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Singularity Fitting in Hydrodynamical Calculations II

by

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SINGULARITY FITTING IN HYDRODYNAMICAL CALCULATIONS II

by

R.D. Richtmyer and R.B. Lazarus

ABSTRACT

This is the second report in a series on the development of techniques for the proper handling of singularities in fluid-dynamical calculations; the first was called *Progress Report on the Shock-Fitting Project*. This report contains six main results: (1) derivation of a free-surface condition, which relates the acceleration of the surface with the gradient of the square of the sound speed just behind it; (2) an accurate method for the early and middle stages of the development of a rarefaction wave, two orders of magnitude more accurate than a simple direct method used for comparison; (3) the similarity theory of the collapsing free surface, where it is shown that there is a two-parameter family of self-similar solutions for $\gamma = 3.9$; (4) the similarity theory for the outgoing shock, which takes into account the entropy increase; (5) a "zooming" method for the study of the asymptotic behavior of solutions of the full initial boundary-value problem; (6) comparison of two methods for determining the similarity parameter δ by zooming, which shows that the second method is preferred.

Future reports in the series will contain discussions of the self-similarity theory for this problem, and for that of the collapsing shock, in more detail and for the full range $(1, \infty)$ of γ ; the values of certain integrals related to neutronic and thermonuclear rates near collapse; and methods for fitting shocks, contact discontinuities, interfaces, and free surfaces in two-dimensional flows.

I. INTRODUCTION

Shock-fitting methods were developed in Los Alamos in 1944 for one-dimensional problems with spherical symmetry, for the special case in which there is just one primary shock, whose position and velocity are known at $t = 0$, and which runs into previously undisturbed material. In spite of the simplifications, the method was sufficiently difficult for the early computers that, when the Hippo project was being planned, in 1948, the pseudo-viscosity method was invented to replace shock fitting.¹ When used with a great deal of care and a certain amount of good luck, the viscosity method can give good results, but is quite risky at best² and is

seriously lacking in spatial resolution in multidimensional problems. It has given quite incorrect results in a few cases. A small project was started here in 1974 to develop shock fitting further in one dimension and to extend it to two dimensions; Ref. 3 is a preliminary report on that work.

In the course of the shock-fitting studies, it became apparent that there are other singularities of flows which also ought to be treated by special methods, which will be called generally *fitting methods*. They include interfaces, contact discontinuities, free surfaces, rarefaction heads and tails, shock interactions, corners, centers of symmetric collapse, and the like. That has led us to the following *working principle* as a basis for study: The finite

difference methods ought to be used only for the smooth parts of the flow, where the differential equations are strictly valid, and all other parts ought to be specially treated by whatever mixture of analytic and numerical methods can be devised.

One advantage of that principle is that it gives one a lot of freedom in the choice of the finite difference method to be used in each of the smooth parts into which the flow is divided by the singularities. (The freedom is made use of in the particular problem to which this report is devoted by the choice of special dependent variables suited to the development of a rarefaction wave.) In particular, the degree of dissipativity of the difference equations can be chosen to satisfy Kreiss's theorem⁴ rather than with any idea of smearing out shocks.

Another advantage of that principle is that it removes the most serious disadvantage of the Eulerian formalism, namely the loss of precise location for material interfaces and other discrete surfaces.

It is recalled that one purpose of the Lax-Wendroff method was to fill a gap in the discussion of the viscosity method in Ref. 1. That discussion showed that the correct description of a flow with shocks is obtainable, at least in principle, by first letting the spacing Δx of the computation net tend to zero and then letting the viscosity coefficient, or equivalently the shock thickness d , tend to zero subsequently. That sidesteps the question of what happens if the limits are taken simultaneously, so that Δx and d remain of the same order of magnitude, as they always are in practical calculations. The same question is unanswered in nearly all the modern versions of the viscosity method.

The Lax-Wendroff method fills that gap in the discussion by conserving mass, momentum, and energy exactly, in a certain sense⁴ already for finite Δx , not merely in the limit as $\Delta x \rightarrow 0$. Since the Rankine-Hugoniot jump conditions for a shock are based on the conservation laws, they also hold, in a sense, for finite Δx .

Another purpose of the Lax-Wendroff method (possibly seen most clearly in retrospect) was to clear up the confusion that existed concerning the dissipativity of a difference scheme and the dissipative terms in such a scheme. Difference schemes are usually analyzed, following von Neumann, by first linearizing, then treating the coefficients as constants (at least in small neighborhoods), and then expanding the solution in a Fourier series or Fourier integral. The time dependence of the Fourier coefficients is then determined by the difference equations. If the absolute value of every Fourier coefficient remains constant in time, as is the case for the differential equations when similarly treated,

the scheme is called *nondissipative*. For any reasonable difference scheme, that must be approximately true for the long and medium wavelength components, but the short wavelength ones are often significantly damped, as t increases, in which case the scheme is called *dissipative*.

It was formerly felt that difference methods for fluid dynamics ought to be nondissipative, because the differential equations are. However, it can be shown that any finite difference scheme necessarily falsifies the phases of the short-wave components; hence, it is pointless to maintain their amplitudes, from the viewpoint of accuracy. That it is also pointless to maintain their amplitudes from the viewpoint of the conservation laws is shown by the Lax-Wendroff equations, which conserve mass, momentum, and energy exactly, but are dissipative. Lastly, Kreiss's theorem shows that a suitable degree of dissipativity, corresponding to a given degree of accuracy, guarantees stability against variability of the coefficients, when the von Neumann condition is satisfied.

When, as in the present work, difference equations are used only for the smooth part of the flow, the two main functions of the Lax-Wendroff method, conservation and dissipation, can be separated. We are interested only in the latter, hence need not require the equations to be in conservation-law form, but can apply the two-step Lax-Wendroff method of differencing directly to equations of the form

$$\frac{\partial U}{\partial t} + A \frac{\partial U}{\partial x} = 0,$$

for it is known that such differencing gives the amount of dissipation required by Kreiss's theorem.

II. THE RAREFACTION (CAVITY COLLAPSE) PROBLEM

At time $t = 0$, a γ -law fluid is at rest under constant pressure in the region of space outside an empty spherical cavity, i.e., in the region $R > R_0$, where R is the Eulerian radial coordinate. See Fig. 2-1. (The computer code was originally planned to handle also the corresponding plane and cylindrical problems, but so far only the spherical one has been studied.) For $0 < t < t_0$, where t_0 is the instant of collapse of the cavity, there is a rarefaction wave between an inward moving free surface at $R = \xi(t)$ and an outward moving head at $R = \eta(t)$, where $0 < \xi(t) < R_0$ and where $\eta(t) = R_0 + c_0 t$, c_0 being the sound speed in the initial state. For $t \ll t_0$, the spherical shell $\xi(t) < R < \eta(t)$ occupied by the wave is very thin, and the rarefaction is approximately a plane simple wave⁵ in which the sound speed c and the material speed u

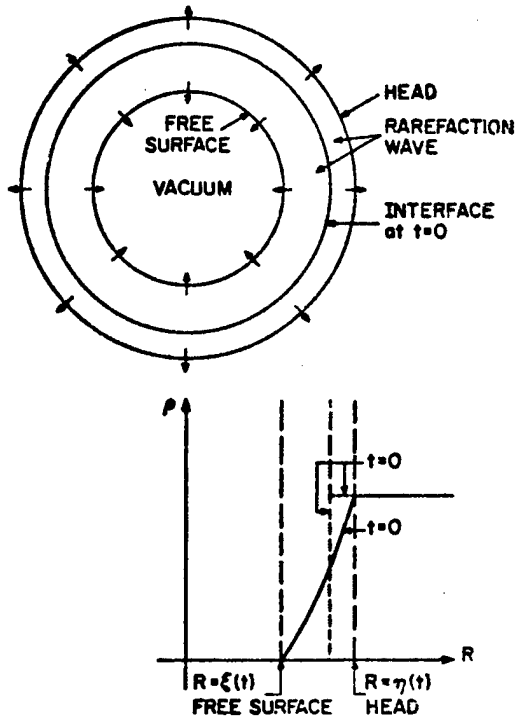


Fig. 2-1.

