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**VIBRATION AND STRESS ANALYSIS
OF
SOFT-BONDED SHUTTLE INSULATION TILES**

Modal Analysis With Compact Widely-Spaced Stringers

GRUMMAN



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OF
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Modal Analysis With Compact Widely-Spaced Stringers

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FOREWORD

The work reported herein was performed by the Grumman Aerospace Corporation under the NASA/Langley Master Agreement and Contract No. NAS 1-10635 for the Development and Implementation of Space Shuttle Structural Dynamics Modeling Technology. The Work Statement of Task Order No. 17, "Development of an Analytical Program to Analyze Reusable Surface Insulation for Shuttle", authorized and specified the tasks to be performed in this study. The period of performance was for 15 months starting in June of 1973.

The overall supervision of programs under the Master Agreement is provided by Mr. E.F. Baird, Master Agreement Program Manager. The Task Order No. 17 Project Manager was Dr. I. U. Ojalvo. Many individuals at Grumman contributed to the work reported here. However, the authors wish to specifically acknowledge the efforts of Mrs. Patricia Ogilvie for greatly assisting in the development of the associated computer program.

ABSTRACT

This report describes an efficient iterative procedure for the vibration and modal stress analysis of Reusable Surface Insulation (RSI) of multi-tiled Space Shuttle panels. The method, which is quite general, is rapidly convergent and highly useful for this application. A user-oriented computer program based upon this procedure and titled RESIST (REusable Surface Insulation Stresses) has been prepared for the analysis of compact, widely spaced, stringer-stiffened panels. RESIST, which uses finite element methods, obtains three dimensional tile stresses in the isolator, arrestor (if any), and RSI materials. Two-dimensional stresses are obtained in the tile coating and the stringer-stiffened primary-structure plate. A special feature of the program is that all the usual detailed finite element grid data is generated internally from a minimum of input data. The program can accommodate tile idealizations with up to 850 nodes (2550 degrees-of-freedom) and primary structure idealizations with a maximum of 10,000 degrees-of-freedom. The primary structure vibration capability is achieved through the development of a new rapid eigenvalue program named ALARM (Automatic LArge Reduction of Matrices to tridiagonal form).

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I. INTRODUCTION

The integrity of a reusable space shuttle system is strongly dependent upon protecting the orbiting vehicle from reentry heating in a manner which does not require significant refurbishment between missions. The thermal protection system (TPS) selected to fulfill this need is a non-structural reusable surface insulation (RSI) material shaped into individual tiles which cover almost all of the orbiter surface area. Most of the tiles are square in planform and measure 15.3 or 20.4 cm (6 or 8 inches) on a side, with thicknesses which vary roughly between 1.3 and 7.6 cm (0.5 and 3 inches). Since loss of a single tile could be catastrophic, RSI tile stresses must be accurately determined for the anticipated static, dynamic and thermal environments.

Recognizing that a system's natural vibration modes may serve as basic building blocks for predicting the response to acoustic excitation, this report describes a new iterative procedure for accurately determining tile stresses associated with typical shuttle-panel lower frequency modes. The results of this work must be combined with a dynamic response program, to obtain realistic forced response results.

A user-oriented computer program based upon the present method of analysis was developed. The program, which is titled RESIST (for RReusable Surface Insulation STresses), is capable of computing undamped vibration mode shapes and frequencies of elastically supported multi-tiled panels and determining associated normalized RSI stresses. Typical numerical results from this computer program are presented. A user's manual to facilitate its implementation is presented as Appendix B.

Because of the complex geometry, nonuniform anisotropic material properties, and detailed three-dimensional stress states, the TPS was idealized by finite element assemblages with up to 2500 degrees of freedom per tile. Since a number of tiles affixed to a given structural panel will, in general, interact with one another, application of the standard direct stiffness method would require simultaneous equation systems involving excessive numbers of unknowns. The present iteration scheme overcomes this problem by treating each tile separately. An important byproduct of this approach is that it avoids conditioning problems associated with combining low-stiffness tile isolation ele-

ments ($E \approx 50$ psi) with high-stiffness primary structure elements ($E \approx 10 \times 10^6$ psi). Typical results from the associated computer program reveal a rapid rate of convergence. In addition, a related effort for obtaining tile-stresses associated with statically loaded and heated shuttle panels is presented in Reference 1.

II. TECHNICAL APPROACH

A. STRUCTURAL CONFIGURATION

The design configuration for which an analysis procedure is presented is shown in Figure 1. It consists of a stringer-stiffened flat rectangular panel which supports a nonstructural thermal protection system (TPS).

The TPS is composed of a series of rigidized (RSI) tiles. The tiles are undercut on all four sides to accommodate "filler-strips". The purpose of the nonrigidized filler-strip insulation is to prevent severe heat penetration through the gaps between adjacent tiles.

The RSI material is not bonded directly to the primary structure, but to a thin, stiff, strain arrestor plate first, and then, in turn, to a soft strain isolator material. The function of both these items is to help isolate the primary-structure strains from the low-strength RSI tiles, for the wide range of loading environments which the vehicle must sustain.

B. GENERAL SOLUTION PROCEDURE

Because of the detailed complexity of the structural configuration and the dependence of material properties upon temperature, an analysis technique based upon finite element methods was selected. However, direct application of the standard direct stiffness finite element procedure, to obtain accurate three-dimensional tile stresses, would require equation systems involving excessive numbers of unknowns. The reason for this is that many tiles affixed to a given structural panel may interact. Since each one is a three-dimensional body requiring a detailed structural idealization, their simultaneous consideration requires the solution of large systems of equations. To overcome this problem a rapidly convergent iteration scheme, which treats each tile separately, was developed. The logic flow for an efficient computer program which employs this procedure is presented in Figure 2. The basis for the method is that the TPS is nonstructural but its stress levels, which are critical, must be computed. Thus, it becomes possible to neglect the stiffness of the TPS initially, but not its mass, to determine approximate primary structure modal deflections associated with the vibrations of the overall system. In these initial calculations, the tiles are assumed rigid in shear and thickness-stretching for the purpose of computing their kinetic energies.

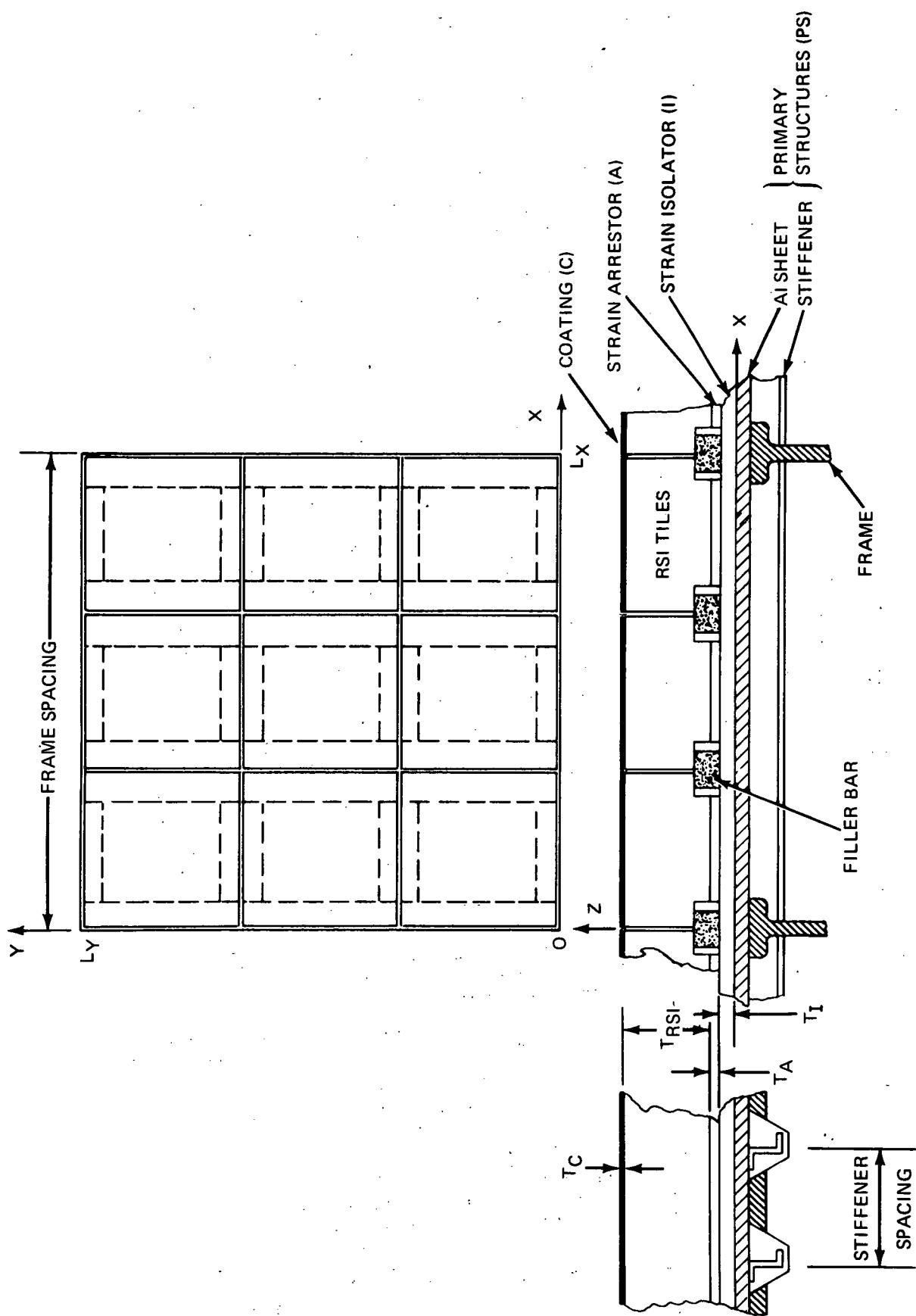


Figure 1. Typical Design Configuration for Shuttle Thermal Protection System

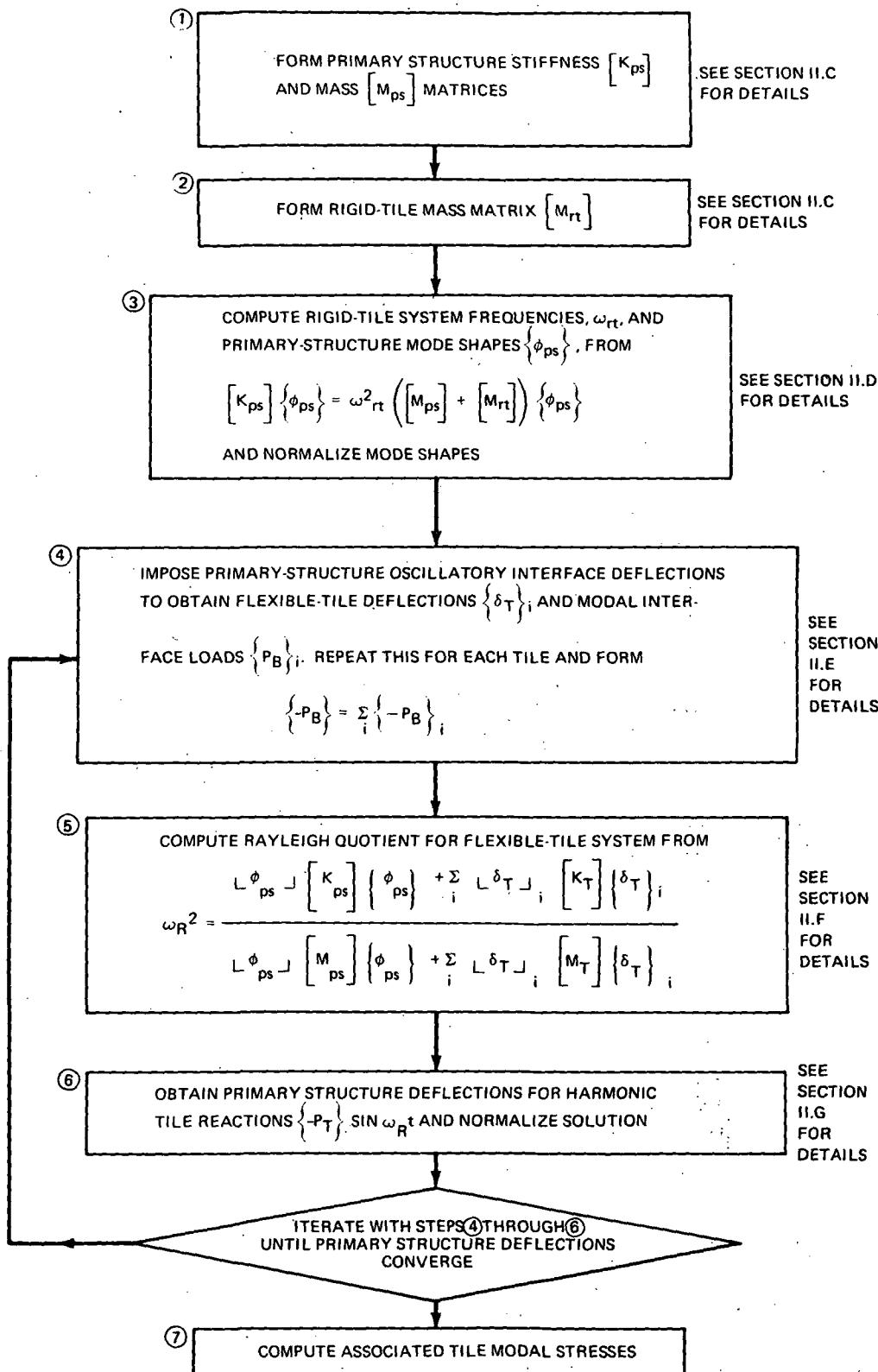


Figure 2. Logic Diagram for Solution Procedure

An interative procedure is then performed where, for each step, the primary structure oscillatory modal deflections are imposed individually upon each tile at the tile/primary-structure interface, and the tile deflections and interface boundary loads are obtained. The frequency is then updated by computing a Rayleigh Quotient, using the latest non-rigid tile displacements in addition to the corresponding primary structure displacements. The individual tile boundary loads obtained are then assembled and their reactions applied to the primary structure. New primary structure deflections are obtained and compared to the previous set. This process is repeated until convergence is obtained. By this method each tile is temporarily assumed uncoupled from all others. Although this is not strictly true, the coupling involved is sufficiently weak so as to ensure accurate approximate results and rapid convergence as well.

It should be noted that an important byproduct of the present tile-by-tile solution procedure is that, unlike the direct stiffness method, it avoids possible numerical precision problems associated with directly combining a low finite element stiffness ($E = 50$ psi) with a high primary structure stiffness element ($E = 10^7$ psi).

C. STRUCTURAL IDEALIZATION

A pictorial representation of the four finite element types incorporated in the RESIST computer program is presented in Appendix C. A brief description of these elements and how each type is used in the overall structural idealization is presented below.

Primary Structure

The primary structure stiffness idealization is based upon two finite elements contained in the Grumman program library described in Ref 2. The panel surface is represented by flat rectangular membrane-bending elements. The elements possess three deflections and two rotations at each node, for a total of 20 independent nodal degrees of freedom per element. The stiffeners are represented by beam elements which bend, stretch and twist. These elements possess 6 degrees of freedom per node, for a total of 12 degrees of freedom per element.

In general, the stiffeners do not attach to the plate nodes at the beam centroids. The stringer attachment points actually used coincide with the plate node which is closest to the centroid of each beam element. The assumption associated with this connection is that the attachment point and beam centroid

move as two points on the same rigid body. Furthermore, the principal axes of the beam cross section may make an arbitrary angle with the plate surface. Refer to Appendix C for further details on the manner of attachment of beam elements to plate nodes.

The translational mass properties of each rectangular plate element are divided equally and lumped at their middle surface node points. The stiffener translation mass properties of each beam element are also divided equally and lumped at their two centroidal nodes. The rotational inertias of the primary structure beam and plate elements are ignored. The assembled structure is rectangular in planform and is assumed to be statically supported with uniform, but different, arbitrary boundary conditions along each edge.

The above idealizations and assumptions appear consistent with the degree of detail required for obtaining accurate dynamic responses of typical aerospace stiffened panel construction.

Thermal Protection System (TPS) Tiles

The basic stiffness element used for the RSI tiles and strain isolator is the anisotropic hexahedron, as defined in References 3 and 4. The version of this element which is being used contains 3 deflection degrees of freedom per node and 8 nodes per element (the more general version which is available permits up to 20 nodes per element). A typical undercut tile with element and node numbering is shown in Figures 3 & 4. The top layer of each tile also has a coating material which is idealized by rectangular membrane elements. The bottom two layers of each tile comprise the strain isolator and strain arrestor, respectively. The undercut regions of each tile are filled with non-rigidized insulation material. This contributes a small amount of mass to the structure but negligible stiffness and negligible effect upon the stresses within the rigidized tile material. Thus, for purposes of computing the tile deflections and stresses, the undercut regions are treated as empty voids.

Mass Matrix Calculations

Two different mass matrices for the TPS must be computed. The first mass matrix required for the TPS "HEX" element and used in Steps 4 and 5 of the logic flow diagram (Figure 2), is based upon lumping 1/8 of each element's mass at each of its 8 corner nodes for x, y and z translational motion. The second type mass matrix, required in Step 2 of Figure 2, is necessary for computing the ap-

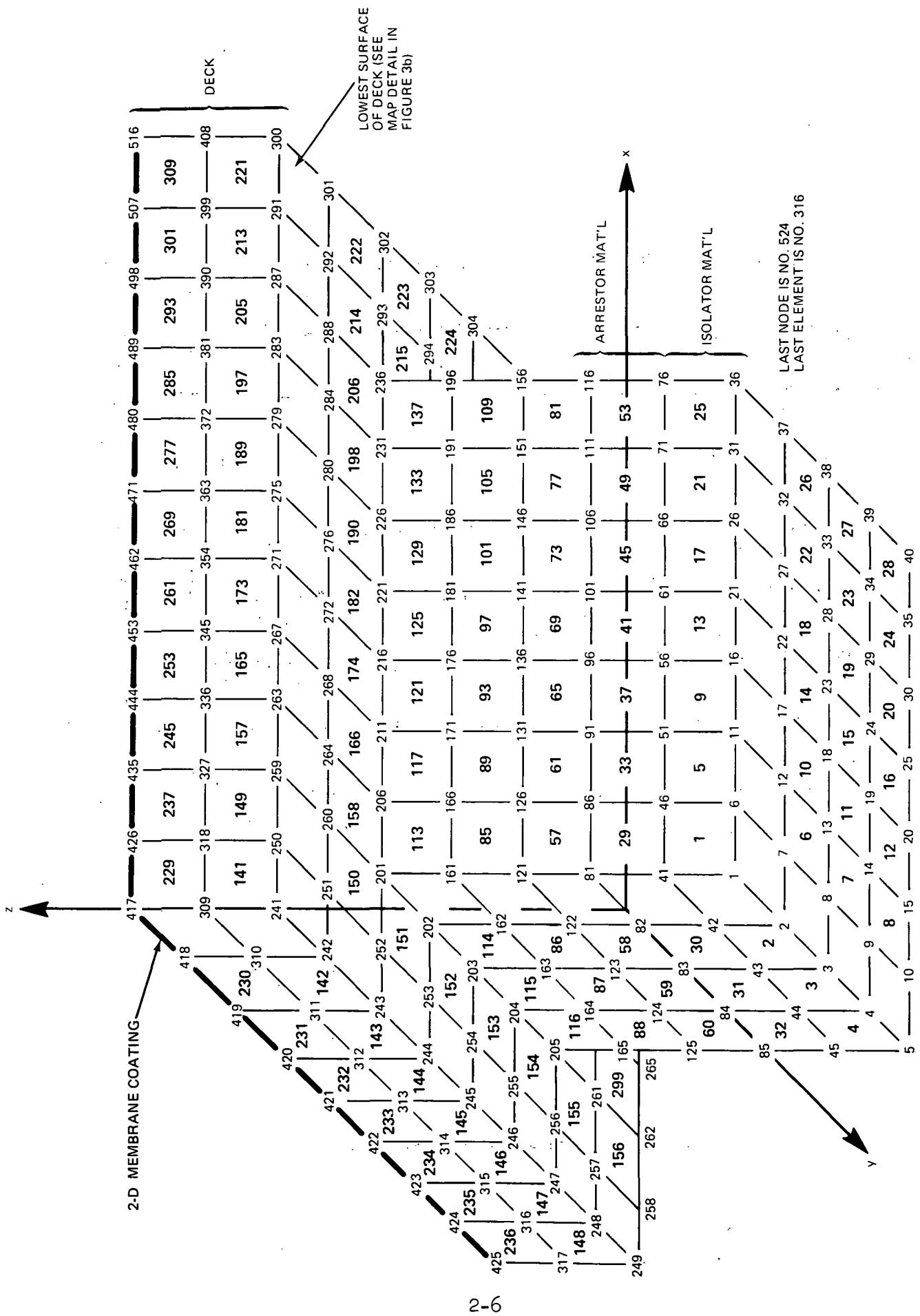


Figure 3. Typical Finite Element Idealization of Shuttle RSI Tile

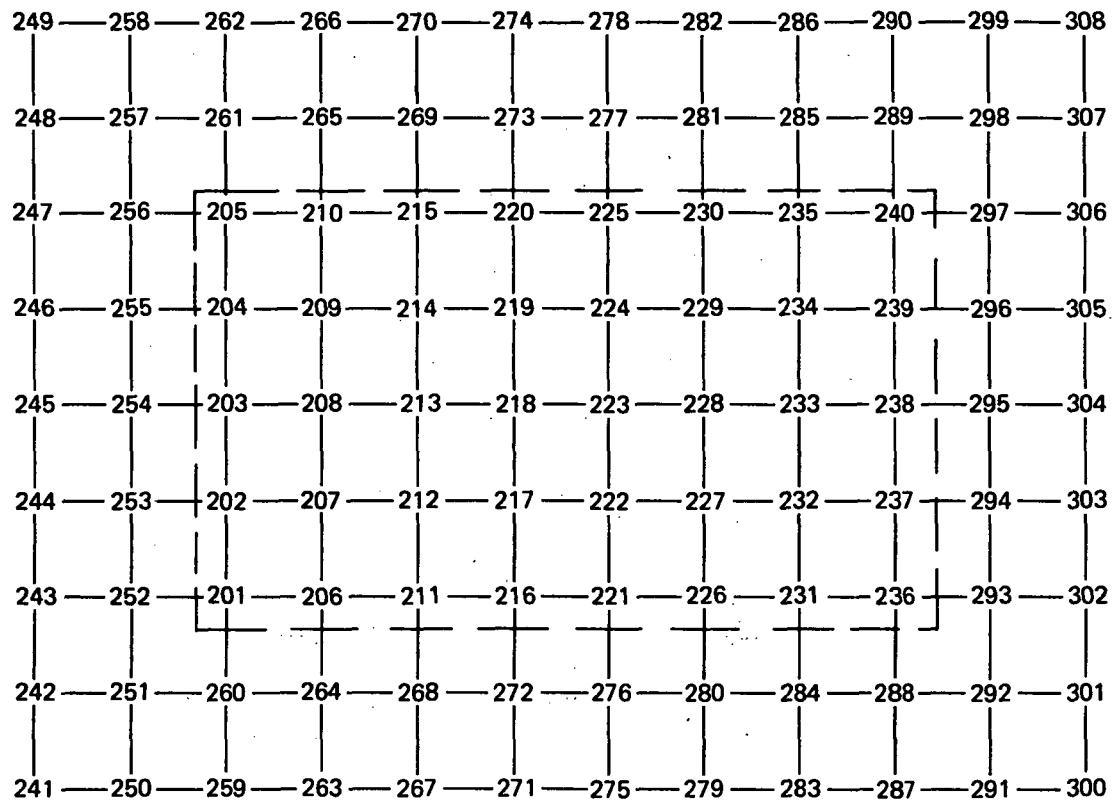


Figure 4. Top View of Lowest Surface of Deck

proximate mode shapes and frequencies (Step 3 of Figure 2) based upon certain assumptions. These assumptions are that the TPS behaves as a series of rigidly attached masses which contribute inertia, but afford no structural stiffness to the primary structure. Analytical details associated with the latter procedure are described as follows.

Let u , v and w be the displacements of a generic TPS point above a corresponding point on the primary structure plate (p) with displacements u_p , v_p and w_p . Then under the assumption of a rigid-in-shear and thickness-stretch TPS we obtain

$$u = u_p - z \frac{\partial w_p}{\partial x}$$

$$v = v_p - z \frac{\partial w_p}{\partial y}$$

$$w = w_p$$

(i.e., normals to the plate middle surface remain normal).

For a section of TPS above a typical plate element, the kinetic energy is given by

$$\begin{aligned} T &= 1/2 \int_A \int_z \rho (\dot{u}^2 + \dot{v}^2 + \dot{w}^2) dz dA \\ &= 1/2 \int_A \int_z \rho \left\{ \dot{u}_p^2 - z \dot{u}_p \frac{\partial \dot{w}_p}{\partial x} + z^2 \left(\frac{\partial \dot{w}_p}{\partial x} \right)^2 \right. \\ &\quad \left. + \dot{v}_p^2 - 2z \dot{v}_p \frac{\partial \dot{w}_p}{\partial y} + z^2 \left(\frac{\partial \dot{w}_p}{\partial y} \right)^2 \right. \\ &\quad \left. + \dot{w}_p^2 \right\} dz dA \end{aligned}$$

where A is the planform area surrounding a node.

