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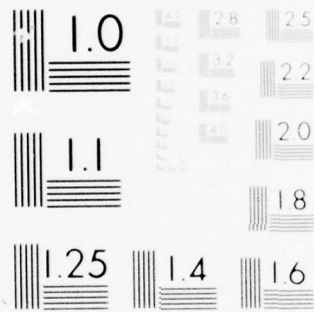
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AIRESEARCH REPORT 21-2548

AN EQUATION SPLITTING METHOD FOR
SHOCK BOUNDARY LAYER INTERACTION

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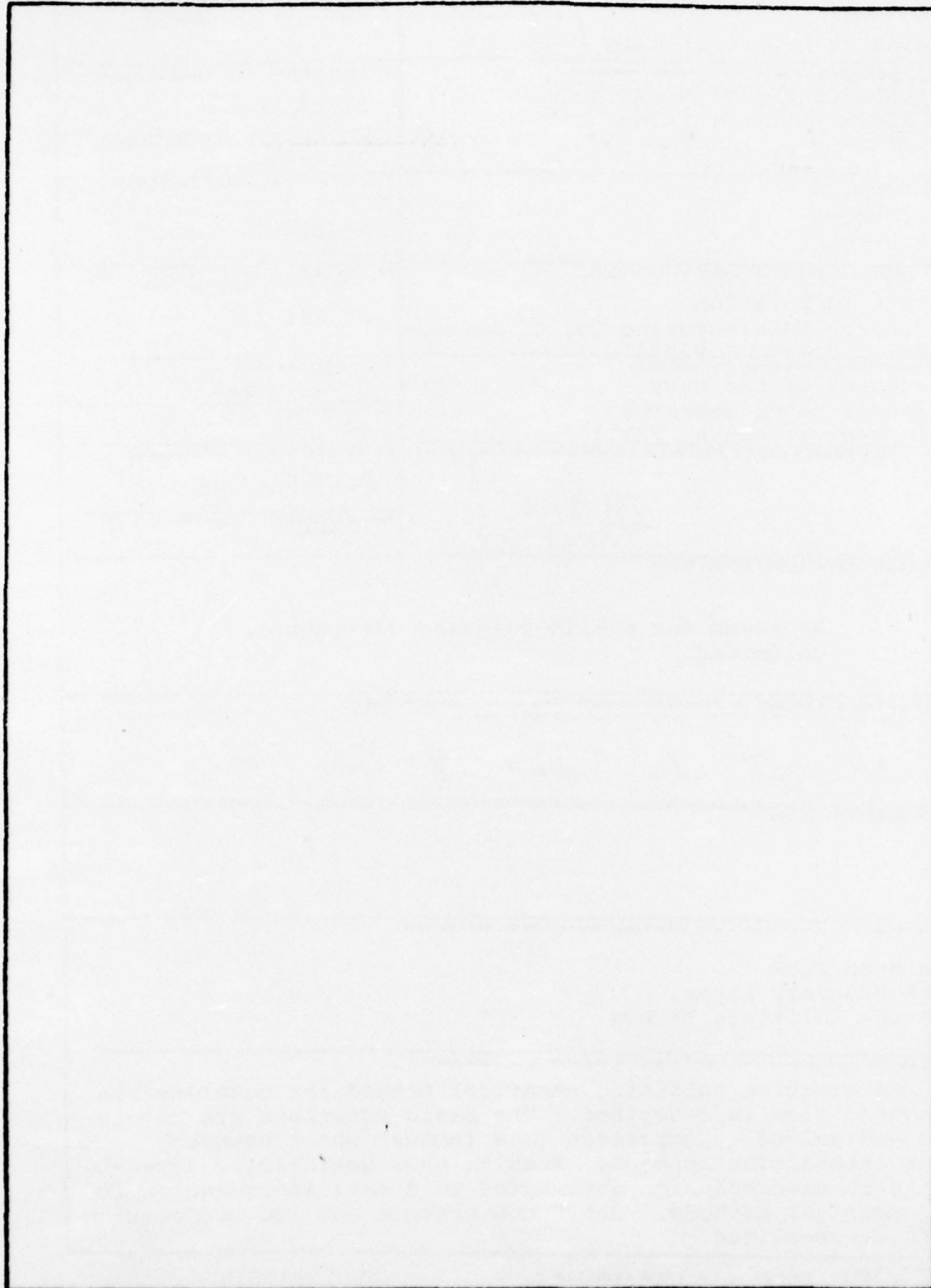
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PREFACE

AiResearch Manufacturing Company of Arizona has conducted a research study for the Office of Naval Research under Contract N00014-74-C-0317, Task NR 061-221, to develop rapid efficient techniques for Navier-Stokes solutions. This document reports the last year's activity. Previous activities have been reported and are referenced herein.

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SECTION I INTRODUCTION

Traditional techniques of solving the Navier-Stokes equations for compressible separated flow have utilized time-dependent methods. These methods have, with extensive development, steadily improved in both accuracy and computational times. Recently, however, Dodge¹ and Dodge and Lieber² have introduced a new steady-state solution method for incompressible flow that splits the pressure and shear forces into two separate equations. One equation is elliptic, and the other nearly parabolic. Contained herein is a description of the first attempt to extend this method to compressible flow. Application is made to the interaction of an oblique shock wave with a laminar boundary layer. The development was guided by the realization that the ultimate value of such a method lies not in the solution of two-dimensional laminar shock boundary layer interactions, but rather in applications to more complete two-dimensional and three-dimensional geometries. Although more development to improve accuracy and further reduce computational times is suggested, the results to date demonstrate the general viability of this numerical method. The following sections describe in detail the basic equation development, numerical techniques, results of the calculations, and suggestions for accuracy improvements.

SECTION II EQUATION DEVELOPMENT

The basis of Dodge's equation-splitting solution method lies in the separation of equations describing pressure and viscous effects. To accomplish this separation, the velocity vector $\vec{W}$ is expressed in the following manner:

$$\vec{W} = \nabla\phi + \vec{U} \quad (1)$$

where ϕ = velocity potential

$\vec{U}$ = velocity vector reflecting viscous effects
(viscous loss velocity vector)

Equations are then required to solve individually for ϕ and $\vec{U}$.

As an initial step, consider the vector form of the momentum equation for steady laminar flow:

$$\rho(\vec{W} \cdot \nabla)\vec{W} = -\nabla P + \nabla(\lambda \nabla \cdot \vec{W}) + \nabla \cdot (\mu \text{def} \vec{W}) \quad (2)$$

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where ρ = density
 P = static pressure
 μ = molecular viscosity
 λ = second coefficient of viscosity

$$\text{def} \vec{W} = \nabla \vec{W} + (\nabla \vec{W})^*$$

$$(\nabla \vec{W})^* = \text{transpose } (\nabla \vec{W})$$

Equation (1) may then be substituted for $\vec{W}$ in Equation (2) as follows:

$$\begin{aligned} \rho \nabla \phi \cdot \nabla (\nabla \phi) + \rho \vec{U} \cdot \nabla (\nabla \phi) + \rho (\vec{W} \cdot \nabla) \vec{U} \\ = - \nabla P + \nabla (\lambda \nabla \cdot \vec{W}) + \nabla \cdot (\mu \text{def} \vec{W}) \end{aligned} \quad (3)$$

To treat the static pressure gradient in Equation (3), consider the density to be described by two terms -- one, ρ^* , related to static pressure and the other, ρ_v , related to viscous effects:

$$\rho = \rho^* + \rho_v \quad (4)$$

The density component ρ^* is defined by

$$\rho^* = \rho_I^* \left(1 - \frac{\nabla \phi \cdot \nabla \phi}{2C_p T_I^*} \right)^{\frac{1}{\gamma-1}} \quad (5)$$

where

ρ_I^* = reference condition of stagnation density
 T_I^* = reference condition of stagnation temperature
 C_p = specific heat at constant pressure
 γ = ratio of specific heats

Further, one can show that all possible static pressure gradients may be represented by

$$-\nabla P = \rho^* \nabla \phi \cdot \nabla (\nabla \phi) \quad (6)$$

where ρ^* is selected so that the right-hand side of Equation (6) is irrotational. Substituting Equation (4) into Equation (3), combining terms, and using Equation (6) yields

$$\begin{aligned} (\rho - \rho^*) \nabla \phi \cdot \nabla (\nabla \phi) + \rho \vec{U} \cdot \nabla (\nabla \phi) + \rho (\vec{W} \cdot \nabla) \vec{U} \\ = \nabla (\lambda \nabla \cdot \vec{W}) + \nabla \cdot (\mu \text{def} \vec{W}) \end{aligned} \quad (7)$$

Equations (6) and (7) could then provide the basis for separate equations for ϕ and $\vec{U}$, respectively. However, when Equation (6) is combined with the continuity equation and the equation of state, it yields a relation that becomes hyperbolic for supersonic conditions based upon potential rather than total velocity. To reconcile this difficulty, a different approach was taken in the development of the potential equation.