The cavity collapse program at, and shortly after, $t = 0$.

vary linearly with R across the shell. In that stage, the velocity of the free surface is given approximately by $\dot{\xi} (= d\xi/dt) = -2c_0/(\gamma-1)$. At later times, the free surface accelerates inward until collapse occurs at $t = t_0$ [$\xi(t_0) = 0$], at which time the pressure and density are instantaneously infinite at $R = 0$. For $t > t_0$, a shock front, whose position is denoted by $R = \zeta(t)$, progresses outward from the center.

The flow has singularities as follows:

(1) $t \approx 0$, $R \approx R_0$. At $t = 0$, the flow quantities are discontinuous across $R = R_0$. At early times thereafter, there is a thin centered rarefaction wave in an interval $[\xi, \eta]$ around R_0 . Although the theory of such waves has been well known in the plane approximation since the time of Riemann, the wave has to be regarded as a singularity from the viewpoint of the finite difference equations, so long as its thickness is comparable with or less than the spacing ΔR of the computational net.

(2) $t > 0$, $R = \eta(t) = R_0 + c_0 t$. At the head of the rarefaction wave, the first derivatives of the flow quantities are discontinuous. A special fitting technique is used; it is described in Sec. V.

(3) $0 < t < t_0$, $R = \xi(t)$. At the free surface, a special boundary condition is needed in addition to fitting. It is described in Sec. IV.

(4) $t \approx t_0$, but $t < t_0$, $R \approx 0$. The collapsing flow may be representable asymptotically by a self-similar flow of the kind described by Hunter⁶ — see also Ref. 7. This question is debated in Sec. X:

(5) $t \approx t_0$, but $t > t_0$, $R \approx 0$. The outgoing shock may also be representable asymptotically by a self-similar flow.

(6) $t > t_0$, $R = \zeta(t)$. Shock fitting, as described in Ref. 3, is required at the outgoing shock.

The main computational problem, to which the latter part of this report is addressed, is how to turn the flow around, at $t \approx t_0$, and get the outgoing shock properly started.

III. DIFFERENCE EQUATIONS FOR THE SMOOTH PART OF THE FLOW

At early times, owing to the approximate plane simple-wave character of the solution, it is natural to take the entropy S per unit mass, the Riemann variable $\rho [= 2c/(\gamma-1)$ for a γ -law gas], and the fluid velocity u , as the dependent variables, because those quantities vary linearly with R in the simple wave, whereas the pressure p and the density ρ vary with a higher power of the distance from the free surface (see Note following Eq. 3.8). The differential equations for S , ρ , and u will next be derived from the Eulerian equations, which, for plane ($\alpha=1$), cylindrical ($\alpha=2$), and spherical ($\alpha=3$) symmetry, are

$$\left. \begin{aligned} \left(\frac{\partial}{\partial t} + u \frac{\partial}{\partial R} \right) \rho &= -\rho \frac{1}{R^{\alpha-1}} \frac{\partial}{\partial R} \left(R^{\alpha-1} u \right), & (a) \\ \left(\frac{\partial}{\partial t} + u \frac{\partial}{\partial R} \right) u &= -\frac{\partial p}{\partial R}, & (b) \\ \left(\frac{\partial}{\partial t} + u \frac{\partial}{\partial R} \right) E &= -p \frac{1}{R^{\alpha-1}} \frac{\partial}{\partial R} \left(R^{\alpha-1} u \right) & (c) \end{aligned} \right\} (3.1)$$

\mathcal{E} is the internal energy per unit mass and is related to p and ρ by an equation of state. A consequence of Eq. 3.1 is the entropy condition

$$\left(\frac{\partial}{\partial t} + u \frac{\partial}{\partial R} \right) S = 0$$

If \tilde{S} is any function of S (not depending on any other variables), then

$$\left(\frac{\partial}{\partial t} + u \frac{\partial}{\partial R}\right) \xi = 0. \quad (3.2)$$

We introduce as further thermodynamic quantity the Reimann variable.

$$\sigma = \int \frac{dp}{\rho c} \Big|_{S=\text{const}} = \sigma(\tilde{S}, \rho). \quad (3.3)$$

For a γ -law gas, we can take

$$p = p(\tilde{S}, \rho) = \tilde{S} \rho^\gamma; \quad (3.4)$$

then

$$\sigma = \sigma(\tilde{S}, \rho) = 2c/(\gamma-1),$$

where

$$c = c(\tilde{S}, \rho) = \sqrt{\gamma \tilde{S} \rho^{\gamma-1}}. \quad (3.5)$$

The program is to take \tilde{S} , σ , and u as dependent variables, define a vector

$$U = (\tilde{S}, \rho, u), \quad (3.6)$$

and derive the differential equations for \tilde{S} , ρ , u in the form

$$\partial U / \partial t + A \partial U / \partial R = g, \quad (3.7)$$

where A is a 3×3 matrix. The first equation is already at hand: it is (3.2). A short calculation gives the system

$$\left. \begin{aligned} \left(\frac{\partial}{\partial t} + u \frac{\partial}{\partial R}\right) \xi &= 0 & (a) \\ \left(\frac{\partial}{\partial t} + u \frac{\partial}{\partial R}\right) \sigma + c \frac{\partial u}{\partial R} &= -uc \frac{\alpha-1}{R} & (b) \\ \left(\frac{\partial}{\partial t} + u \frac{\partial}{\partial R}\right) u + c \frac{\partial \sigma}{\partial R} + \left(\frac{1}{\rho} \frac{\partial p}{\partial \xi} - c \frac{\partial \sigma}{\partial \xi}\right) \frac{\partial \xi}{\partial R} & & (c) \end{aligned} \right\} = 0. \quad (3.8)$$

In the coefficient of $\partial \tilde{S} / \partial R$ in Eq. (3.8c), the differentiations with respect to \tilde{S} are understood to be for fixed ρ .

Note: For a γ -law fluid, either of the variables σ and c can be eliminated by use of the equation $\sigma = 2c/(\gamma-1)$. For a fluid with a non- γ -law equation of state, neither σ nor c varies exactly linearly with distance across a plane simple rarefaction wave, but σ

varies more nearly linearly than c , hence Eq. (3.8) is the preferred form of the equations, in which c may be thought of as a function of σ and \tilde{S} .

As stated in the Introduction, we wish to use dissipative difference equations of second-order accuracy. An easy way of obtaining them is the Lax-Wendroff two-step method of differencing (Ref. 4, pp. 300-306). For equations of the form (3.7) we have, in the usual notation,

Step 1:

$$U_{j+\frac{1}{2}}^{n+\frac{1}{2}} = \frac{1}{2} (U_{j+1}^n + U_j^n) - \frac{\Delta t}{2\Delta R} \bar{A}^n (U_{j+1}^n - U_j^n) + \Delta t \bar{g}^n, \quad (3.9a)$$

Step 2:

$$U_j^{n+1} = U_j^n - \frac{\Delta t}{\Delta R} \bar{A}^{n+\frac{1}{2}} (U_{j+\frac{1}{2}}^{n+\frac{1}{2}} - U_{j-\frac{1}{2}}^{n+\frac{1}{2}}) + \Delta t \bar{g}^{n+\frac{1}{2}}, \quad (3.9b)$$

where the overbars denote the appropriate spatial averages of the matrix A and the vector g . In the code, the following minor modification of these equations is used: In the system (3.8), g appears only in the second equation and can be included by rewriting the last term of the first member of that equation (for $\alpha = 3$) as

$$\frac{c}{R^2} \frac{\partial (R^2 u)}{\partial R} = 3c \frac{\partial (R^2 u)}{\partial (R^3)}.$$

This term is differenced as

$$3c \frac{R_{j+\frac{1}{2}}^2 u_{j+\frac{1}{2}}^{n+\frac{1}{2}} - R_{j-\frac{1}{2}}^2 u_{j-\frac{1}{2}}^{n+\frac{1}{2}}}{R_{j+\frac{1}{2}}^3 - R_{j-\frac{1}{2}}^3}$$

in step 2, and in a similar way in step 1. Special treatment of points near the free surface, near the head of the rarefaction wave, and (after collapse) near the center and near the shock front is described in the following sections.

