IF(INTR1.EQ.0) INT1=1
GO TO 405
404 U1=-UTD(I,J)
V1= VTD(I,J)

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	E1= ETD(I,J)	II-1	24
	DM1=0.	II-1	25
	INT1=1	II-1	26
405	IF(J.EQ.1) GO TO 406	II-1	27
	U2=UTD(I,J-1)	II-1	28
	V2=VTD(I,J-1)	II-1	29
	E2=ETD(I,J-1)	II-1	30
	DM2=DMUL(I)	II-1	31
	INT2=0	II-1	32
	IF(DM2.GE.0. ) INT2=1	II-1	33
	GO TO 407	II-1	34
406	U2= UTD(I,J)	II-1	35
	V2=VTD(I,J)	II-1	36
	E2=ETD(I,J)	II-1	37
	DM2=0.	II-1	38
	INT2=0	II-1	39
407	IF(1.EQ.IN.OK.(I.EQ.IN1.AND.J.GT.MEMR)) GO TO 410	II-1	40
	U3=UTD(I+1,J)	II-1	41
	V3=VTD(I+1,J)	II-1	42
	E3=ETD(I+1,J)	II-1	43
	IF(UTD(I,J)+UTD(I+1,J).LT.0. ) GO TO 408	II-1	44
	DM3=ROU(I,J)*(UTD(I,J)+UTD(I+1,J))*S1*DT/2.	II-1	45
	INT3=0	II-1	46
	GO TO 409	II-1	47
408	DM3=ROU(I+1,J)*(UTD(I,J)+UTD(I+1,J))*S1*DT/2.	II-1	48
	INT3=1	II-1	49
409	DMR1=DM3	II-1	50
	INTR1=INT3	II-1	51
	GO TO 411	II-1	52
410	IF(1.EQ.IN) GO TO 460	II-1	53
	U3=-UTD(I,J)	II-1	531
	V3= VTD(I,J)	II-1	54
	E3= ETD(I,J)	II-1	55
	DM3=0.	II-1	56
	INT3=0	II-1	57
	GO TO 411	II-1	571
460	U3=UTD(I,J)	II-1	572
	V3=VTD(I,J)	II-1	573
	E3=ETD(I,J)	II-1	574
	DM3=ROU(I,J)*UTD(I,J)*S1*DT	II-1	575
	INT3=0	II-1	576
	IF(UTD(I,J).LT.0. ) INT3=1	II-1	577
	ETH=ETH-DM3*(ETH(I,J)+(GAM-1. ) *EI(I,J))	II-1	578
	CMTH=CMTH-DM3	II-1	579
411	IF(J.EQ.JN) GO TO 414	II-1	58
	U4=UTD(I,J+1)	II-1	59
	V4=VTD(I,J+1)	II-1	60
	E4=ETD(I,J+1)	II-1	61
	IF(VTD(I,J)+VTD(I,J+1).LT.0. ) GO TO 412	II-1	62
	DM4=ROU(I,J)*(VTD(I,J)+VTD(I,J+1))*S4*DT/2.	II-1	63
	INT4=0	II-1	64
	GO TO 413	II-1	65
412	DM4=ROU(I,J+1)*(VTD(I,J)+VTD(I,J+1))*S4*DT/2.	II-1	66
	INT4=1	II-1	67
413	DMUL(I)=DM4	II-1	68
	GO TO 415	II-1	69
414	U4= UTD(I,J)	II-1	70

V4=-VTD(I,J)	II-1	71
E4= ETD(I,J)	II-1	72
D4=0.	II-1	73
INT4=0	II-1	74
415 ROUNP1=ROU(I,J)+(DM1+DM2-DM3-DM4)/VOL	II-1	75
AIN1=INT1	II-1	76
AIN2=INT2	II-1	77
AIN3=INT3	II-1	78
AIN4=INT4	II-1	79
BIN1=1-INT1	II-1	80
BIN2=1-INT2	II-1	81
BIN3=1-INT3	II-1	82
BIN4=1-INT4	II-1	83
U(I,J)=(AIN1*U1*DM1+AIN2*U2*DM2-AIN3*U3*DM3-AIN4*U4*DM4+UTD(I,J)*	II-1	84
1(ROU(I,J)*VOL+BIN1*DM1+BIN2*DM2-BIN3*DM3-BIN4*DM4))/(VOL*ROUNP1)	II-1	85
V(I,J)=(AIN1*V1*DM1+AIN2*V2*DM2-AIN3*V3*DM3-AIN4*V4*DM4+VTD(I,J)*	II-1	86
1(ROU(I,J)*VOL+BIN1*DM1+BIN2*DM2-BIN3*DM3-BIN4*DM4))/(VOL*ROUNP1)	II-1	87
E(I,J)=(AIN1*E1*DM1+AIN2*E2*DM2-AIN3*E3*DM3-AIN4*E4*DM4+ETD(I,J)*	II-1	88
1(ROU(I,J)*VOL+BIN1*DM1+BIN2*DM2-BIN3*DM3-BIN4*DM4))/(VOL*ROUNP1)	II-1	89
EI(I,J)=E(I,J)-(U(I,J)**2+V(I,J)**2)/2.	II-1	90
ROU(I,J)=ROUNP1	II-1	91
IF(EI(I,J).LT.0. ) GO TO 440	II-1	92
IF(I.GE.ISTP(J)) GO TO 470	II-1	93M
I=I+1	II-1	94
GO TO 403	II-1	95
416 IF(J.EQ.JN) GO TO 500	II-1	96
J=J+1	II-1	97
GO TO 401	II-1	98
440 WRITE(6,450) I,J,ITIME,EI(I,J)	II-1	99
450 FORMAT(1H0,38HINTERNAL ENERGY BECOMES NEGATIVE AT I= 14,3H J= 14,	II-1	100
1 2OH AT TIME CYCLE NO.= 18, 9H EI(I,J)= E15.7)	II-1	101
GO TO 60	II-1	102
470 IF(ISTP(J).GE.IN) GO TO 472	II-1	103M
ISTP1=ISTP(J)	II-1	104M
DO 471 I=ISTP1+1,IN	II-1	105M
E(I,J)=EINF	II-1	106M
471 CONTINUE	II-1	107M
472 GO TO 416	II-1	108M
C ***		
C ***		
C *** II-2 THE TOTAL MASS AND ENERGY CALCULATION	II-2	1
C ***		
C ***		
500 J=1	II-2	2
SIGMM=0.	II-2	3
SIGME=0.	II-2	4
VOL=DR*DX	II-2	5
501 IF(ITYPE.EQ.2 ) VOL=2. *PI*(FLOAT(J)-.5)*DX*DR**2	II-2	6
502 I=1	II-2	7
503 SIGMM=SIGMM+ROU(I,J)*VOL	II-1	8
SIGME=SIGME+ROU(I,J)*VOL*E(I,J)	II-1	9
IF(I.EQ.IN) GO TO 505	II-1	10
504 I=I+1	II-1	11
GO TO 503	II-1	12
505 IF(J.EQ.JN) GO TO 600	II-1	13
J=J+1	II-1	14
GO TO 501	II-1	15