The pressure gradient may be expressed in terms of the equation of state, as

$$\nabla P = a^2 \nabla \rho + a^2 \rho \left(\frac{\gamma - 1}{\gamma R} \right) \nabla S \quad (8)$$

where

a = local speed of sound at static conditions

R = gas constant

S = entropy

Replacing the pressure gradient term in the momentum equation, Equation (2), with Equation (8), and dotting the resulting expression with $\vec{W}$, yields Equation (9).

$$\begin{aligned} \vec{W} \cdot [\rho (\vec{W} \cdot \nabla) \vec{W}] &= -a^2 (\vec{W} \cdot \nabla \rho) - a^2 \rho \left(\frac{\gamma - 1}{\gamma R} \right) \vec{W} \cdot \nabla S \\ &+ \vec{W} \cdot \nabla (\lambda \nabla \cdot \vec{W}) + \vec{W} \cdot [\nabla \cdot (\mu \text{def} \vec{W})] \end{aligned} \quad (9)$$

The vector form of the compressible continuity equation, given by

$$\nabla \cdot (\rho \vec{W}) = 0 \quad (10)$$

may be expanded and re-expressed as

$$\vec{W} \cdot \nabla \rho = - \rho \nabla \cdot \vec{W} \quad (11)$$

Substituting Equation (11) into the first term of the right-hand side of Equation (9), rearranging, and dividing by ρ , yields Equation (12).

$$\begin{aligned} \vec{W} \cdot [(\vec{W} \cdot \nabla) \vec{W}] - a^2 (\nabla \cdot \vec{W}) &= - a^2 \left(\frac{\gamma - 1}{\gamma R} \right) \vec{W} \cdot \nabla S \\ &+ \vec{W} \cdot \left[\frac{1}{\rho} \nabla (\lambda \nabla \cdot \vec{W}) \right] + \vec{W} \cdot \left[\frac{1}{\rho} \nabla \cdot (\mu \text{def} \vec{W}) \right] \end{aligned} \quad (12)$$

However, Equation (12) may be recast by replacing $\vec{W}$ in the left-hand terms of Equation (12) with the velocity components of Equation (1). Then, placing all terms involving potential on the left-hand side of the resulting equation, and expanding the last two terms, yields the following:

$$\begin{aligned} \vec{W} \cdot [(\vec{W} \cdot \nabla) \nabla \phi] - a^2 \nabla^2 \phi &= a^2 \nabla \cdot \vec{U} - \vec{W} \cdot [(\vec{W} \cdot \nabla) \vec{U}] \\ &- a^2 \left(\frac{\gamma - 1}{\gamma R} \right) \vec{W} \cdot \nabla S + \vec{W} \cdot \left[\frac{1}{\rho} \nabla (\lambda \nabla \cdot \vec{W}) \right] \\ &+ \vec{W} \cdot \left\{ \frac{\mu}{\rho} [\nabla (\nabla \cdot \vec{W}) + \nabla^2 \vec{W}] + (\text{def} \vec{W}) \frac{\nabla \mu}{\rho} \right\} \end{aligned} \quad (13)$$

Equation (13) thus describes pressure effects in terms of the velocity potential, ϕ . The equation will be elliptic or hyperbolic, depending upon the local Mach number, defined in terms of the velocity, W .

The equation for the viscous loss velocity vector is developed from the momentum equation, as presented in Equation (7). The terms on the right-hand side of Equation (7) may be expanded in the same manner as for the incompressible case presented by Dodge¹. Then, rearranging, and dividing the resulting equation by ρ , yields

$$\begin{aligned} \vec{U} \cdot \nabla (\nabla \phi) + (\vec{W} \cdot \nabla) \vec{U} &= \left(\frac{\rho^*}{\rho} - 1 \right) \nabla \phi \cdot \nabla (\nabla \phi) \\ &+ \frac{1}{\rho} \{ (\nabla^2 \phi) \nabla \lambda + (\nabla \cdot \vec{U}) \nabla \lambda + (2\mu + \lambda) \nabla (\nabla^2 \phi) \\ &+ (\mu + \lambda) \nabla (\nabla \cdot \vec{U}) + \mu \nabla^2 \vec{U} + 2 [\nabla (\nabla \phi)]^* \nabla \mu \\ &+ (\text{def} \vec{U}) \nabla \mu \} \end{aligned} \quad (14)$$

When expanded, Equation (14) will produce scalar equations for the viscous loss velocity components, U_i . These equations may be parabolized by treating the streamwise diffusion term as a known quantity from the previous iteration, as discussed by Dodge¹ and Dodge and Lieber².

Because the flow is compressible, an energy equation is also required. In terms of static enthalpy, the laminar steady-flow energy equation may be given as

$$\begin{aligned} \rho (\vec{W} \cdot \nabla) \left(h + \frac{1}{2} W^2 \right) &= \nabla \cdot (\lambda \vec{W} \nabla \cdot \vec{W}) \\ &+ \nabla \cdot (\vec{W} \cdot \mu \text{def} \vec{W}) + \nabla \cdot (k \nabla T) \end{aligned} \quad (15)$$

where

h = enthalpy at static conditions

k = conductivity of the fluid

T = static temperature

Expressing enthalpy in terms of static temperature, i.e.,

$$h = C_p T \quad (16)$$

and expanding the left-hand side of Equation (15), yields Equation (17)

$$\begin{aligned} \rho (\vec{W} \cdot \nabla) (C_p T) + \frac{1}{2} \rho (\vec{W} \cdot \nabla) W^2 &= \nabla \cdot (\lambda \vec{W} \nabla \cdot \vec{W}) \\ &+ \nabla \cdot (\vec{W} \cdot \mu \text{def} \vec{W}) + \nabla \cdot (k \nabla T) \end{aligned} \quad (17)$$

Assuming a constant Prandtl number, Pr , given by

$$Pr = \frac{\mu C_p}{k} \quad (18)$$

then conductivity may be expressed as

$$k = \frac{\mu C_p}{Pr} \quad (19)$$

In addition, specific heat is assumed to be constant. Thus, substituting Equation (19) into Equation (17), rearranging, and dividing by ρC_p , yields

$$\begin{aligned} (\vec{W} \cdot \nabla) T - \frac{1}{\rho C_p} \nabla \cdot \left[\left(\frac{\mu C_p}{Pr} \right) \nabla T \right] \\ = \frac{1}{\rho C_p} \left[-\frac{1}{2} \rho (\vec{W} \cdot \nabla) W^2 \right. \\ \left. + \nabla \cdot (\lambda \vec{W} \nabla \cdot \vec{W}) + \nabla \cdot (\vec{W} \cdot \mu \text{def} \vec{W}) \right] \end{aligned} \quad (20)$$

Expanding the second term on the left-hand side of Equation (20) results in the following:

$$\begin{aligned} (\vec{W} \cdot \nabla) T - \frac{1}{\rho Pr} [\nabla \mu \cdot \nabla T + \mu \nabla^2 T] \\ = \frac{1}{\rho C_p} \left[-\frac{\rho}{2} (\vec{W} \cdot \nabla) W^2 + \nabla \cdot (\lambda \vec{W} \nabla \cdot \vec{W}) \right. \\ \left. + \nabla \cdot (\vec{W} \cdot \mu \text{def} \vec{W}) \right] \end{aligned} \quad (21)$$

Expanding the last two terms of Equation (21) yields

$$\begin{aligned} (\vec{W} \cdot \nabla) T - \frac{1}{\rho Pr} (\nabla \mu \cdot \nabla T) - \frac{1}{\rho Pr} (\mu \nabla^2 T) \\ = \frac{1}{\rho C_p} \left\{ -\frac{\rho}{2} (\vec{W} \cdot \nabla) W^2 + \lambda (\nabla \cdot \vec{W})^2 \right. \\ \left. + \mu [\nabla \cdot (\vec{W} \cdot \text{def} \vec{W}) - \vec{W} \cdot (\nabla \cdot \text{def} \vec{W})] \right. \\ \left. + \vec{W} \cdot [\nabla (\lambda \nabla \cdot \vec{W}) + \nabla \cdot (\mu \text{def} \vec{W})] \right\} \end{aligned} \quad (22)$$

Substituting the momentum equation in the form of Equation (7) for the last two terms in Equation (22) gives the final form of the laminar, steady-flow energy equation in terms of static temperature:

$$\begin{aligned} (\vec{W} \cdot \nabla) T - \frac{1}{\rho Pr} (\nabla \mu \cdot \nabla T) - \frac{1}{\rho Pr} (\mu \nabla^2 T) \\ = \frac{1}{\rho C_p} \left\{ -\frac{\rho}{2} (\vec{W} \cdot \nabla) W^2 + \lambda (\nabla \cdot \vec{W})^2 \right. \\ \left. + \mu [\nabla \cdot (\vec{W} \cdot \text{def} \vec{W}) - \vec{W} \cdot (\nabla \cdot \text{def} \vec{W})] \right. \\ \left. + \rho \vec{W} \cdot [\vec{U} \cdot \nabla (\nabla \phi) + (\vec{W} \cdot \nabla) \vec{U}] \right. \\ \left. + (1 - \frac{\rho^*}{\rho}) \nabla \phi \cdot \nabla (\nabla \phi) \right\} \end{aligned} \quad (23)$$

As with the equation for $\vec{U}$, Equation (23) may be parabolized by treating the diffusion term as a known quantity from the previous iteration.

Equations (13), (14), and (23) comprise the system of equations that must be solved to yield ϕ , $\vec{U}$, and T . In conjunction with these equations, a set of boundary conditions is required. Figures 1 and 2 illustrate the boundary conditions on the solution space for the elliptic/hyperbolic and parabolic equations.

In Figure 1, the potential velocities V_I , V_M , and V_E are determined, based upon the oblique shock equations. The value of ϕ_E is determined by integrating the values of $\partial\phi/\partial x_1$ to the downstream boundary.

The upstream values of $\vec{U}$ and T in Figure 2 are dependent upon the corresponding value of $\vec{W}$, which is initialized using an arbitrary boundary layer profile similar to a Pohlhausen profile. An isothermal plate is assumed, with plate temperature equal to the reference stagnation temperature.