IV. THE FREE-SURFACE CONDITION

It is assumed that $\sigma(R,t)$ and $u(R,t)$ are smooth, for $R \geq \xi(t)$, so that the entropy law $(\partial/\partial t + u\partial/\partial R)\tilde{S} = 0$ holds near the free surface. Then, for problems in which \tilde{S} is constant initially, \tilde{S} is constant throughout the flow until shocks form. (The present discussion needs modification for problems in which $\tilde{S} \neq$ constant initially.) We set

$$\sigma(R,t) = \sum_{p=0}^{\infty} \sigma_p(t) [R - \xi(t)]^{\alpha+p}, \quad (4.1)$$

$$u(R,t) = \dot{\xi}(t) + \sum_{p=0}^{\infty} u_p(t) [R - \xi(t)]^{\beta+p}, \quad (4.2)$$

where α and β are positive constants to be determined later. If we substitute (4.1) and (4.2) into (3.8b) and (3.8c), with $\partial\tilde{S}/\partial R = 0$, then, writing only the lowest order terms and replacing the others by dots, we have

$$\begin{aligned} & \ddot{\sigma}_0 (R - \xi)^\alpha + \dots \\ & + u_0 \sigma_0 \left(\alpha + \frac{\gamma-1}{2}\beta \right) (R - \xi)^{\beta+\alpha-1} + \dots = \dots \end{aligned} \quad (4.3)$$

$$\begin{aligned} & \ddot{\xi} + \dots + \dot{u}_0 (R - \xi)^\beta + \dots + u_0^2 \beta (R - \xi)^{2\beta-1} + \dots \\ & + \frac{\gamma-1}{2} \sigma_0^2 \alpha (R - \xi)^{2\alpha-1} + \dots = 0 \end{aligned} \quad (4.4)$$

In order to achieve cancellation of the lowest order terms in (4.3), β must be =1; then, the second and third terms indicated in (4.4) are of higher order, and to achieve cancellation of the first and fourth terms, α must be =1/2, and we find

$$\ddot{\xi} + \frac{\gamma-1}{4} \sigma_0^2 = 0.$$

With $\alpha = 1/2$, (4.1) shows that $\sigma^2 [= \sigma(R,t)^2]$ is a power series in $R - \xi$, starting with $\sigma_0^2 (R - \xi)$; hence,

$$\ddot{\xi} = - \frac{\gamma-1}{4} \left. \frac{\partial(\sigma^2)}{\partial R} \right|_{R=\xi}. \quad (4.5)$$

That is the boundary condition at the free surface; it connects the inward acceleration of the free surface with the rate at which $\sigma^2 \rightarrow 0$ as the free surface is approached. In the code, $\partial(\sigma^2)/\partial R$ is approximated by two difference quotients containing values of σ^2 at ξ (where $\sigma^2 = 0$) and at nearby net points and is

then extrapolated back to $R = \xi$. Equation (4.5) is implied by Eq. (3.6) of Ref. 7.

When σ has been calculated at all regular net points at time $t = t^{n+1}$ by the method of Sec. III, Eq. (4.5), for $t = t^{n+1}$, contains two unknowns ξ^{n+1} and $\dot{\xi}^{n+1}$, because ξ^{n+1} enters into the approximation to $\partial(\sigma^2)/\partial R$. That equation is solved, together with the two additional equations

$$\begin{aligned} \xi^{n+1} &= \xi^n + \Delta t/2(\dot{\xi}^n + \dot{\xi}^{n+1}) \\ & \quad \quad \quad (4.6) \end{aligned}$$

$$\dot{\xi}^{n+1} = \dot{\xi}^{n+1} = \dot{\xi}^n + \Delta t/2(\ddot{\xi}^n + \ddot{\xi}^{n+1})$$

for the unknowns ξ^{n+1} , $\dot{\xi}^{n+1}$, and $\ddot{\xi}^{n+1}$ by Newton's iterative procedure. (Only one or two iterations are required at each time step.)

It is easily verified that the boundary condition (4.5) at the free surface is satisfied by the self-similar solutions of the flow equations discussed in Sec. VII, by virtue of the first equation of the pair (7.9).

The free-surface boundary condition on the variable u has already been incorporated in (4.2); it is

$$u(\xi,t) = \dot{\xi}(t). \quad (4.7)$$

V. FITTING THE HEAD OF THE WAVE AND THE FREE SURFACE

The singularity at $R = \eta(t)$ is of a very mild kind, and the only boundary condition or joining condition needed for the differential equations is the continuity of the function values. Rather little harm is done if this singularity is ignored completely in the calculation; the main effect of so doing is loss of the second-order accuracy of the finite difference equations for that interval that contains the front, at which the second derivatives are infinite. However, it is very easy to treat the singularity correctly, and to do so costs almost nothing computationally — it costs less than nothing in the present problem because it obviates the necessity of any computations whatever for $R > \eta(t)$.

A portion of the computational net near the rarefaction head is shown in Fig. 5-1. Whenever the path of the rarefaction head is in either of the positions indicated as (a) and (b) in the figure, that is, lies to the right of a centered point like x_2 but cuts through the net rectangle that contains that point, values of the flow quantities at x_2 are obtained by the following special procedure in order that they may then be used in the regular Lax-Wendroff step 2 to yield the values at the point indicated by the circle on the line $n+1$: A special Lax-Wendroff step 1 is

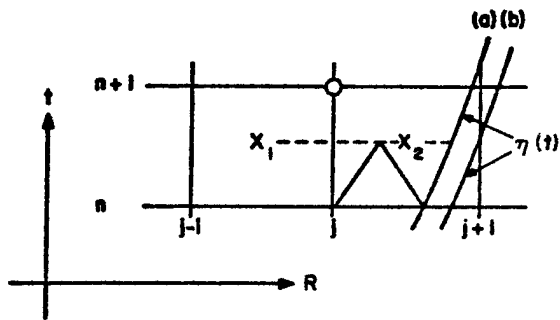


Fig. 5-1.

Treatment of the (mild) singularity at the head of the rarefaction.

performed using the triangle indicated; it is the same as the normal Lax-Wendroff step 1 except that the base of the triangle has been reduced from ΔR to $\eta(t^n) - R_j$. This step gives values of the flow quantities at the tip of the triangle; they are then obtained at x_2 by quadratic interpolation on R using also the known values at x_1 and at the rarefaction head itself.

If the head lies in a position like (b), new values of the flow quantities are needed also at the net point $j+1$ that has just been uncovered by the motion of the head; those values are obtained by linear or quadratic interpolation on R at time t^{n+1} .

The procedure for determining the flow quantities near the free surface is similar, once the motion of the free surface is known. The sequence is: The modified Lax-Wendroff step 1 uses the old position $\xi(t^n)$ of the free surface; the interpolation at $t^{n+1/2}$ uses an extrapolated value of $\xi(t^{n+1/2})$; the Lax-Wendroff step 2 can then be performed at all ordinary net points, and therefrom the final position $\xi(t^{n+1})$ of the free surface is determined as described in the preceding section, followed by interpolation of the flow quantities, at t^{n+1} , if needed for an uncovered net point.