Defining the integrals in the kinetic energy equation as follows (refer to Figure 5):

$$\int_z \rho dz = \sum_{i=1}^3 \rho_i (z_{i+1} - z_i) \equiv \overline{\rho h_1}$$

$$\int_z \rho z dz = \sum_{i=1}^3 \rho_i \left(\frac{z_{i+1}^2 - z_i^2}{2} \right) \equiv \overline{\rho h_2}$$

$$\int_z \rho z^2 dz = \sum_{i=1}^3 \rho_i \left(\frac{z_{i+1}^3 - z_i^3}{3} \right) \equiv \overline{\rho h_3}$$

and computing

$$\frac{d}{dt} \left\{ \frac{\partial T}{\partial (\dot{u}_p, \dot{v}_p, \dot{w}_p, \frac{\partial w}{\partial x}, \frac{\partial w}{\partial y})} \right\}$$

then yields for the TPS mass matrix, per plate node

$$A \begin{bmatrix} \overline{\rho h_1} & 0 & 0 & -\overline{\rho h_2} & 0 \\ \overline{\rho h_1} & 0 & 0 & -\overline{\rho h_2} & 0 \\ \overline{\rho h_1} & 0 & 0 & 0 & 0 \\ \overline{\rho h_3} & 0 & & & \\ \overline{\rho h_3} & & & & \end{bmatrix} \begin{Bmatrix} \dot{u}_p \\ \dot{v}_p \\ \dot{w}_p \\ \frac{\partial \dot{w}}{\partial x} \\ \frac{\partial \dot{w}}{\partial y} \end{Bmatrix} \equiv \begin{bmatrix} \bar{M} \end{bmatrix} \begin{Bmatrix} \ddot{\delta}_p \end{Bmatrix}$$

symmetric

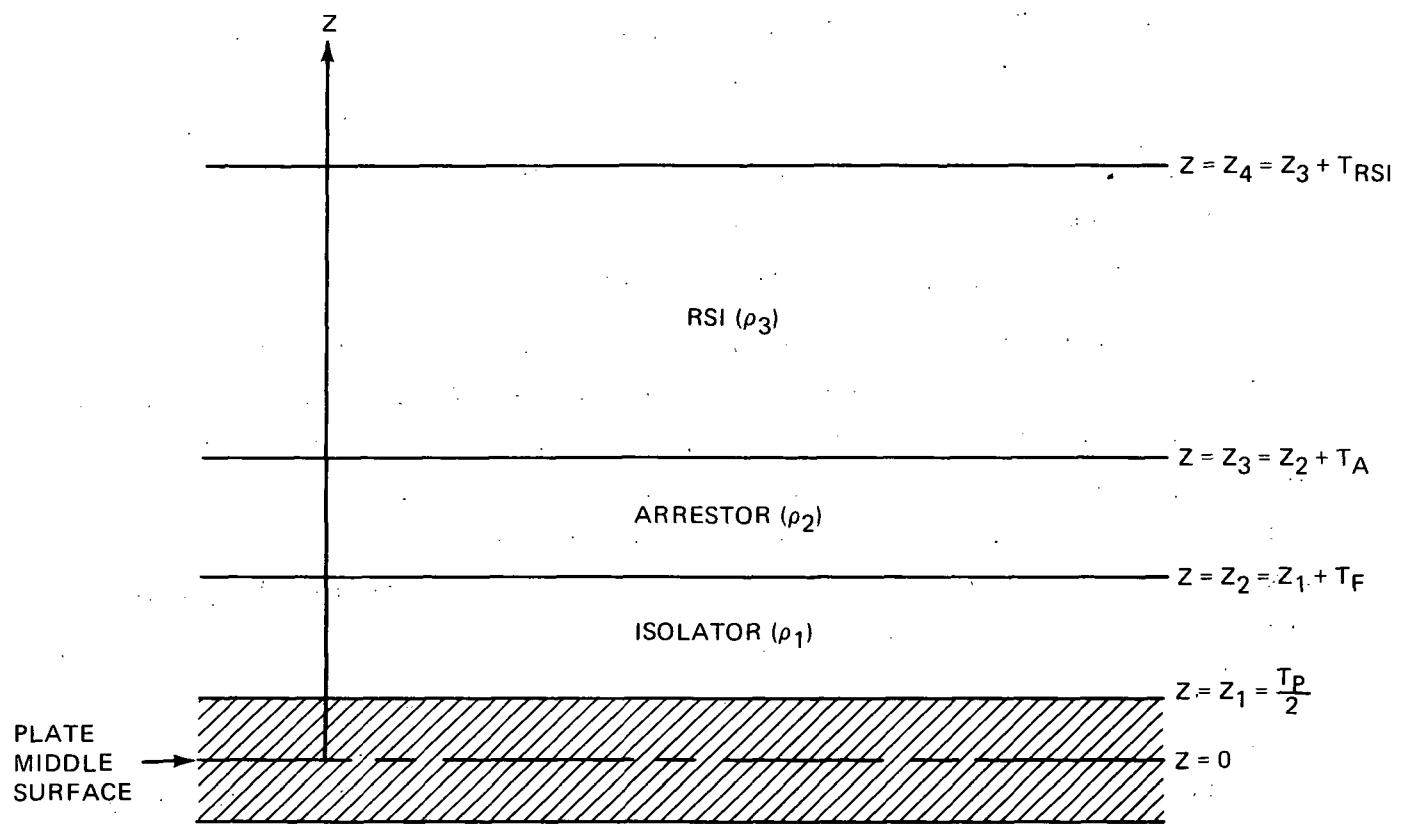


Figure 5. TPS Z-Coordinates for Computation of Rigid Mass Matrix

Automatic Grid Generators

To reduce the quantity of input data significantly, two automatic mesh generating schemes were developed. The first procedure generates the primary structure geometry, node and element numbering system, material properties, and boundary conditions; while the second does the exact same thing for the TPS. Each requires a minimum of information, such as overall dimensions stringer pitch, material properties (as a function of temperature) and TPS material temperatures and element properties from which data may be interpolated (see Appendix B for details).

D. APPROXIMATE FREQUENCIES AND MODE SHAPES

This phase employs the primary-structure mass and stiffness matrices (the latter being in lower triangular decomposed form L , where $K = LL^T$) as well as the rigid-tile approximate mass matrix, just described, to obtain approximate mode shapes and frequencies. Since the primary structure idealizations are quite refined, the number of degrees of freedom will be in the thousands with 10,000 as an upper limit. Thus, an efficient and reliable eigenvalue reduction scheme is required.

The reduction procedure employed was suggested by Crandall⁽⁵⁾ and is based upon an algorithm for accommodating the full matrix due to Lanczos⁽⁶⁾. However, the basic scheme is numerically unstable. The essential improvements necessary to correct this weakness were made by Ojalvo and Newman⁽⁷⁾, who were the first to develop a successful reduction scheme for large scale problems. The work was further streamlined by Newman and Pipano⁽⁸⁾ through the use of banded matrix packing techniques and the computation of convenient error bounds on frequency squared.

The present program, which is a synthesis of these works, is called ALARM (Automatic Large Reduction of Matrices). It is designed to operate efficiently through use of the Grumman Comprehensive Matrix Algebra Procedures (COMAP) interface programs developed by C. Wilkie and F. Nolan.

A mathematical statement of the original vibration problem considered is

$$[K] \{\phi\} = \omega^2 [M] \{\phi\} \quad (1)$$

where $n \times n$ stiffness matrices $[M]$ and $[K]$, which are generally highly banded and sparse, may be singular but non-negative, with the only restriction being $n \leq 10,000$. The associated reduced eigenvalue problem is symmetric, tri-diagonal and of the form⁽⁷⁾

$$\begin{bmatrix} \alpha_1 & \beta_1 & & & \\ \alpha_2 & \beta_2 & & & \\ \cdot & \cdot & \ddots & & \\ \cdot & \cdot & \cdot & \ddots & \\ \alpha_{m-1} & \beta_{m-1} & & & \\ \alpha_m & & & & \end{bmatrix} \begin{Bmatrix} y \end{Bmatrix} = \lambda \begin{Bmatrix} y \end{Bmatrix} \quad (2)$$

where m is typically several orders of magnitude smaller than n , and $\lambda = 1/\omega^2$.

The algorithm used for obtaining the α_i and β_i of Eq. (2), as will be shown, is based upon the power method⁽⁹⁾. Unlike the power method, however, in which each vector is discarded as a new one is generated, the present procedure retains each vector V_i after orthogonalizing and normalizing it to all previous vectors such that $\langle V_i | V_j \rangle = \delta_{ij}$, where δ_{ij} is the Kronecker delta. Thus, the involvement of the original system's lower modes is strengthened. A Rayleigh-Ritz approximation, using the $n \times m$ matrix $[V]$ whose columns are the V_i , follows from the substitution of

$$[L]^T \{\phi\} \equiv [V] \{y\} \quad (3)$$

into Eq. (1) where

$$[K] = [L] [L]^T \quad (4)$$

Premultiplication of the resulting equation by $\lambda [V] [L]^{-1}$ then yields Eq. (2). This result is now solved (in core) by Sturm-Sequencing for the λ 's and power iterations (as described on page 622 of Reference 9) for the eigenvectors.

Experience has shown⁽⁷⁾ that over half the number of modes, $\frac{m}{2}$, obtained by this method are "exact" with a sharp drop-off in accuracy somewhere around $\frac{2m}{3}$; thus, the program usually sets m to $2q+1$, where q is the number of desired modes. Newman-Pipano error bounds⁽⁸⁾ are useful indicators of where a loss of accuracy occurs. The computational steps for obtaining the approximate frequencies and mode shapes are presented in the Appendix A.

E. FLEXIBLE TILE SOLUTIONS

The tiles are originally treated as nonstructural mass items. However, since their modal stress states are desired, it is necessary to eventually account for their flexibility. This is achieved by imposing the primary-structure/TPS interface deflections $\{\delta_B\}_i \sin \omega t$, to each tile and computing the associated tile deflections, $\{\delta_A\}_i \sin \omega t$, to obtain tile stresses and reactions, $\{-P_B\} \sin \omega t$. As described earlier, this is repeatedly done in an iterative manner until convergence is obtained.

Referring to Figure 6, the associated steady-state partitioned matrix equations governing each tile are

$$\begin{bmatrix} K_{AA} - \omega^2 M_{AA} & K_{AB} \\ K_{BA} & K_{BB} - \omega^2 M_{BB} \end{bmatrix} \begin{Bmatrix} \delta_A \\ \delta_B \end{Bmatrix}_i = \begin{Bmatrix} 0 \\ P_B \end{Bmatrix}_i \quad (5)$$

where use has been made of the fact that the tile mass matrix is diagonal.

This equation is then decomposed into

$$[K_{AA} - \omega^2 M_{AA}] \{\delta_A\}_i = -[K_{AB}] \{\delta_B\}_i \quad (6)$$

and

$$-[K_{BA}] \{\delta_A\}_i - [K_{BB} - \omega^2 M_{BB}] \{\delta_B\}_i = \{-P_B\}_i \quad (7)$$

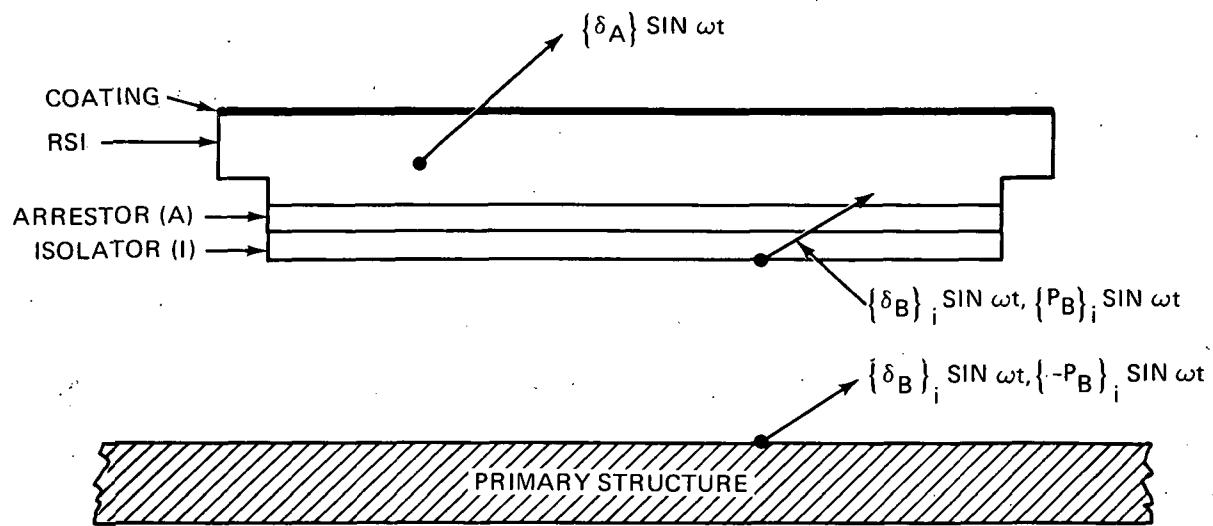


Figure 6. Notation for Flexible Tile Solution Approach

Prior to solution of Eq. (6) for $\{\delta_A\}_i$, it is noted that $[K_{AA} - \omega^2 M_{AA}]$ is not positive definite if any of the eigenvalues, λ_A , of

$$\lambda_A [K_{AA}] \{\delta_A\} = [M_{AA}] \{\delta_A\}$$

satisfy the equation

$$\lambda_A \leq \frac{1}{\omega^2} \quad (8)$$

where ω is the current (approximate) frequency value.

Since the efficient solution routine⁽⁹⁾ that is used for large banded symmetric equations, such as Eq. (2), requires positive definiteness, Eq. (6) is premultiplied by $(K_{AA} - \omega^2 M_{AA})$ to insure this property. Thus, the equation treated to obtain $\{\delta_A\}_i$ is

$$[K_{AA} - \omega^2 M_{AA}]^2 \{\delta_A\}_i = [K_{AA} - \omega^2 M_{AA}] \{-K_{AB} \delta_B\}_i \quad (9)$$

Following this, Eq. (7) is used to compute $\{-P_B\}_i$. Once convergence of the primary structure deflections has been achieved, the corresponding tile modal stresses are computed from the hexahedron stress recovery equations (3).

F. FREQUENCY UPDATE

Since the modes and frequencies obtained originally ignored the tile stiffness and true displacements, the resulting frequencies are only approximate. To correct these, the frequencies are updated through use of a Rayleigh Quotient calculation in which the latest tile and primary structure deflections are used in the equation

$$\omega_R^2 = \frac{\sum_i [\delta_T]_i \left[K_T \right] \left\{ \delta_T \right\}_i + [\delta_{ps}]_i \left[K_{ps} \right] \left\{ \delta_{ps} \right\}_i}{\sum_i [\delta_T]_i \left[M_T \right] \left\{ \delta_T \right\}_i + [\delta_{ps}]_i \left[M_{ps} \right] \left\{ \delta_{ps} \right\}_i} \quad (10)$$

where the subscripts "T" and "ps" denote tile and primary structure, respectively. Thus, $\{\delta_{ps}\}$ are the latest primary structure deflections, which

originally are the $\{\delta\}$ of Eq. (1), $[K_{ps}]$ and $[M_{ps}]$ are the banded stiffness and mass matrices, respectively, of the primary structure and tile. Thus, the $[K_T]$ and $[M_T]$ matrices of Eq. (10) are related to the stiffness and mass matrices of Eq. (5) as follows:

$$[K_T] = \begin{bmatrix} K_{AA} & K_{AB} \\ K_{BA} & K_{BB} \end{bmatrix} \quad [M_T] = \begin{bmatrix} M_{AA} & 0 \\ 0 & M_{BB} \end{bmatrix}$$

G. PRIMARY-STRUCTURE DEFLECTION UPDATE

The original primary-structure modal deflections are approximate, should be checked and, if necessary, corrected. Employing the latest compatible set of tile and primary-structure deflections, the frequencies are recomputed, to give $\omega_{Rayleigh}$, by the method described in the previous subsection. The steady oscillating tile boundary reactions $\{-P_B\} \sin \omega_R t$ are collected for all the tiles and applied to the primary-structure, to yield new modal deflections, as follows:

$$[K_{ps} - \omega_R^2 M_{ps}] \quad \{\delta_{ps}\} = \{-P_B\} \quad (11)$$

where $\{-P_B\} \sin \omega_R t$ is the assembled load reaction of all the tiles acting upon the primary-structure.

Once again, since the solution routine⁽²⁾ that is used for large, banded symmetric equations requires the positive-definiteness property, Eq. (11) is premultiplied by $[K_{ps} - \omega_R^2 M_{ps}]$ prior to solution for the $\{\delta_{ps}\}$, i.e., the actual equation solved for $\{\delta_{ps}\}$ is:

$$[K_{ps} - \omega_R^2 M_{ps}]^2 \quad \{\delta_{ps}\} = [K_{ps} - \omega_R^2 M_{ps}] \{-P_{ps}\} \quad (12)$$

For consistency, the deflections $\{\delta_{ps}\}$ are then normalized such that

$$\|\delta_{ps}\| \quad \{\delta_{ps}\} = 1$$

prior to comparison to the previous set, to establish convergence. The convergence test employed consists in satisfying

$$\text{max element of } \left\{ \begin{vmatrix} \delta_{\text{ps}} & -\delta_{\text{ps}} \\ \text{present} & \text{previous} \end{vmatrix} \right\} \leq \epsilon \left| \begin{matrix} \text{largest element of } \delta_{\text{ps}} \\ \text{present} \end{matrix} \right|$$

where ϵ is an empirically determined positive input quantity.

III. NUMERICAL RESULTS

A. TILED CONFIGURATIONS

To demonstrate the numerical convergence properties of the proposed iteration scheme, two examples are presented here. The material properties and loading conditions used, while not precise in a specific design sense, are representative of the problem parameters the method is intended to accommodate.

Figure 7 shows the configuration, finite element idealization, material properties, and boundary conditions considered for Example 1. The fundamental frequency, largest primary-structure normalized deflection and average stress components for certain critical RSI elements are presented as a function of iteration number in Table 1. Besides revealing a rapid convergence rate, the results presented indicate a high level of accuracy for the first iterate in that the maximum frequency, deflection and σ_x errors are less than 2%. The errors in σ_z were higher for the first iteration (approximately 20%), but these settled down to under 4% in the second iteration.

Similar observations were made for the more realistic shuttle tile configuration of Example 2 shown in Figure 8. This problem consists of three 6 x 6 inch tiles, each of which is two inches thick and has 1 x 1/2 inch edge undercuts. The primary structure is a .041 inch aluminum plate with offset stringers spaced half an inch apart. The two opposite shorter boundaries are clamped while the longer ends are free.

The fundamental frequency and peak normalized deflection, as a function of iteration number, are shown in Table 2. The primary-structure mode-shape is plotted in Figure 9 for iterations 1, 2, and 4. The corresponding stresses for iterations 1 and 4 are shown in Figures 10 through 12. Running time for this realistic type problem on an IBM 370 averaged 1 1/3 CPU minutes per tile plus a fixed time of 13 minutes per iteration. Thus, four iterations of the three tiled configuration required approximately 68 CPU minutes.

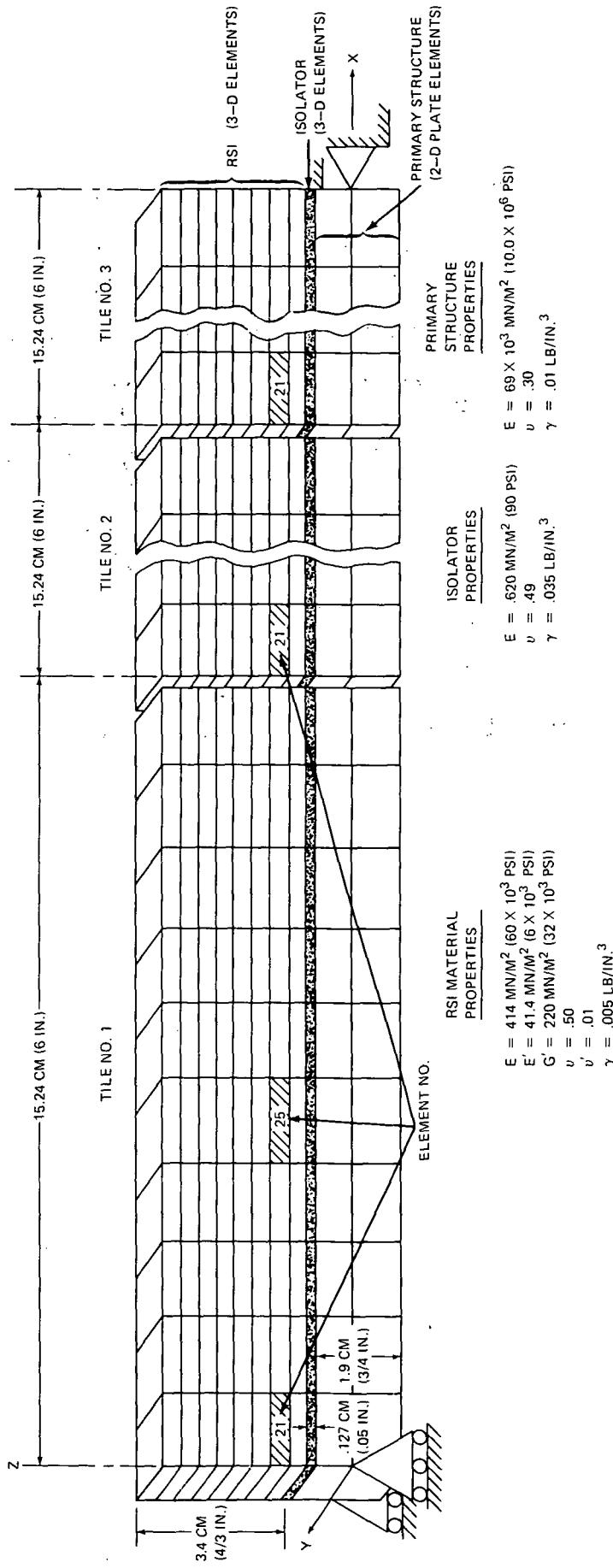


Figure 7. Example 1: Sample Problem Used to Demonstrate Numerical Convergence of Iteration Scheme

TABLE 1
 FUNDAMENTAL FREQUENCY, MAXIMUM DEFLECTION, AND CRITICAL
 TILE STRESSES VS. ITERATION NUMBER (Refer to Figure 7)

ITERATION NO.	ω_1 (Hz)	MAXIMUM NORMALIZED DEFLECTION (INCHES)	RSI STRESSES					
			σ_x (psi) ELEMENT NO. 25			σ_z (psi) ELEMENT NO. 21		
			TILE NO. 1	TILE NO. 2	TILE NO. 3	TILE NO. 1	TILE NO. 2	TILE NO. 3
0	510	0.180	-	-	-	-	-	-
1	425	0.181	-43	-88	-46	22	57	35
2	427	0.181	-40	-86	-43	22	49	28
3	425	0.181	-40	-86	-46	21	48	28
4	425	0.181	-40	-86	-46	22	49	28

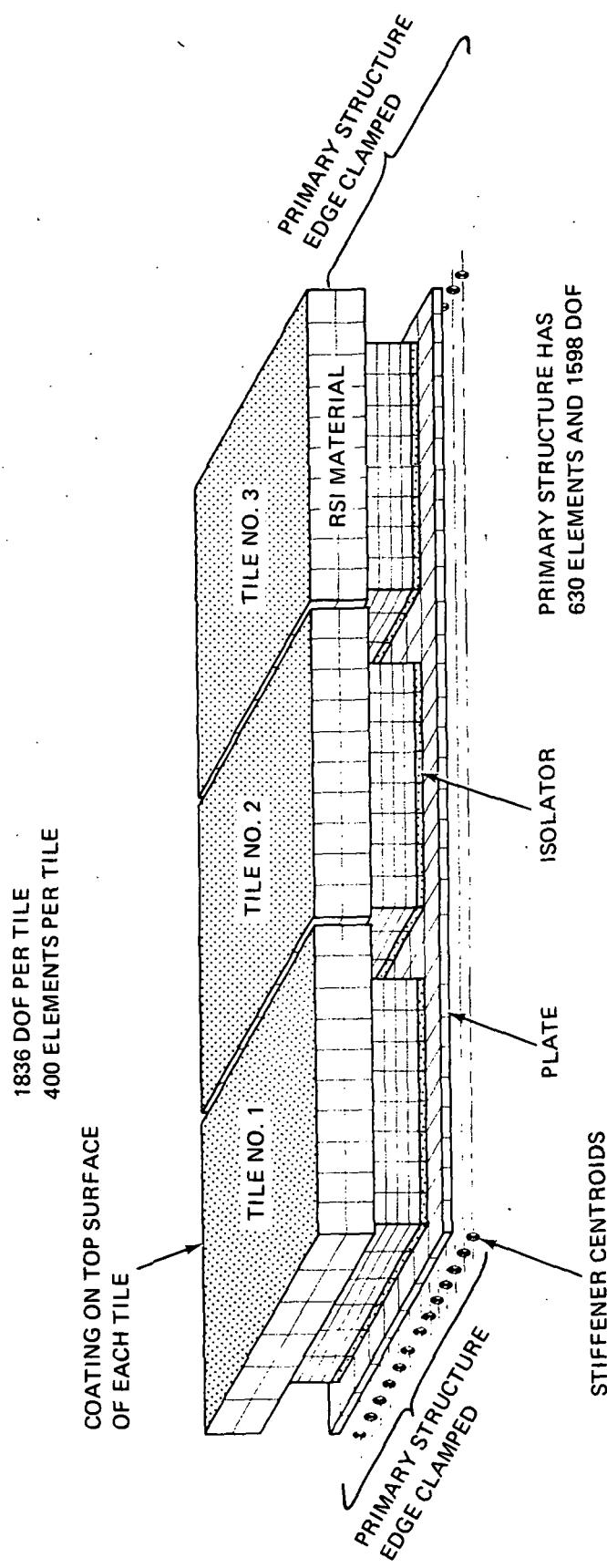


Figure 8. Example 2: Typical Configuration That Computer Program is Capable of Analyzing

TABLE 2

FUNDAMENTAL FREQUENCY AND MAXIMUM PRIMARY STRUCTURE
DEFLECTION AS A FUNCTION OF ITERATION NUMBER FOR
EXAMPLE 2

ITERATION NUMBER	FUNDAMENTAL FREQUENCY (Hz)	NORMALIZED MAX PRIM STRUCT DEFLECTION (inches)
0	146	0.111
1	153	0.108
2	153	0.111
3	154	0.110
4	153	0.110