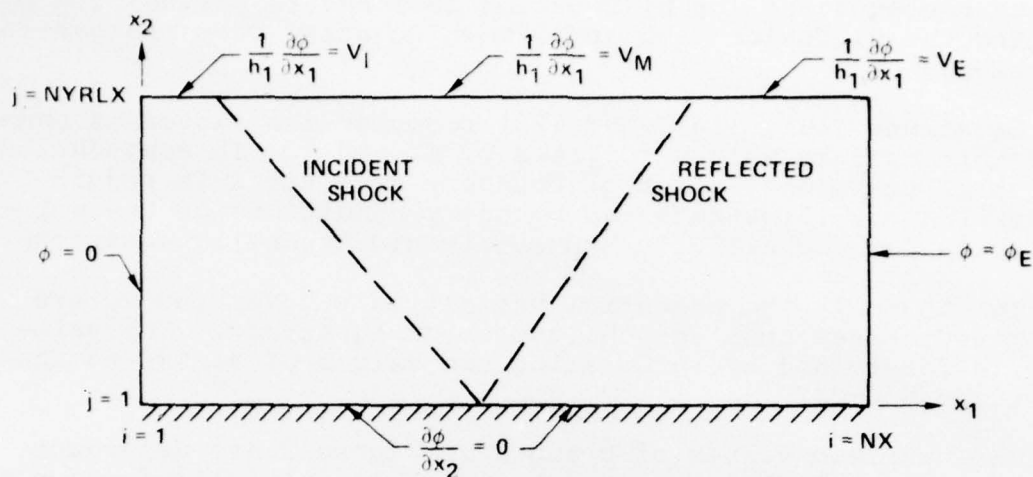


Figure 1. Boundary conditions on ϕ for elliptic/hyperbolic equation.

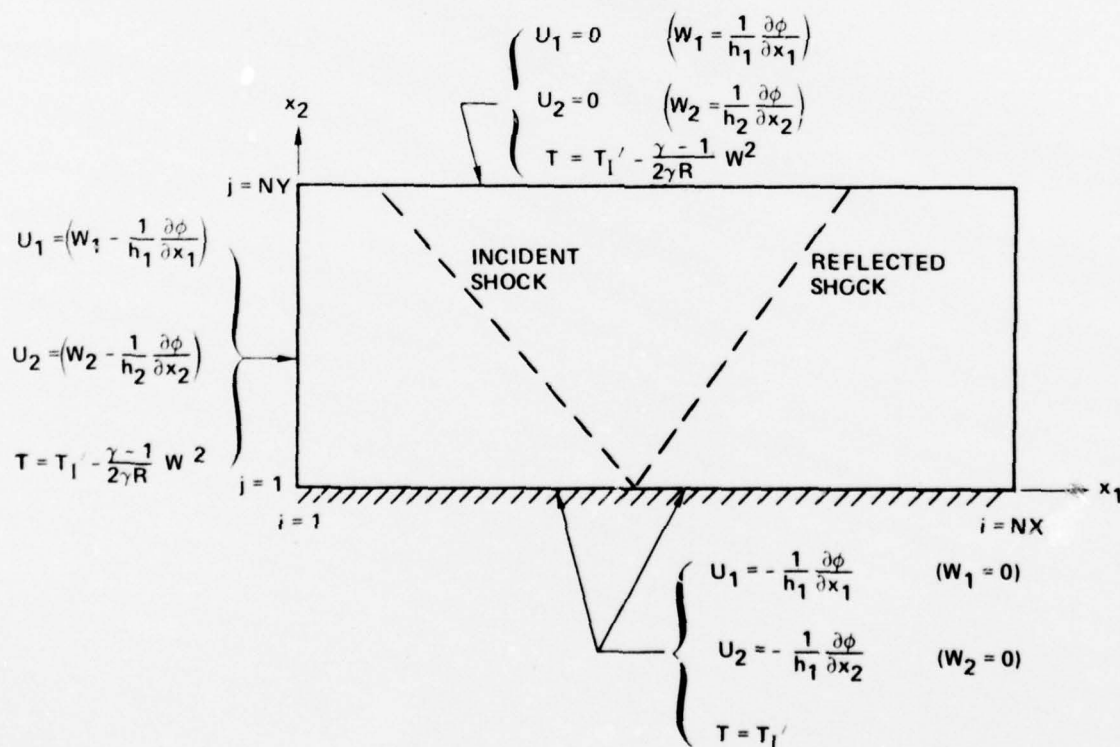


Figure 2. Boundary conditions on U_1 , U_2 , T for parabolic equation.

SECTION III NUMERICAL METHODS

Two major numerical solutions are required to implement the above algorithm: a solution to a mixed elliptic/hyperbolic potential equation, and a solution to a nearly parabolic viscous-shear equation. Murman and Cole³, in a pioneering paper, developed a technique for modifying a difference star with Mach number, and then using conventional relaxation to solve a similar equation. Modifications to the Murman and Cole method have been necessary to improve its accuracy, and allow its use in totally supersonic flow. Dodge^{4,5} describes such a method and its application to supersonic inlet, subsonic exit compressor blades. Murman⁶, in a subsequent paper, reported further improvements in the basic methods to account for differencing across discontinuities. From these works, a list of ideal characteristics can be assembled for the supersonic portion of any such algorithm.

- o Region of influence of difference star matches precisely the region of influence of the differential equation.
- o Across a potential slope discontinuity, the method must preserve the proper jump conditions.
- o The method remains integrally conservative in the non-discontinuous portion of the flow field.

Unfortunately, such a numerical method does not now exist.

To assist in the evaluation of the importance of each of the above criteria, the one-dimensional potential Equation (24) was studied.

$$(v^2 - a^2) \frac{d^2 \phi}{dx^2} = 0 \quad (24)$$

where

a = sound speed

$v = \frac{d\phi}{dx}$

ϕ = velocity potential

For the solution to be other than trivial, the flow must enter supersonically, undergo a jump, and exit subsonically. Boundary conditions then consist of an inlet Mach number and an exit potential. By integrating Equation (24) and assuming an ideal gas, a jump condition can be derived that relates downstream velocity to inlet velocity.

$$\begin{aligned} V_1 (a'^2 - \frac{\gamma + 1}{6} V_1^2) \\ = V_2 (a'^2 - \frac{\gamma + 1}{6} V_2^2) \end{aligned} \quad (25)$$

where

a' = stagnation speed of sound
 γ = ratio of specific heats
 V_1 = velocity ahead of the jump
 V_2 = velocity behind the jump

The analytical solution ahead of and behind the jump is obviously that of constant velocity. Jump location is determined by matching to the exit potential boundary condition using Equation (26).

$$x_J = \frac{\phi_{EXIT} - V_2 L}{V_1 - V_2} \quad (26)$$

where

x_J = position of the jump
 L = total length of the solution region ($X = 0$ is the inlet)
 ϕ_{EXIT} = exit potential

Consider what happens when the original Murman and Cole method is applied. Upstream of the assumed jump location, the difference star is cast in a backwards direction.

$$(\phi_i - 2\phi_{i-1} + \phi_{i-2}) = 0 \quad (27)$$

where

ϕ_i = potential at the i th node

Note that the value of $v^2 - a^2$ is immaterial since it is a non-zero constant at any particular node, and thus divides out. Downstream of the assumed jump location, the equation is elliptic.

$$\phi_{i+1} - 2\phi_i + \phi_{i-1} = 0 \quad (28)$$

The solution to both Equations (27) and (28) is a linear variation in ϕ . An initial guess for potential would be made. Equation (27) is solved starting at node 3, and simply broadcasts forward the initial velocity gradient until a node is reached, which was initially assumed subsonic. At this point, the subsonic solution proceeds as a boundary value problem with ϕ at the change-over (supersonic to subsonic) point established by Equation (27). There is, however, no communication between upstream and downstream. Thus, the location of the jump never changes, and the value of the velocity at the jump is only controlled by the initial guess of jump location and potential variation. No aspect of the physical problem appears in the numerical solution at the jump. To overcome this, special conditions at the jump must be invoked, as pointed out by Murman⁶. It is, however, immaterial what form is used away from the jump. A conservative form results in a solution that is no different from a nonconservative form.

When actually applying methods to introduce upstream-downstream communication, two classes result. The first class, into which Murman's method falls, allows the downstream flow to be adjusted based on upstream conditions. This class allows only for jump propagation in the forward direction. The second class, permits the jump to retreat by allowing the upstream flow to be affected by downstream conditions. Whether the jump needs to retreat or go forward is only dependent on the initial guess, and thus, for complete generality, both classes must be combined into a single method. For one-dimensional flow, such techniques are readily generated, the results of which are shown in Figure 3 for a forward propagating jump and Figure 4 for a retreating jump. The jump-node region is treated by the following algorithm

$$D = r\phi_{i+1} + \phi_{i-1} - (1+r)\phi_i = 0$$

applied to the node ahead of the jump if $D \geq 0$,

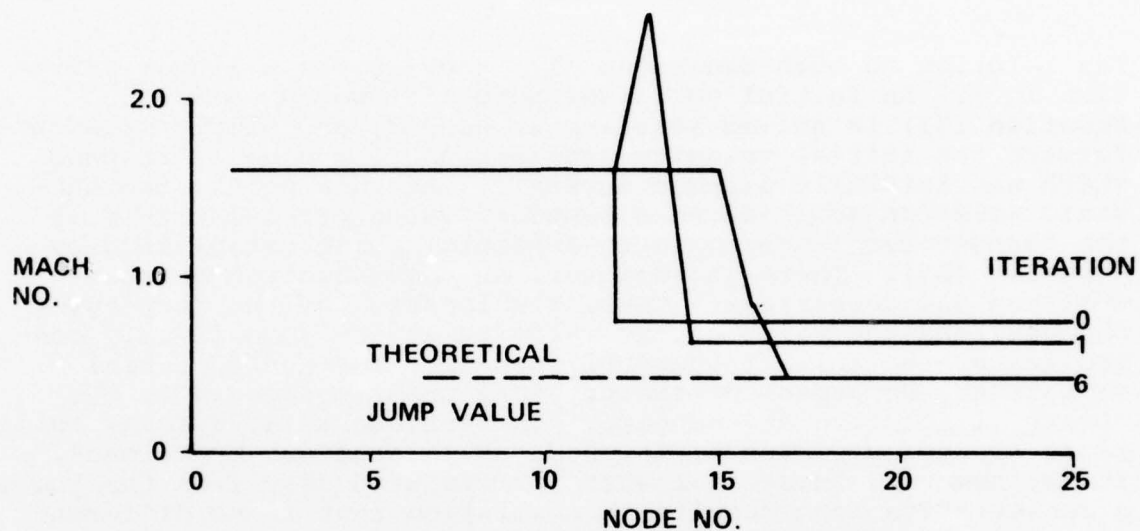


Figure 3. Forward propagating one-dimensional jump.