VI. SOME NUMERICAL RESULTS FOR THE RAREFACTION PHASE PRIOR TO COLLAPSE

Several calculations were made, for $\gamma = 3.0$, using the methods described in the three preceding sections. Curves of c vs R at four values of t are given, for a calculation in which ΔR was $= 0.02$, in Fig. 6-1. Except where shown, the calculated points lie on the curves drawn within the accuracy of the drawing. The steepening of the gradient at the free surface, which causes the inward acceleration, is clearly visible.

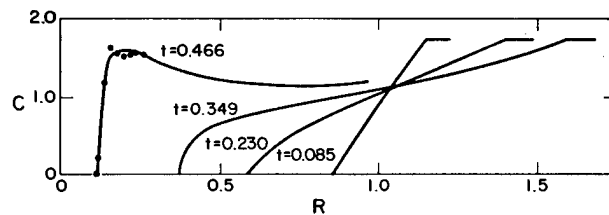


Fig. 6-1.

Progress of the cavity collapse wave, for $\gamma = 3$. Except for those shown, calculated points are within the curves as drawn.

As a provisional measure of overall accuracy, the total energy, kinetic plus internal, of the fluid between $R = \xi(t)$ and $R = \eta(t)$ was computed and compared with the initial energy $[4\pi/3(\eta^3 - \xi^3)\rho_0/\gamma - 1]$ of that same fluid. The percentage errors are given in Table I (together with the corresponding errors for a calculation with $\Delta R = 0.0025$). A comparison calculation was made with the standard Eulerian equations, with no fitting or boundary condition on the free surface, but with the cavity initially filled, as is often done, with fluid at a very low density and pressure ($= 2 \times 10^{-4}$ and 8×10^{-12} , in units of the initial density and pressure outside the cavity). Comparison of the second and third columns of the table shows that, by this particular measure of the accuracy, the errors are reduced by a factor 50-200 by the methods described in the preceding sections. Most of the improvement results from the proper treatment of the free surface; the special choice of dependent variables given in Sec. III gives only an additional improvement by a factor 2-3.

VII. SOME SELF-SIMILAR SOLUTIONS OF THE BOUNDARY-VALUE PROBLEM

If the initial conditions at $t = 0$ are ignored, one has a boundary-value problem consisting of the partial differential equations (3.8) for $R > \xi(t)$ together with the free-surface condition at $R = \xi(t)$. For a γ -law fluid, for which σ and c differ merely by a numerical constant, that problem contains no characteristic length or time, hence is likely to have self-similar solutions, i.e., solutions such that the functions of R obtained at any distinct instants t_1 and t_2 become identical if all lengths, velocities, and other quantities are suitably rescaled between the two instants. (Such solutions can be obtained by solving ordinary differential equations, as shown below.) In the literature, one often sees arguments claiming to show that the solution of the full initial boundary-value problem, though not self-similar

TABLE I
 ERROR OF TOTAL ENERGY OF RAREFACTION FAN
 $\gamma = 3.0$

Time	Comparison Calculation: Eulerian Variables, No Free-Surface Fitting		Riemannian Variables, Free-Surface Fitting	
	$\Delta R = 0.02$	$\Delta R = 0.02$	$\Delta R = 0.02$	$\Delta R = 0.0025$
t = 0.0846			0.22%	
0.1401	6.99%			
0.1646			0.13%	
0.1993	7.29%			
0.2302				0.016%
0.2399			0.10%	
0.3097			0.075%	
0.3104	6.93%			
0.3495				0.008%
0.3726			0.64%	
0.4254			0.39%	
0.4550				0.018%
0.4657			0.033%	
0.4923			0.035%	
0.5009	6.00%			

(the initial conditions introduce a characteristic length R_0), is asymptotically self-similar near collapse, i.e., for $t_0 - t \ll t_0$ and $R \ll R_0$. Although we feel there is reason to doubt the general validity of those arguments, we have investigated the self-similar solutions of the boundary-value problem, and we have attempted to examine numerically the solution of the full problem asymptotically near collapse to see whether it has a similarity property. The results so far are inconclusive, but further work is planned. The goal of such work is to be able to patch together numerical calculations just before and just after collapse by solutions obtained from the similarity theory.

For application to the cavity-collapse problem, the class of self-similar solutions is further restricted by the assumption of isentropy, which introduces a characteristic constant entropy S_0 . (That is to be contrasted with the assumption made in the Guderley-Butler shock-collapse problem, in which there is a characteristic constant density ρ_0 — the density of the stationary fluid into which the collapsing shock is running. The two sets of assumptions are contrasted in Table II — they lead to different systems of ordinary differential equations.)

The similarity assumptions are these: First, there are positive constants A and δ such that the position of the free surface is given by

$$\xi(t) = A(t_0 - t)^\delta \quad (7.1)$$

In the discussion of the self-similar solutions, t_0 could be taken $= 0$, but we retain the notation of the preceding sections in the interest of returning later to the full initial boundary-value problem.) Then, a dimensionless variable η is defined:

$$\eta = \frac{R}{A(t_0 - t)^\delta} \quad (7.2)$$

the region of interest is $\eta \geq 1$; $\eta = 1$ corresponds to the free surface. It is then assumed that the dependent variables depend on R and t only through η , after suitable scaling.

The dependent variables could be chosen as σ and u , or as c and u , since, for a γ -law fluid, σ and c differ merely by a constant. We make the latter choice, in conformity with most of the literature. Since c and u have the dimensions of velocity, they can be written as

$$c(R, T) = RC(\eta)/(T_0 - t) = A(t_0 - t)^{\delta-1} \eta C(\eta) \quad (7.3)$$

$$u(R, t) = RV(\eta)/(t_0 - t) = A(t_0 - t)^{\delta-1} \eta V(\eta),$$

where C and V are dimensionless functions, called the *reduced* sound and material speeds. When the expressions (7.3) are substituted into (3.8), with $\tilde{S} = \text{const}$, the explicit dependence on R and t cancels out, and the following system of ordinary differential equations is obtained:

TABLE II
SIMILARITY ASSUMPTIONS

$$\eta = \frac{R}{A(t_0 - t)^\delta}$$

Collapsing Shock
(Guderley, Butler, etc.)

$$c = A(t_0 - t)^{\delta-1} C(\eta)$$

$$u = A(t_0 - t)^{\delta-1} V(\eta)$$

$$\rho = P(\eta) \left[\text{e.g., } P(1) = \frac{\gamma+1}{\gamma-1} \rho_0 \right]$$

Consequence:

$$s = A^2 (t_0 - t)^{2\delta-2} S(\eta)$$

Collapsing Cavity
(Hunter, Clarke, etc.)

$$c = A(t_0 - t)^{\delta-1} C(\eta)$$

$$u = A(t_0 - t)^{\delta-1} V(\eta)$$

$$s = S(\eta) \quad (\text{In fact, } s \text{ is independent of both } R \text{ and } t \text{ until collapse.})$$

Consequence:

$$\rho = A^{\frac{2}{\gamma-1}} (t_0 - t)^{\frac{2\delta-1}{\gamma-1}} P(\eta)$$

$$\eta C' + C$$

$$= \frac{(\delta-1) \left\{ (V+\delta)C - \frac{\gamma-1}{2} CV \right\} - (\gamma-1)(V+\delta)CV}{\Delta}$$

$$\eta V' + V = \frac{(\delta-1) \left\{ (V+\delta)V - \frac{2}{\gamma-1} C^2 \right\} + 2C^2V}{\Delta}$$

(7.4)

where

$$\Delta = \Delta(C, V) = (V+\delta)^2 - C^2. \quad (7.5)$$

It is noted that the quantity of $dC/dV = C'/V'$ is a function of C and V and does not depend explicitly on η . An ordinary differential equation system of the general form (7.4) having that property is called *autonomous*; it has the advantage for interpretation and presentation of the solutions that a solution curve in the three-dimensional space C, V, η is uniquely determined by its projection on the C, V plane. (Even the explicit appearance of η in the left members of (7.4) can be eliminated by taking $\tau = \log \eta$ as the independent variable.)