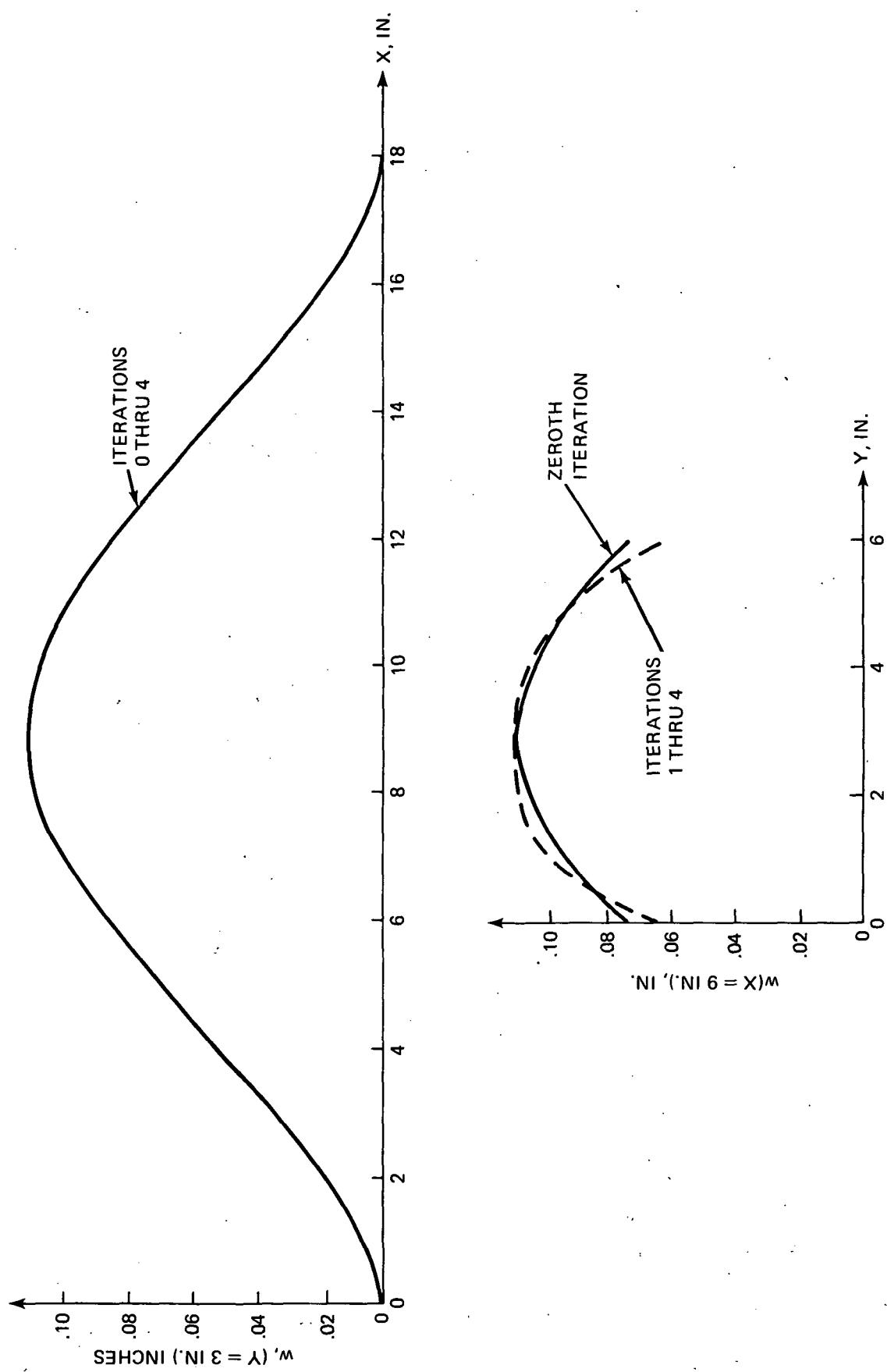


Figure 9. Example 2: Plate Structure Normal Deflections of First Mode

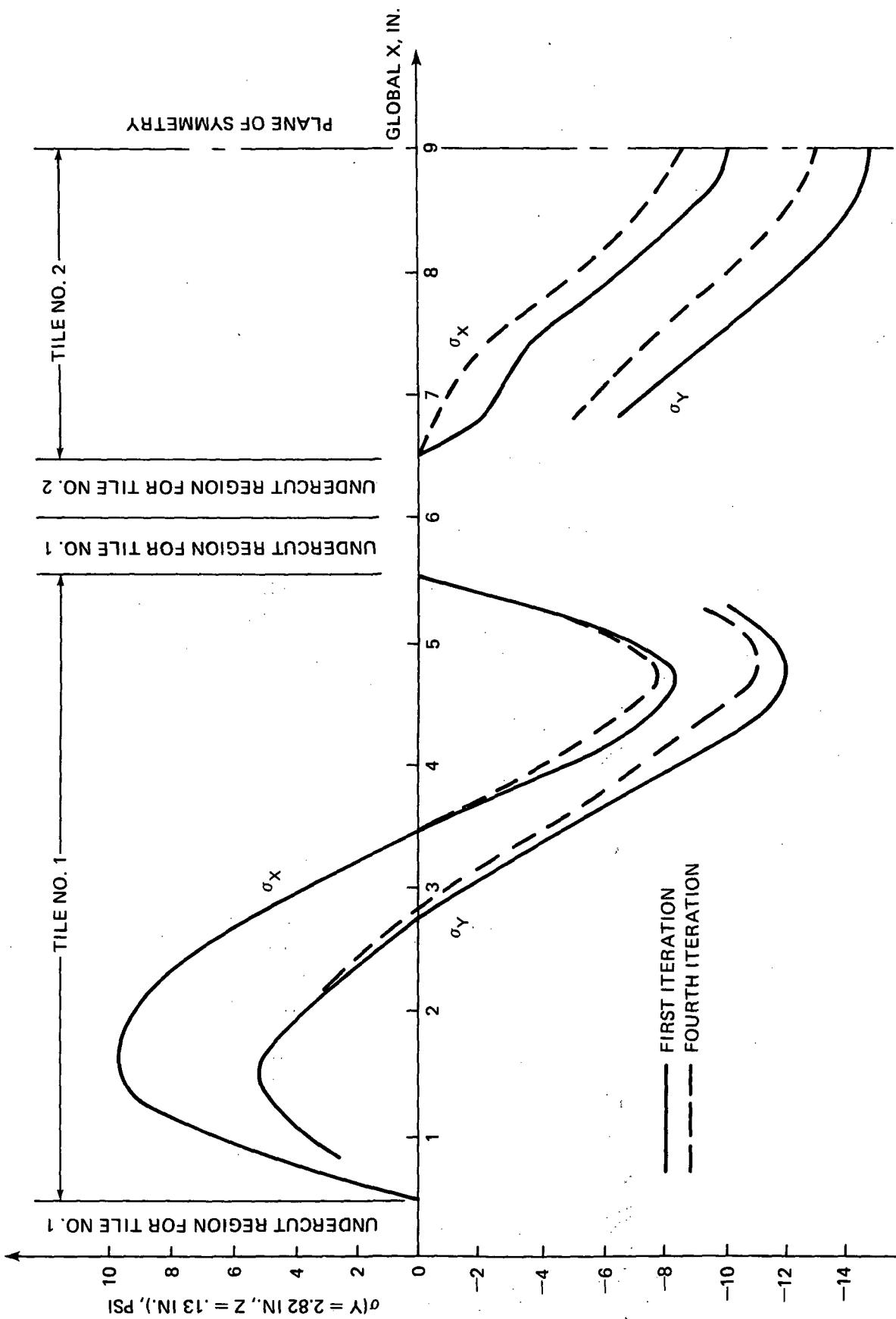
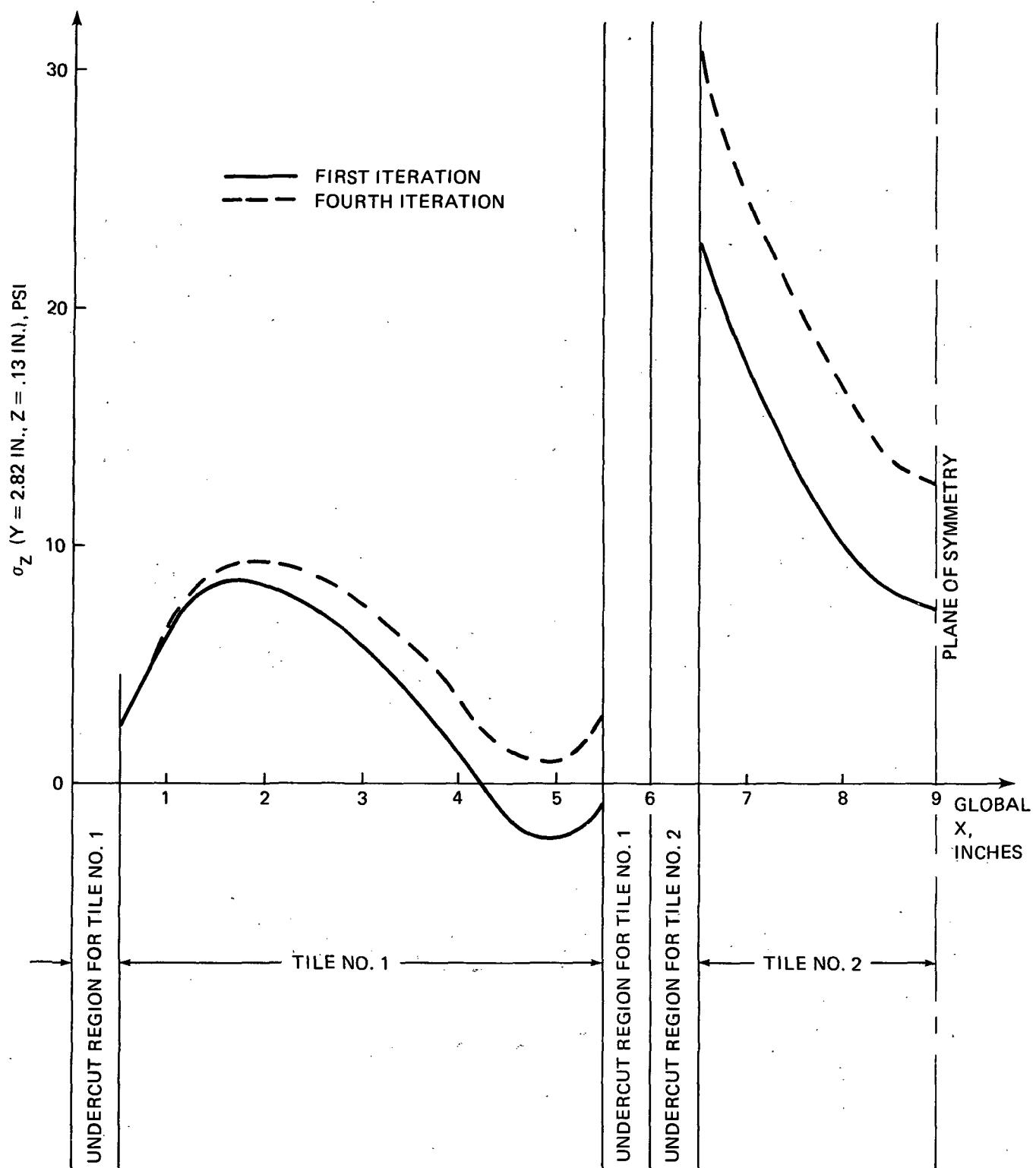


Figure 10. Example 2: Normalized First Mode Direct Stresses for Tiles 1 and 2 in Planer Directions as a Function of Spanwise Coordinate



**Figure 11. Example 2: Normalized First Mode Direct Stresses for Tiles 1 and 2
in Planer Directions as a Function of Spanwise Coordinate**

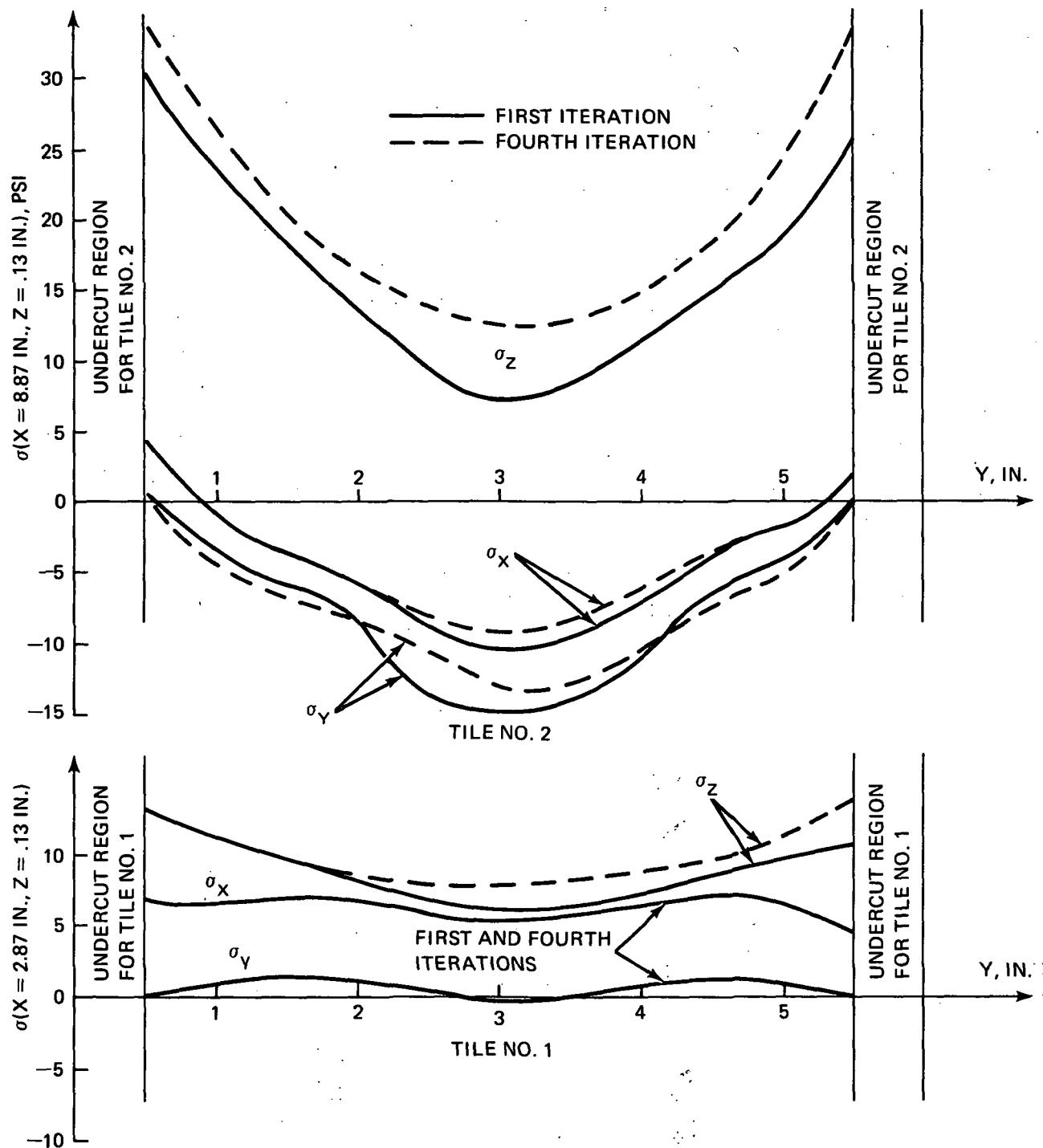


Figure 12. Example 2: Normalized First Mode Direct Stresses for Tiles 1 and 2 as a Function of Chordwise Coordinate

B. UNTILED PRIMARY STRUCTURE

Figure 13a is representative of a typical stiffened untiled "shuttle panel" undergoing dynamic testing at the Langley Research Center. Figure 13b shows the finite element idealization that the RESIST computer program generated for this panel. The numerical results obtained (see Table 3) reveal a tendency for the frequencies to occur in groups of 7. The geometric significance of this clustering is that the modes in each group are related to the physical presence of 7 stringers, which subdivide the panel into 6 bays.

The idealized stringers are assumed to be attached at only a single rivet line. However, in the configuration of Figure 13a, each stringer is attached by two rivet lines to form a closed section torque-tube. Thus, the idealization of Figure 13b is inadequate for predicting the higher cross-stringer modes (i.e. $m > 0$) when the stiffened plate-pitch, a , is not small compared to the unstiffened plate-pitch, b . Since a is of the order of b for realistic shuttle panels, the idealization capability within RESIST should be extended to accommodate closed-section stringers with narrow spacing in addition to the present capability which treats widely spaced, or open-section, stringers.

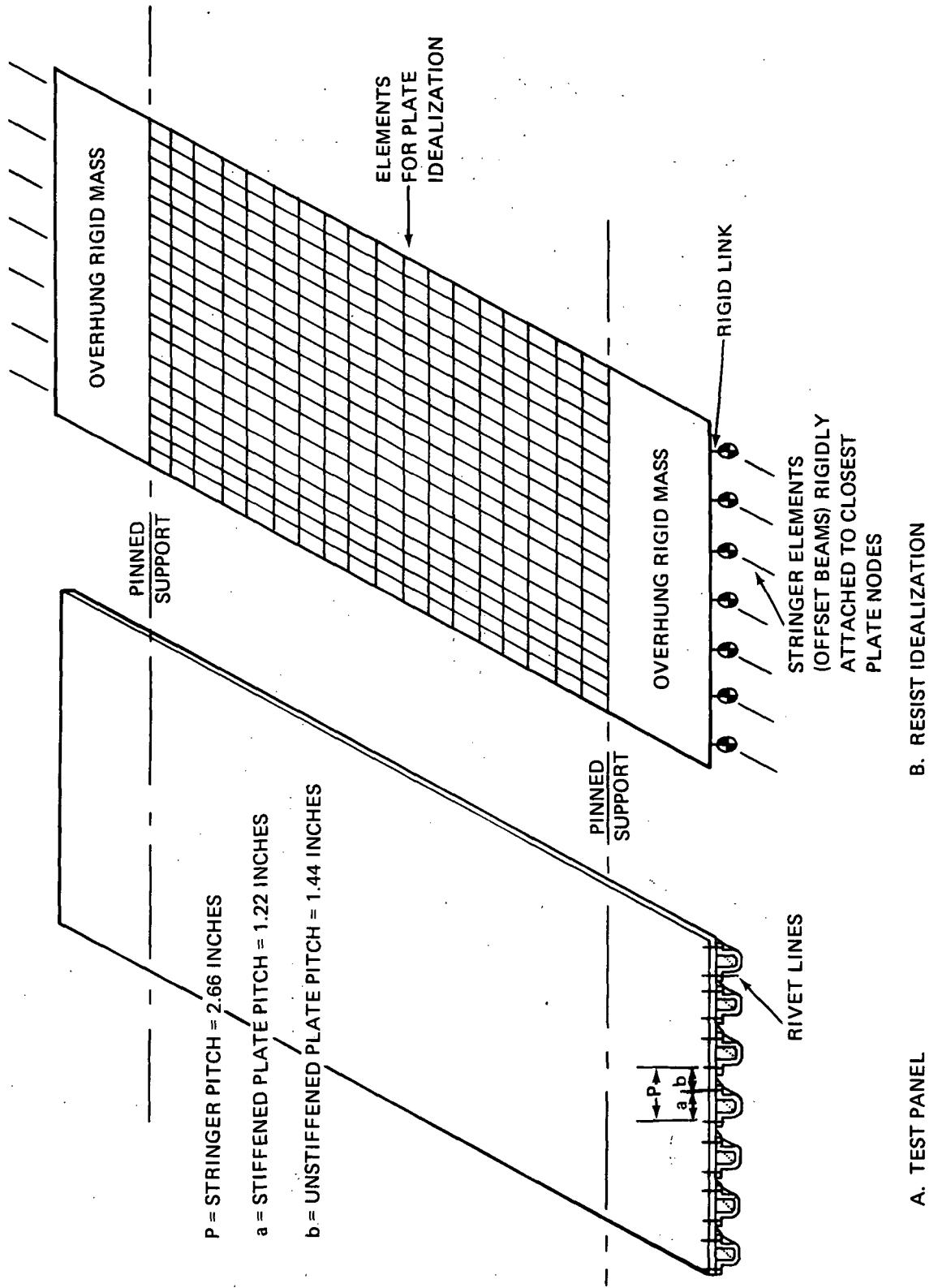


Figure 13. RESIST Idealization of Shuttle Panel for Test at Langley

TABLE 3
NATURAL FREQUENCIES FOR RESIST IDEALIZATION OF
SHUTTLE PANEL FOR TEST AT Langley

Mode Shapes		RESIST Finite Element Analysis
n*	m**	
1	0	106
	1	108
	2	126
	3	136
	4	150
	5	183
	6	203
2	0	286
	1	288
	2	302
	3	315
	4	328
	5	343
	6	346

* n = number of 1/2 sine waves between spanwise supports

** m = number of nodes in cross-stringer direction

IV. CONCLUSIONS AND RECOMMENDATIONS

An iterative procedure for the vibration stress analysis of RSI multi-tiled shuttle panels has been developed. The method, which is quite general, is rapidly convergent and highly useful for this application.

A user-oriented computer program based upon this procedure and titled RESIST has been coded. RESIST, which uses finite element methods, obtains three dimensional tile stresses in the isolator, arrestor (if any) and RSI materials. Two dimensional stresses are obtained in the tile coating and the stringer-stiffened primary structure plate. A special feature of the program is that all the usual detailed finite element grid data is generated internally from a minimum of input data.

The program may be used in an iterative mode to obtain detailed results. However, for parametric design studies, reasonably accurate results may be obtained by using only one iteration at a significant savings in running time. This is achieved by having the program analyze certain specific tiles of a multi-tiled panel only once, while ignoring less critically stressed tiles.

At present the program can accommodate tile idealizations with up to 850 nodes (2,550 degrees-of-freedom) and primary structure idealizations with a maximum of 10,000 degrees-of-freedom. In addition, the tile pattern must begin and end at a structural frame and the isolator material must be isotropic. Should such restrictions, or any similar ones, require alteration as the shuttle TPS design changes, it would appear logical to update RESIST.

To enhance the usefulness of the present work, it should be extended to accommodate closely spaced, closed-cell stiffeners (which are more typical of Space Shuttle panel designs) and to perform acoustic response analyses as well.

V. NOMENCIATURE

A	Plate area associated with a given primary-structure node
[A]	Matrix used in large eigenvalue algorithm ($= [L]^{-1} [M] [L]^{-T}$)
[G]	Matrix used in large eigenvalue algorithm
[K]	Stiffness matrix used in large eigenvalue algorithm
[K _T]	Tile stiffness matrix
[K _{ps}]	Primary structure stiffness matrix
[K _{AA}], [K _{AB}],	Matrix partitions of [K _T]
[K _{BA}], [K _{BB}]	
[L]	Lower triangular decomposition of stiffness matrix
[M _{ps}]	Diagonal primary structure mass matrix
[M _{rt}]	Mass matrix associated with rigid tile
[M _T]	Tile mass matrix
[\bar{M}]	Approximate tile mass matrix associated with a typical plate node
[M]	Mass matrix used in large eigenvalue algorithm
[M _{AA}], [M _{BB}]	Diagonal partition matrices of tile mass matrix
{-P _B }	Reaction loads caused by all tiles acting upon primary structure
{P _B } _i	Load acting upon i th tile caused by primary structure
T	Approximate kinetic energy associated with a typical node
{ $\bar{v}_i^{(s)}$ }	Unnormalized and unrefined Lanczos vector
{ \bar{v}_i }	Refined but unnormalized Lanczos vector
{v _i }	Normalized and refined Lanczos vector

z_i	Thickness coordinates. Refer to Figure 5.
$\{b_i\}$	Vector used in eigenvalue algorithm
$\{g_i\}$	Vector used in eigenvalue algorithm
m	Order of reduced eigenvalue problem
n	Order of unreduced eigenvalue problem
\bar{n}	Number of dynamic degrees of freedom
q	Number of desired approximate system modes
t	Time
u, v, w	Approximate displacement components of arbitrary point within tile
u_p, v_p, w_p	Displacement components of primary structure
x, y, z	Spacial coordinates
$\{y\}, \{y_i\}$	Eigenvectors of reduced eigenvalue problem
y_i^m	Last element of $\{y_i\}$
α_i, β_i	Diagonal and superdiagonal elements of tridiagonal reduced eigenvalue matrix
$\{\delta_{ps}\}$	Primary structure deflections
$\{\delta_A\}_i, \{\delta_B\}_i$	Deflections of i^{th} tile at node points not in contact (A) and in contact (B) with plate
$\{\delta_T\}_i$	Deflections of i^{th} tile
ϵ	Primary structure deflection convergence parameter
λ	Eigenvalue of reduced problem

λ_A	Eigenvalue problem associated with "A" degrees-of-freedom of tile
ρ	Mass density
$\bar{\rho}_i^h$	Integrated mass density parameter
$\{\phi_{ps}\}$	Approximate primary structure mode shape
ω	Frequency of large eigenvalue problem
ω_R^2	Rayleigh Quotient
ω_{rt}	Approximate rigid-tile frequency

VI. REFERENCES

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APPENDIX A

LARGE EIGENVALUE SOLUTION ALGORITHM

The algorithm employed to initially determine the approximate mode shapes and frequencies of the system^(7,8) is outlined below:

1. Initialization: Set $i = 0$ and determine m , the order of the reduced problem as follows:

For $\bar{n} \leq 7$ set $m = \bar{n}$

For $\bar{n} > 7$ set $m = \bar{n}$ if $2q + 1 \geq \bar{n}$

$m = 2q + 1$ if $2q + 1 < \bar{n}$ and $q > 3$

$m = 7$ if $q \leq 3$

"Reverse", transpose, and "reverse" $[L]$ to obtain

$[L]^{RTR}$ and save result for repeated use in Step 6
(see discussion in Step 6).

2. Generate $\{V_0\}$: Use a random number generator for each element.
Go to 6.

3. Compute $\theta_{i-1} = [\bar{V}_i] \downarrow \{\bar{V}_i\}$

4. Test $i \geq m+1$: if no--continue; if yes--go to 15

5. Compute $\{V_i\} = \frac{1}{\theta_{i-1}} \{\bar{V}_i\}$: This step normalizes the vectors such that

$$[V_i] \downarrow \{V_i\} = 1$$

6. Solve for $\{g_i\}$ and save it if $i \geq 1$. This is the solution vector of

$$[L]^T \{g_i\} = \{V_i\}$$

The usual Grumman routine used for solutions of this type equation is titled QBAC (see Reference 10). However, it is not too efficient when there is only one right-hand side, as is our case.

A more efficient procedure is to manipulate $[L]^T$, which is an upper triangular matrix, into lower triangular form, and then to use QFOR to solve the equation

$$[L]^{RTR} \{g_i\}^R = \{V_i\}^R, \text{ where } \{g_i\}^R \text{ and } \{V_i\}^R \text{ correspond to } \{g_i\} \text{ and } \{V_i\}, \text{ respectively, each with the order of its elements reversed (e.g., the first element of } \{g_i\} \text{ is the last element } \{g_i\}^R \text{ and vice versa). The manipulation of } [L]^T \text{ into } [L]^{RTR} \text{ requires two matrix row reversals (COMAP routine REVERS*) and one transpose (TRAN*) operation in the order REVERS-TRAN -REVERS. The most time consuming routine of these three is TRAN, which is even slower than QFOR*. However, since it is only necessary to compute } [L]^{RTR} \text{ once (see Step 1), each solution for } \{g_i\}, \text{ by the present procedure, requires far less running time than by the direct method (e.g., 1/30th the CPU time for a 3,000 DOF system).}$$

7. Compute $\{b_i\}$ from $\{b_i\} = [M] \{g_i\}$. This is actually accomplished by forming $([g_i] [M])^T$ which is considerably more efficient using the COMAP associated subroutines.
8. Solve for $\{A V_i\}$ from $[L] \{A V_i\} = \{b_i\}$ using QFOR.
Note that $[A] = [L]^{-1} [M] [L]^T$. Thus $\{A V_i\}$ is one power iteration of V_i and so accentuates its lower modal content.
9. Test i:
if $i = 0$, set $\bar{V}_1 = A V_0$ and return to step 3.
if $i \neq 0$, continue.

*See Reference 10 for descriptions of these subroutines.

10. Compute $\alpha_i = [V_i] \{A V_i\}$ and step up i by 1.

This step computes the Rayleigh Quotient associated with the vector $\{V_i\}$.

11. For $i = 2$, compute $\left\{ \bar{V}_2^{(1)} \right\} = \{A V_1\} - \alpha_1 \{V_1\}$

12. For $i \geq 3$, compute $\left\{ \bar{V}_1^{(1)} \right\} = \left\{ A V_{i-1} \right\} - \alpha_{i-1} \left\{ V_{i-1} \right\} - \alpha_{i-2} \left\{ V_{i-2} \right\}$
and set $s = 1$.

This step theoretically (7) orthogonalizes $\{\bar{V}_i\}$.

However, any slight numerical inprecision tends to bias the resulting vectors unfavorably through the present iterative generation scheme. The suggested correction to this numerical stability problem is contained in the following two steps.