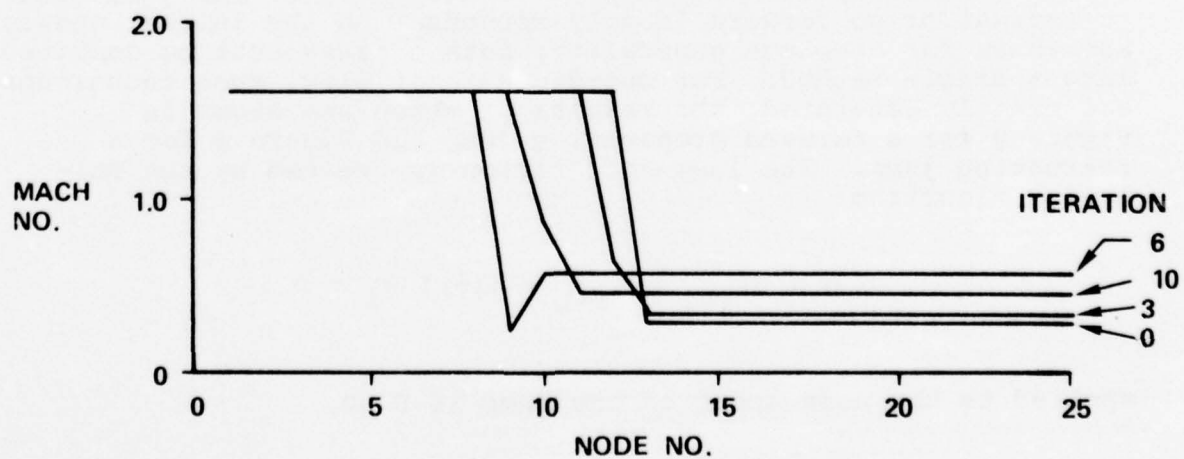


Figure 4. Retreating one-dimensional jump.

applied to the node behind the jump if $D < 0$

$$r = \frac{\alpha_{\text{DOWN}}}{\alpha_{\text{UP}}}$$

$$\alpha = V(a'^2 - \frac{\gamma+1}{6} V^2)$$

where DOWN and UP subscripts refer to downstream and upstream of the jump.

It is necessary to evaluate the velocity away from the jump, since values at the jump, particularly for intermediate iterations, are in significant error.

When the flow field is multidimensional and jumps are other than normal, the proper procedure is complicated by the lack of accepted numerical solutions. Murman's method assumes an oblique jump passing through the grid system at an arbitrary angle. However, this derivation loses sight of the fact that the differential equation is left behind when solutions are made to the finite-difference grid. Discontinuities on a fixed orthogonal finite-difference system can only have one of three forms, illustrated by Figure 5. A physically oblique wave must be made up of discrete components of the three jumps. The normal jump causes little or no problems, but the other two result in differencing across a discontinuity. To be totally correct, as Murman points out, any difference scheme applied to any combination of the above discontinuities must be integrally (summation in difference system) conservative of the desired jump condition. Even if such a system existed, however, practical problems, such as detecting a jump and determining proper upstream and downstream conditions away from the perturbed region of the jump, intervene to prevent reliable treatment of jump conditions.

The method by Dodge⁴ avoids some of the difficulty by only differencing along characteristic lines, and not differencing across oblique jumps, which may explain some of its good results. However, it requires a nonstationary grid system that is difficult to implement in two dimensions, and is nearly impossible in three dimensions. Thus, after comparing other methods, such as the two by Murman, to a range of six supersonic and supercritical cases, it was decided to mix a nonorthogonal and orthogonal difference system, staying with a fixed orthogonal grid system.



Figure 5. Possible jump nodes.

The degree of mixing was determined as a function of local Mach number, based upon a criterion developed from a stability analysis of the basic relaxation difference equation. Figure 6 illustrates the difference stars utilized. In simplified form, the mixing of difference operators is accomplished in the following following manner:

$$C_1 \frac{\partial^2 \phi}{\partial x^2} + C_2 \frac{\partial^2 \phi}{\partial y^2} = [C_1 \frac{\partial^2 \phi}{\partial x^2} + C_2 \frac{\partial^2 \phi}{\partial y^2}] (f_H) \\ + [C_1 \frac{\partial^2 \phi}{\partial x^2} + C_2 \frac{\partial^2 \phi}{\partial y^2}] (1 - f_H)$$

where f_H represents the fraction of hyperbolic difference operator utilized, and $(1-f_H)$ represents the fraction of single-spaced implicit difference operator. The local value of f_H was determined for supersonic flow to be

$$f_H = -C_2/C_1 \text{ for } (-C_2/C_1) < 1.0$$

or

$$f_H = 1.0 \text{ for } (-C_2/C_1) \geq 1.0$$

based upon a stability analysis.

If the variable supersonic difference star is examined in relation to the location of local characteristic lines, it may be seen that, as Mach number increases, the difference star becomes weighted to more closely reflect the region of influence of the characteristic lines (refer to Figure 7), eventually becoming totally hyperbolic.

Admittedly, such a system does not meet all the criteria for a potential equation solver set out above, but represents a workable compromise between complete agreement with the above criterion, and the practical constraints of computer code development.

The viscous equation, on the other hand, represents a much easier form to implement. Dodge¹ and Dodge and Lieber² have discussed many of the difficulties associated with marching equation solution. For this program, the streamwise diffusion term was determined from the previous iteration. In separated flow regions, the difference star for the convection term was turned upstream, utilizing the values established on a previous iteration. With these adjustments, the solution to the viscous equation proceeds as a normal marching solution to a parabolic equation. The component of velocity normal to the wall in the

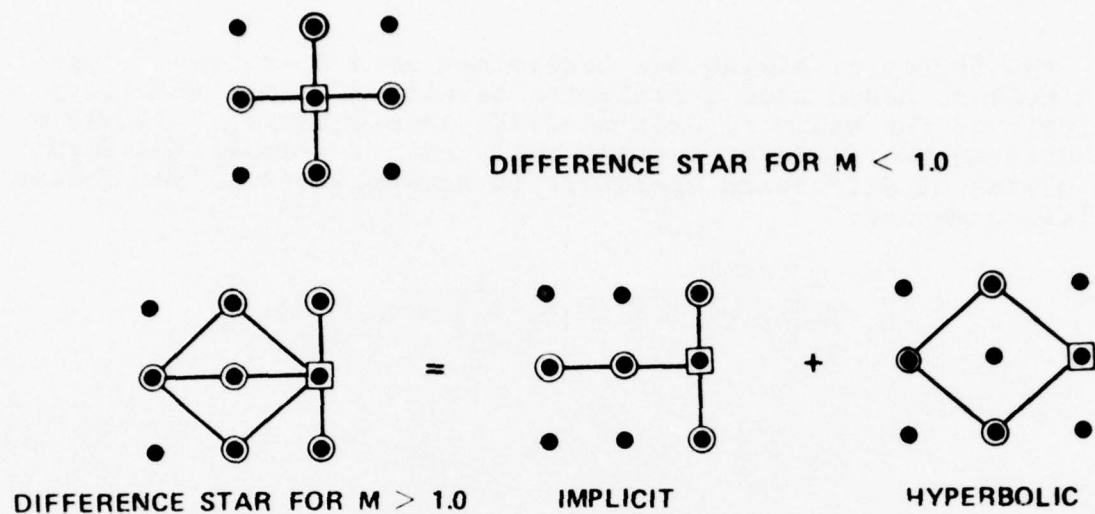


Figure 6. Transonic difference stars.

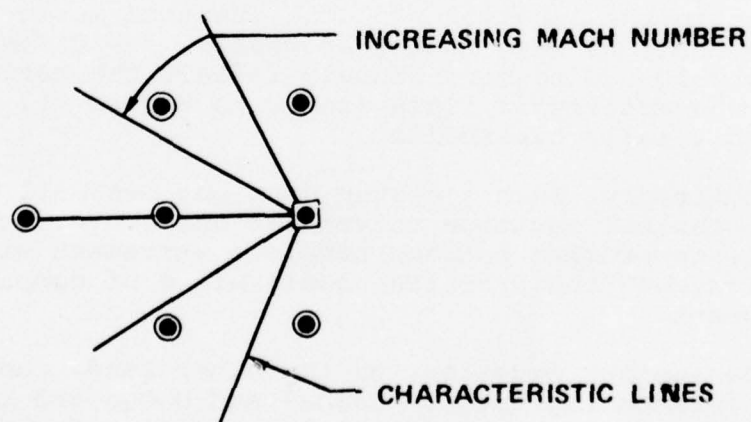


Figure 7. Supersonic difference star.

region between the wall node and the first relaxation node was, however, determined from the continuity equation. This was found in past studies by Dodge¹ to produce somewhat more accurate results.

The program developed to implement this numerical solution method consisted of a modularized code written in a structured FORTRAN language. A basic logic diagram for the program may be found in Figure 8.