It should be noted that the autonomous property of (7.4) follows from the particular substitution (7.3) and does not follow, for example, from the dimen-

sionally equivalent substitution $c(R, t) = \xi C_1(\eta)$, $u(R, t) = \xi V_1(\eta)$. The possibility of putting the equations into autonomous form is a consequence of the dimensional properties of the fluid-dynamical equations and can be seen as follows: First, we take $A=1$ and $t_0=0$ in this paragraph, so that $\eta = R(-t)^{-\delta}$. We write

$$\frac{C'(\eta)}{V'(\eta)} = \Psi(C(\eta), V(\eta), \eta). \quad (7.6)$$

Let a be any positive constant, and let $b = a^{1-\delta}$. Now,

$$\frac{C'(b\eta)}{V'(b\eta)} = \Psi(C(b\eta), V(b\eta), b\eta). \quad (7.7)$$

On the other hand, for $\tilde{S} = \text{const}$, it is seen that if $c(R, t)$ and $u(R, t)$ satisfy (3.8), then the functions

$$\hat{c}(R, t) = c(aR, at)$$

$$\hat{u}(R, t) = u(aR, at)$$

also satisfy (3.8), hence the functions

$$\hat{C}(\eta) = C\left(\frac{aR}{(-at)^\delta}\right) = C(b\eta)$$

$$\hat{V}(\eta) = V \left(\frac{aR}{(-at)^\delta} \right) = V(b\eta)$$

also satisfy (7.6); that is,

$$\frac{C'(b\eta)}{V'(b\eta)} = \Psi(C(b\eta), V(b\eta), \eta).$$

Since b is arbitrary (if $\delta \neq 1$), comparison with (7.7) shows that Ψ does not depend explicitly on η . (The argument must be modified for $\delta = 1$.)

Equation (7.4) can be written as

$$\eta C'(\eta) = \frac{F(C, V)}{\Delta(C, V)} \quad (7.8)$$

$$\eta V'(\eta) = \frac{G(C, V)}{\Delta(C, V)}$$

where F and G , like Δ , are simply polynomials in C and V . The boundary conditions at the free surface ($\eta = 1$) are

$$C(1) = 0 \text{ (from vanishing pressure),}$$

$$V(1) = -\delta \text{ (from } u = \xi \text{ at } R = \xi). \quad (7.9)$$

The boundary conditions for $\eta \rightarrow \infty$ are discussed below.

In principle, the system (7.8) can be integrated by Runge-Kutta with (7.9) as starting values. However, there is a difficulty that the point with C and V given by (7.9) is a critical point, i.e., a point where the polynomials F , G , and Δ all vanish. A standard analysis shows that, near the critical point, V has the form of a power series in $\eta - 1$, while C has the form $\sqrt{\eta - 1}$ times a power series. For that and other reasons, the numerical work was done with functions $q(\eta)$ and $v(\eta)$ given by

$$q(\eta) = \frac{2}{\gamma - 1} \frac{\eta}{\sqrt{\eta - 1}} C(\eta) \quad (7.10)$$

$$v(\eta) = \eta V(\eta)$$

(then, the system no longer has the autonomous form), but the results will be described in the C, V plane. The complete set of critical points is described below and in Fig. 7-1.

The boundary conditions for $\eta \rightarrow \infty$ come from the requirement that $c(R, t)$ and $u(R, t)$ have definite limiting values, as $t \rightarrow t_0$ for fixed R , i.e., as $\eta \rightarrow \infty$.

From the definition (7.2) of η , it is seen that (7.3) can be rewritten as

$$c(R, t) = \left(\frac{R}{(t_0 - t)^\delta} \right)^{1/\delta} R^{1-1/\delta} C(\eta)$$

$$= (A\eta)^{1/\delta} R^{1-1/\delta} C(\eta)$$

$$u(R, t) = \dots$$

$$= (A\eta)^{1/\delta} R^{1-1/\delta} V(\eta)$$

Hence, as $\eta \rightarrow \infty$,

$$C(\eta), V(\eta) \sim \eta^{-1/\delta}. \quad (7.11)$$

Examination of (7.4) shows that if C and $V \rightarrow 0$, as $\eta \rightarrow \infty$, then they behave asymptotically according to (7.11); hence the boundary condition is that C and $V \rightarrow 0$.

For most values of γ and δ in the relevant ranges, the system (7.4) has nine critical points, of which six are shown in Fig. 7-1 for $\gamma = 3$, $\delta = 0.7$ and the remaining three are obtained by reflection in the V axis. For the general classification of critical points, the reader is referred to Birkhoff and Rota,⁸ p. 130 ff. A little algebra shows that, in the present problem, if either F or G vanishes along with Δ , then the other vanishes too. That simplifies locating and classifying the critical points.

Factorization of the denominator $\Delta(C, V)$, (7.5), of the expressions (7.8) shows that, along any solution curve in the C, V plane, the direction of change of η reverses, as the curve crosses either of the lines $V + \delta = \pm C$, one of which is shown dashed in Fig. 7-1, unless F and G also vanish at the same time. According

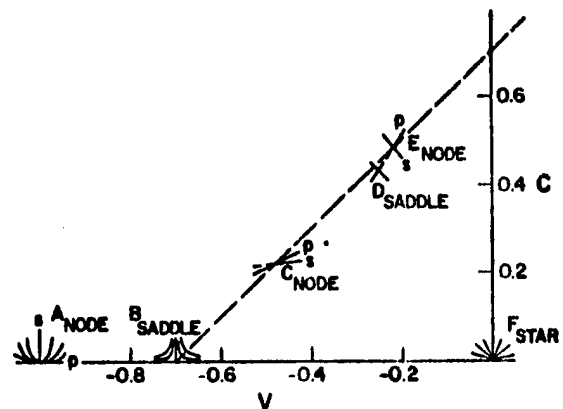


Fig. 7-1.

Critical points of $C = C(V)$ in the upper half-plane, for $\gamma = 3$.

to the boundary condition given above, the solution curve starts from the saddle point B in an upward direction and terminates, for $\eta \rightarrow \infty$, at the star point F; hence, in order that C and V be single-valued functions of η , the curve must cross the dashed line either at the node E or the node C. Numerical integration with third-order Runge-Kutta for various values of δ shows that, when a complete solution is obtained at all, it goes through point C, not E.*

The situation is in marked contrast with that for the corresponding collapsing shock problem studied by Guderley and Butler, where the point corresponding to C is a saddle point, instead of a node. (Qualification is necessary, because our studies have been made for $\gamma = 3.0$, while those of Guderley and Butler were for $1 < \gamma < 2$; detailed studies of both problems for general γ will be presented elsewhere.) In the case of a node, all solution curves that come reasonably close are drawn into the node, and all, with one exception, pass through it in the direction indicated by the letter "p" (for "primary") in Fig. 7-1; the one exception takes the direction "s" (for "secondary"); whereas, in the case of a saddle point, only the curve that is aimed in precisely the right way passes through the point at all; all other solutions deviate either to the right or to the left. Consequently, in order to get a solution at all, for the collapsing shock problem, the value of δ has to be precisely chosen, whereas there is range of admissible δ 's for the collapsing cavity problem. In either case, once the dashed line has been crossed, the solution is drawn automatically to the star point F at the origin, and the terminal boundary condition is satisfied.

The results described from now on are all for $\gamma = 3.0$.