13. Test $s > \bar{s}$: if no, continue; if yes, set $\{V_i\} = \left\{ \bar{V}_i^{(s)} \right\}$ and return to 3.

14. Replace $\left\{ \bar{V}_i^{(s)} \right\}$ by $\left\{ \bar{V}_i^{(s)} \right\} - \left([V_j] \left\{ \bar{V}_i^{(s)} \right\} \right) \{V_j\}$ for $j = 1$
Continue correcting $\left\{ V_i^{(s)} \right\}$ by repeating the above step for $j = 2, 3, \dots, i-1$. Step up s by 1 and return to 13.

15. Solve for the lowest $3/4^{\text{ths}}$ (rounded to an integer) of the frequencies of Eq. 2 using Sturm sequencing (9).

16. Solve for the eigenvectors associated with the highest "q" eigenvalues λ , each normalized such that

$$[y] \{y\} = 1$$

17. Compute and print $\omega_i = \lambda_i^{-1/2}$ and $f_i = \frac{\omega_i}{2\pi}$

18. Compute $\left| \frac{e_m y_i^m}{\lambda_i} \right|$, where y_i^m is the m^{th} (last) component of $\{y_i\}$.

This term is an error bound (9) on

$$\left| \frac{\frac{\omega_i^2}{2} - 1}{\omega_i \text{ exact}} \right|$$

19. Compute $[G] \{y_i\}$ for $i = 1, 2, \dots, q$ where the columns of $[G]$ are the vectors $\{g_i\}$. These are the vibration mode shapes $\{\phi_i\}$ associated with the lower q eigenvalues.

These are then normalized such that

$$\|\phi_i\|_2 = 1$$

APPENDIX B

USER'S MANUAL FOR

RE*S*I*ST

(STATIC AND DYNAMIC REUSABLE SURFACE INSULATION STRESS PROGRAM)

A. INTRODUCTION

This Appendix describes the use of a finite element based structural computer program for determining the static response and natural vibrations of TPS protected shuttle panels. The program is titled "RESIST" for static and dynamic RReusable Surface Insulation Stresses. The logic flow for RESIST is presented in Figure B-1.

The basis for the method is that the TPS is nonstructural but its stress levels, which are critical, must be computed. Thus, it becomes possible to neglect the stiffness of the TPS initially, but not its mass in the vibration, to determine approximate primary structure deflections.

An iterative procedure is then performed where, for each step, the primary structure deflections are imposed individually upon each tile at the tile/primary-structure interface, and the tile deflections and interface boundary loads are obtained. For the vibration option, the frequency is updated by computing a Rayleigh Quotient, using the latest non-rigid tile displacements in addition to the corresponding primary structure displacements. The individual tile boundary loads obtained are then assembled and their reactions are applied to the primary structure. New Primary structure deflections are obtained and compared to the previous set. This process is repeated until convergence is obtained.

B. PROGRAM LIMITATIONS

The usual assumptions for programs based upon the linear elastic finite element method are applicable to RESIST. However, to facilitate the preparation of program input, a number of simplifications regarding the configuration and loadings have been made. Thus, the generation of a voluminous quantity of finite element input data has been greatly reduced by inclusion of a series of data preprocessing subroutines within RESIST. The restrictions upon which these subroutines are based follows:

1. Boundary conditions and edge loadings are assumed uniform along the four rectangular plate edges defined by $x = 0$, L_x and $y = 0$, L_y .
2. The primary structure plate temperature and properties are all uniform.

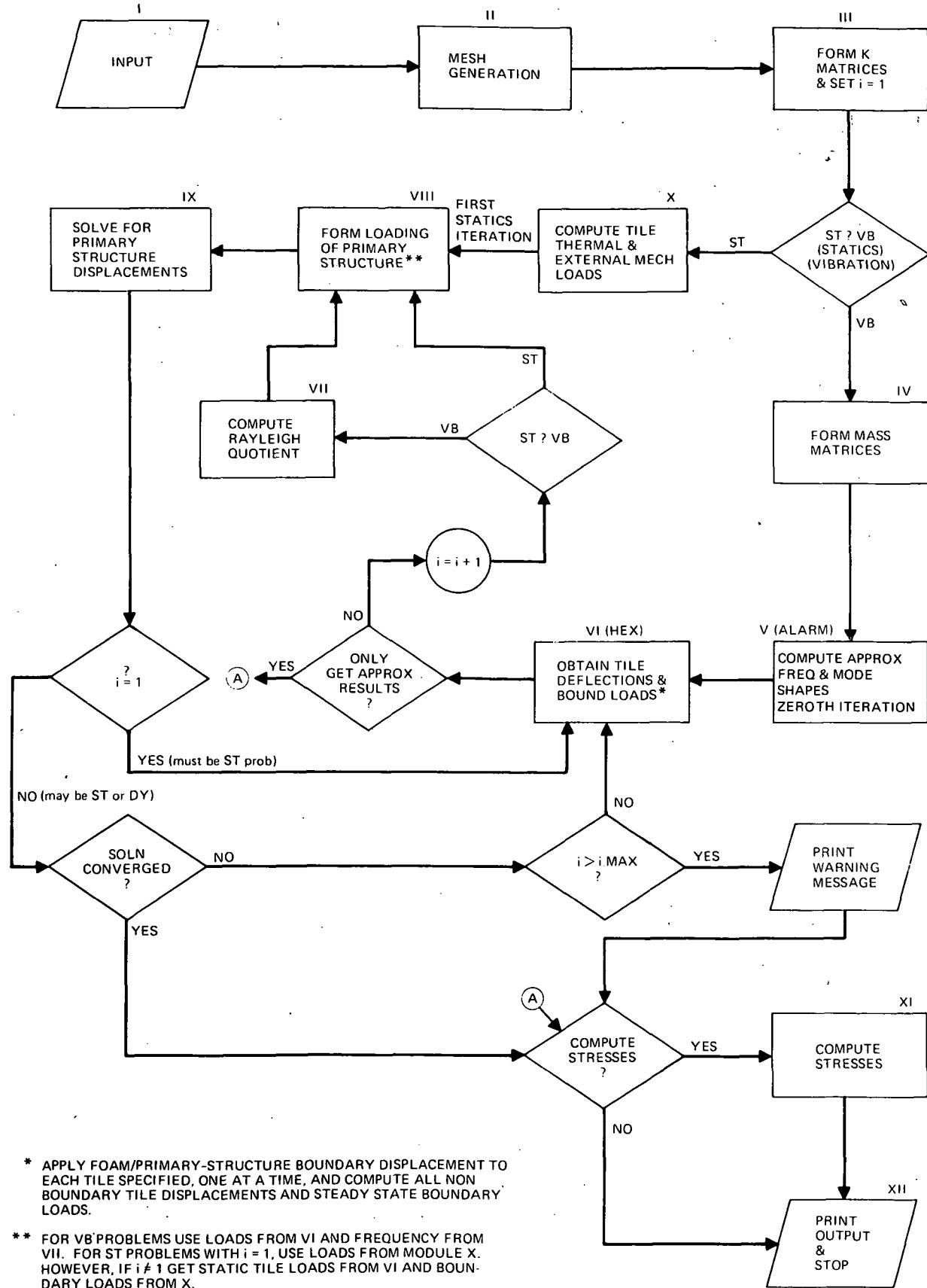


Figure B-1 Flow Chart for RSI Stress Analysis Program "Resist"

3. The stringers are equally spaced with temperatures and properties which are all uniform.
4. All tiles are geometrically identical as are their temperature distributions and uniform pressure loadings.
5. The boundary conditions must be selected such that the primary structure is statically stable.

The remaining limitations are primarily concerned with the program's capacity and should be adhered to by the user. These limitations are as follows:

6. Maximum number of nodes in a tile = 850.
7. Maximum number of finite elements running in any one direction in a tile = 20.
8. Maximum number of nodes in primary structure = 2500.
9. Maximum number of primary structure nodes along x or y direction = 1,000.
10. Maximum number of degrees of freedom in primary structure

$$= \begin{cases} 10,000 & \text{for vibration option.} \\ 15,000 & \text{for statics option.} \end{cases}$$
11. Maximum number of natural mode shapes = 50.
12. Maximum number of stringers = 15.
13. To avoid a singular stiffness matrix, I_z , and $\sin \beta$ must not both be zero for a given stringer.

A violation of restrictions 6-13, inclusive, will cause the program to stop and an appropriate warning message to appear.

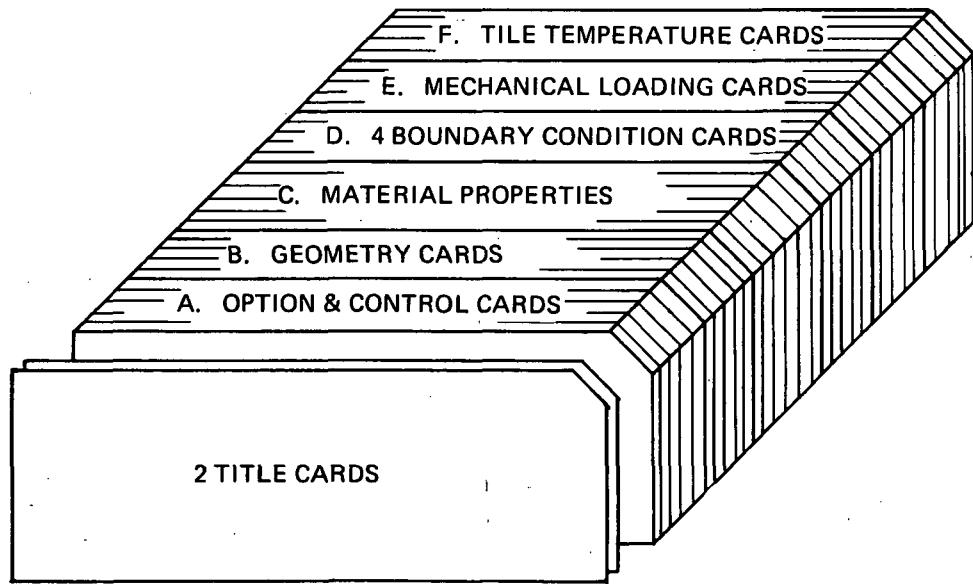
To insure symmetry of solutions for panels which are symmetric with regard to stringer locations about $y = L_y/2$, care should be taken with the input data to see that the plate nodes associated with the stringers are symmetric about $y = L_y/2$. Thus, the number of primary structure finite elements in the y direction should not be odd if the number of panel stringers is even.

C. INPUT INSTRUCTIONS

A description of the card input for the IBM 370 and CDC 6600 versions of this program is presented in this section.

In addition to the first two input cards which contain literal data, such as special program title and date, in columns 1 through 80, inclusive; there are six groups of input cards containing the following information:

- Group A - Instructions regarding the type of problem being performed, number of iterations desired, and type of output information.
- Group B - Details of the geometric configuration and finite element mesh of the primary structure and tiles. (Card B.4 is omitted if there are no tiles)
- Group C - Defines the primary structure and RSI temperature dependent material properties. If there is no TPS, cards C.3 through C.11 are omitted.
- Group D - Specifies the primary structure boundary conditions
- Group E - Describes the mechanical loading upon the primary structure as well as its temperature. These cards are omitted when the vibration option is used
- Group F - Defines the RSI temperature distribution. These cards are omitted if there is no TPS.

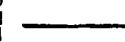


CARD(S)	COL(S)	FORMAT	SYMBOLS	UNITS	DESCRIPTION
A.1	1-5	I5	-	-	1 in col. 5 denotes that a statics problem is being treated. Skip cols. 6-25 in such cases
6-10	I5	N _D	-	-	2 in col. 5 denotes that a natural vibration problem is being treated.
16-20	I5	- S	-	-	Number of desired mode shapes (50 is the maximum permitted).
					Number of reorthogonalizations for eigenvalue algorithm. A min of 2 and a max. of 5 is suggested with 3 as an adequate compromise for most problems. The run should be repeated with greater values for S or N _D if the frequency error bound of a desired mode is greater than 1%.
21-25	I5	-	-	-	Vibration mode number for which tile modes are desired.
26-30	I5	i _{max}	{ in. or rad.	-	Maximum number of iterations
31-40	E10.0	ε	-	-	Convergence parameter. Maximum primary structure deflection or rotation difference between iterations divided by magnitude of largest element.
46-50	I5	-	-	-	0 in col. 50 indicates that <u>primary structure stresses and strains are not required.</u>
					1 in col. 50 indicates that only <u>midplate strains and stresses of primary structure are required.</u>
					2 in col. 50 indicates that only <u>top of plate strains and stresses of primary structure are required.</u>
					3 in col. 50 indicates that only <u>bottom of plate strains and stresses of primary structure are required.</u>
					4 in col. 50 indicates that only <u>mid and top of plate strains and stresses of primary structure are required.</u>
					5 in col. 50 indicates that only <u>mid and bottom of plate strains and stresses of primary structure are required.</u>
					6 in col. 50 indicates that only <u>top and bottom of plate strains and stresses of primary structure are required.</u>
					7 in col. 50 indicates that <u>top, bottom, and mid plate strains and stresses of primary structure are required.</u>
51-60	E10.0	-	-	lb-in- sec ²	Overhung rotatory mass inertia ε associated with each stringer. Used if plate overhang x = 0 and x = L _x boundaries.

CARD(S)	COL(S)	FORMAT	SYMBOLS	UNITS	DESCRIPTION
A.2	1-5	I5	-	-	0 in col. 5 indicates that tile stresses are <u>not</u> required. 1 in col. 5 indicates that tile stresses are to be computed after each iteration is performed.
	6-10	I5	-	-	2 in col. 5 indicates that tile stresses are to be computed only after last iteration is performed or only after convergence is obtained.
					0 in col. 10 if primary structure stresses and strains were not requested in column 50 of Card A.1.
					1 in col. 10 indicates that primary structure stresses and strains are required after each iteration.
					2 in col. 10 indicates that primary structure stresses and strains are required only after last iteration or, only after convergence.
					O in col. 15 indicates no tiles on the primary structure. Skip card A.3*
					1 in col. 15 indicates that there are tiles on the primary structure.
					1 in col. 20 indicates tile node map printout desired. 0 = no node map printout.
					1 in col. 25 indicates tile element map printout desired. 0 = no element map printout.
					1 in col. 30 indicates tile nodal coordinate, temp. and nodes per element printout.
					O in col. 30 indicates suppression of this printout.
					1 in col. 35 indicates printout of element stiffness matrices. 0 = no element stiffness matrices.
					1 in col. 40 indicates printout of assembled stiffness matrices and ALARM reorthog. info.
					O in col. 40 indicates suppression of this printout.
					1 in col. 45 indicates printout of unit no., file no., and matrix storage info. for program debugging.
					O in col. 45 indicates suppression of this printout.
A.3	1-4, 5-8, 9-12, etc.	I4 I4 I4	- - -	-	This card is used to indicate which tile stress states are desired. User may specifically request up to 20 tile stress states (see Figure A.2 for tile numbering scheme). A zero in col. 4 indicates that stress states for all tiles are desired.

* If there are no tiles then \bar{n}_x and \bar{n}_y , together with n_{B2} and n_{D2} , are still required since they determine the primary structure finite element grid. In analyzing panels without tiles, leave out cards B.4, C.3 through C.10 and all "F" cards.

B. GEOMETRIC CONFIGURATION – Sheet 1 of 2 (See Figure' B-2)

CARD(S)	COL(S)	FORMAT	SYMBOLS	UNITS	DESCRIPTION
B.1	1-10 11-20 21-30	2E10.0 	L_x L_y t_p	inches inches inches	Panel dimension Panel dimension Panel thickness
B.2	1-10 11-20 21-30 31-40 41-50 51-60 61-70 71-80	8E10.0 	Y_1 Z_s Y_s A_s I_y' I_z' $J_{x'}$ β_s	inches inches inches in. in. in. in. Degrees	Position of first stringer centroid. If there are no stringers, set $Y_1 > L_y$ and skip to next card. Distance of stringer centroid below plate middle surface Discrete stiffener spacing Stringer cross sectional area Stiffener principal mom. of inertia about y' axis Stiffener principal mom. of inertia about z' axis Stiffener twisting stiffness geometric parameter Angle between z and z' axis measured positive clockwise along x
B.3	1-10 11-20	110 110	\bar{n}_x \bar{n}_y	- -	Integer number of tiles between $x = 0$ and L_x^* Integer number of tiles between $y = 0$ and L_y^*

* If there are no tiles then \bar{n}_x and \bar{n}_y , together with n_{B2} and n_{D2} , are still required since they determine the primary structure finite element grid. In analyzing panels without tiles, leave out cards B.4, C.3 through C.10 and all "F" cards.

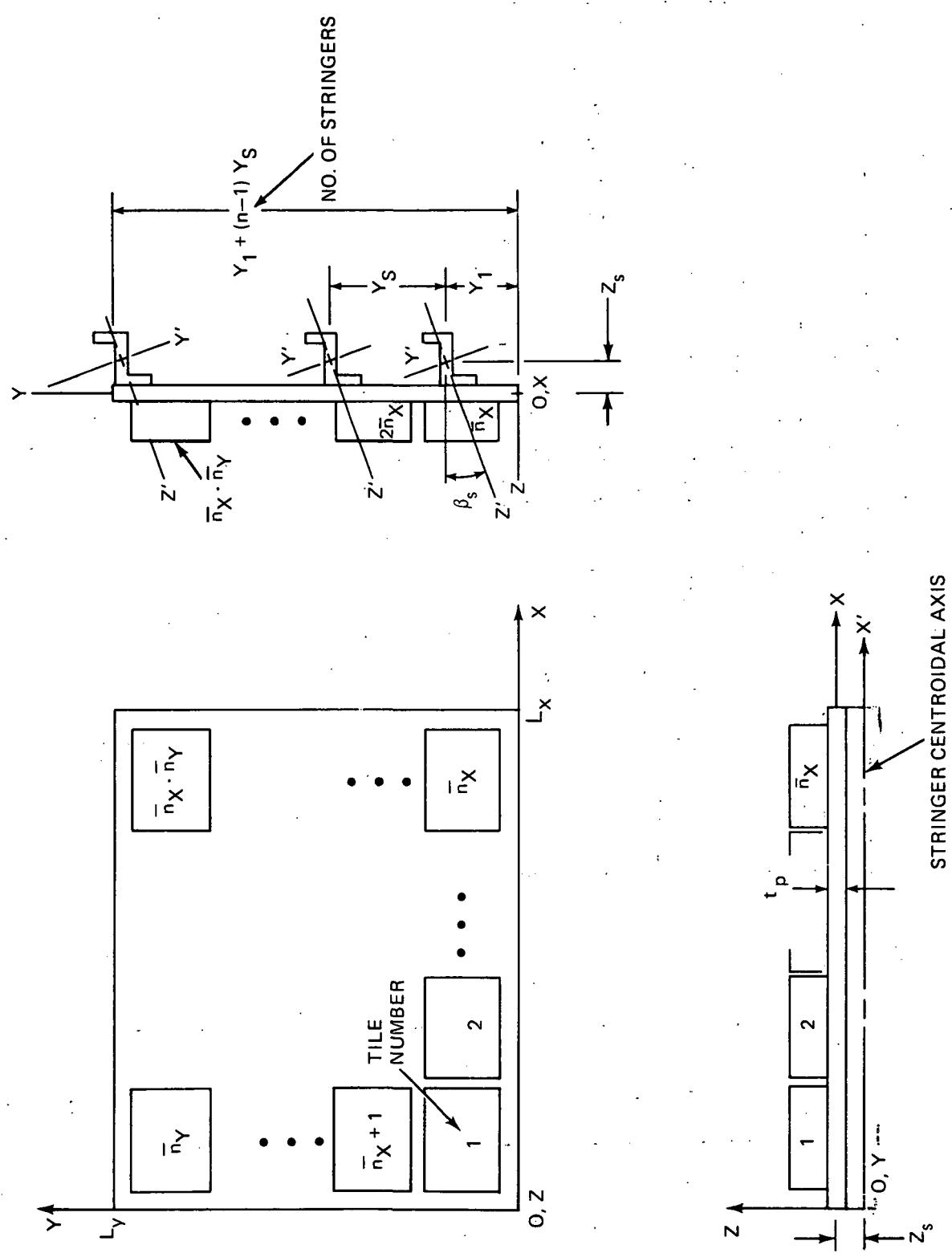
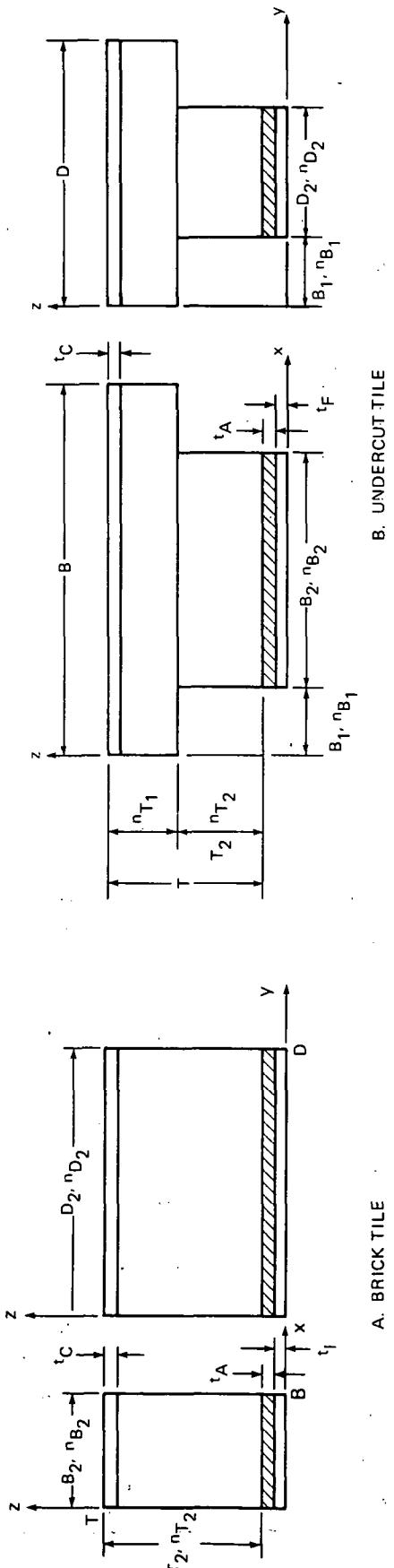


Figure B-2 TPS Configuration on Stiffened Primary Structure

B. GEOMETRIC CONFIGURATION - Sheet 2 of 2 (See Figure B-3)

CARD (S)	COL(S)	FORMAT	SYMBOLS	UNITS	DESCRIPTION
B.4	1-10	6E10.0	T B ₁	inches	Undercut RSI tile thickness. Leave blank if tile is brick shaped or if there are no tiles.
	11-20			inches	Tile undercut dimension. Leave blank if tile is brick shaped.
	21-30		T ₂	inches	Tile undercut dimension or, height of brick shaped tile.
	31-40		t _A	inches	Strain arrestor plate (SAP) thickness. May replace with layer of isolator, RSI or bond material if no SAP.
	41-50		t _T	inches	Strain isolator thickness (SIP)
	51-60		t _c	inches	Coating thickness. Leave blank if no tile coating.
B.5	1-5	T ₅	n _{B₁}	-	Number of elements along B ₁ . Leave blank if tile is brick shaped or if there are no tiles.
	6-10		n _{B₂}	-	Number of elements along B ₂ .
	11-15		n _{D₂}	-	Number of elements along D ₂ .
	16-20		n _{T₁}	-	Number of elements along T-T ₂ . Leave blank if tile is brick shaped.
	21-25		n _{T₂}	-	Number of elements along T ₂ .



NOTE: SUBSCRIPTED SYMBOLS BEGINNING WITH "n" ARE THE NUMBER OF ELEMENTS WHICH SUBDIVIDE THE INDICATED SPAN.
THE OTHER SYMBOLS ARE DIMENSIONS.

Figure B-3 RSI Tile Parameters

C. MATERIAL PROPERTIES – Sheet 1 of 2 (See Figure B-4)

CARD (S)	COL(S)	FORMAT	SYMBOLS	UNITS	DESCRIPTION
C.1	1-10 11-20 21-30 31-40	4E10.0 →	E_p ν_p γ_p α_p	psi - lb/in ³ σ_F^{-1}	Isotropic plate modulus of elasticity Poisson's ratio for plate Weight density for plate Coefficient of thermal expansion for plate
C.2	1-10 11-20 21-30 31-40	4E10.0 →	E_s ν_s γ_s α_s	psi - lb/in ³ σ_F^{-1}	Stringer modulus of elasticity (enter zero if no stringers) Poisson's ratio for stringer Weight density for stringer Coefficient of thermal expansion for stringer
C.3	1-10 11-20 21-30 31-40 41-50 51-60	6E10.0 →	E_x E_y E_z ν_{xy} ν_{yz} ν_{zx}	psi psi psi - - -	Arrestor x direction orthotropic stiffness Arrestor y direction orthotropic stiffness Arrestor z direction orthotropic stiffness See Figure A.4 See Figure A.4 See Figure A.4
C.4	1-10 11-20 21-30 31-40 41-50 51-60 61-70	7E10.0 →	G_{xy} G_{yz} G_{zx} γ_A α_{AX} α_{AY} α_{AZ}	psi psi psi σ_F^{-1} σ_F^{-1} σ_F^{-1}	See Figure A.4 See Figure A.4 See Figure A.4 Weight density for arrestor X coefficient of thermal expansion for arrestor Y coefficient of thermal expansion for arrestor Z coefficient of thermal expansion for arrestor

NOTE: This matrix is symmetric;
thus, the program insures
that

$$\begin{bmatrix}
 \frac{1}{E_x} & -\frac{\nu_{xy}}{E_y} & -\frac{\nu_{xz}}{E_z} \\
 -\frac{\nu_{yx}}{E_x} & \frac{1}{E_y} & -\frac{\nu_{yz}}{E_z} \\
 -\frac{\nu_{zx}}{E_x} & -\frac{\nu_{zy}}{E_y} & \frac{1}{E_z}
 \end{bmatrix}
 = \begin{bmatrix}
 \epsilon_x & \epsilon_y & \epsilon_z \\
 \epsilon_y & \frac{1}{G_{xy}} & 0 \\
 \epsilon_z & 0 & \frac{1}{G_{yz}}
 \end{bmatrix}
 \begin{bmatrix}
 \sigma_x & \sigma_y & \sigma_z \\
 \sigma_y & \frac{\nu_{xy}}{E_x} & \sigma_{xy} \\
 \sigma_z & \frac{\nu_{xz}}{E_x} & \frac{\nu_{yz}}{E_y}
 \end{bmatrix}
 \begin{bmatrix}
 \gamma_{xy} & \gamma_{yz} & \gamma_{zx} \\
 \gamma_{yz} & \frac{\nu_{zy}}{E_z} & 0 \\
 \gamma_{zx} & 0 & \frac{1}{G_{zx}}
 \end{bmatrix}$$

Figure B-4 Orthotropic Stress-Strain Law for 3-Dimensional Elements

C. MATERIAL PROPERTIES - Sheet 2 of 2

CARD(S)	COL(S)	FORMAT	SYMBOLS		UNITS
C.5	1-10 11-20 21-30 31-40	4E10.0 4E10.0	E _I ν _I γ _I α _I	psi - lb/in ³ °F ⁻¹	Isolator modulus of elasticity Poisson's ratio for isolator (Note: max ν _I = .499) Weight density for isolator Coefficient of thermal expansion for isolator
C.6	1-10 11-20 21-30	E10.0 E10.0 E10.0	γ _R α _y /α _x α _z /α _x	lb/in ³ - -	Weight density of RSI material RSI coefficient of thermal expansion in y direction (°F ⁻¹) divided by coefficient of thermal expansion in x direction (α _x) RSI coefficient of thermal expansion ratio in z vs. x direction

C. TEMPERATURE DEPENDENT MATERIAL PROPERTIES, sheet 1 of 2

CARD(S)	COL(S)	FORMAT	SYMBOLS	UNITS	DESCRIPTION
C. 7.1	1-5	I5	-	-	Number of entry sets in the following table of E_R vs. temperature (°F)
C. 8.1	1-10	E10.0	T_1	°F	Temperature (absolute, not relative) corresponding to following value of E_R
	11-20	E10.0	$E_R(T_1)$	psi	Value of E_R (RTI modulus - refer to equations below*) associated with previous temp.
	31-30 etc.	E10.0 etc.	T_2 etc.	°F etc.	Repeat above set of data as often as necessary, 4 sets to a card.
					Program uses closest 3 data pts. for 2nd order Langrangian interpolation of properties if element temp. is within data specified temp. range and at least 3 data-points are input. Program uses closest data-point properties for element temp. outside range. Uniform property value is used for any given property if only one value of that property is specified. Thus, program requires a minimum of 1 or 3 value(s) per property for proper execution.