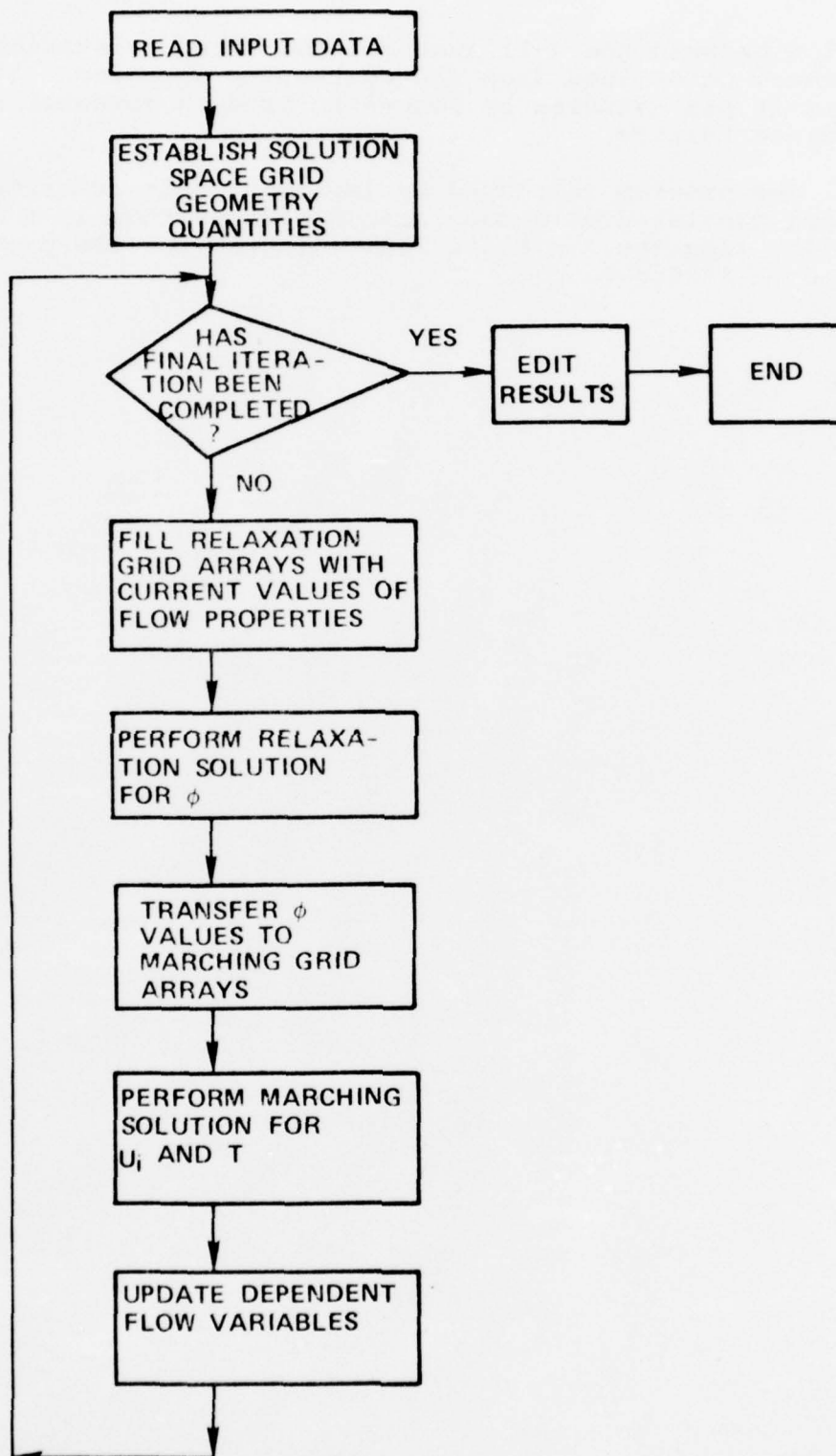


Figure 8. Program logic diagram.

SECTION IV CALCULATION RESULTS

The numerical method discussed herein has been utilized to predict separation of a two-dimensional laminar flat-plate shock boundary layer interaction flow. The test case used was that of Hakkinen, et al⁷, consisting of a laminar flat-plate flow with an upstream Mach number $M_o = 2.0$, and a Reynolds number $Re_L = 2.96 \times 10^5$, where

$$Re_L = \frac{W_o L}{\nu_o}$$

and

W_o = freestream inlet velocity
($W_o = 1664$ ft/sec)

ν_o = freestream inlet value of kinematic viscosity at static conditions

L = distance from leading edge to shock intersection with plate
($L = 0.1624$ ft)

An oblique shock wave, with shock angle $\theta = 32.59^\circ$, intersects the flat plate as indicated in Figure 9.

The grid system employed in the numerical solution consisted of a linear relaxation grid and a marching grid composed of the relaxation nodes, with an additional nonlinear grid segment near the wall. This scheme permits a coarse grid to be used for the relaxation solution, and yet provides a more finely-spaced grid within the boundary layer for the viscous marching equation solution. Figure 10 illustrates the relative spacing of the relaxation and marching grid nodes within the boundary layer.

Results of the present numerical method were compared with both the experimental data of Hakkinen, and results of the time-dependent scheme of MacCormack⁸.

Figure 11 compares surface pressure distribution for the three sets of data. Clearly, the pressure plateau in the region of the shock/plate intersection is not indicated by the present method. A much sharper pressure rise is predicted than evidenced by experiment.

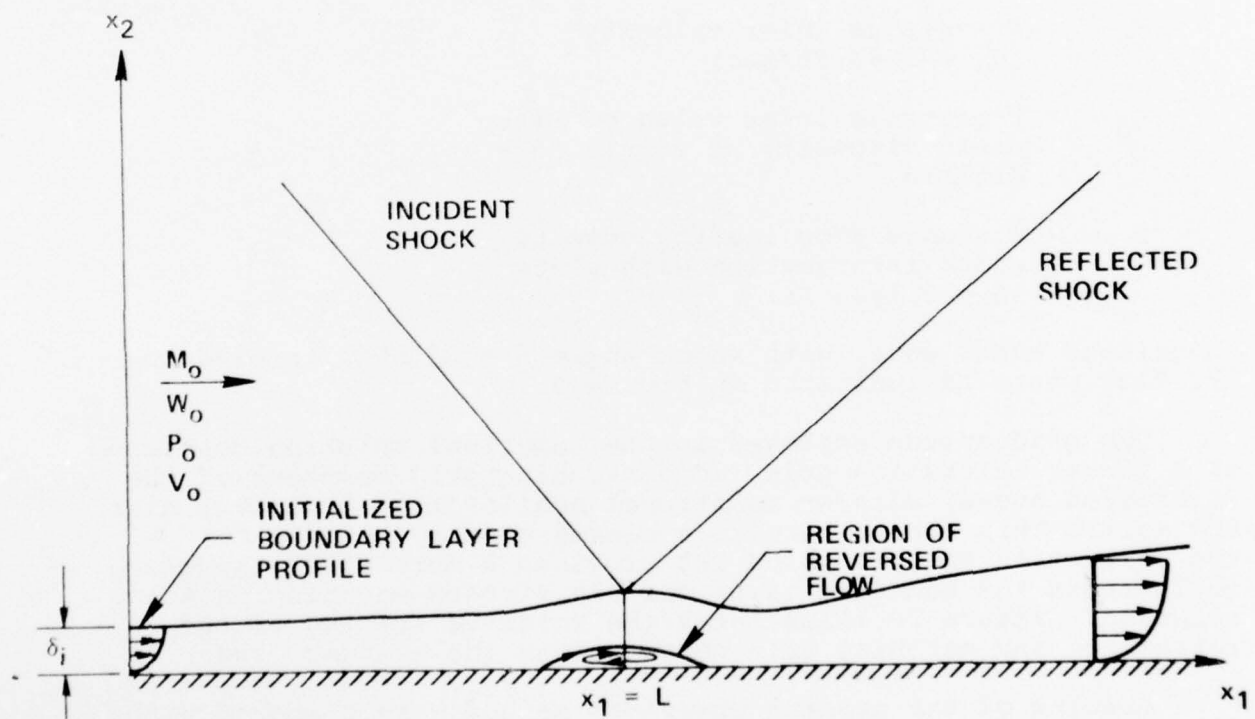


Figure 9. Generalized two-dimensional shock boundary layer interaction flow field.