For δ in the interval $0.666667 < \delta < 0.708542$, the solution curve arrives at the node C along the primary direction, but there are many curves leaving C to the right: one starting in the secondary direction, as seen most clearly in Fig. 7-2, for $\delta = 0.68$, and a one-parameter family of curves starting in the primary direction, as illustrated by the upper three curves of Fig. 7-3, for $\delta = 0.70$. (The arrows indicate the direction in which the curves were calculated; in particular, the lowest one was calculated both

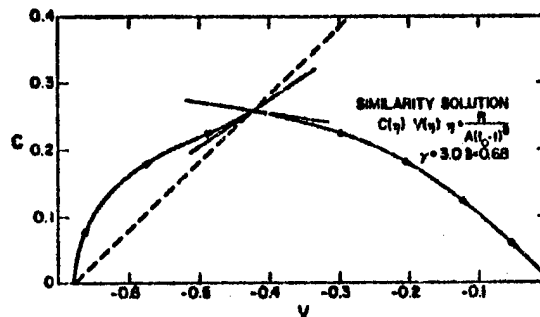


Fig. 7-2.

A similarity solution for $\delta = 0.68$, $\gamma = 3$.

forward and backward, by reversing the sign of $\Delta\eta$ in the Runge-Kutta method; the direction of increase of η is from left to right along all the curves.) The curves of the one-parameter family (for given δ) pass through C with change of direction, but generally without discontinuity of higher derivatives.

In the special case $\delta = 0.708542$, the curve comes to C along the secondary direction and, if continued so as to leave in the same direction, passes through C with continuity of all derivatives and in fact analytically. For certain special values δ in (0.666667, 0.708542), one curve of the one-parameter family referred to is also analytic at C. That appears to be a rather intricate Diophantine affair and is discussed in detail by Hunter,⁶ who, however, rejects those solutions for reasons that are rather hard to follow.

An argument given by Hunter claims to show that only those solutions analytic at C can appear asymptotically in a physical problem. That argument seems doubtful to us for the following reason: In the first place, any of the curves in the C, V plane discussed above, including those of the two-parameter family (where δ is now regarded as the second parameter) lead to an acceptable solution of the partial differential equations of fluid dynamics, by

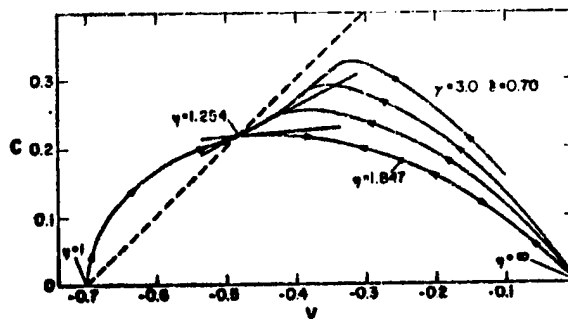


Fig. 7-3.

Some of the similarity solutions for $\delta = 0.70$, $\gamma = 3$.

*At least for $\gamma = 3$. For the collapsing shock problem (to be published), the situation reverses at a critical value of γ (near 1.9, but different for cylindrical and spherical geometry). It remains to be investigated whether the same is true for the collapsing cavity problem, and, if so, whether the critical values of γ are the same.

means of Eqs. (7.2) and (7.3). Since the fluid-dynamical equations are hyperbolic, there is no requirement of analyticity; in fact, jumps of the various derivatives of the flow quantities can be propagated along the characteristics; the first derivatives are discontinuous at the head of a rarefaction wave, and the functions themselves can be discontinuous for a weak solution. Now, the solution of the full initial boundary-value problem of the cavity-collapse problem is indeed analytic; it can be shown (again by consideration of the characteristics), but it does not follow that the self-similar solution to which it converges must be analytic.

If η_0 is the value of η at the node C, then the curve $\eta = \eta_0$ in the R,t plane, namely the curve

$$R = A\eta_0(T_0 - t)^\delta,$$

is characteristic; it is the path of an incoming spherical sound wave that just catches up with the free surface at collapse. For determining how the free surface itself collapses, nothing outside that sound wave has any influence, hence the part of the solution curve to the right of the dashed line in the figures has no effect, until after collapse.

The properties of the critical point C as a function of δ in (0,1) (still for $\gamma = 3.0$) can now be summarized. (It always lies on the dashed line in the figures.) For $0.708542 < 0.711405$, the solution curve comes to C along the primary direction from above and to the right and crosses the dashed line first, hence is unacceptable. For $\delta > 0.711405$, C is a spiral point, hence again the solution curve crosses the dashed line (in fact infinitely often) before arriving at C. For $\delta < 0.666667$, the point C disappears (it merges with F, then becomes complex, as δ decreases). Hence, we are left with the interval (0.666667, 0.708542).

VIII. THE SIMILARITY SOLUTION AFTER COLLAPSE; THE OUTGOING SHOCK

At the instant of collapse ($t = t_0$), according to (7.8) and the equations just preceding it and the relations $p = s\rho^\gamma$, $c^2 = \gamma p/\rho$, the flow quantities vary as inverse powers of R, namely,

$$\begin{aligned} c(R, t_0) &= c_{00} R^{-\frac{1-\delta}{\delta}} \\ u(R, t_0) &= u_{00} R^{-\frac{1-\delta}{\delta}} \end{aligned}$$

$$\begin{aligned} \rho(R, t_0) &= \rho_{00} R^{-\frac{2(1-\delta)}{\delta(\gamma-1)}} \\ p(R, t_0) &= p_{00} R^{-\frac{2\gamma(1-\delta)}{\delta(\gamma-1)}} \end{aligned} \quad (8.1)$$

The infinite pressure gradient at $R = 0$ starts an outgoing shock, and we look for a self-similar solution for that shock and the flow behind it. The problem was considered by Hunter,⁶ but only in the isentropic approximation.

As in Sec. VII, the flow quantities are written as functions of a similarity variable, each multiplied by a power of $t - t_0$. As similarity variable, we choose

$$\hat{\eta} = \frac{R}{A(t-t_0)^\delta} = e^{-i\pi\delta} \eta \quad (8.2)$$

(This choice makes $\hat{\eta}$ real, for $t \geq t_0$. Since only $\log \eta$ appears in the ordinary differential equations, the extra factor $e^{-i\pi\delta}$ is irrelevant.) The flow ahead of the shock, for $t \geq t_0$, is simply the smooth continuation of the flow found in Sec. VII; hence, the value of δ has to be the same as in that section, to give the same behavior (8.1) for $t = t_0$. It then follows that the power of $t - t_0$ appearing in each flow quantity must be the same as in Sec. VII, because the compression ratio ρ_2/ρ_1 across the shock, the corresponding pressure ratio p_2/p_1 , velocity ratio u_2/u_1 , and "entropy" ratio s_2/s_1 (where $s = p\rho^{-\gamma}$), and so on, are all independent of t for a self-similar solution. Therefore, we write [compare with (7.3)]:

$$\begin{aligned} c(R, t) &= \frac{R}{t_0 - t} C(\hat{\eta}) = -A(t-t_0)^{\delta-1} \hat{\eta} C(\hat{\eta}), \\ u(R, t) &= \frac{R}{t_0 - t} V(\hat{\eta}) = -A(t-t_0)^{\delta-1} \hat{\eta} V(\hat{\eta}), \\ s(R, t) &= S(\hat{\eta}), \end{aligned} \quad (8.3)$$

for $t > t_0$. The functions C, V, and S can have jumps at the point $\hat{\eta} = \hat{\eta}^*$ where the shock occurs. (Note that c and C have opposite signs, as do u and V; that is unfortunate, but it obviated recoding some of the computer programs.)

In the similarity variables, the entropy equation

$$\frac{\partial s}{\partial t} + u \frac{\partial s}{\partial R} = 0$$

takes the form

$$[\delta + V(\hat{\eta})]S'(\hat{\eta}) = 0.$$

It can be shown that $V(\hat{\eta})$ cannot be $\equiv -\delta$. In fact, $V(\hat{\eta})$ cannot be $= -\delta$ immediately behind the shock, at $\hat{\eta} = \hat{\eta}^*$, no matter what value $\hat{\eta}^*$ has; that follows from the numerical values of the self-similar solution already obtained for the flow ahead and the Rankine-Hugoniot shock conditions. Therefore, $S'(\hat{\eta}) = 0$, and $s(R,t)$ has a constant value s_2 in the flow behind the shock, which is, of course, not necessarily equal to the constant value s_1 ($= 1$ in Sec. VII) ahead of the shock; in fact, $s_2 > s_1$ because a shock always increases the entropy.