C. TEMPERATURE DEPENDENT MATERIAL PROPERTIES — Sheet 2 of 2

CARD(S)	COL(S)	FORMAT	SYMBOLS	UNITS	DESCRIPTION
C.7.2 & C.8.2					Repeat above two card sets for E'_R *
C.7.3 & C.8.3 through C.7.6 & C.8.6					Repeat above card sets for remaining RSI properties in following order: G'_R , ν_R , ν'_R , and α'_x where α_x = RSI coefficient of thermal expansion in x direction.
C.9.1 & C.10.1 through C.9.3 & C.10.3					Repeat above card sets for coating properties in following order: E_c , ν_c , α'_c

* For RSI

(refer to Figure B-4)

$$G_{xy} = G_{yx} = \frac{E_R}{2(1 + \nu_R)}$$

$$G_{yz} = G_{zy} = G_{zx} = G_{xz} = \frac{G'_R}{E'_R}$$

$$\nu_{xy} = \nu_{yx} = \nu_R$$

$$\nu_{xz} = \nu_{yz} = \nu'_R$$

$$\nu_{zx} = \nu_{zy} = \frac{E_R}{E'_R} \nu'_R$$

D. BOUNDARY CONDITIONS - Sheet 2 of 3 (See Figure B-5)

CARD(S)	COL(S)	FORMAT	SYMBOLS	UNITS	DESCRIPTION
D.1-D-4, (continued)	31	A1	-	-	<p>0 denotes edge is <u>not held</u> from <u>in-plane deflections</u></p> <p>1 denotes edge is <u>held</u> from <u>in-plane deflections</u></p> <p>2 denotes edge is not held for y deflection, but is held for x deflection (PARTIALLY HELD)</p> <p>4 denotes edge is not held for x deflection, but is held for y deflection (PARTIALLY HELD)</p> <p>3 denotes edge is <u>flexibly held</u> for <u>in-plane deflections</u></p>
					<p>NOTE: For non-vibratory heated or cooled primary structure problems, refer to special instructions on bottom of page</p>
32-40	E9.0		K_{uu}	lb/in^2	In-plane x force per unit length on an edge caused by in-plane x direction unit deflection
41-49			K_{uv} or K_{vu}	lb/in^2	In-plane x force per unit length on an edge caused by in-plane y direction unit deflection or In-plane y force per unit length on an edge caused by in-plane x direction unit deflection
50-58				lb/in^2	In-plane y force per unit length on an edge caused by in-plane y direction unit deflection

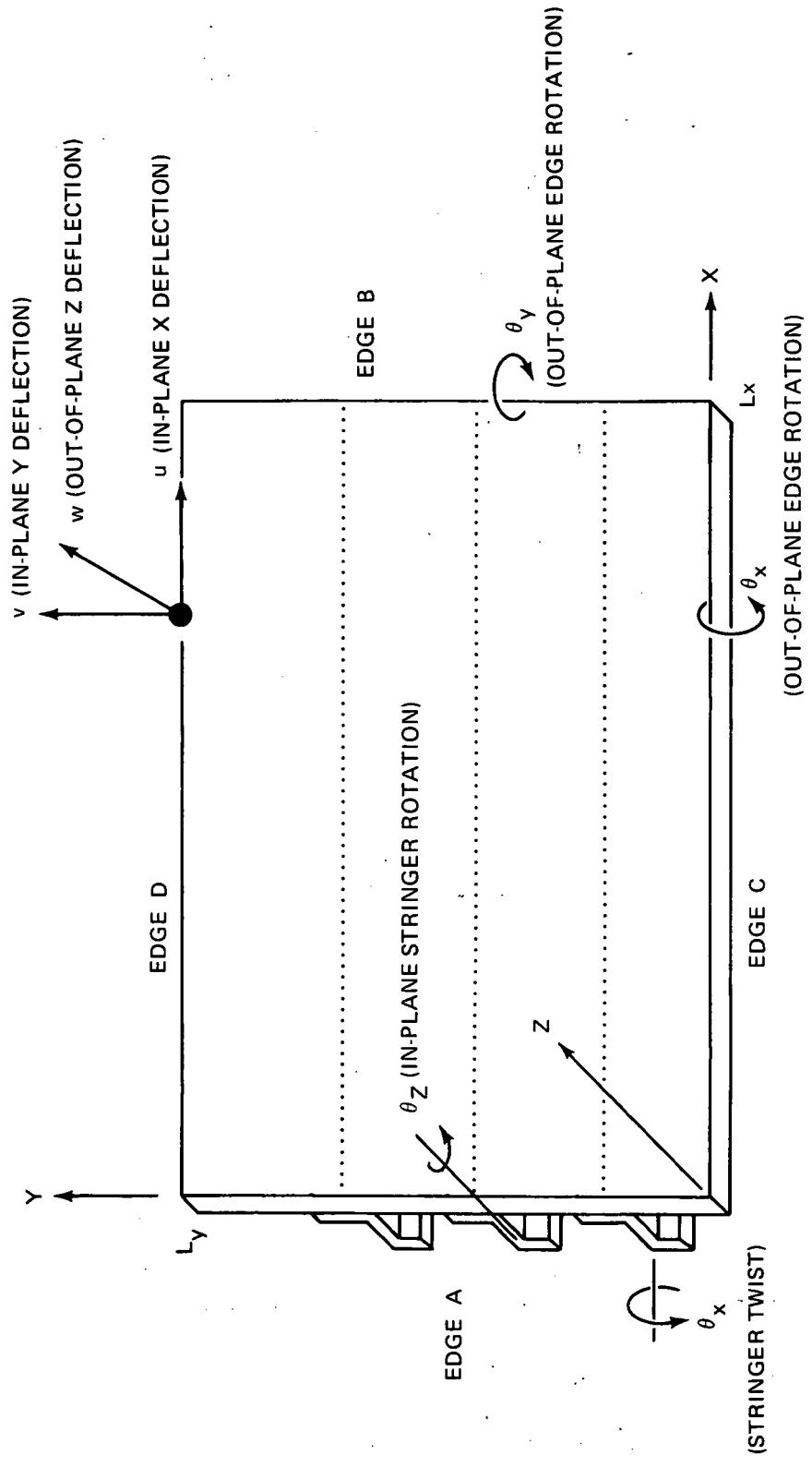


Figure B-5 Primary Structure Boundary Condition Notation

D. BOUNDARY CONDITIONS – Sheet 1 of 3 (See Figure B-5)

CARD(S)	COL(S)	FORMAT	SYMBOLS	UNITS	DESCRIPTION
D.1, D.2, D.3, D.4	1	A1	-	-	<p>B denotes the $x=L_x$ edge of the plate (CARD D.2)</p> <p>C denotes the $y=0$ edge of the plate (CARD D.3)</p> <p>D denotes the $y=L_y$ edge of the plate (CARD D.4)</p>
These are four simi- lar bound- ary condi- tion cards	2	A1	-	-	<p>0 indicates that the plate edge is <u>free</u> to deflect and rotate <u>out</u> of the $z=0$ plane (FREE)</p> <p>1 indicates that the plate edge is <u>not free</u> to deflect or rotate <u>out</u> of the $z=0$ plane (CLAMPED)</p> <p>2 indicates that the plate edge is <u>not free</u> to deflect but is <u>free to rotate</u> <u>out</u> of the $z=0$ plane (PINNED)</p> <p>3 indicates that the plate edge is <u>flexibly held</u> with regard to <u>out</u> of plane motion</p>
3-11	E9.0		K_{WW}	lb/in^2	Out-of-plane force per unit edge-length caused by out-of-plane unit deflection
12-20			$K_{W\theta}$ or $K_{\theta W}$	lb/in	Out-of-plane force per unit edge-length caused by out-of-plane unit rotation or Out-of-plane moment per unit edge-length caused by out-of-plane unit deflection
21-29			$K_{\theta\theta}$	1lb.	Out-of-plane moment per unit edge-length caused by out-of-plane unit rotation

D. BOUNDARY CONDITIONS — Sheet 1 of 3 (See Figure B-5)

CARD(S)	COL(S)	FORMAT	SYMBOLS	UNITS	DESCRIPTION
Add'l info. for cards D1 and D2 only.	60	A1	-	-	O denotes stringer edge not held for in-plane rotation (θ_z) 1 denotes stringer edge held for in-plane rotation ($\theta_z = 0$) 3 denotes stringer edge flexibly held for in-plane rotation Note A _z is not a primary structure degree of freedom unless a stringer element is also present at a particular plate node
61-69	E9.0	$K_s\theta_z$	in-lb.	In-plane stringer edge moment produced by unit rotation θ_z	
71	A1	-	-	O denotes stringer edge free to twist ($\theta_x = 0$) 1 denotes stringer edge not free to twist ($\theta_x = 0$) 3 denotes stringer edge flexibly held against twist	
72-80	E9.0	$K_s\theta_x$	in-lb.	Twist moment on end of stringer for a unit twist-rotation	

Special Instructions for running a thermal stress problem when the primary structure is at a uniform temperature other than the reference temperature are required to permit free in-plane thermal straining; e.g.:

1. Permit the $x=0$ boundary to move freely or be elastically held in-plane elastically held if elastically held along $x=0$.
2. Permit the $x=L_x$ boundary to move freely in the y direction but not the x direction if free, or be elastically held if elastically held in-plane.
3. Permit the $y=0$ boundary to move freely in the x direction but not the y direction if free, or be elastically held if elastically held along $y=0$.
4. Permit the $y=L_y$ boundary to move freely in the x direction but not the y direction if free, or be elastically held if elastically held in-plane.

— See pages B-53 and B-57 for a typical example of the above instructions.

E. PRIMARY STRUCTURE LOADING (See Figure B-6)

CARD(S)	COL(S)	FORMAT	SYMBOLS	UNITS	DESCRIPTION
E.1	1-10	E10.0	N_x	lb/in	Uniform, direct cover-plate running load in x direction on $x = 0$ edge (see Figure 7)
	11-20		N_y	lb/in	Uniform, direct cover-plate running load in y direction on $y = 0$ edge (see Figure 7)
	21-30		N_{xy}	lb/in	Uniform, shearing cover-plate running load on $x = 0$ edge (see Figure 7)
	31-40		N_{yx}	lb/in	Uniform, shearing cover-plate running load on $y = 0$ edge (see Figure 7)
	41-50		P_z	psi	Uniform external normal pressure acting upon tiles
	51-60		T	lb	Tension force acting upon centroid of each stiffener at $x = 0$
	61-70		M	in-lb	Out-of-plane bending moment acting upon each stiffener
	71-80		V	lb	Shear load acting upon each stiffener
					Note, boundary conditions for B and D edges should be selected carefully to produce desired effect, e.g. to produce a uniform primary structure tension in the x direction, $\sigma_x = \bar{\sigma}$ and $\sigma_y = 0$; set $N_x = t_p \bar{\sigma}$, $N_y = A_s \bar{\sigma}$, and P_z , M and V all equal to zero; then hold plate edge B from in plane x, but not y, motion and also hold the stringers at edge B; next, make edges C and D free for X motion, and only hold one of these edges against y motion.
E.2	1-10	E10.0	ΔT_p	F^o	Temperature difference of plate from T_{Ref}
	11-20	E10.0	ΔT_s	F^o	Temperature difference of stringers from T_{Ref}

Note: Leave out cards E.1 and E.2 if vibration option is used

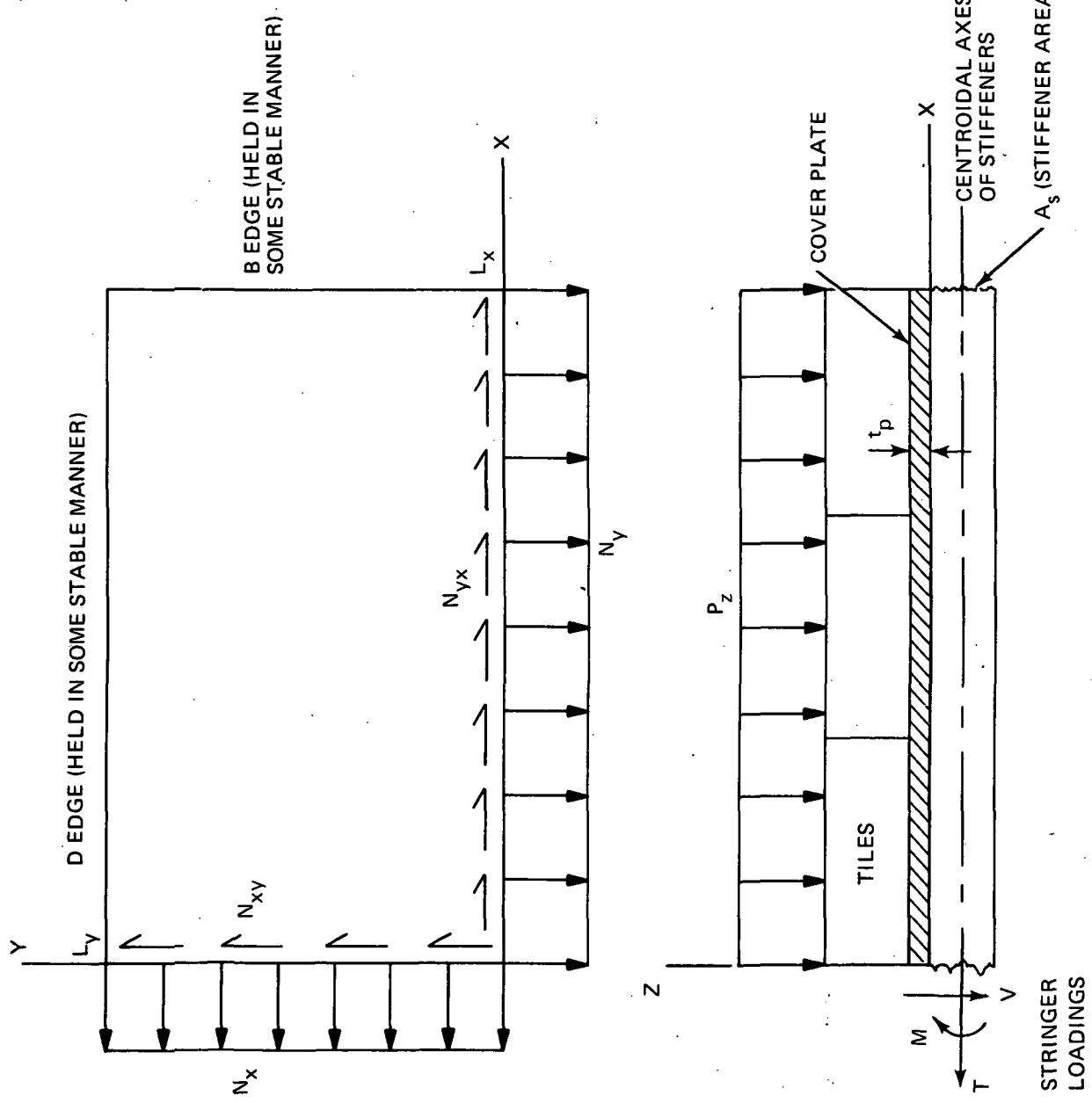


Figure B-6 Possible Static Mechanical Loadings Upon Panel

F. TILE TEMPERATURE DISTRIBUTIONS - Sheet 1 of 2

Each tile is assumed to have the same temperature distribution. There are 3 temperature distribution options, each of which is considered separately below. Tile temperature differences, rather than absolute tile temperatures, are required for each of these options (since thermal strains depend upon temperature differences). However, since temperature-dependent material property data are presented in terms of absolute temperature scales, a reference temperature (which is also input) is added to the differences to obtain absolute temperatures for internally computing material properties.

CARD(S)	COL(S)	FORMAT	SYMBOLS	UNITS	DESCRIPTION
F.1	5	I1	-	-	O in col. 5 of this card indicates no thermal static loading effects will be considered. But material properties used in forming the TPS stiffness properties will be based upon the specified temperature distribution. 1 in col. 5 indicates that thermal static loading will be considered in the analysis. In such cases, refer to bottom of following page for special instructions regarding boundary condition cards (D.1 through D.4).
	10	I1	-	-	1 in col. 10 indicates that each tile is at the same uniform temperature.
	E10.0	T Ref	O_F		2 in col. 5 indicates that each tile temperature distribution is governed by Lagrangian interpolation formulas.
	11-20				3 in col. 5 indicates that each tile temperature distribution is input by consecutive finite element node-temperature differences from the reference temperature.
					Panel reference temperature (added to temp. differences when obtaining mat'l. properties)

F. TITLE TEMPERATURE DISTRIBUTIONS - Sheet 2 of 2

CARD(S)	COL(S)	FORMAT	SYMBOLS	UNITS	DESCRIPTION
UNIFORM TEMPERATURE OPTION (1)					
F.2	1-10	E10.1	ΔT_u	$^{\circ}F$	Uniform temperature difference from T_{Ref}
or LAGRANGIAN INTERPOLATION TEMPERATURE OPTION (2)					
F.2.1	1-5	215	-	-	Number of x coordinates through which temperature differences will be interpolated.
	6-10		-	-	Order of Lagrangian interpolation polynomial in x direction. Must be at least 1 less than number of coords given in col. 5.
F.3.1	1-10	E10.0	x_1	inches	The local x coordinates used in the x direction temperature difference interpolation. Eight to a card until all are accounted for.
F.2.2 - 3.2					Repeat card types F.2.1 and F.3.1 for the y coordinates
F.2.3 - 3.3					Repeat card types F.2.1 and F.3.1 for the z coordinates
or ELEMENT NODE TEMPERATURE OPTION (3)					
F.2	1-10 11-20 etc.	E10.0	ΔT_R	$^{\circ}F$	Temperature differences above reference temperatures, node by node, in consecutive order. Seven temperature differences to a card until all nodes are accounted for. Cols. 71-80 of each card are reserved for user's card identification.

D. DESCRIPTION OF OUTPUT

Output from a typical run of the RESIST computer program is explained below in outline form. References in parentheses refer to pages in this Appendix.

1. Program title and date indicating latest update of program version which was run.

INPUT INFORMATION

2. Listing of input cards, the first two of which are the title assigned to any given run by the user.
3. User selected input options are listed.
4. Plate, stringer and tile geometry and specification of finite element grids for primary structure and tiles (pp. B.8 - B.11).
5. Plate, stringer, strain isolator and arrestor material properties (p. B.12). Note, if there is no strain arrestor, RSI or isolator material properties may be used for the arrestor. If this is done, the thickness dimension of the usual isolator or RSI should be appropriately reduced to compensate for this addition.
6. Temperature-dependent RSI material property data used for generating curves used internally by program to compute RSI average finite element properties.
7. Plate and stringer boundary conditions (pp. B.16-B.21).
8. Applied primary structure static mechanical and thermal loading if not a vibration problem.
9. RSI temperature distribution input data. Used for property data (item 6 above) and thermal loading if a statics problem.

OUTPUT INFORMATION

10. Map showing typical tiles three dimensional finite element ordering, by layers. Top, or first layer also corresponds to two-dimensional tile coating elements as well.

11. Map showing ordering of a typical tiles finite element nodes by layers.
12. Position and temperatures for a typical tile in a local coordinate system (reference Figures 3.a &.b).
13. Global geometry of primary structure nodes and plate nodal degree-of-freedom numbering . D_x , D_y and D_z refer to nodal deflections, and R_x , R_y and R_z are the nodal rotations. Nodes with no degrees-of-freedom are used to define the stringer centroids and axes.
- 14.a. Statics Option: Primary structure nodal deflections by iteration number. Nodes with the same x coordinate are grouped together. These groups are separated with dashed lines.
- 14.b. Vibration Option: Mode numbers, approximate frequencies and corresponding modal error bound (which should be less than 2% to be a reliable approximate mode). This is followed by the primary structure mode shapes with a similar nodal deflection format as for the Statics Option.
15. If requested by the user, the computed convergence parameter is printed out along with the input quantity it was tested against. This is done for each iteration after the first for the Static Option. The primary structure degree-of-freedom with the largest change from the previous iteration is also identified.
16. Tile nodal displacements by tile and iteration number. For a vibration option, this calculation and the subsequent ones are performed only for the user-specified vibration mode.
17. Three dimensional tile stresses and strains for the bottom two layers of elements by element number. These quantities are computed at each element's 8 Gauss integration points. Gauss point stresses are believed to be more accurate than nodal values and provide more detail than simply the elements' average stresses.

18. Three dimensional element average stresses and strains (by tile and iteration number).
19. Two-dimensional element average coating stresses. Coating element numbers correspond to three dimensional element numbering directly below them.
- 20.a. Statics Option: Repeat of items 16-19 for each tile. Repeat of item 14.a and 15 for each iteration.
- 20.b. Vibrations Option: Computation of Rayleigh Quotient (OMEGA SQUARED) if all tiles have been treated. Repeat of items 16-19 for each iteration. Repeat of items 14.b, 15 and Rayleigh Quotient until convergence or last iteration is performed.
21. Plate element stresses and strains for mid and/or top and/or bottom surfaces. This computation is done after each iteration if requested by the user. Otherwise, it is computed only after convergence or the last iteration is performed.

E. SAMPLE PROBLEMS

Output for three sample problems, one vibration case and two statics cases, are presented in the remaining pages of this report. Only portions of the output for each problem are shown. However, the pages presented are representative of the types of information, and their respective formats, which the RESIST Program can deliver.

STATIC AND DYNAMIC

RELIABILITY SURFACE INSULATION STRENGTH

VERSION DATE

卷之三

PREPARED BY

I. OJALVO, P. OGILVIE, A. LEVY AND F. AUSTIN

© GRUMMAN AEROSPACE CORPORATION

FOR

THE LANGLEY RESEARCH CENTER

PROGRAM LISTING OF INPUT DATA CARDS

SAMPLE PROBLEM 1 - VIBRATION OF THREE SIMPLE TIES

123456789012345678901234567890123456789012345678901234567890

O P T I O N S

FREE VIBRATION MODES

NO. DESIRED MODES = 5 NO. RENORTHOGONALIZATIONS = 2 MODE NO. = 1

MAXIMUM NO. ITERATIONS = 4

OVERHUNG ROTATORY MASS INERTIA ASSOCIATED WITH EACH STRINGER = 0.0

PRIMARY STRUCTURE STRESSES PRESENTED AFTER LAST ITERATION AT PLATE MID, TOP AND BOTTOM SURFACES

TITLE FOR PRIMARY STRUCTURE

TITLE STRESSES PRESENTED AFTER EACH ITERATION

TITLE NODE MAP OR CHIRP

TITLE ELEMENT AND DEGREE

TITLE VOLUME CONCENTRATES PRECULLED

DO NOT PRINT REFORMATTED SURFACE SURFACES

DO NOT PRINT ASSEMBLED SURFACE MATRICES

DO NOT PRINT TITLE NUMBERING FOR THIS COMPUTATION

TRANSIENT STRESSES & TITLE TITLES

GEOMETRY

PLATE LX = 1.80000F 01 LY = 3.33330E-01 LP = 7.50000E-01
STRINGERS Y1 = 1.00000E 00 ZS = 0.0 YS = 0.0 AS = 0.0
 YY' = 0.0 YZ' = 0.0 JX' = 0.0 BETA S = 0.0

TILES NXP = 3 NYR = 1
 T = 0.0 R1 = 0.0 T2 = 1.16667E 00
 TA = 1.666670F-01 TI = 5.00000E-02 TC = 0.0
 NT1 = 0 NQ2 = 10 ND2 = 1
 NT2 = 0 NT3 = 7

MATERIAL PROPERTIES

PLATE P1 = 1.00000E 07 NU P = 3.00000E-01 GAMMA D = 1.00000E-02 ALPHA P = 0.0
STRINGERS PS = 0.0 NU S = 0.0 GAMMA S = 0.0 ALPHA P = 0.0
 FX = 1.00000E 04 FY = 6.00000F 04 FZ = 6.00000F 03
 NU XY = 5.00000E-01 NU YZ = 1.00000E-01 NU ZX = 1.00000E-02
 GXY = 2.00000E 04 GYZ = 2.20000E 04 GZX = 3.20000E 04
 GAMMA A = 5.00000E-01 ALPHA Y = 0.0 ALPHA Z = 0.0
 ALPHA X = 0.0 ALPHA Y = 0.0 ALPHA Z = 0.0

ISPLAT12 RT = 9.00000E 01 NU I = 4.00000E-01 GAMMA I = 3.50000E-02 ALPHA I = 0.0
 GAMMA ? = 5.00000E-01 ALPHA BY / ALPHA BX = 0.0 ALPHA RZ / ALPHA RX = 0.0

TEMPERATURE DEPENDENT MATERIAL PROPERTIES

	PROPERTY	TEMPERATURE	PROPERTY	TEMPERATURE	PROPERTY	TEMPERATURE
1	FR	ALL	6.000E-04			
1	ER	ALL	6.000E-03			
1	GR	ALL	3.200E-04			
1	NJ_R	ALL	5.000E-01			
1	NJ_R'	ALL	1.000E-02			
1	ALPHA_R	ALL	0.0			
1	FC	ALL	0.0			
1	NU_C	ALL	0.0			
1	ALPHA_C	ALL	0.0			

BOUNDRAY CONDITIONS

EDGE PLATE OUT OF PLANE

STRINGERS

EDGE

PINNED

FREE

FREE

PINNED

U HELD, V FREE

FREE

FREE

FREE

FREE

V HELD, U FREE

FREE

FREE

RESISTIVE TEMPERATURES

NO STATIC THERMAL LOADING

UNIFORM TEMPERATURE OPTION

T REFERENCE = 7.0000E 01

DEFL T U = T - T REF = 0.0

NODE MAP

SURFACE 1

200	202	204	206	208	210	212	214	216	218	220
199	201	203	205	207	209	211	213	215	217	219