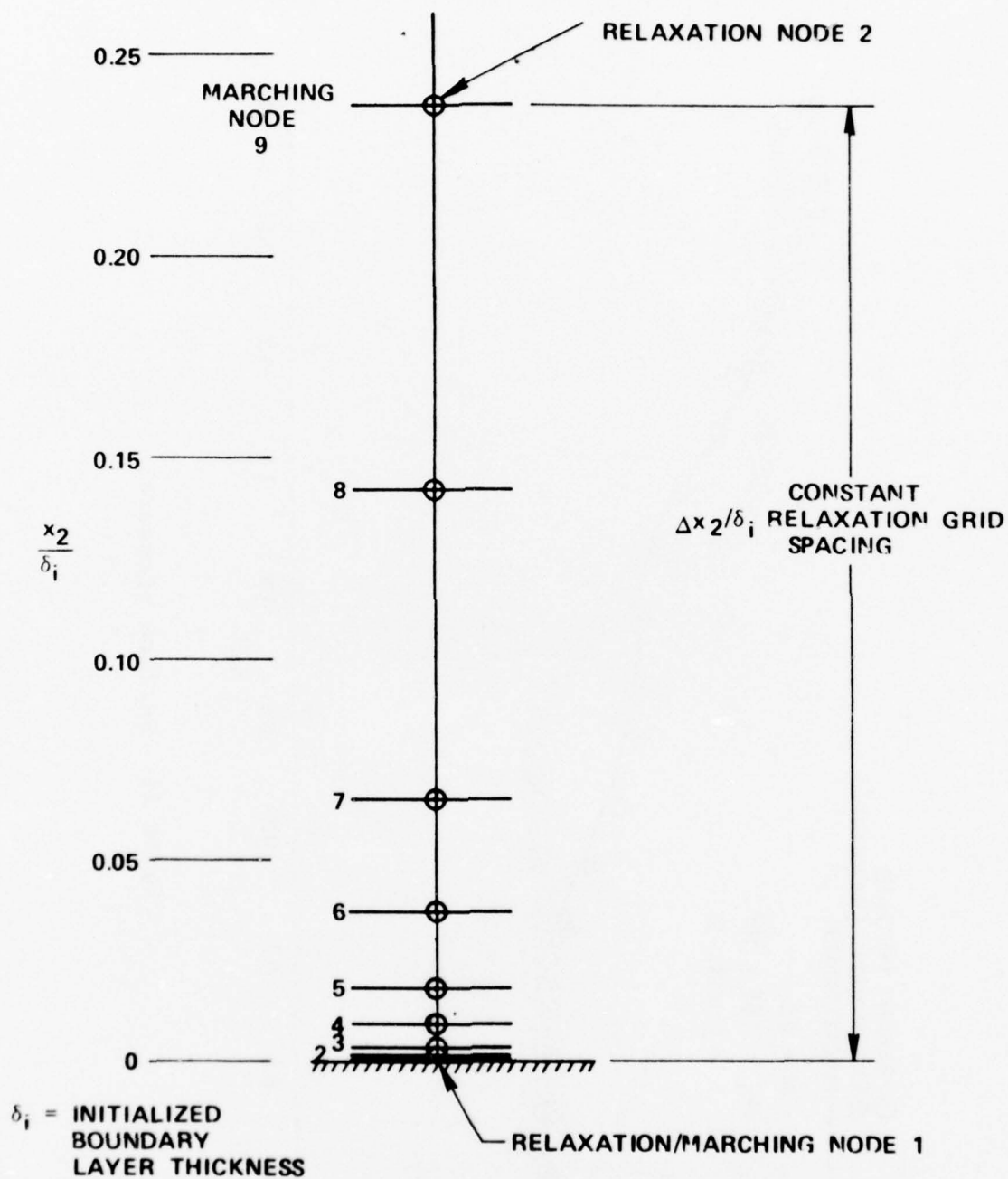


Figure 10. Relative relaxation and marching grid spacing.

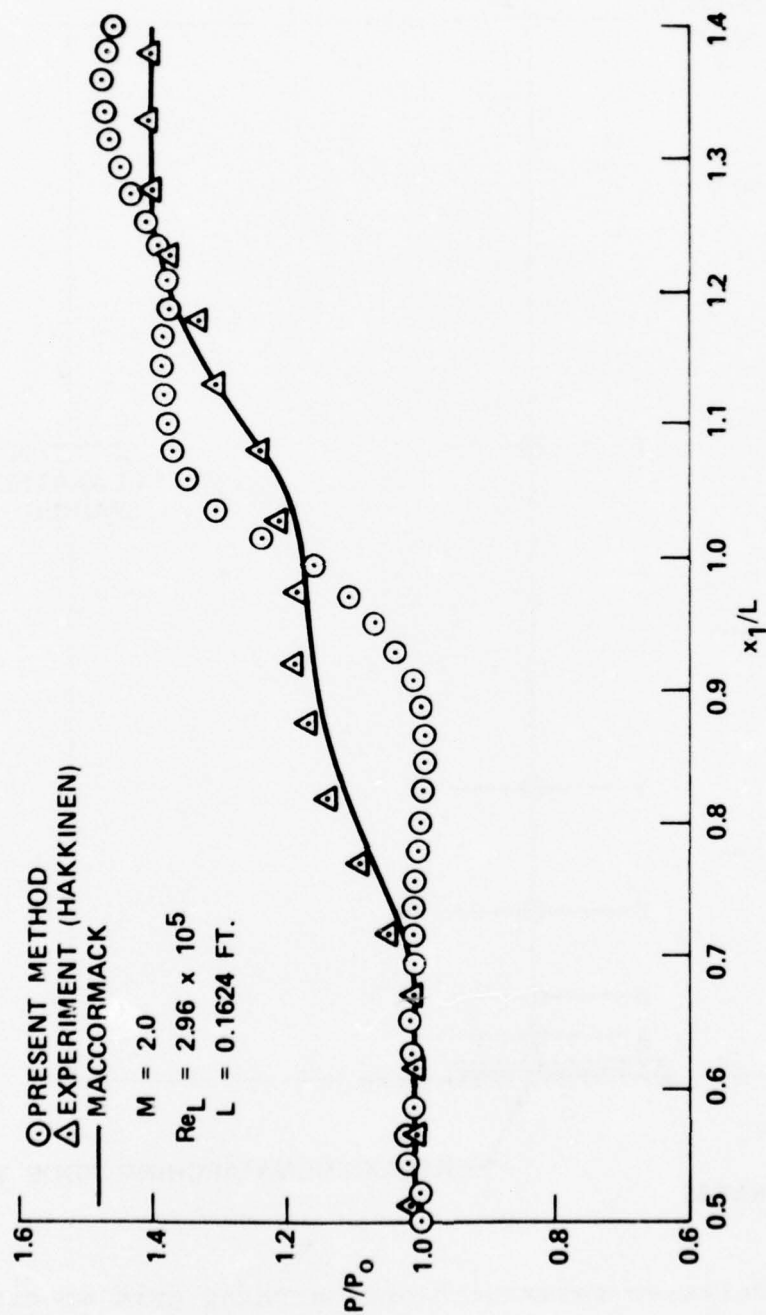


Figure 11. Surface pressure.

As a consequence of this too-sudden rise in static pressure, the predicted values of skin friction coefficient, C_f , shown in Figure 12, fail to decrease as far upstream as the experimental data. The predicted region of reversed flow is also much smaller than that indicated by Hakkinen's data.

The predicted values of C_f are presented as average values within a bandwidth of error representing the standard deviation of C_f as calculated at as many as 8 nodes within the constant pressure and density portion of the boundary layer. The skin friction coefficient is defined by

$$C_f = \frac{\rho_w v_w \frac{1}{h_2} \left. \frac{\partial W_1}{\partial x_2} \right|_w}{\frac{1}{2} \rho_\infty W_\infty^2}$$

where the subscript w refers to condition at the wall, and the subscript ∞ refers to the freestream. A considerable fluctuation in values of $\partial W_1 / \partial x_2$ takes place, apparently due to the nonlinear character of the grid system near the plate.

Figures 13 through 15 are comparisons of boundary layer velocity profiles between the present steady-flow method and MacCormack's technique. In general, agreement is fair; however, once again it is apparent that the present method predicts a smaller region of separated flow, and generally larger displacement thicknesses. Discrepancies in the assumed inlet boundary layer profile are also apparent. Note that dimensions in the x_2 direction have been normalized by H_f (where $H_f = 0.003$ ft) -- a grid parameter used in MacCormack's data presentation.

Figure 16 presents static pressure contours over the solution space, based upon the results of the steady-flow method discussed here. The incident and reflected shocks are apparent, along with a spurious downstream reflection from the upper boundary of the region. The effect of the reflection is clearly visible in the downstream portion of the skin friction and surface pressure curves.

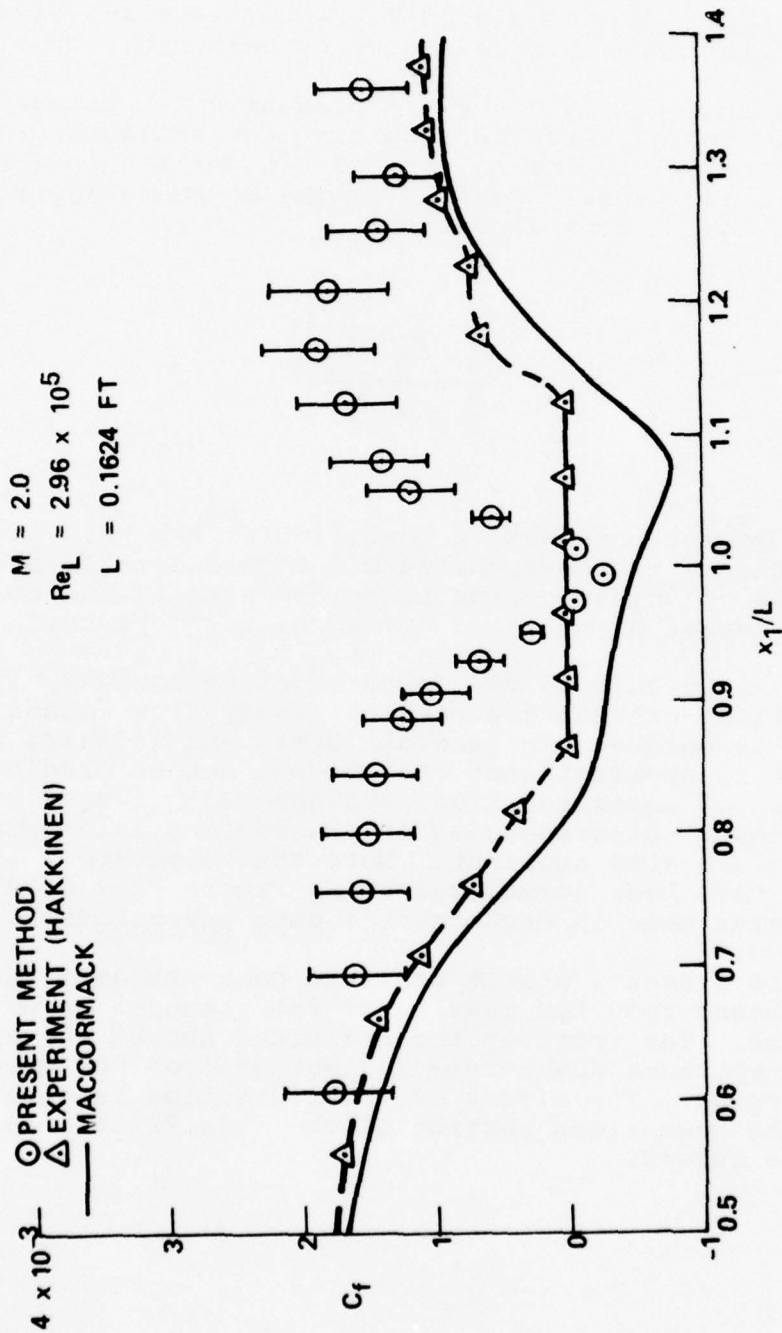


Figure 12. Skin friction coefficient.