Since the entropy is constant, the ordinary differential equations (7.4) and (7.8) hold also behind the shock. In Fig. 10-4, two solution curves are shown in the C, V plane: the curve C , consisting of the parts C_1, C_2 , and C_3 , which describe the flow ahead of the shock, and the curve C' , which describes the flow behind it. The shock is a jump from a point P_0 on the first curve to a point P'_0 on the second.

The boundary conditions at $\hat{\eta} = 0$ for the curve C' come from conditions at the center of the system. For $R = 0$, the velocity u vanishes by symmetry, while the sound speed c assumes a positive value. Hence, by (7.3), $V(0)$ is finite, while $C(\hat{\eta}) \rightarrow \infty$, as $\hat{\eta} \rightarrow 0$. By letting $C(\hat{\eta}) \rightarrow \infty$ in the second differential equation (7.4), we find that

$$V(0) = -2(1-\delta)/3(\gamma-1). \quad (8.4)$$

That suffices to start the curve C' at very small $\hat{\eta}$, hence very large negative $C(\hat{\eta})$.

Let $\hat{\eta}$ be so normalized on C' (it can be multiplied by an arbitrary constant), that it has the value $\hat{\eta}^*$ at P' . Then, the constant A in (8.2) has the same value on both curves, because the coordinate R of the shock is the same when viewed from in front of the shock or behind it; the shock's position is

$$R_{sh} = A\hat{\eta}^*(t-t_0)^\delta,$$

and its speed is

$$\dot{R}_{sh} = \delta A\hat{\eta}^*(t-t_0)^{\delta-1}. \quad (8.5)$$

For the Rankine-Hugoniot jump conditions across the shock, let subscripts 1 and 2 denote values just before and just behind the shock, respectively. The conditions are

$$(\dot{R}_{sh} - u_1)\rho_1 = (\dot{R}_{sh} - u_2)\rho_2 = \sqrt{\frac{p_2 - p_1}{\frac{1}{\rho_1} - \frac{1}{\rho_2}}}, \quad (8.6)$$

$$\frac{p_2}{p_1} = \frac{\theta x - 1}{\theta - x}, \text{ where } x = \frac{\rho_2}{\rho_1}, \theta = \frac{\gamma + 1}{\gamma - 1}. \quad (8.7)$$

In the similarity variables, these equations take the form

$$x = \frac{\gamma\phi_0}{\delta\phi_0 + \theta + 1}, \text{ where } \phi_0 = \frac{(V_1 + \delta)^2}{C_1^2} \quad (8.8)$$

$$C_2 = C_1 \sqrt{\frac{\theta x - 1}{(\theta - x)x}} \quad (8.9)$$

$$V_2 = \frac{V_1 + \delta}{x} - \delta. \quad (8.10)$$

The numerical procedure for locating the jump is this: For each point P on C_3 (where C_1 and V_1 are known), C_2 and V_2 are determined from these equations as target values to be attained by the jump, if the jump were to occur at point P . Then, for each P' on C' , the value V_2 of V at P' fixes the point P on C_3 from which the jump would have to start; P' is moved along C' (upward in the figure) and P correspondingly along C_3 , until condition (8.9) is also satisfied; $\hat{\eta}^*$ is then the value of $\hat{\eta}$ at the point P on C_3 , and the shock is completely determined.

Each of the solutions, for $t < t_0$, of the two-parameter family discussed in Sec. VII can be completed for $t > t_0$, in the way described here, by a self-similar solution containing an emerging shock. An example is shown in Fig. 10-4.

IX. THE FLUID-DYNAMICAL EQUATIONS IN SIMILARITY VARIABLES

To follow the motion of the free surface and the fluid just behind it near collapse, a special computational net and a special set of difference equations are used for the region $t \approx t_0, R \approx 0$. The independent variables are η and t , instead of R and t , where

$$\eta = R/\xi(t), \quad (9.1)$$

$\xi(t)$ being the position of the free surface, as determined by the calculation itself. Relative to the R grid, the η grid represents a moving and shrinking frame of reference. The two calculations are coupled by interpolation at intermediate radii, and the whole is called *similarity-fitting*. The dependent variables are $Q(\eta, t)$ and $V(\eta, t)$ and are related to the Reimann variable σ and the fluid speed u by the equations

$$\begin{aligned}\sigma(R, t) &= \dot{\xi}(t)\sqrt{\eta-1} Q(\eta, t) \\ u(R, t) &= \dot{\xi}(t)V(\eta, t).\end{aligned}\quad (9.2)$$

The factors $\dot{\xi}(t)$ are included to make Q and V dimensionless. (*Note 1:* It would have been more reasonable to include a minus sign in those definitions, because $\dot{\xi} < 0$, but that was not done. *Note 2:* The entropy equation was also carried along in the calculation, but vacuously, because only the isentropic case was computed.)

The partial differential equations result from substituting (9.2) into (3.8) and dropping the terms in \tilde{S} ; they are:

$$\begin{aligned}\dot{Q} + \frac{\ddot{\xi}}{\xi} Q + \frac{\dot{\xi}}{\xi} [(V-\eta) \\ \cdot (Q' + \frac{Q}{2(\eta-1)}) + \frac{\gamma-1}{2} Q \frac{1}{2}(\eta^2 V)'] &= 0, \\ \dot{V} + \frac{\ddot{\xi}}{\xi} V + \frac{\dot{\xi}}{\xi} [(V-\eta)V' \\ + \frac{\gamma-1}{2} Q\{(\eta-1)Q' + \frac{1}{2}Q\}] &= 0,\end{aligned}\quad (9.3)$$

where the dot denotes $\partial/\partial t$ and the prime $\partial/\partial \eta$. The system is of the form (3.7) and was differenced by the Lax-Wendroff two-step procedure (3.9).

If the flow is asymptotically self-similar, then $Q(\eta, t)$ and $V(\eta, t)$ should become independent of t , as $t \rightarrow t_0$.

The free-surface boundary conditions come from (4.5) and (4.7); they are

$$\left. \begin{aligned}\frac{\ddot{\xi}}{\xi} &= -\frac{\gamma-1}{4} \frac{\dot{\xi}}{\xi} Q^2 \\ v &= 1\end{aligned} \right\} \text{at } \eta = 1. \quad (9.4)$$

Note: The symbol V is used differently here and in Sec. VII. To get the quantities C and V of that section, the present quantities $\sqrt{\eta-1} Q$ and V must be multiplied by δ/η .

A consequence of the transformation (9.2) of the dependent variables is that, whereas σ and u were both known in advance at the free surface for the R grid (in particular, $\sigma = 0$), Q is now unknown there, and an additional equation is needed; it is obtained by setting $\eta = 1$, $V = 1$ in the first equation of (9.3) and evaluating $(V - \eta)/(\eta - 1)$ by L'Hôpital's rule [setting $\eta = 1$, $V = 1$ in the second equation of (9.3) merely gives the first of (9.4) again]. We find

$$\dot{Q} + \frac{\ddot{\xi}}{\xi} Q + \frac{\dot{\xi}}{\xi} \frac{Q}{2} [\gamma(V' + 2) - 3] = 0.$$

This equation is used (in effect as a special step 2 of the Lax-Wendroff) to advance Q in time at $\eta = 1$. The derivative V' is obtained at $t^{n+1/2}$ from the results of step 1 at $\eta - 1 = \Delta\eta/2$ and $3\Delta\eta/2$, together with the value $V = 1$ at $\eta = 1$, by differencing the extrapolation. Since Q' has disappeared, the above equation is not coupled to the equations for advancing Q at the regular net points $\eta - 1 = k\Delta\eta$ ($k = 1, 2, \dots$). A more careful treatment provides such coupling. Instead of merely evaluating the differential equation at $\eta = 1$, we average it over the interval $(1, 1 + \Delta\eta)$ with a weight that decreases linearly from 1 to 0 across the interval. (The corresponding effect is achieved in the normal Lax-Wendroff by the averaging that takes place in step 1.) Then, the above equation has an extra term and takes the following form:

$$\dot{Q} + \frac{\ddot{\xi}}{\xi} Q + \frac{\dot{\xi}}{\xi} \left\{ \frac{Q}{2} [\gamma(V' + 2) - 3] + \frac{1}{3}(V_1 - \eta_1)Q' \right\} = 0, \quad (9.5)$$

where $\eta_1 = 1 + \Delta\eta$ and $V_1 = V(\eta_1)$. Since the last term is of smaller order than the others, it is adequate to evaluate Q' to first order from the values at t^n .