NODE MAP

SURFACE 2

178	150	182	184	186	188	190	192	194	196	198
177	179	181	183	185	187	189	191	193	195	197

ELEMENT MAP

LAYER 1
RSI

ELEMENT MAP

LAYER 9
ISOLATOR

1 2 3 4 5 6 7 8 9 10

... NO. 33111 C. M. E. S. H. 111

1	2	24	23
3	4	26	25
5	6	26	25
7	8	28	27
9	10	30	29
11	12	32	31
13	14	34	33
15	16	36	35
17	18	38	37
19	20	40	39
21	22	42	41
23	24	44	43
25	26	46	45
27	28	48	47
29	30	50	49
31	32	52	51
33	34	54	53
35	36	56	55
37	38	58	57
39	40	60	59
41	42	62	61

LOCAL TIME COORDINATES

TEMPERATURE

NODE	X	Y	Z	0.0	0.0	0.0	0.0	0.0
1	0.0	0.0	0.0	3.33330E-01	0.0	0.0	0.0	0.0
2	6.00000E-01	0.0	0.0	3.33330E-01	0.0	0.0	0.0	0.0
3	6.00000E-01	3.33330E-01	0.0	3.33330E-01	0.0	0.0	0.0	0.0
4	1.20000E-01	0.0	0.0	3.33330E-01	0.0	0.0	0.0	0.0
5	1.20000E-01	0.0	0.0	3.33330E-01	0.0	0.0	0.0	0.0
6	1.20000E-01	0.0	0.0	3.33330E-01	0.0	0.0	0.0	0.0
7	1.80000E-01	0.0	0.0	3.33330E-01	0.0	0.0	0.0	0.0
8	1.80000E-01	0.0	0.0	3.33330E-01	0.0	0.0	0.0	0.0
9	2.40000E-01	0.0	0.0	3.33330E-01	0.0	0.0	0.0	0.0
10	2.40000E-01	0.0	0.0	3.33330E-01	0.0	0.0	0.0	0.0
11	3.00000E-01	0.0	0.0	3.33330E-01	0.0	0.0	0.0	0.0
12	3.00000E-01	0.0	0.0	3.33330E-01	0.0	0.0	0.0	0.0
13	3.60000E-01	0.0	0.0	3.33330E-01	0.0	0.0	0.0	0.0
14	3.60000E-01	0.0	0.0	3.33330E-01	0.0	0.0	0.0	0.0
15	4.20000E-01	0.0	0.0	3.33330E-01	0.0	0.0	0.0	0.0
16	4.20000E-01	0.0	0.0	3.33330E-01	0.0	0.0	0.0	0.0
17	4.80000E-01	0.0	0.0	3.33330E-01	0.0	0.0	0.0	0.0
18	4.80000E-01	0.0	0.0	3.33330E-01	0.0	0.0	0.0	0.0
19	5.40000E-01	0.0	0.0	3.33330E-01	0.0	0.0	0.0	0.0
20	5.40000E-01	0.0	0.0	3.33330E-01	0.0	0.0	0.0	0.0
21	6.00000E-01	0.0	0.0	3.33330E-01	0.0	0.0	0.0	0.0
22	6.00000E-01	0.0	0.0	3.33330E-01	0.0	0.0	0.0	0.0
23	0.0	0.0	0.0	5.00000F-02	0.0	0.0	0.0	0.0
24	0.0	0.0	0.0	5.00000F-02	0.0	0.0	0.0	0.0
25	6.00000E-01	0.0	0.0	5.00000F-02	0.0	0.0	0.0	0.0
26	6.00000E-01	0.0	0.0	5.00000F-02	0.0	0.0	0.0	0.0
27	1.20000E-01	0.0	0.0	5.00000F-02	0.0	0.0	0.0	0.0
28	1.20000E-01	0.0	0.0	5.00000F-02	0.0	0.0	0.0	0.0
29	1.80000E-01	0.0	0.0	5.00000F-02	0.0	0.0	0.0	0.0
30	1.80000E-01	0.0	0.0	5.00000F-02	0.0	0.0	0.0	0.0
31	2.40000E-01	0.0	0.0	5.00000F-02	0.0	0.0	0.0	0.0
32	2.40000E-01	0.0	0.0	5.00000F-02	0.0	0.0	0.0	0.0
33	3.00000E-01	0.0	0.0	5.00000F-02	0.0	0.0	0.0	0.0
34	3.00000E-01	0.0	0.0	5.00000F-02	0.0	0.0	0.0	0.0
35	3.60000E-01	0.0	0.0	5.00000F-02	0.0	0.0	0.0	0.0
36	3.60000E-01	0.0	0.0	5.00000F-02	0.0	0.0	0.0	0.0
37	4.20000E-01	0.0	0.0	5.00000F-02	0.0	0.0	0.0	0.0
38	4.20000E-01	0.0	0.0	5.00000F-02	0.0	0.0	0.0	0.0
39	4.80000E-01	0.0	0.0	5.00000F-02	0.0	0.0	0.0	0.0
40	4.80000E-01	0.0	0.0	5.00000F-02	0.0	0.0	0.0	0.0
41	5.40000E-01	0.0	0.0	5.00000F-02	0.0	0.0	0.0	0.0
42	5.40000E-01	0.0	0.0	5.00000F-02	0.0	0.0	0.0	0.0
43	6.00000E-01	0.0	0.0	5.00000F-02	0.0	0.0	0.0	0.0
44	6.00000E-01	0.0	0.0	5.00000F-02	0.0	0.0	0.0	0.0
45	7.20000E-01	0.0	0.0	5.00000F-02	0.0	0.0	0.0	0.0
46	0.0	0.0	0.0	2.16670E-01	0.0	0.0	0.0	0.0
47	6.00000E-01	0.0	0.0	2.16670E-01	0.0	0.0	0.0	0.0
48	6.00000E-01	0.0	0.0	2.16670E-01	0.0	0.0	0.0	0.0
49	1.20000E-01	0.0	0.0	2.16670E-01	0.0	0.0	0.0	0.0
50	1.20000E-01	0.0	0.0	2.16670E-01	0.0	0.0	0.0	0.0

NODE	PRIMARY STRUCTURE			DEGREES OF FREEDOM						
	X	Y	Z	DX	DY	DZ	RX	RY	RZ	
1	0.5	0.0	0.0	1	2	0	3	4	0	
2	0.0	3.333300E-01	0.0	5	0	0	6	7	0	
3	6.000000E-01	0.0	0.0	8	9	10	11	12	0	
4	6.000000F-01	3.333300E-01	0.0	13	0	14	15	16	0	
5	1.200000E-00	0.0	0.0	17	18	19	20	21	0	
6	1.200000E-00	3.333300F-01	0.0	22	0	23	24	25	0	
7	1.799999E-00	0.0	0.0	26	27	28	29	30	0	
8	1.799999E-00	3.333300F-01	0.0	31	0	32	33	34	0	
9	2.400000E-00	0.0	0.0	35	36	37	38	39	0	
10	2.400000E-00	3.333300F-01	0.0	40	0	41	42	43	0	
11	2.999999E-00	0.0	0.0	44	45	46	47	48	0	
12	2.999999E-00	3.333300E-01	0.0	49	0	50	51	52	0	
13	3.599999E-00	0.0	0.0	53	54	55	56	57	0	
14	3.599999E-00	3.333300E-01	0.0	58	0	59	60	61	0	
15	4.199999E-00	0.0	0.0	62	63	64	65	66	0	
16	4.199999E-00	3.333300F-01	0.0	67	0	68	69	70	0	
17	4.799999E-00	0.0	0.0	71	72	73	74	75	0	
18	4.799999E-00	3.333300F-01	0.0	76	0	77	78	79	0	
19	5.400000E-00	0.0	0.0	80	81	82	83	84	0	
20	5.400000F-00	0.0	3.333300F-01	0.0	85	0	86	87	88	0
21	5.999999E-00	0.0	0.0	89	90	91	92	93	0	
22	6.975999E-00	0.0	3.333300F-01	0.0	94	0	95	96	97	0
23	6.976999E-00	0.0	0.0	98	99	100	101	102	0	
24	7.745594E-00	3.333300E-01	0.0	103	0	104	105	106	0	
25	7.160999E-00	0.0	0.0	107	108	109	110	111	0	
26	7.195999E-00	3.333300E-01	0.0	112	0	113	114	115	0	
27	7.799999E-00	0.0	0.0	116	117	118	119	120	0	
28	7.7245594E-00	3.333300F-01	0.0	121	0	122	123	124	0	
29	8.458595E-00	0.0	0.0	125	126	127	128	129	0	
30	8.194669E-00	0.0	3.333300F-01	0.0	130	0	131	132	133	0
31	8.690967E-00	0.0	0.0	134	135	136	137	138	0	
32	8.699959E-00	3.333300F-01	0.0	139	0	140	141	142	0	
33	9.595999E-00	0.0	0.0	143	144	145	146	147	0	
34	1.029469E-00	3.333300E-01	0.0	148	0	149	150	151	0	
35	1.020000E-01	0.0	0.0	152	153	154	155	156	0	
36	1.020000F-01	3.333300E-01	0.0	157	0	158	159	160	0	
37	1.040000E-01	0.0	0.0	161	162	163	164	165	0	
38	1.060000E-01	3.333300F-01	0.0	166	0	167	168	169	0	
39	1.140000E-01	0.0	0.0	170	171	172	173	174	0	
40	1.140000F-01	3.333300F-01	0.0	175	0	176	177	178	0	
41	1.200000E-01	0.0	0.0	179	180	181	182	183	0	
42	1.200000F-01	3.333300F-01	0.0	184	0	185	186	187	0	
43	1.260000E-01	0.0	0.0	188	189	190	191	192	0	
44	1.260000F-01	3.333300E-01	0.0	193	0	194	195	196	0	
45	1.320000E-01	0.0	0.0	197	198	199	200	201	0	
46	1.320000F-01	3.333300E-01	0.0	202	0	203	204	205	0	
47	1.380000E-01	0.0	0.0	206	207	208	209	210	0	
48	1.380000F-01	3.333300E-01	0.0	211	0	212	213	214	0	
49	1.440000E-01	0.0	0.0	215	216	217	218	219	0	
50	1.440000F-01	3.333300E-01	0.0	220	0	221	222	223	0	
51	1.500000E-01	0.0	0.0	224	225	226	227	228	0	

ORDER OF UNREFINED PROBLEM = 273
NUMBER OF DESIRED MODES = 5
NUMBER OF REORTHONALIZATIONS = 2

MODE NUMBER	FREQUENCY (RAD / SEC)	FREQUENCY (HERTZ)	FRFQ. SQRD. ERROR BOUND (PERCENT)
1	3.204729E .03	5.100483E .02	7.609798E-14
2	1.096593E .04	1.745266E .03	2.610086E-13
3	2.262648E .04	3.601117E .03	4.191322E-06
4	2.400171E .04	3.819991E .03	9.051610E-06
5	3.946659E .04	6.281301E .03	1.497426E-01
6	4.502263E .04	7.165570E .03	9.145379E-02
7	6.271878E .04	9.982004E .03	4.056820E .00
8	7.672000E .04	1.221937E .04	7.066679E .01

PRIMARY STRUCTURE MODE SHAPE 1

ITÉRATION NO..0

NODE	DX	DY	DZ	RX	RY	RZ
1	-2.372577E-07	4.973345E-08	-	-	1.429838E-06	-3.161962E-02
2	-2.371776E-07	-	-	-	-1.400851E-06	-3.162404E-02
3	-2.104755E-07	4.944270E-08	-	1.893834E-02	-2.142981E-05	-3.144493E-02
4	-2.104563E-07	-	1.894102E-02	3.746948E-05	-3.144924E-02	-
5	-1.819400E-07	4.864863F-08	3.766652E-02	-4.241391E-05	-3.092045E-02	-
6	-1.838776E-07	-	3.767179E-02	7.399049E-05	-3.092459E-02	-
7	-1.578119E-07	4.731215E-08	5.597616E-02	-6.385770E-05	-3.005014E-02	-
8	-1.577289E-07	-	5.598381F-02	1.099205E-04	-3.005387E-02	-
9	-1.323493E-07	4.545071E-08	7.366067E-02	-8.621626E-05	-2.883692E-02	-
10	-1.322388E-07	-	7.367039F-02	1.444587F-04	-2.884041E-02	-
11	-1.077374E-07	4.307981E-08	9.051722F-02	-1.059153E-04	-2.729557E-02	-
12	-1.076069F-07	-	9.052908E-02	1.770441E-04	-2.729934E-02	-
13	-6.422490E-08	4.024713F-08	1.063567F-01	-1.229580F-04	-2.545035E-02	-
14	-8.406460E-08	-	1.063711E-01	2.091781F-04	-2.545403E-02	-
15	-6.143574F-08	3.657092F-08	1.210015E-01	-1.404429F-04	-2.331475F-02	-
16	-6.175902E-08	-	1.210190F-01	?392771F-04	-2.331808E-02	-
17	-4.137149F-09	3.328104F-09	-	-	-	-
18	-4.085524F-08	-	1.342934F-01	-1.547958F-04	-2.091386F-02	-
19	-2.176171F-08	2.925146F-08	1.4670547F-01	-1.638929F-04	-1.828836E-02	-
20	-2.165723F-08	-	1.466758F-01	?906900E-04	-1.829213F-02	-
21	-6.124917F-09	2.454605F-03	1.561939F-01	-1.707078F-04	-1.548051E-02	-
22	-3.875034F-09	-	1.562172E-01	3.108693F-04	-1.548352E-02	-
23	1.174755F-08	2.042313E-09	1.546020F-01	-1.776290F-04	-1.252230E-02	-
24	1.293594F-08	-	1.646270F-01	?271538E-04	-1.252472E-02	-
25	2.537176F-08	1.570584F-09	1.711965F-01	-1.823899F-04	-9.440076E-03	-
26	2.615315F-08	-	1.712228F-01	3.403833E-04	-9.441860E-03	-
27	2.614896F-08	-	-	-	-	-
28	3.915121F-08	-	-	-	-	-
29	4.851621F-08	5.017776F-09	1.759129F-01	-1.952148F-04	-6.268606F-03	-
30	4.874832F-08	-	1.759400F-01	3.493910E-04	-6.269876F-03	-
31	5.697772F-08	9.227483E-17	1.767075F-01	-1.877103F-04	-3.038223F-03	-
32	5.714250E-08	-	1.787353F-01	3.546392F-04	-3.039023F-03	-
33	-	-	1.795799F-01	3.578277F-04	2.286215E-04	-
34	-	-	-	-	?281381F-04	-

PRIMARY STRUCTURE NODES ASSOCIATED WITH TILE NO. 1

1	3	5	7	9	11	13	15	17	19	21
2	4	6	8	10	12	14	16	18	20	22

TOPS DISPLACEMENTS FOR TILE NO. 1 AND ITERATION NO. 1

MDF	Y COMPONENT(U)	Y COMPONENT(V)	Z COMPONENT(W)
1	-1.1857588E-02	-4.8645558E-07	0.0
2	-1.1859249E-02	5.2531914E-07	0.0
3	-1.1792056E-02	8.08561189E-06	1.8938444E-02
4	-1.1793669E-02	-1.4051055E-05	1.8941015E-02
5	-1.1595350E-02	1.5953861E-05	3.7666518E-02
6	-1.1569703E-02	-2.7746423E-05	3.7671786E-02
7	-1.1268958E-02	2.3993940E-05	5.997156E-02
8	-1.1270355E-02	-4.1220163E-05	5.5983808E-02
9	-1.0813974E-02	3.2376542E-05	7.3660672E-02
10	-1.0815281E-02	-5.4172007F-05	7.3670387E-02
11	-1.0237359E-02	3.9761304E-05	9.0517223E-02
12	-1.0237359E-02	-6.6391527E-05	9.0529084E-02
13	-9.5439628E-03	4.6149493E-05	1.0635674E-01
14	-9.5453411E-03	-7.8441793E-05	1.0637110E-01
15	-8.7430837F-03	5.2703050E-05	1.200154E-01
16	-8.7442367E-03	-8.9728899E-05	1.201799E-01
17	-7.8427382E-03	5.8081714E-05	1.3428342E-01
18	-7.8441054E-03	-9.9733385E-05	1.3430196E-01
19	-6.8581515E-03	6.1489074E-05	1.4605474E-01
20	-6.8595633E-03	-1.090875E-04	1.4607584E-01
21	-5.8051944E-03	6.4040360E-05	1.5619385E-01
22	-5.8063194E-03	-1.1657560F-04	1.56221716E-01
23	-4.4547564E-02	-6.3910964E-04	1.2398793F-03
24	-1.4194820E-02	-8.7081082E-04	1.48911145E-03
25	-1.4516903E-02	-5.0962073E-05	1.95530174E-02
26	-1.4168964E-02	-1.9416506E-04	1.9783155E-02
27	1.4426317E-02	4.8901746E-04	3.8002007E-02
28	1.4112670E-02	4.4811233F-04	3.9206004E-02
29	1.4063956E-02	9.4013456E-04	5.6169845E-02
30	1.3832748E-02	1.0023819E-03	5.3762256E-02
31	1.3425775E-02	1.2302840E-03	7.3761225E-02
32	1.3346210E-02	1.3594626E-03	7.3977792E-02
33	1.2545496E-02	1.2185425E-03	9.0547393E-02
34	1.2709964E-02	1.4096978E-03	9.0835929E-02
35	1.1494334E-02	7.3654414E-04	1.7633683E-01
36	1.1976253E-02	9.4510523E-04	1.7675156E-01
37	1.03376227E-02	-3.6612675E-04	1.2097675E-01
38	1.12502797E-02	-1.3315647E-04	1.2157196E-01
39	9.3781725E-03	-2.1529541E-03	1.3437933F-01
40	1.1673909E-02	-2.0691941E-03	1.3517231E-01
41	8.6540478E-02	-4.5677647E-03	1.46522973F-01
42	1.1241371E-02	-6.6531380E-03	1.4751375E-01
43	8.3560693E-03	-7.3943339E-03	1.5772129E-01
44	1.2027479E-02	-7.7313706E-03	1.5879726F-01
45	9.5742457E-03	-8.5340976E-04	2.1167265E-03
46	9.2171517E-03	-9.480652E-04	2.4439321E-03
47	9.5377430E-03	-2.2878307E-04	2.0108391E-02
48	6.1875307E-03	-2.9676839E-04	2.0416472F-02
49	6.4466721E-03	3.5079474E-04	3.9066460E-02
50	9.1335963E-03	3.47779799E-04	3.959393HF-02

STRESSES FOR ISOLATOR AND ARRESTOR FOR TILE NO. 1 AND ITERATION NO. 1

	LOCAL COORDINATES			STRESSES					
	X	Y	Z	XX	YY	ZZ	XY	YZ	ZX
ELEMENT NUMBER									
1	4.7320E-01	2.6289E-01	3.9434E-02	2.6906E 01	2.6883E 01	2.8029E 01	8.8016E-06	-1.6162E-01	1.6666E 01
2	4.7320E-01	7.0441E-02	3.9434E-02	2.2899E 01	2.2876E 01	2.3860E 01	-1.0059E-03	-1.1118E-01	1.6789E 01
3	1.2679E-01	2.6289E-01	3.9434E-02	3.7726E 01	3.7696E 01	3.9298E 01	8.0522E-04	-3.9381E-01	1.6698E 01
4	1.2679E-01	7.0441E-02	3.9434E-02	3.3534E 01	3.3504E 01	3.4934E 01	-1.0808E-03	-3.2088E-01	1.6821E 01
5	4.7320E-01	2.6289E-01	1.0566E-02	2.7413E 01	2.7399E 01	2.8531E 01	-2.2875E-04	-1.7388E-01	1.6685E 01
6	4.7320E-01	7.0441E-02	1.0566E-02	2.3408E 01	2.3395E 01	2.4363E 01	-2.4156E-04	-1.2343E-01	1.6807E 01
7	1.2679E-01	2.6289E-01	1.0566E-02	3.8398E 01	3.8385E 01	3.9965E 01	-2.5049E-04	-6.0664E-01	1.6717E 01
8	1.2679E-01	7.0441E-02	1.0566E-02	3.4207E 01	3.4194E 01	3.5603E 01	-2.6328E-04	-3.3371E-01	1.6839E 01
ELEMENT NUMBER									
1	1.0732E 00	2.6289E-01	3.9434E-02	1.6174E 01	1.6165E 01	1.6847E 01	1.5363E-03	2.2179E-01	1.6533E 01
2	1.0732E 00	7.0441E-02	3.9434E-02	1.2621E 01	1.2613F 01	1.3152E 01	-6.7633E-04	2.2994E-01	1.6655E 01
3	7.2679E-01	2.6289E-01	3.9434E-02	2.1127E 01	2.1109E 01	2.2009E 01	-3.2834E-06	3.251E-03	1.6624E 01
4	7.2679E-01	7.0441E-02	3.9434E-02	1.7262E 01	1.7246E 01	1.7990E 01	-2.1545E-03	2.6417E-02	1.6744E 01
5	1.0732E 00	2.6289E-01	1.0566E-02	1.5621E 01	1.5599E 01	1.7278E 01	-2.8970E-05	2.1105E 01	1.6542E 01
6	1.0732E 00	7.0441E-02	1.0566E-02	1.3099E 01	1.3077E 01	1.3614E 01	-1.6293E-04	2.1920E-01	1.6654E 01
7	7.2679E-01	2.6289E-01	1.0566E-02	2.1757E 01	2.1734E 01	2.2624E 01	-4.4361E-04	8.3737E-03	1.6633E 01
8	7.2679E-01	7.0441E-02	1.0566E-02	1.7924E 01	1.7901E 01	1.8635E 01	-5.7781E-04	2.4718E-02	1.6752E 01
ELEMENT NUMBER									
1	1.6732E 00	2.6289E-01	3.9434E-02	1.0661E 01	1.0681F 01	1.1127E 01	3.2492E-03	5.5905E-01	1.6188E 01
2	1.6732E 00	7.0441E-02	3.9434E-02	7.3571E 00	7.3810E 00	7.6956E 00	1.0171E-03	5.2378E-01	1.6274E 01
3	1.3268E 00	2.6289E-01	3.9434E-02	1.3076E 01	1.3038E 01	1.3648F 01	-1.5320E-04	3.7069E-01	1.6405E 01
4	1.3268E 00	7.0441E-02	3.9434E-02	9.6333E 00	9.6495E 00	1.0073E 01	-2.3954E-03	3.6053E-01	1.6508E 01
5	1.6732E 00	2.6289E-01	1.0566E-02	1.1327E 01	1.1300E 01	1.1756E 01	4.4718E-04	5.4923E-01	1.6192E 01
6	1.6732E 00	7.0441E-02	1.0566E-02	8.0979E 00	8.0767E 00	8.3925E 00	3.0674E-04	5.1397E-01	1.6278E 01
7	1.3268E 00	2.6289E-01	1.0566E-02	1.3925E 01	1.3898E 01	1.4461E 01	-4.7055E-04	3.6045E-01	1.6410E 01
8	1.3268E 00	7.0441E-02	1.0566E-02	1.0554E 01	1.0529E 01	1.0954E 01	-6.1104E-04	3.5029E-01	1.6512E 01
ELEMENT NUMBER									
1	2.2732E 00	2.6289E-01	3.9434E-02	7.8193F 00	7.8658F 00	8.1889E 00	8.8366E-03	8.0050E-01	1.5616E 01
2	2.2732E 00	7.0441E-02	3.9434E-02	4.1944F 00	4.2473F 00	4.4296F 00	4.2046E-03	7.2903E-01	1.5654E 01
3	1.9264E 00	2.6289E-01	3.9434E-02	8.9758E 00	9.0158F 00	9.3979E 00	-4.2068E-04	6.7421E-01	1.5955E 01
4	1.9264E 00	7.0441E-02	3.9434E-02	5.5634E 00	5.6104E 00	5.8604E 00	-2.0530E-03	6.2295E-01	1.6024E 01
5	2.2732E 00	2.6289E-01	1.0566E-02	6.7375F 00	6.7054F 00	9.0500F 00	1.1650F-03	7.8995E-01	1.5618E 01
6	2.2732E 00	7.0441E-02	1.0566E-02	5.241HE 00	5.2115F 00	5.4154F 00	1.1814F-03	7.1847E-01	1.5656E 01
7	1.9264E 00	2.6289E-01	1.0566E-02	1.0139E 01	1.0307E 01	1.0404F 01	-5.1494F-04	6.6431E-01	1.5957F 01
8	1.9264E 00	7.0441E-02	1.0566E-02	6.7561F 00	6.7260E 00	6.9915F 00	-4.9975E-04	6.1262E-01	1.6026E 01
ELEMENT NUMBER									
1	2.8772E 00	2.6289E-01	3.9434E-02	6.9036E 00	6.9647F 00	7.2467F 00	9.1163E-03	8.6363E-01	1.4849E 01
2	2.8772E 00	7.0441E-02	3.9434E-02	2.1431E 00	2.2185F 00	2.3172F 00	8.0451F-03	7.7022E-01	1.4808E 01
3	2.5268E 00	2.6289E-01	3.9434E-02	7.0614E 00	7.1217F 00	7.4130E 00	-9.5026E-04	8.5259E-01	1.5291E 01
4	2.5268E 00	7.0441E-02	3.9434E-02	2.9963E 00	3.0677F 00	3.2094F 00	-2.0213E-03	7.6860E-01	1.5798E 01
5	2.8772E 00	2.6289E-01	1.0566E-02	8.0533F 00	8.0143E 00	8.3259F 00	2.0526E-03	8.5493E-01	1.4850E 01
6	2.8772E 00	7.0441E-02	1.0566E-02	3.5047F 00	3.4688E 00	3.5978E 00	2.1829E-03	7.5652F-01	1.4809E 01
7	2.5268E 00	2.6289E-01	1.0566E-02	9.3207E 00	9.2815E 00	9.6030E 00	-6.4047E-04	8.4102E-01	1.5291E 01
8	2.5268E 00	7.0441E-02	1.0566E-02	4.4638E 00	4.4283F 00	4.5955E 00	-5.1222E-04	7.5703E-01	1.5300E 01