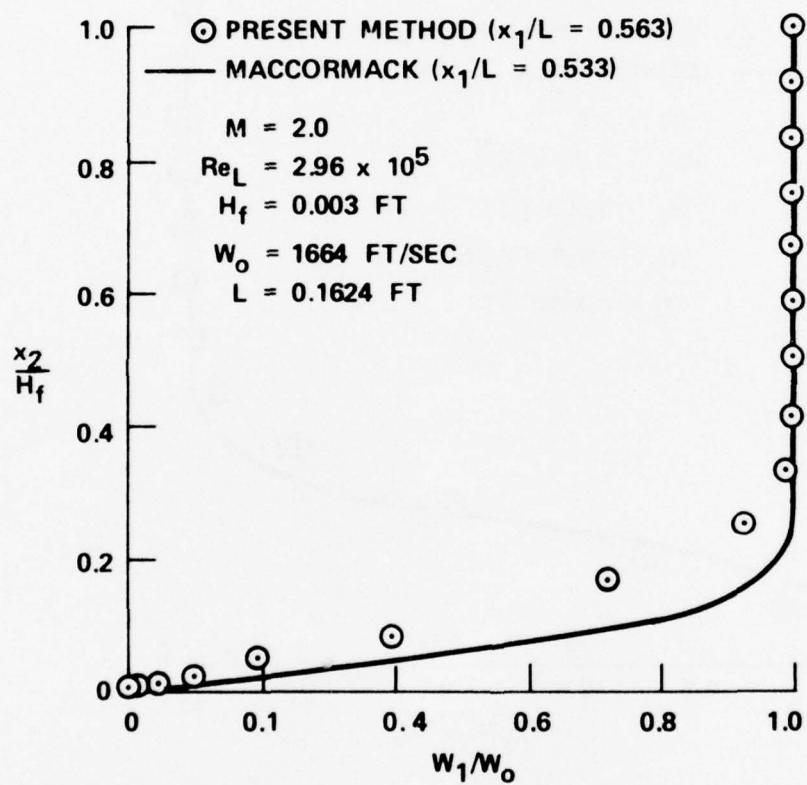


Figure 13. Velocity profile upstream of separation.

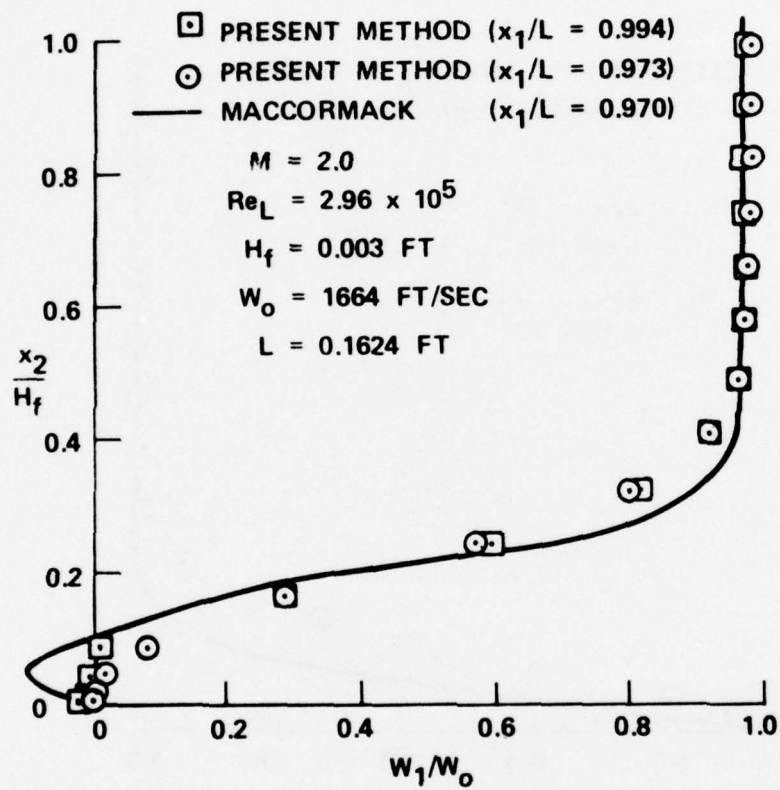


Figure 14. Velocity profile in region of shock wave.

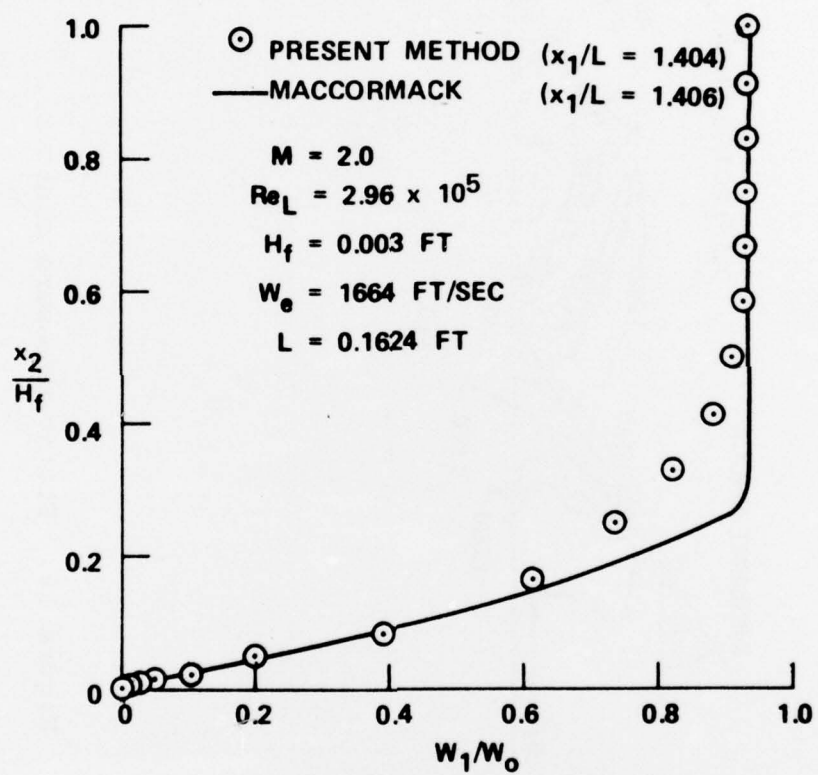


Figure 15. Velocity profile downstream of reattachment.

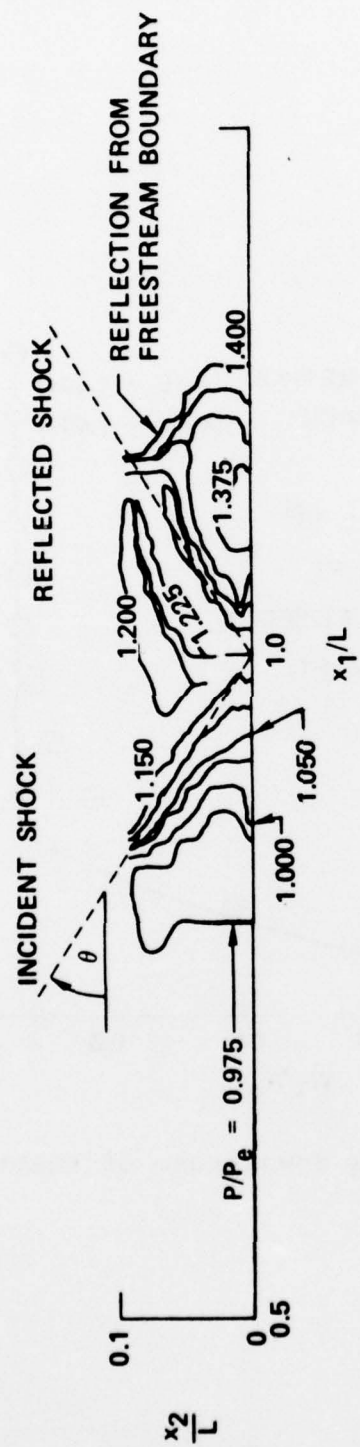


Figure 16. Static pressure contours.

SECTION V HOW TO IMPROVE ACCURACY

Because of time and fiscal constraints, the development of this method halted with the results shown above. It is the opinion of the authors, based on an analysis of the computed results to date, that considerable improvements in accuracy are possible by improving the accuracy of both the relaxation and marching numerical methods. The relaxation nodes need only be a subset of the total grid. However, relaxation is responsible for satisfying the continuity equation. The effects of the increased displacement thickness downstream of the shock are propagated forward through the subsonic region of the flow field, creating the shock spreading observed above. It is of primary importance that this effect be properly accounted for. Figures 17 and 18 show the results of the above calculation on the mass balance of each grid element. In the relaxation grid, agreement is very good; but inside the marching-only region, substantial error occurs. This error comes about from several sources associated with the nonlinearity between the last relaxation node and the wall. A finite-difference approach to numerical method derivation assumes that a match of appropriate derivatives at a point yields a reasonable estimate of the founding equation in the vicinity of the match point. This is only true when variations in dependent variables are nearly linear, or at most quadratic. In a boundary layer, such as shown above, this assumption is in doubt. Substantial errors can occur in the numerical grid transformation, the evaluation of the source term for the relaxation equation, the evaluation of derivatives across the jump, and the like. The continuity error shown in Figures 17 and 18 is the result of the sum total of these effects. Undoubtly, a similar momentum and energy error also occur. The cure is to first produce a marching system that is derived from an integral equation. Thus, it can be integrally conservative through rapid changes in dependent variables. Secondly, an integral average for a source term of the potential equation must be established. As an example, for the position shown in Figure 17, $\nabla \cdot \vec{U}$, part of the source term, differs by a factor of 9 between the wall and the first relaxation node. The net effects of the source are not then well modeled by utilizing its value at the relaxation node, but rather a weighted average across the span would be needed. A grid system that is more nearly uniform would also help. It is obvious when examining the skin friction coefficient, as an example, that for W to deviate so far from linear very near the wall, there must be substantial error in the computation of the numerical transformation functions. The superposition of two fixed-spaced grids, as utilized by MacCormack, would substantially alleviate this difficulty.