The application of (9.4) and (9.5) is similar to the fitting procedure at the free surface described for the R net in Secs. IV and V, but with the following differences: (a) there is no need for a special Lax-Wendroff step 1, because the free surface is always at the net point $k = 0$; (b) after step 1 at $k = 1/2, 3/2, \dots$ and the approximation to $\ddot{\xi}$, $\dot{\xi}$, and ξ at $t^{n+1/2}$ that result therefrom, and after the regular step 2 at $k = 1, 2, \dots$, Eq. (9.5) is then used to advance Q at $\eta = 1$ as described above, and the values of $\ddot{\xi}$, $\dot{\xi}$ and ξ at t^{n+1} are obtained from (9.4); (c) there is no need for interpolations at t^{n+1} , because no points of the η net are uncovered by the motion of the free surface.

When the method of this section is combined with that of Sec. III, involving a standard $R - t$ net, and the two calculations are coupled by mutual interpolations at intermediate radii, the effect is to

provide a refinement of the $R - t$ net near $t = t_0$, $R = 0$, the degree of refinement increasing without limit as collapse is approached. The overall procedure is called *zooming*, to borrow a term from photography.

The resulting algorithm was tested by a calculation in which the initial values of Q and V were taken, for an interval $1 = \eta_0 \leq \eta \leq \eta_k (= 2.0 \text{ or } 5.0)$, from the similarity theory of Sec. VII, for $\delta = 0.708542$. The computed Q and V were very nearly independent of t over a very large number of cycles, as they ought to be, while $\xi(t)$ and $\dot{\xi}(t)$ varied as shown in the log-log plot of Fig. 9-1. It is seen that the values lie on a straight line with slope $-(1 - \delta)/\delta$, as they should.

For $\eta > \eta_c$ (η_c is the value of η at the critical point through which the solution of the ordinary differential equation passes; $\eta_c = 1.22195$ for $\delta = 0.708542$), both characteristics of the system (9.3) slope to the right; they both represent signals moving away from the region near the free surface. One consequence is that in the test calculation just referred to, one must not impose a boundary condition at $\eta = \eta_k$; the boundary value must float freely with the solution of the differential equation. To achieve that, we calculated the boundary values by one step of the so-called Courant-Isaacson-Rees method, which consists of writing the equation in characteristic form, following each characteristic back from $\eta = \eta_k$, $t = t^{n+1}$ to a point between η_{k-1} and η_k at $t = t^n$, obtaining a value of the Riemann invariant at that point by linear interpolation on η , and then using that value of the invariant at η_k , t^{n+1} .

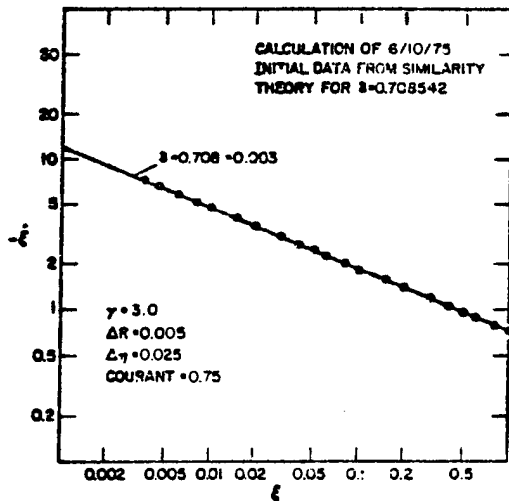


Fig. 9-1.

Agreement of the solution of the partial differential equations with the similarity solution, when the initial data are taken from the similarity solution.

X. THE ASYMPTOTIC BEHAVIOR OF THE CAVITY-COLLAPSE PROBLEM

In view of the large number (in fact, two-parameter family) of self-similar solutions found in Secs. VII and VIII, the question arises, to which of them the solution of the full cavity-collapse problem is asymptotic, and indeed whether the solution of that problem is asymptotically self-similar at all. The same questions arise about other problems of cavity collapse, such as ones in which the fluid is initially moving inward, ones with spherical layers of different material, and so on. To begin investigating those questions, a few calculations have been made with the zooming method described in the preceding section, for the present problem as described in Sec. II. One such calculation, for $\gamma = 3.0$, which used 400 points in the R net and 700 in the η net, will now be briefly analyzed, and will be referred to simply as the *full calculation*.

One method, in principle, for testing the asymptotic self-similarity of the solution obtained by the full calculation is to plot $\dot{\xi} \log\log$ vs ξ ; the graph should be asymptotically a straight line with slope $-(1 - \delta)/\delta$. If the solution were truly self-similar, that would be as good as any other method of test. In practice, it is unsatisfactory, because, as stated in Secs. VII and IX, the motion of that part of the fluid corresponding to the interval $(1, \eta_c)$ of the variable η , a part having vanishing mass in the limit $t = t_0$, is unaffected by the motion of all the fluid outside, until after collapse; hence, its motion is not necessarily representative of that of the bulk of the fluid. The log-log plot of $\dot{\xi}$ vs ξ is given in Fig. 10-1, and it is

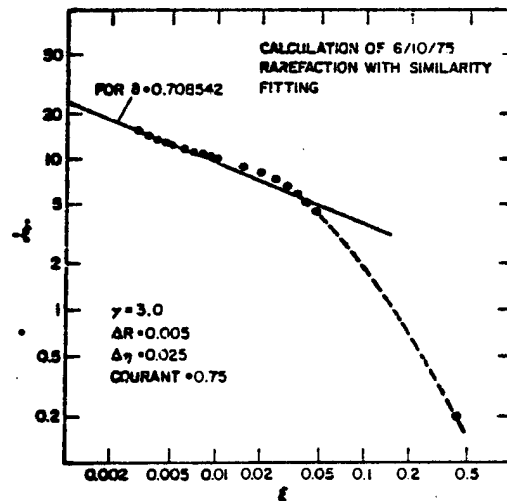


Fig. 10-1.

Weak approach of $\dot{\xi}$ vs ξ , from solution of the partial differential equations, to the similarity solution.

XI. CONCLUSIONS

Satisfactory numerical methods have been developed for handling singularities of the first three kinds listed in Sec. II, namely the free surface, the rarefaction head, and the region in between at early times. The methods are described in Secs. III-V and were shown to give about two orders of magnitude greater accuracy than a comparison calculation by conventional methods. The methods that have been developed for singularities of the fourth and fifth kind, namely the cavity collapse at late times and the emergent shock at early times, are described in Secs. VII-X; they are somewhat less satisfactory from a theoretical point of view, owing to unresolved questions whether the solution is asymptotically self-similar, and, if so, what the values of the appropriate parameters are.

A method of testing the asymptotic self-similarity of the solution of the full problem is given, which is superior to the perhaps more obvious method of a log-log plot of ξ vs ξ .

The similarity theory has been developed in some detail. The present work goes beyond that of Hunter⁶ in that the entropy increase at the shock is taken into account. It is shown that there is a two-parameter family of self-similar solutions that contrasts with the collapsing shock problem of Guderley⁹ and Butler,¹⁰ in which there is only one.

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