STRESSES AND DIRECT STRAINS FOR FILE NO. 1 AND ITERATION NO. 1

MFM	TEMP	LOCAL COORDINATES	STRAINS			STRESSES			STRAINS			STRESSES		
			X	Y	Z	XX	YY	ZZ	XY	YZ	ZX	XY	YY	ZZ
1	0.	0.30	0.17	0.02	3.234E-05	-2.970F-04	2.091E-02	3.056F-01	3.054E-01	3.182E-01	-1.733E-04	-2.531E-01	1.675E-01	1.675E-01
2	0.	0.90	0.17	0.02	1.015E-04	-1.875F-04	1.162E-02	1.706F-01	1.706F-01	1.777E-01	-3.217E-04	1.173E-01	1.664E-01	1.664E-01
3	0.	1.50	0.17	0.02	4.441F-06	-6.568F-05	7.209E-03	1.058E-01	1.057E-01	1.101E-01	1.742E-04	4.485E-01	1.635E-01	1.635E-01
4	0.	2.10	0.17	0.02	-8.944E-05	3.724E-05	4.898E-03	7.166F-00	7.173E-00	7.447E-00	1.112E-03	7.014E-01	1.581E-01	1.581E-01
5	0.	2.70	0.17	0.02	-1.502F-04	1.032E-04	3.722E-03	5.430F-00	5.445E-00	5.664E-00	2.159E-03	8.087E-01	1.506E-01	1.506E-01
6	0.	3.30	0.17	0.02	-1.275F-04	1.275E-04	3.488E-03	5.107E-00	5.124E-00	5.327E-00	3.080E-03	6.764E-01	1.413E-01	1.413E-01
7	0.	3.90	0.17	0.02	-8.861E-05	9.413E-05	4.449E-03	6.587E-00	6.598E-00	6.861E-00	3.630E-03	6.052E-01	1.308E-01	1.308E-01
8	0.	4.50	0.17	0.02	7.243F-05	-1.763F-05	4.477F-03	1.115E-01	1.115E-01	1.160E-01	3.314E-03	-6.779E-01	1.199E-01	1.199E-01
9	0.	5.10	0.17	0.02	3.389E-04	-2.475E-04	1.440E-02	2.146E-01	2.143E-01	2.231E-01	2.243E-03	-1.978E-00	1.093E-01	1.093E-01
10	0.	5.70	0.17	0.02	6.551F-04	-5.952E-04	3.010E-02	4.469F-01	4.462F-01	4.647E-01	-4.048E-04	-3.614E-00	9.993E-00	9.993E-00
11	0.	6.30	0.17	0.13	-4.985E-05	-3.958F-04	4.535E-03	-1.038E-00	3.234E-00	2.753E-01	8.851E-02	-1.630E-01	1.170E-01	1.170E-01
12	0.	6.90	0.17	0.13	-1.229E-04	-1.840F-04	2.811F-03	-5.491E-00	3.418F-00	1.720E-01	2.288E-02	1.013E-01	2.029E-01	2.029E-01
13	0.	7.50	0.17	0.13	-4.461E-04	6.061E-05	6.627F-03	1.664E-01	1.664E-01	1.722E-00	5.416E-01	3.229E-01	1.916E-01	1.916E-01
14	0.	8.10	0.17	0.13	-7.918F-04	2.861F-04	9.667E-04	-4.907E-01	1.244E-00	5.628E-00	1.432E-00	4.998E-01	1.798E-01	1.798E-01
15	0.	8.70	0.17	0.13	-1.072F-03	4.517F-04	6.049F-04	-6.545F-01	2.283E-00	3.336E-00	2.375E-00	5.264E-01	1.622E-01	1.622E-01
16	0.	9.30	0.17	0.13	-1.241F-03	5.412F-04	4.851F-04	-7.593E-01	2.955E-00	2.539E-00	3.089E-00	3.452E-01	1.390E-01	1.390E-01
17	0.	9.90	0.17	0.13	-1.259F-03	5.281F-04	6.761F-04	-7.714E-01	3.226E-00	3.657E-00	3.282E-00	-8.202E-02	1.107E-01	1.107E-01
18	0.	4.50	0.17	0.13	-1.461E-04	3.784F-04	1.402F-03	-6.698E-01	2.716E-00	2.644E-00	2.644E-00	-7.877E-01	7.748E-00	7.748E-00
19	0.	5.10	0.17	0.13	-7.355F-04	3.815F-05	3.189E-03	-4.440E-01	9.126E-01	1.907E-01	1.313E-00	-1.602E-00	3.380E-00	3.380E-00
20	0.	5.70	0.17	0.14	-3.018E-04	-4.679F-04	6.799E-03	-1.473E-01	5.936E-00	4.137E-01	-1.089E-00	-2.552E-00	-1.602E-01	-1.602E-01
21	0.	6.30	0.17	0.14	-6.129F-05	-1.187E-04	3.741E-03	-5.198E-00	7.484F-00	2.732F-01	2.389E-01	-1.448E-01	1.003E-01	1.003E-01
22	0.	6.90	0.17	0.14	-1.123F-04	-2.359F-05	2.569E-03	-7.267E-00	-4.119E-00	1.530E-01	4.555E-01	5.484E-02	2.092E-01	2.092E-01
23	0.	7.50	0.17	0.14	-2.8679F-04	9.123E-05	1.546F-03	-1.749F-01	-2.363E-00	9.080E-03	7.577E-01	1.650E-01	2.282E-01	2.282E-01
24	0.	8.10	0.17	0.14	-5.154F-04	2.763F-04	8.677F-04	-3.120E-01	-1.536E-00	4.879F-00	1.205E-00	2.657E-01	2.101E-01	2.101E-01
25	0.	8.70	0.17	0.14	-7.053E-04	3.332E-04	4.927E-04	-4.927E-01	-1.054E-00	2.520E-00	1.556E-00	4.460E-01	1.784E-01	1.784E-01
26	0.	9.30	0.17	0.14	-8.065E-04	3.879F-04	2.527E-04	-4.868F-01	-9.017E-01	1.620F-00	1.566E-00	7.629E-02	1.338E-01	1.338E-01
27	0.	9.90	0.17	0.14	-7.354E-04	3.714F-04	6.143F-04	-4.745F-01	-1.180F-01	2.599E-00	1.068F-00	-2.490E-01	7.850E-00	7.850E-00
28	0.	4.50	0.17	0.14	-2.8679F-04	2.682E-04	1.226F-03	-3.787E-01	-2.363E-00	9.080E-03	7.577E-01	1.650E-01	2.282E-01	2.282E-01
29	0.	5.10	0.17	0.14	-5.154F-04	2.763F-04	8.677F-04	-3.120E-01	-1.536E-00	4.879F-00	1.205E-00	2.657E-01	2.101E-01	2.101E-01
30	0.	5.70	0.17	0.14	-7.053E-04	3.332E-04	4.927E-04	-4.927E-01	-1.054E-00	2.520E-00	1.556E-00	4.460E-01	1.784E-01	1.784E-01
31	0.	6.30	0.17	0.14	-8.065E-04	3.879F-04	2.527E-04	-4.868F-01	-9.017E-01	1.620F-00	1.566E-00	7.629E-02	1.338E-01	1.338E-01
32	0.	6.90	0.17	0.14	-6.240E-04	2.905E-03	2.905E-03	-3.857E-00	3.206E-01	1.740E-01	2.590E-01	1.068F-00	-2.490E-01	7.850E-00
33	0.	7.50	0.17	0.14	-2.8679F-04	2.682E-04	1.226F-03	-3.787E-01	-2.363E-00	9.080E-03	7.577E-01	1.650E-01	2.282E-01	2.282E-01
34	0.	8.10	0.17	0.14	-5.154F-04	2.763F-04	8.677F-04	-3.120E-01	-1.536E-00	4.879F-00	1.205E-00	2.657E-01	2.101E-01	2.101E-01
35	0.	8.70	0.17	0.14	-7.053E-04	3.332E-04	4.927E-04	-4.927E-01	-1.054E-00	2.520E-00	1.556E-00	4.460E-01	1.784E-01	1.784E-01
36	0.	9.30	0.17	0.14	-8.065E-04	3.879F-04	2.527E-04	-4.868F-01	-9.017E-01	1.620F-00	1.566E-00	7.629E-02	1.338E-01	1.338E-01
37	0.	9.90	0.17	0.14	-6.240E-04	2.905E-03	2.905E-03	-3.857E-00	3.206E-01	1.740E-01	2.590E-01	1.068F-00	-2.490E-01	7.850E-00
38	0.	4.50	0.17	0.14	-2.8679F-04	2.682E-04	1.226F-03	-3.787E-01	-2.363E-00	9.080E-03	7.577E-01	1.650E-01	2.282E-01	2.282E-01
39	0.	5.10	0.17	0.14	-5.154F-04	2.763F-04	8.677F-04	-3.120E-01	-1.536E-00	4.879F-00	1.205E-00	2.657E-01	2.101E-01	2.101E-01
40	0.	5.70	0.17	0.14	-7.053E-04	3.332E-04	4.927E-04	-4.927E-01	-1.054E-00	2.520E-00	1.556E-00	4.460E-01	1.784E-01	1.784E-01
41	0.	6.30	0.17	0.14	-8.065E-04	3.879F-04	2.527E-04	-4.868F-01	-9.017E-01	1.620F-00	1.566E-00	7.629E-02	1.338E-01	1.338E-01
42	0.	6.90	0.17	0.14	-6.240E-04	2.905E-03	2.905E-03	-3.857E-00	3.206E-01	1.740E-01	2.590E-01	1.068F-00	-2.490E-01	7.850E-00
43	0.	7.50	0.17	0.14	-2.8679F-04	2.682E-04	1.226F-03	-3.787E-01	-2.363E-00	9.080E-03	7.577E-01	1.650E-01	2.282E-01	2.282E-01
44	0.	8.10	0.17	0.14	-5.154F-04	2.763F-04	8.677F-04	-3.120E-01	-1.536E-00	4.879F-00	1.205E-00	2.657E-01	2.101E-01	2.101E-01
45	0.	8.70	0.17	0.14	-7.053E-04	3.332E-04	4.927E-04	-4.927E-01	-1.054E-00	2.520E-00	1.556E-00	4.460E-01	1.784E-01	1.784E-01
46	0.	9.30	0.17	0.14	-8.065E-04	3.879F-04	2.527E-04	-4.868F-01	-9.017E-01	1.620F-00	1.566E-00	7.629E-02	1.338E-01	1.338E-01
47	0.	9.90	0.17	0.14	-6.240E-04	2.905E-03	2.905E-03	-3.857E-00	3.206E-01	1.740E-01	2.590E-01	1.068F-00	-2.490E-01	7.850E-00
48	0.	4.50	0.17	0.14	-2.8679F-04	2.682E-04	1.226F-03	-3.787E-01	-2.363E-00	9.080E-03	7.577E-01	1.650E-01	2.282E-01	2.282E-01
49	0.	5.10	0.17	0.14	-5.154F-04	2.763F-04	8.677F-04	-3.120E-01	-1.536E-00	4.879F-00	1.205E-00	2.657E-01	2.101E-01	2.101E-01
50	0.	5.70	0.17	0.14	-7.053E-04	3.332E-04	4.927E-04	-4.927E-01	-1.054E-00	2.520E-00	1.556E-00	4.460E-01	1.784E-01	1.784E-01
51	0.	6.30	0.17	0.14	-8.065E-04	3.879F-04	2.527E-04	-4.868F-01	-9.017E-01	1.620F-00	1.566E-00	7.629E-02	1.338E-01	1.338E-01
52	0.	6.90	0.17	0.14	-6.240E-04	2.905E-03	2.905E-03	-3.857E-00	3.206E-01	1.740E-01	2.590E-01	1.068F-00	-2.490E-01	7.850E-00
53	0.	7.50	0.17	0.14	-2.8679F-04	2.682E-04	1.226F-03	-3.787E-01	-2.363E-00	9.080E-03	7.577E-01	1.650E-01	2.282E-01	2.282E-01
54	0.	8.10	0.17	0.14	-5.154F-04	2.763F-04	8.677F-04	-3.120E-01	-1.536E-00	4.879F-00	1.205E-00	2.657E-01	2.101E-01	2.101E-01
55	0.	8.70	0.17	0.14	-7.053E-04	3.332E-04	4.927E-04	-4.927E-01	-1.054E-00	2.520E-00	1.556E-00	4.460E-01	1.784E-01	1.784E-01
56	0.	9.30	0.17	0.14	-8.065E-04	3.879F-04	2.527E-04	-4.868F-01	-9.017E-01	1.620F-00	1.566E-00	7.629E-02	1.338E-01	1.338E-01
57	0.	9.90	0.17	0.14	-6.240E-04	2.905E-03	2.905E-03	-3.857E-00	3.206E-01	1.740E-01	2.590E-01	1.068F-00	-2.490E-01	7.850E-00

SUM T1E DEN = 2.18633D-06

P.S. DEN = 1.37088D-06

SUM DEN = 4.06620D-06

DNMGA = 2.64955E-03
DNMGA_SQUARE(D) = 7.12655E-06

PRIMARY STRUCTURE DEFLECTIONS FOR ITERATION NO. 1

NODE	DX	DY	DZ	RX	RY	RZ
33	-	-	-	-	-	-
34	-	6.681679E-05	-1.673227E-06	1.785551E-01	-5.587914E-05	3.189787E-03
	-	6.6930365E-05	-	1.786259E-01	4.809038E-04	3.191435E-03
35	-	-	-	-	-	-
36	-	5.671559E-05	-1.613901E-06	1.756771E-01	-5.725912E-05	6.399900E-03
	-	5.673004E-05	-	1.757467E-01	4.751717E-04	6.402336E-03
37	-	-	-	-	-	-
38	-	4.701459E-05	-1.5366437E-06	1.708924E-01	-5.781297E-05	9.573631E-03
	-	4.699403E-05	-	1.709508E-01	4.673784E-04	9.576578E-03
39	-	-	-	-	-	-
40	-	3.7779292E-05	-1.429681E-06	1.642013E-01	-5.394027E-05	1.267944E-02
	-	3.7756843E-05	-	1.642683E-01	4.556030E-04	1.268287E-02
41	-	-	-	-	-	-
42	-	2.907454E-05	-1.320172F-06	1.556897E-01	-4.895801F-05	1.566771E-02
	-	2.905251E-05	-	1.557546E-01	4.377079E-04	1.567218E-02
43	-	-	-	-	-	-
44	-	2.120507E-05	-1.232098E-06	1.454340E-01	-3.918960E-05	1.847837F-02
	-	2.116353E-05	-	1.454960E-01	4.109822E-04	1.848471E-02
45	-	-	-	-	-	-
46	-	1.428868E-05	-1.367070F-06	1.335610F-01	-3.102203E-05	2.105218E-02
	-	1.423488E-05	-	1.336186F-01	3.767714E-04	2.106069E-02
47	-	-	-	-	-	-
48	-	8.277575E-06	-8.915556E-07	1.202210F-01	-2.576198F-05	2.336588E-02
	-	8.214299F-07	-	1.202730F-01	3.377073F-04	2.337612E-02
49	-	-	-	-	-	-
50	-	2.616151E-06	-7.026549F-07	1.055738E-01	-2.197572F-05	2.540788E-02
	-	2.477469E-06	-	1.056194E-01	2.953170E-04	2.541899E-02
51	-	-	-	-	-	-
52	-	-1.711983F-07	-5.027475E-07	8.979516E-02	-1.980127E-05	2.716987E-02
	-	-2.311528F-07	-	8.982366E-02	2.503702E-04	2.718170E-02
53	-	-	-	-	-	-
54	-	-2.632518E-06	-2.924773F-07	7.302600F-02	-1.593972E-05	2.863894F-02
	-	-2.742852E-06	-	7.305735E-02	2.037658E-04	2.865112E-02
55	-	-	-	-	-	-
56	-	-3.945375E-06	-7.611419E-08	5.547747E-02	-1.253435F-05	2.979947E-02
	-	-4.037647E-06	-	5.551126F-02	1.549640E-04	2.981209E-02
57	-	-	-	-	-	-
58	-	-2.917401E-06	-1.474601F-07	3.732769E-02	-9.897214E-06	3.064273E-02
	-	-2.954422E-06	-	3.734367F-02	1.049516E-04	3.065582E-02
59	-	-	-	-	-	-
60	-	-2.677151E-06	-2.967929F-07	1.476934E-02	-5.443655E-06	3.16091E-02
	-	-2.645435E-06	-	1.477641F-02	5.394142E-05	3.117428E-02
61	-	-	-	-	-	-
62	-	-5.462543E-07	-	-	1.710423E-06	3.133634E-02
	-	-	-	-	-	3.134975E-02
63	-	-	-	-	-	-

MAXIMUM DEFLECTION:

= 1.79574E-31

FOR DDF 140

MAXIMUM DEFLECTION DIFFERENCE = 7.54654E-34 FOR DDF 194

MID-POINT PLATE MEMBER STRAINS AND STRESSES FOR ITERATION NO. 1

MEMBER	COORDINATES		STRAINS		STRESSES	
	X	Y	EPS X	EPS Y	SIG X	SIG Y
1	-	3.0000E-01	1.6666E-01	-	-1.1270F-06	2.8437E-07
2	-	9.0000E-01	1.6666E-01	-	-3.0212F-06	8.1109E-07
3	-	1.5000E-00	1.6666E-01	-	-4.9522F-06	1.4223E-06
4	-	2.1000E-00	1.6666E-01	-	-6.8412F-06	1.9951E-06
5	-	2.7000E-00	1.6666E-01	-	-8.6617F-06	2.5449F-06
6	-	3.3000E-00	1.6666E-01	-	-1.0399E-05	3.0682E-06
7	-	3.9000E-00	1.6666E-01	-	-1.2010E-05	3.5613E-06
8	-	4.5000E-00	1.5666E-01	-	-1.3517E-05	4.0194E-06
9	-	5.1000E-00	1.6666E-01	-	-1.4923E-05	4.4374E-06
10	-	5.7000E-00	1.6666E-01	-	-1.6246E-05	4.7443E-06
11	-	6.3000E-00	1.5666E-01	-	-1.7193E-05	4.9652E-06
12	-	6.9000E-00	1.6666E-01	-	-1.7465E-05	5.1359E-06
13	-	7.5000E-00	1.6666E-01	-	-1.7671E-05	5.2154E-06
14	-	8.1000E-00	1.5666E-01	-	-1.7738E-05	5.2275E-06
15	-	8.7000E-00	1.6666E-01	-	-1.7630E-05	5.1921E-06
16	-	9.3000E-00	1.5666E-01	-	-1.7330E-05	5.0826E-06
17	-	9.9000E-00	1.6666E-01	-	-1.6337E-05	4.9377E-06
18	-	1.0500E-01	1.5666E-01	-	-1.6173E-05	4.7256E-06
19	-	1.1100E-01	1.6666E-01	-	-1.5373E-05	4.4492E-06
20	-	1.1700E-01	1.6666E-01	-	-1.4524E-05	4.1243E-06
21	-	1.2300E-01	1.5666E-01	-	-1.3132E-05	3.8294E-06
22	-	1.2900E-01	1.6666E-01	-	-1.1539E-05	3.4483E-06
23	-	1.3500E-01	1.5666E-01	-	-9.8535E-06	2.9333E-06
24	-	1.4100E-01	1.6666E-01	-	-8.0711E-06	2.3913E-06
25	-	1.4700E-01	1.5666E-01	-	-6.1804E-06	1.3081E-06

PROGRAM LISTING OF INPUT DATA CARDS

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O P T I O N S

STATICS PROBLEM

MAXIMUM NO. ITERATIONS = 3

CONVERGENCE PARAMETER = 0.0

PRIMARY STRUCTURE STRESSES PRESENTED AFTFR EACH ITERATION AT PLATE MID, TOP AND BOTTOM SURFACES

TILFS ON PRIMARY STRUCTURE

TILF STRESSES PRESENTED AFTER LAST ITERATION

TILF NODE MAP NOT REQUESTED

TILF ELEMENT MAP REQUESTED

TILF NODE COORDINATES REQUESTED

DO NOT PRINT ELEMENT STIFFNESS MATRICES

DO NOT PRINT ASSEMBLED STIFFNESS MATRICES

DO NOT PRINT TILF REQUESTED INFORMATION

COMPUTE STRESSES FOR ALL TILFS

GEOMETRY

PLATE LX = 6.66667E 00 LY = 3.33330E-01 TP = 7.50000E-01
STOPGERS Y1 = 4.00000E-01 TS = 0.0 VS = 0.0 AS = 0.0
 Y1 = 0.0 T1 = 0.0 JX1 = 0.0 BETAS = 0.0

TILES NX0 = 1 NYR = 1 NR = 1 R1 = 0.0 R2 = 1.16667E 00
 TA = 1.66670E-01 TI = 5.00000E-02 TC = 0.0
 NR1 = 0 NB2 = 10 ND2 = 1
 NT1 = 0 NT2 = 7

HATF01

HATF01 LINES FEATURES
PLATE ED = 1.00000E 07 NJD = 3.00000E-01 GAMMA P = 0.0 ALPHA P = 1.31000E-05
STOPGERS ES = 0.0 NS = 0.0 GAMMA S = 0.0 ALPHA S = 0.0
ADFSY100 IX = 5.00000E 04 CY = 5.00000E 04 F7 = 6.00000E 03
OR_PCF IY XY = 5.00000E-01 NU Y7 = 1.00000CF-01 NU ZX = 1.00000E-02
 CY = 3.00000E 04 CY7 = 3.02000E 04 G7X = 3.20000E 04
 GAMMA A = 0.0
 ALPHA V = 9.999999E-07 ALPHA Y = 9.999999E-07 ALPHA Z = 9.999999E-07
 IY = 4.00000E 01 NJI I = 4.00000E-01 GAMMA I = 0.0 ALPHA I = 2.71000E-04
 GAMMA Q = 0.0

251 GAMMA Q = 0.0 ALPHA QY / ALPHA RX = 1.00000E 00 ALPHA RZ / ALPHA RX = 1.00000E 00

TEMPERATURE DEPENDENT MATERIAL PROPERTIES

		PROPERTY	TEMPERATURE	PROPERTY	TEMPERATURE	PROPERTY	TEMPERATURE
1	EQ	ALL	6.000E-04				
1	ER	ALL	6.000E-03				
1	Ga	ALL	3.200E-04				
1	NU_R	ALL	5.000E-01				
1	NU_R	ALL	1.000E-02				
1	ALPHA_R	ALL	3.470E-07				
1	FC	ALL	0.0				
1	NU_C	ALL	0.0				
1	ALPHA_C	ALL	0.0				

B O U N D A R Y C O N D I T I O N S

EDGE	PLATE OUT OF PLANE	PLATE IN PLANE	STRINGERS
A	PINNED	FREE	FREE
B	PINNED	U HELD, V FREE	FREE
C	FREE	FREE	FREE
D	FREE	V HELD, U FREE	FREE
E	FREE	FREE	FREE
F	FREE	FREE	FREE
G	FREE	FREE	FREE
H	FREE	FREE	FREE
I	FREE	FREE	FREE
J	FREE	FREE	FREE
K	FREE	FREE	FREE
L	FREE	FREE	FREE
M	FREE	FREE	FREE
N	FREE	FREE	FREE
O	FREE	FREE	FREE
P	FREE	FREE	FREE
Q	FREE	FREE	FREE
R	FREE	FREE	FREE
S	FREE	FREE	FREE
T	FREE	FREE	FREE
U	FREE	FREE	FREE
V	FREE	FREE	FREE
W	FREE	FREE	FREE
X	FREE	FREE	FREE
Y	FREE	FREE	FREE
Z	FREE	FREE	FREE

R S I - T E M P E R A T U R E S

STATIC THERMAL LOADING

UNIFORM TEMPERATURE OPTION

T REFERENCE = 7.0000E 01

DEL T U = T - T REF = -3.2500E 02

FILEMENT MAP

**LAYER 1
RSI**

81 82 83 84 85 86 87 88 89 90

TITLE MESH
ELEMENT

		TITLE NODES	
	ELEMENT	1	2
1	1	3	4
2	2	5	6
3	3	7	8
4	4	9	10
5	5	11	12
6	6	13	14
7	7	15	16
8	8	17	18
9	9	19	20
10	10	21	22
11	11	23	24
12	12	25	26
13	13	27	28
14	14	29	30
15	15	31	32
16	16	33	34
17	17	35	36
18	18	37	38
19	19	39	40
20	20	41	42
21	21	43	44
22	22	45	46
23	23	47	48
24	24	49	50
25	25	51	52
26	26	53	54
27	27	55	56
28	28	57	58
29	29	59	60
30	30	61	62
31	31	63	64
32	32	65	66
33	33	67	68
34	34	69	70
35	35	71	72
36	36	73	74
37	37	75	76
38	38	77	78
39	39	79	80
40	40	81	82
41	41	83	84

NODE

LPCAL TIME COORDINATES

TEMPERATUR

	X	Y	Z	TIME COORDINATES	TEMPERATUR
1	0.0	0.0	0.0	3.33330E-01	-3.25000E 02
2	0.0	0.0	0.0	3.33330F-01	-3.25000E 02
3	6.66667E-01	0.0	0.0	3.33330F-01	-3.25000E 02
4	6.66667E-01	0.0	0.0	3.33330F-01	-3.25000E 02
5	1.33333E-01	0.0	0.0	3.33330F-01	-3.25000E 02
6	1.33333E-01	0.0	0.0	3.33330F-01	-3.25000E 02
7	2.00000C 00	0.0	0.0	3.33330F-01	-3.25000E 02
8	2.02000F 00	0.0	0.0	3.32230F-01	-3.25000F 02
9	2.66667E 00	0.0	0.0	3.33330F-01	-3.25000F 02
10	2.66667E 00	0.0	0.0	2.333330F-01	-3.25000E 02
11	3.33333E 00	0.0	0.0	2.33333E 00	-3.25000E 02
12	3.33333E 00	0.0	0.0	2.333330F-01	-3.25000F 02
13	4.00000C 00	0.0	0.0	2.333330F-01	-3.25000F 02
14	4.66667E 00	0.0	0.0	2.333330F-01	-3.25000E 02
15	4.66667E 00	0.0	0.0	2.333330F-01	-3.25000F 02
16	5.33333E 00	0.0	0.0	3.333330F-01	-3.25000E 02
17	5.33333E 00	0.0	0.0	3.333330F-01	-3.25000F 02
18	5.99999C 00	0.0	0.0	3.333330F-01	-3.25000F 02
19	6.66667E 00	0.0	0.0	3.333330F-01	-3.25000E 02
20	6.66667E 00	0.0	0.0	3.332230F-01	-3.25000F 02
21	6.66667E 00	0.0	0.0	3.332230F-01	-3.25000E 02
22	6.66667E 00	0.0	0.0	3.332230F-01	-3.25000F 02
23	7.0	0.0	0.0	3.332230F-01	-3.25000F 02
24	7.0	0.0	0.0	3.332230F-01	-3.25000F 02
25	6.66667E-01	0.0	0.0	3.332230F-01	-3.25000F 02
26	6.66667E-01	0.0	0.0	3.332230F-01	-3.25000F 02
27	1.33333E 00	0.0	0.0	3.332230F-01	-3.25000E 02
28	1.33333E 00	0.0	0.0	3.332230F-01	-3.25000E 02
29	2.00000C 00	0.0	0.0	3.332230F-01	-3.25000F 02
30	2.66667E 00	0.0	0.0	3.332230F-01	-3.25000E 02
31	2.66667E 00	0.0	0.0	3.332230F-01	-3.25000F 02
32	3.33333E 00	0.0	0.0	3.332230F-01	-3.25000E 02
33	3.33333E 00	0.0	0.0	3.332230F-01	-3.25000F 02
34	4.00000C 00	0.0	0.0	3.332230F-01	-3.25000E 02
35	4.66667E 00	0.0	0.0	3.332230F-01	-3.25000F 02
36	4.66667E 00	0.0	0.0	3.332230F-01	-3.25000E 02
37	5.33333E 00	0.0	0.0	3.332230F-01	-3.25000F 02
38	6.00000C 00	0.0	0.0	3.332230F-01	-3.25000E 02
39	6.66667E 00	0.0	0.0	3.332230F-01	-3.25000F 02
40	7.0	0.0	0.0	3.332230F-01	-3.25000E 02
41	7.0	0.0	0.0	3.332230F-01	-3.25000F 02
42	7.66667E 00	0.0	0.0	3.332230F-01	-3.25000E 02
43	8.33333E 00	0.0	0.0	3.332230F-01	-3.25000F 02
44	8.33333E 00	0.0	0.0	3.332230F-01	-3.25000E 02
45	9.0	0.0	0.0	3.332230F-01	-3.25000F 02
46	9.0	0.0	0.0	3.332230F-01	-3.25000F 02
47	9.66667E 00	0.0	0.0	3.332230F-01	-3.25000E 02
48	10.33333E 00	0.0	0.0	3.332230F-01	-3.25000F 02
49	11.0	0.0	0.0	3.332230F-01	-3.25000E 02
50	11.66667E 00	0.0	0.0	3.332230F-01	-3.25000F 02