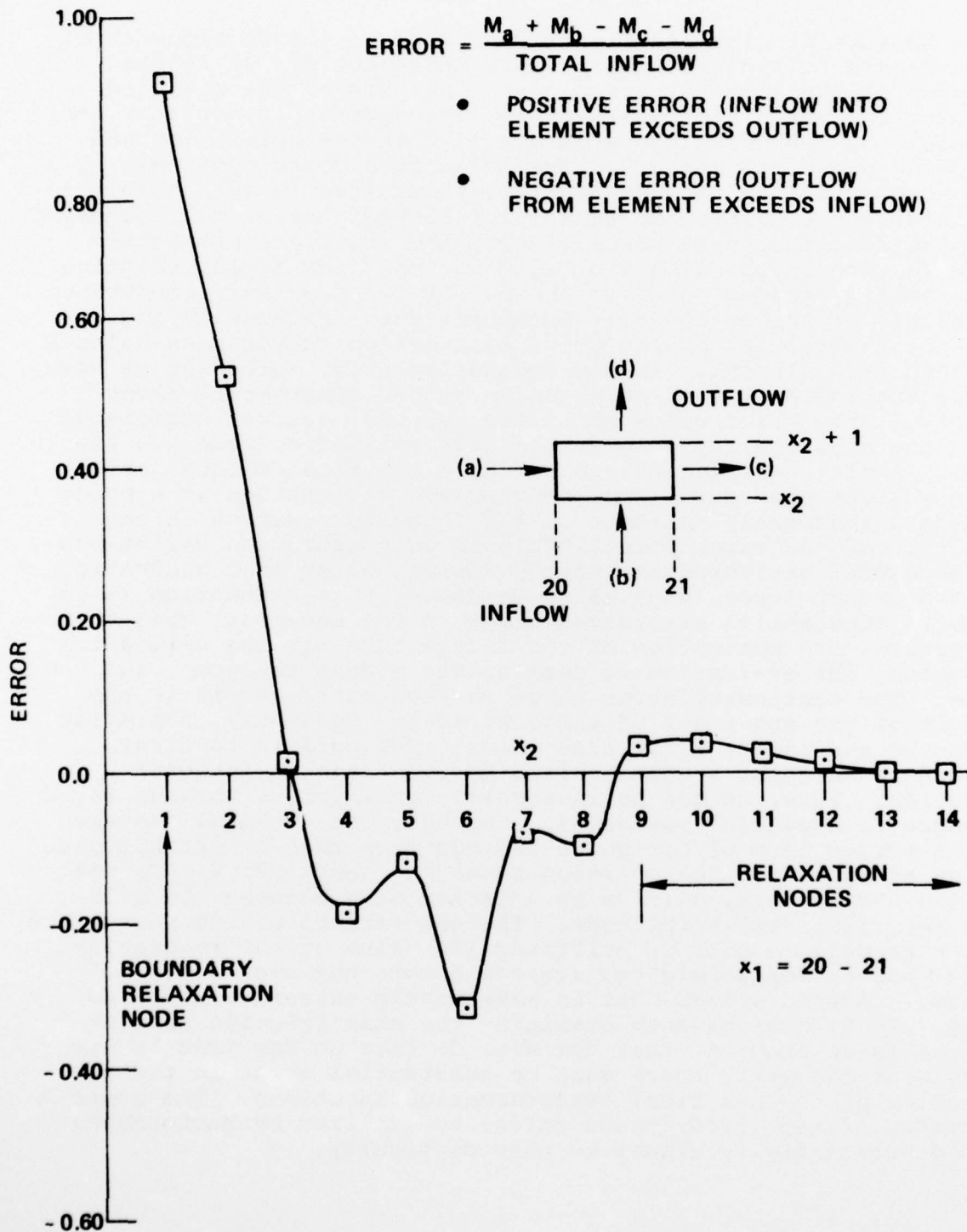


Figure 17. Element continuity error at $x/L = 0.908$.

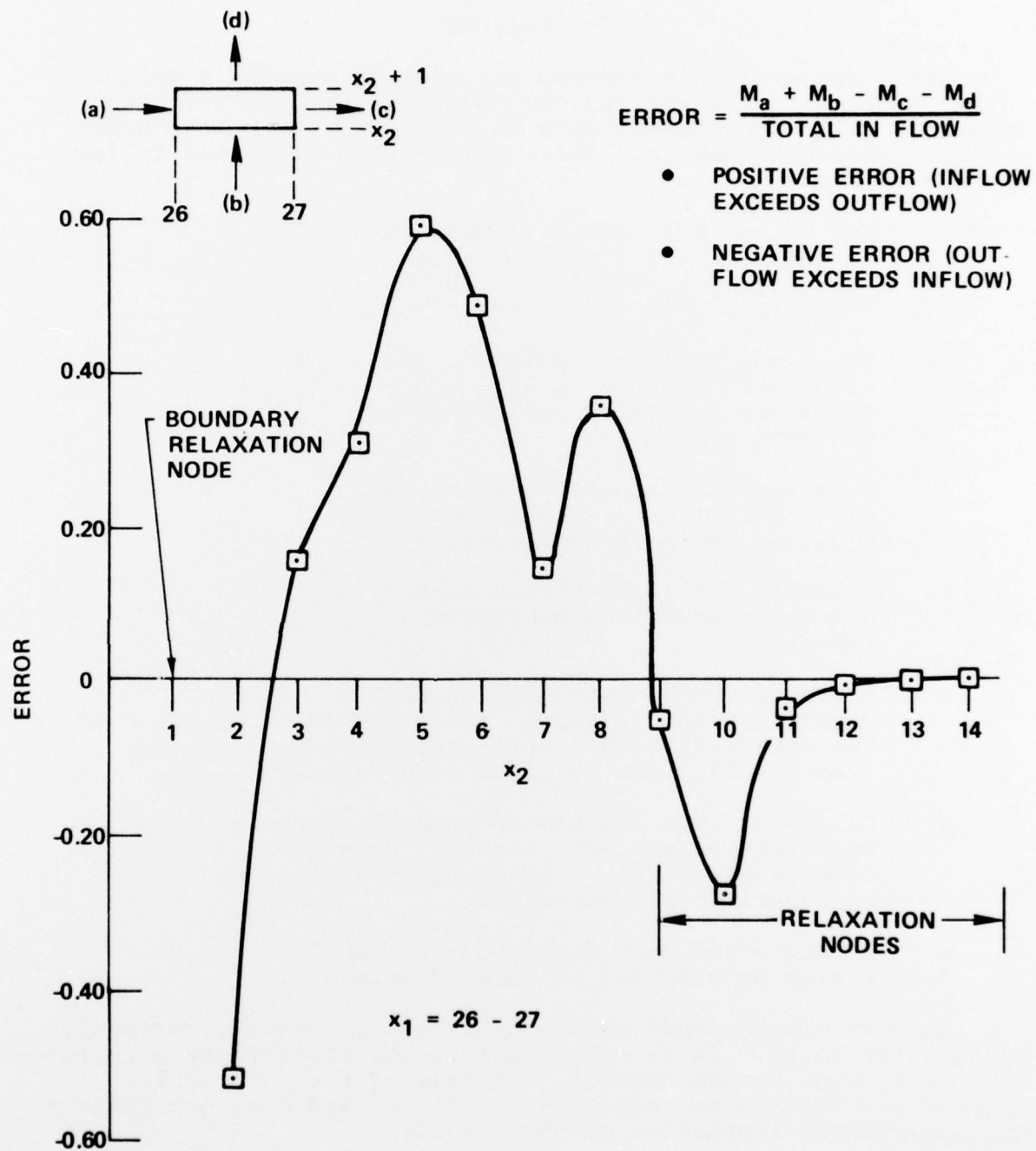


Figure 18. Element continuity error at $x/L = 1.038$.

SECTION VI CONCLUSIONS

The purpose of this program has been to compare a unique method of numerically solving the steady-state Navier-Stokes equations with experimental data unconfused by turbulent model choice. This goal has been met. On the plus side, the following can be stated:

- o The method has met its rapid speed potential, with run times in the vicinity of 1 minute on a CDC 7600 for a 3060 node problem.

Run times could be improved substantially by more careful programming, and better selection of the grid system.
- o A suitable method for solving the mixed elliptic-hyperbolic potential equation has been developed.
- o Iteration between marching and potential equations converged rapidly.
- o A simple method for handling both separation and diffusion terms within the framework of a single-pass marching equation has been utilized successfully.
- o Calculated results are in general agreement with experiment, although pressure rise is not spread over as great a distance as experimentally observed.
- o A method has been developed that could easily be extended to three-dimensions.

Agreement with experimental data could, however, certainly stand improvement. It is expected that the route to such improvement is through a more careful treatment of the numerical details involved in the solution technique, and does not reflect any fundamental limitation of the method.

SECTION VII
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