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C *** ***
C *** ***
C *** III AUXILIARY CALCULATION AND DATA PLOT III 1
C *** ***
C *** ***
600 ETOT=SIGME III 2
DE=ETOT-ETH III 3
IF(ABS(DE/ETH).GT.EPSE) GO TO 603 III 4
601 CMTOT=SIGMM III 5
DM=CMTOT-CMTH III 6
IF(ABS(DM/CMTH).GT.EPSM) GO TO 604 III 7
602 ETH=ETOT III 8
CMTH=CMTOT III 9
T=FLOAT(ITIME-ITOFF)*DT III 10
IF(ITIME.GE.IDT) T=TOFF+T III 10A
IF(MOD(ITIME,IPRINT).EQ.0) GO TO 700 III 11M
GO TO 102 III 12
603 WRITE(6,1111) III 13
WRITE(6,1000) III 14
1111 FORMAT(1H0,15HENERGY CHECK) III 15
1000 FORMAT(1H0,19X ,5H ETH=E15.7,6H ETOT=E15.7,4H DE=E15.7) III 16
GO TO 601 III 17
604 WRITE(6,2222) III 18
WRITE(6,1001)CMTH,CMTOT,DM III 19
2222 FORMAT(1H0,15HMASS CHECK) III 20
1001 FORMAT(1H0,4HCMTH=E15.7,6H CMTOT=E15.7,4H DM=E15.7) III 21
GO TO 602 III 22
605 WRITE(6,15) T,ITIME,MEMR,ITYPE III 23
DO 606 J=1,JN III 25
WRITE(6,608) J III 26
WRITE(6,3333) III 24
ISTPPI=ISTPP(J) III 26.1
DO 607 I=1,ISTPPI III 27
ROU=ROU(I,J) III 28
EIM=EI(I,J) III 29
ROU(I,J)=ROU(I,J)/ROINF III 30
EI(I,J)=EI(I,J)/EIMF III 31
IF(MOD(I,2).EQ.0) GO TO 609 III 32
WRITE(6,1003)I, ROU(I,J),U(I,J),V(I,J),EI(I,J) III 33
GO TO 610 III 34
609 WRITE(6,1004)I, ROU(I,J),U(I,J),V(I,J),EI(I,J) III 35
610 ROU(I,J)=ROUM III 36
EI(I,J)=EIM III 37
607 CONTINUE III 38
606 CONTINUE III 39
15 FORMAT(1H1,2HT=F12.6,16H TIME CYCLE NO.=18,6H MEMR=18,5H TYPE18) III 40
3333 FORMAT(1H0,2(60H I ROU(I,J) U(I,J) V(I,J) III 41
1 EI(I,J) )) III 42
608 FORMAT(1H0,2HJ=15) III 43
1003 FORMAT(1H ,18,4X,4F12.6) III 44
1004 FORMAT(1H+,60X,18,4X,4F12.6) III 45
IF(1TAPR.NE.1) GO TO 102 III 45A
WRITE(3) ITIME,IPRINT,CMTH,ETH,DT III 45B
WRITE(3) ((ROU(I,J),U(I,J),V(I,J),EI(I,J),I=1,IN),J=1,JN) III 45C
GO TO 102 III 46
C *** ***
C *** ***

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C	III2 PRINT PAGE CONTROL ROUTINE	III2	1
C ***	***		
C ***	***		
700	J=1	III2	2
701	I=1	III2	3
702	IF(ROU(I,J)-ROINF.LT.1.E-6) GO TO 703	III2	4
	IF(I.GE.IN) GO TO 703	III2	5
	I=I+1	III2	6
	GO TO 702	III2	7
703	ISTPP(J)=1	III2	8
	IF(J.GE.JN) GO TO 605	III2	9
	J=J+1	III2	10
	GO TO 701	III2	11
	END		