NODE	X	Y	Z	PRIMARY STRUCTURE						SECONDARY STRUCTURE						DEGREES OF FREEDOM					
				DX	DY	DZ	RX	RY	RZ	DX	DY	DZ	RX	RY	RZ	DX	DY	DZ	RX	RY	RZ
1	0.0	0.0	0.0	0.0	0.0	0.0	0.0	0.0	0.0	1	2	0	3	4	0	1	2	0	6	7	0
2	0.0	6.666670E-01	0.0	3.333300E-01	0.0	0.0	0.0	0.0	0.0	5	0	0	0	0	0	10	11	12	0	0	0
3	6.666670E-01	0.0	3.333300F-01	0.0	0.0	0.0	0.0	0.0	0.0	8	9	0	0	0	0	14	15	16	0	0	0
4	6.666670E-01	3.333300F-01	0.0	0.0	0.0	0.0	0.0	0.0	0.0	13	0	0	0	0	0	17	18	19	20	21	0
5	1.33334F 00	0.0	0.0	0.0	0.0	0.0	0.0	0.0	0.0	17	18	0	0	0	0	20	21	22	23	24	0
6	1.333334E 00	3.333300E-01	0.0	0.0	0.0	0.0	0.0	0.0	0.0	22	0	0	0	0	0	23	24	25	26	27	0
7	2.0000001E 00	0.0	0.0	0.0	0.0	0.0	0.0	0.0	0.0	26	27	0	0	0	0	28	29	30	31	32	0
8	2.0000001E 00	3.333300E-01	0.0	0.0	0.0	0.0	0.0	0.0	0.0	31	0	0	0	0	0	32	33	34	35	36	0
9	2.666668E 00	0.0	0.0	0.0	0.0	0.0	0.0	0.0	0.0	35	36	0	0	0	0	37	38	39	39	39	0
10	2.666668F 00	3.333300F-01	0.0	0.0	0.0	0.0	0.0	0.0	0.0	40	0	0	0	0	0	41	42	43	43	43	0
11	3.333335E 00	0.0	0.0	0.0	0.0	0.0	0.0	0.0	0.0	44	45	0	0	0	0	46	47	48	48	48	0
12	3.333335F 00	3.333300F-01	0.0	0.0	0.0	0.0	0.0	0.0	0.0	49	50	0	0	0	0	51	52	52	53	54	0
13	4.0000002F 00	0.0	0.0	0.0	0.0	0.0	0.0	0.0	0.0	53	54	0	0	0	0	55	56	57	57	57	0
14	4.0000002E 00	3.333300E-01	0.0	0.0	0.0	0.0	0.0	0.0	0.0	58	0	0	0	0	0	59	60	61	61	61	0
15	4.666669E 00	0.0	0.0	0.0	0.0	0.0	0.0	0.0	0.0	62	63	0	0	0	0	64	65	66	66	66	0
16	4.666669F 00	3.333300F-01	0.0	0.0	0.0	0.0	0.0	0.0	0.0	67	0	0	0	0	0	68	69	70	70	70	0
17	5.333336E 00	0.0	0.0	0.0	0.0	0.0	0.0	0.0	0.0	71	72	0	0	0	0	73	74	75	75	75	0
18	5.333336F 00	3.333300F-01	0.0	0.0	0.0	0.0	0.0	0.0	0.0	76	0	0	0	0	0	77	78	79	79	79	0
19	6.0000003F 00	0.0	0.0	0.0	0.0	0.0	0.0	0.0	0.0	80	81	0	0	0	0	82	83	84	84	84	0
20	6.0000003E 00	3.333300E-01	0.0	0.0	0.0	0.0	0.0	0.0	0.0	85	0	0	0	0	0	86	87	88	88	88	0
21	6.6666670E 00	0.0	0.0	0.0	0.0	0.0	0.0	0.0	0.0	99	0	0	0	0	0	90	91	92	93	93	0
22	6.6666670F 00	3.333300F-01	0.0	0.0	0.0	0.0	0.0	0.0	0.0	0	0	0	0	0	0	92	93	93	93	93	0

PRIMARY STRUCTURE DEFLECTIONS FOR ITERATION NO. 1

NODE	DX	DY	DZ	RX	RY	RZ
1	-	-	-	-	-	-
2	2.838170E-02	-	1.419157E-03	-	-	0.0
3	-	-	-	-	-	0.0
4	2.554341E-02	-	1.419158E-03	0.0	-	0.0
5	2.554342E-02	-	-	0.0	-	0.0
6	-	-	-	-	-	-
7	2.270516E-02	-	1.419166E-03	0.0	-	0.0
8	2.270516E-02	-	-	0.0	-	0.0
9	-	-	-	-	-	-
10	1.936694E-02	-	1.419171E-03	0.0	-	0.0
11	1.936694E-02	-	-	0.0	-	0.0
12	-	-	-	-	-	-
13	1.419059E-02	-	1.419178E-03	0.0	-	0.0
14	1.419060E-02	-	-	0.0	-	0.0
15	-	-	-	-	-	-
16	8.514372E-03	-	1.419182E-03	0.0	-	0.0
17	8.514374E-03	-	-	0.0	-	0.0
18	-	-	-	-	-	-
19	2.000105E-03	-	1.419196E-03	0.0	-	0.0
20	2.000106E-03	-	-	0.0	-	0.0
21	-	-	-	-	-	-
22	-	-	-	1.419196E-03	-	0.0
	-	-	-	-	-	0.0

PRIMARY STRUCTURE NODES ASSOCIATED WITH FILE NO. 1

1	3	5	7	9	11	13	15	17	19	21
2	4	6	8	10	12	14	16	18	20	22

MID-POINT PLATE MEMBER STRAINS AND STRESSES FOR ITERATION NO. 1

MEMREP	COORDINATES		STRAINS		STRESSES			
	X	Y	EPS X	EPS Y	EPS XY	SIG X	SIG Y	
1	-3.3333F-01	1.6666E-01	-4.2574E-03	-4.2575E-03	2.5495F-08	6.5625E-01	8.2031E-02	1.8211E-01
2	1.0000F 00	1.6666E-01	-4.2574E-03	-4.2575E-03	1.4232E-08	1.3516E 00	1.6406E-01	1.0166E-01
3	1.5667F 10	1.6666E-01	-4.2573E-03	-4.2575E-03	1.8335F-09	1.8008F 00	8.2031E-02	1.3097E-02
4	2.3333E 00	1.6666E-01	-4.2573E-03	-4.2576F-03	7.0723E-09	2.3750E 00	1.2500E-01	5.0516E-02
5	3.0000F 20	1.6666E-01	-4.2572E-03	-4.2576E-03	1.2049E-08	2.6211E 00	1.2500E-01	8.6065E-02
6	3.6667E 20	1.6666E-01	-4.2572E-03	-4.2575F-03	1.5629E-08	2.9063E 00	8.2031E-02	1.1163E-01
7	4.2223F 30	1.6666E-01	-4.2572E-03	-4.2576F-03	1.6939E-08	3.2344E 00	1.2500E-01	1.2099E-01
8	5.0000F 10	1.6666E-01	-4.2572E-03	-4.2576F-03	9.5170F-09	3.2555E 00	1.2500E-01	6.7979E-02
9	5.6667F 30	1.6666E-01	-4.2572E-03	-4.2576F-03	1.7462F-10	3.5195F 00	8.2031E-02	1.2473E-03
10	6.3333F 20	1.6656E-01	-4.2572F-03	-4.2576F-03	8.7311F-11	3.5195E 00	8.2031E-02	6.2365E-04

PRIMARY STRUCTURE DEFLECTIONS FOR ITERATION NO. 2

NODE	DX	DY	DZ	RX	RZ
1	-	-	-	-	-
2	2.837436E-02	1.419182E-03	-	1.110410E-08	-4.857246E-05
3	2.553628E-02	1.419224E-03	-	-2.494943E-07	-4.696766E-05
4	-	-	-	3.204057E-05	-1.107075E-08
5	2.269858E-02	1.419286E-03	-	3.203745E-05	2.304330E-07
6	2.269857E-02	-	-	6.173004E-05	-6.444173E-07
7	1.986120E-02	1.419325E-03	-	6.172433E-05	6.098011E-07
8	1.935119E-02	-	-	8.592271E-05	-9.869391E-07
9	-	-	-	8.591506E-05	9.408456E-07
10	1.702403E-02	1.419350F-03	-	1.017759E-04	-1.215332E-06
11	1.413696E-02	1.419358E-03	-	1.017165E-04	1.158402E-06
12	1.413696E-02	-	-	1.072070E-04	-1.294960F-06
13	1.134495E-02	1.419356E-03	-	1.016948E-04	-1.215301E-06
14	1.134499E-02	-	-	1.016850E-04	1.155909F-06
15	8.512772E-03	1.419315F-03	-	8.587058E-05	-1.233488E-06
16	8.512784E-03	-	-	1.071968E-04	2.492761E-08
17	5.675440E-02	1.419301E-03	-	-	-
18	5.675441E-02	-	-	-	-
19	2.937912E-02	1.4192743E-03	-	3.200154E-05	-2.472392E-07
20	2.937926E-02	-	-	3.199324E-05	2.270913E-07
21	-	-	-	1.419209F-03	-
22	-	-	-	-	-

MAXIMUM DEFLECTION =

= 2.93744E-02 DEF. UNITS = 5

MAXIMUM DEFLECTION DIFFERENCE = 1.072075E-04 DEF. UNITS = 46

MAXIMUM SURFACE PARAMETER = 3.077830E-02

STRUCTURE HAS NOT CONVERGED

TPS DISPLACEMENTS FOR TIEF NO. 1 AND ITERATION NO. 3

NODNE	X COMPONENT (U)	Y COMPONENT (V)	Z COMPONENT (W)	ITERATION NO. 3	
				1	2
1	2.9356735F-02	1.4191780E-03	0.0	2.9356735F-02	1.4191780E-03
2	2.9356742E-02	4.0737547E-09	0.0	2.9356742E-02	4.0737547E-09
3	2.5519237E-02	1.4193156E-03	3.1108400E-05	2.5519237E-02	1.4193156E-03
4	2.5519237E-02	-8.450016F-08	3.1105301E-05	2.5519237E-02	-8.450016F-08
5	2.2683658E-02	1.4195214F-03	5.9920829E-05	2.2683658E-02	1.4195214F-03
6	2.2683650E-02	-2.2263924F-07	5.9915212F-05	2.2683650E-02	-2.2263924F-07
7	1.9850120F-02	1.4196830F-03	8.3183013E-05	1.9850120F-02	1.4196830F-03
8	1.9850120F-02	-3.42352.00E-07	8.3375431E-05	1.9850120F-02	-3.42352.00E-07
9	1.7018143E-02	1.4197913F-03	9.8699835E-05	1.7018143E-02	1.4197913F-03
10	1.7018139E-02	-4.2083934E-07	9.8690522E-05	1.7018139E-02	-4.2083934E-07
11	1.4186997E-02	1.4198280E-03	1.0401016E-04	1.4186997E-02	1.4198280E-03
12	1.4186997E-02	-4.4781825E-07	1.0400009E-04	1.4186997E-02	-4.4781825E-07
13	1.1355862E-02	1.4197964F-03	9.8668752E-05	1.1355862E-02	1.4197964F-03
14	1.1355866E-02	-4.1985641F-07	9.8659031E-05	1.1355866E-02	-4.1985641F-07
15	8.5239187E-03	1.4196937F-03	8.3330786E-05	8.5239187E-03	1.4196937F-03
16	8.5239336E-03	-3.4791357E-07	8.3322491E-05	8.5239336E-03	-3.4791357E-07
17	5.6904666F-03	1.4195368E-03	5.9863130E-05	5.6904666F-03	1.4195368E-03
18	5.6904779E-03	-2.2099863E-07	5.9857077E-05	5.6904779E-03	-2.2099863E-07
19	2.85550987E-03	1.4193379F-03	3.1069256E-05	2.85550987E-03	1.4193379F-03
20	2.85550987E-03	-8.3249513F-08	3.1069256E-05	2.85550987E-03	-8.3249513F-08
21	1.7662009F-05	1.4192052F-03	0.0	1.7662009F-05	1.4192052F-03
22	1.7662024E-05	4.0019721E-09	0.0	1.7662024E-05	4.0019721E-09
23	1.5587583E-02	7.6064793F-04	-1.29556962E-02	1.5587583E-02	7.6064793F-04
24	1.5587766E-02	6.7323772E-04	-1.2960125E-02	1.5587766E-02	6.7323772E-04
25	1.5332391E-02	7.7265571E-04	-1.2648195E-02	1.5332391E-02	7.7265571E-04
26	1.5332391E-02	6.6036545E-04	-1.2651250E-02	1.5332391E-02	6.6036545E-04
27	1.50757739E-02	7.7950767E-04	-1.2551006E-02	1.50757739E-02	7.7950767E-04
28	1.5057940E-02	6.5354165E-04	-1.2553982E-02	1.5057940E-02	6.5354165E-04
29	1.4790314E-02	7.3239455E-04	-1.2500148F-02	1.4790314E-02	7.3239455E-04
30	1.4790516E-02	6.4307037E-04	-1.2503054E-02	1.4790516E-02	6.4307037E-04
31	1.4522307E-02	7.8311702E-04	-1.2469962F-02	1.4522307E-02	7.8311702E-04
32	1.4522433E-02	6.4752677E-04	-1.2472779E-02	1.4522433E-02	6.4752677E-04
33	1.4252145E-02	7.8328699E-04	-1.2460098E-02	1.4252145E-02	7.8328699E-04
34	1.42522149E-02	6.46976279E-04	-1.2462813F-02	1.42522149E-02	6.46976279E-04
35	1.3691205E-02	7.8261830F-04	-1.2470119F-02	1.3691205E-02	7.8261830F-04
36	1.3691205E-02	6.4715979E-04	-1.2472757E-02	1.3691205E-02	6.4715979E-04
37	1.3713446E-02	7.04135147E-04	-1.2500532E-02	1.3713446E-02	7.04135147E-04
38	1.3713501E-02	6.4794134E-04	-1.2503073E-02	1.3713501E-02	6.4794134E-04
39	1.2445462E-02	7.7666598E-04	-1.2551751E-02	1.2445462E-02	7.7666598E-04
40	1.3445556E-02	6.5199039E-04	-1.2554195E-02	1.3445556E-02	6.5199039E-04
41	1.3163744E-02	7.6977460E-04	-1.2497746E-02	1.3163744E-02	7.6977460E-04
42	1.3164061E-02	6.5793012E-04	-1.2657103E-02	1.3164061E-02	6.5793012E-04
43	1.2713762E-02	7.5655939E-04	-1.2960803E-02	1.2713762E-02	7.5655939E-04
44	1.2714191E-02	6.6170486E-04	-1.2963068E-02	1.2714191E-02	6.6170486E-04
45	1.5435151E-02	7.4291974E-04	-1.3182176E-02	1.5435151E-02	7.4291974E-04
46	1.5435334E-02	6.9463020E-04	-1.3186198E-02	1.5435334E-02	6.9463020E-04
47	1.52469137E-02	7.4337330E-04	-1.2763843E-02	1.52469137E-02	7.4337330E-04
48	1.5246933E-02	6.9313263E-04	-1.277731F-02	1.5246933E-02	6.9313263E-04
49	1.5213333E-02	7.4724360E-04	-1.2557903E-02	1.5213333E-02	7.4724360E-04
50	1.501126533E-02	6.73245333E-04	-1.26011636E-02	1.501126533E-02	6.73245333E-04

STRESSES FOR ISOLATOR AND ARRESTOR FOR TILE NO. 1 AND ITERATION NO. 3

LOCAL COORDINATES			STRESSES		
X	Y	Z	XX	YY	ZZ
1	5.2578F-01	2.6289F-01	3.9434F-02	1.5710E-01	1.5712E-01
2	5.2578F-01	7.0441E-02	3.9434F-02	1.5762F-01	1.5765F-01
3	1.4088F-01	2.6289F-01	3.9434F-02	1.1014E-01	1.1019E-01
4	1.4088F-01	7.0441E-02	3.9434F-02	1.1068F-01	1.1073E-01
5	5.2578F-01	2.6289E-01	1.0566E-02	8.8938F-00	8.3943E-00
6	5.2578E-01	7.0441E-02	1.0566E-02	9.9462F-00	8.9467E-00
7	1.4088E-01	2.6289E-01	1.0566E-02	4.1617E-01	4.1628E-00
8	1.4088F-01	7.0441E-02	1.0566E-02	4.2153F-01	4.2165E-00
			ELEMENT NUMBER		
1	1.1925F-00	2.6269F-01	3.9434F-02	1.9934F-01	1.8937E-01
2	1.1925F-00	7.0441E-02	3.9434F-02	1.8986E-01	1.8988F-01
3	9.0755E-01	2.6289F-01	3.9434E-02	1.7791E-01	1.7795E-01
4	3.0755F-01	7.0441F-02	3.9434E-02	1.7843E-01	1.7846E-01
5	1.1925F-01	2.6289F-01	1.0566F-02	1.2201E-01	1.2201E-01
6	1.1925F-01	7.0441F-02	1.0566F-02	1.2252E-01	1.2253E-01
7	3.0755F-01	2.6289F-01	1.0566F-02	1.1039E-01	1.1040E-01
8	9.0755F-01	7.0441F-02	1.0566F-02	1.1091E-01	1.1092E-01
			ELEMENT NUMBER		
1	1.3551F-01	2.6278F-01	3.9434F-02	1.9995E-01	1.9995E-01
2	1.4591F-01	7.0441F-02	3.9434F-02	2.0445F-01	2.0045F-01
3	1.4742F-01	2.6289F-01	3.9434F-02	1.9544F-01	1.9545E-01
4	1.4742F-01	7.0441F-02	3.9434F-02	1.9594F-01	1.9595E-01
5	1.8591F-01	2.6289F-01	1.0566F-02	1.3727F-01	1.3727F-01
6	1.8591F-01	7.0441F-02	1.0566F-02	1.3320E-01	1.3320E-01
7	1.4742F-01	2.6289F-01	1.0566F-02	1.2897E-01	1.2897E-01
8	1.4742F-01	7.0441F-02	1.0566F-02	1.2957F-01	1.2957F-01
			ELEMENT NUMBER		
1	2.5253F-01	2.6241F-01	3.9434F-02	2.0523F-01	2.0523F-01
2	2.5253F-01	7.0441F-02	3.9434F-02	2.0551F-01	2.0551E-01
3	2.1409F-01	2.6289F-01	3.9434F-02	2.0252E-01	2.0252F-01
4	2.1409F-01	7.0441F-02	3.9434F-02	2.0301F-01	2.0301F-01
5	2.5253F-01	2.6241F-01	1.0566F-02	1.3566F-01	1.3566F-01
6	2.5253F-01	7.0441F-02	1.0566F-02	1.3838F-01	1.3838F-01
7	2.1409F-01	2.6241F-01	1.0566F-02	1.3536F-01	1.3536F-01
8	2.1409F-01	7.0441F-02	1.0566F-02	1.3585F-01	1.3585F-01
			ELEMENT NUMBER		
1	1.1924F-01	2.6241F-01	3.9434F-02	2.0698F-01	2.0697F-01
2	1.1924F-01	7.0441F-02	3.9434F-02	2.0744F-01	2.0744F-01
3	2.4755F-01	2.6289F-01	3.9434F-02	2.0620F-01	2.0620F-01
4	2.3275F-01	7.0441F-02	3.9434F-02	2.0668F-01	2.0667F-01
5	3.1924F-01	2.6241F-01	1.0566F-02	1.3990F-01	1.3990F-01
6	3.1924F-01	7.0441F-02	1.0566F-02	1.4037E-01	1.4037E-01
7	2.9775F-01	2.6289F-01	1.0566F-02	1.3912F-01	1.3912F-01
8	2.9775F-01	7.0441F-02	1.0566F-02	1.3755F-01	1.3755F-01

STRESSES AND DIRECT STRAINS FOR TILE NO. 3

MEM	TEMP	LOCAL COORDINATES			STRAINS			STRESSES			ITERATION NO. 3				
		X	Y	Z	XX	YY	ZZ	XY	YZ	ZX	XX	YY	ZZ		
1	-325.	0.33	0.02	0.17	-2.315E-03	-2.279E-03	-2.564E-01	9.971E-00	-5.376E-00	4.430E-07	4.162E-03	-6.923E-00			
2	-325.	1.00	0.02	0.17	-2.337F-03	-2.307F-03	-2.529E-01	1.502E-01	-1.196E-01	-5.625E-07	3.872E-03	-5.375E-00			
3	-325.	1.67	0.02	0.17	-2.326E-03	-2.324E-03	-2.520F-01	1.643F-01	1.349E-01	-7.735E-08	3.593E-03	-3.830E-00			
4	-325.	2.33	0.02	0.17	-2.326E-03	-2.332E-03	-2.516F-01	1.704E-01	1.990E-00	7.876E-07	3.383E-03	-2.281E-00			
5	-325.	3.00	0.02	0.17	-2.326E-02	-2.334E-03	-2.514E-01	1.733E-01	2.286E-00	-6.610E-07	3.229E-03	-7.337E-01			
6	-325.	3.67	0.02	0.17	-2.326E-03	-2.334E-03	-2.514E-01	1.733E-01	2.286E-00	-2.496E-06	3.092E-03	8.124E-01			
7	-325.	4.33	0.02	0.17	-2.325E-03	-2.332F-03	-2.516F-01	1.704E-01	1.704E-01	-2.890E-06	2.960E-03	2.359E-00			
8	-325.	5.00	0.02	0.17	-2.326E-02	-2.323F-03	-2.520E-01	1.642E-01	1.642E-01	-3.417E-06	2.805E-03	3.908E-00			
9	-325.	5.67	0.02	0.17	-2.318E-02	-2.307E-03	-2.529E-01	1.499E-01	1.500E-01	-1.440E-01	-3.052E-06	2.572E-03	5.453E-00		
10	-325.	6.33	0.02	0.17	-2.315E-03	-2.278E-03	-2.564E-01	9.904E-00	9.907E-00	-5.445E-00	-2.173E-06	2.244E-03	7.001E-00		
11	-325.	0.33	0.17	0.13	-3.264E-04	-2.236E-04	-1.025E-03	1.280E-00	2.893E-00	-3.919E-00	-1.090E-03	4.053E-03	-5.748E-00		
12	-325.	1.00	0.09	0.17	0.13	-2.874E-04	-2.523E-04	-4.899E-04	-2.601E-00	2.297E-00	-7.626E-01	-1.588E-03	3.457E-03	-6.509E-00	
13	-325.	1.67	0.17	0.13	-3.898E-04	-2.692E-04	-2.060E-04	-2.163E-00	-2.163E-00	-6.566E-04	3.457E-03	-3.888E-00			
14	-325.	2.33	0.17	0.13	-7.125E-04	-2.793E-04	-9.189E-05	-2.304E-00	-3.315E-00	1.728E-00	-2.841E-04	3.576E-03	-2.252E-00		
15	-325.	3.00	0.17	0.13	-2.948E-04	-2.820E-04	-4.342E-05	-2.485E-00	-3.354E-00	2.023E-00	-1.690E-03	3.695E-03	-7.064E-01		
16	-325.	3.67	0.17	0.13	-2.450E-04	-2.819E-04	-4.357E-05	-2.495E-00	-3.314E-00	2.022E-00	-3.535E-03	3.338E-03	8.361E-01		
17	-325.	4.33	0.17	0.13	-3.218E-04	-2.790E-04	-4.754E-05	-2.335E-00	-3.314E-00	1.724E-00	-3.544E-03	3.457E-03	-2.380E-00		
18	-325.	5.00	0.17	0.13	-2.996E-04	-2.586E-04	-2.078E-04	-2.213E-00	-3.310E-00	1.032E-00	-3.811E-03	2.742E-03	4.013E-00		
19	-325.	5.67	0.17	0.13	-2.985E-04	-2.514E-04	-4.97E-04	-2.675E-00	-2.287E-00	-7.920E-01	-2.240E-03	1.907E-03	6.642E-00		
20	-325.	6.33	0.17	0.13	-3.272E-04	-2.224E-04	-1.034E-03	1.213E-00	2.787E-00	-3.977E-00	-1.364E-03	2.980E-03	5.857E-00		
21	-325.	0.33	0.17	0.30	-2.380E-04	-1.005E-04	-1.711E-04	-9.988E-00	-4.434E-01	-2.295E-00	-2.021E-03	2.980E-03	-3.986E-00		
22	-325.	1.00	0.20	0.17	0.20	-3.212E-04	-2.611E-04	-2.611E-04	-1.575E-01	-6.259E-20	-1.110E-00	-1.639E-03	3.219E-03	-5.603E-00	
23	-325.	1.67	0.17	0.13	-3.682E-04	-7.308E-04	-1.225E-05	-1.795E-01	1.541E-01	5.053E-01	-7.451E-04	2.980E-03	-3.971E-00		
24	-325.	2.33	0.17	0.13	-3.681E-04	-7.187E-05	-1.521E-04	-1.853E-01	1.675E-01	1.337E-00	-1.495E-03	3.099E-03	-2.166E-00		
25	-325.	3.00	0.17	0.17	0.39	-3.725E-04	-2.224E-04	-1.034E-03	1.213E-00	1.642E-00	-1.937E-03	3.219E-03	-6.657E-01		
26	-325.	3.67	0.17	0.17	0.39	-3.727E-04	-1.005E-04	-1.711E-04	-1.891E-01	1.750E-00	1.641E-01	-2.231E-03	3.099E-03	8.199E-03	
27	-325.	4.33	0.17	0.17	0.30	-3.682E-04	-2.605E-04	-2.611E-04	-1.575E-01	-6.259E-20	-1.110E-00	-1.639E-03	3.219E-03	-5.603E-00	
28	-325.	5.00	0.17	0.17	0.30	-3.681E-04	-7.187E-05	-1.521E-04	-1.853E-01	1.675E-01	1.337E-00	-1.495E-03	3.099E-03	-2.166E-00	
29	-325.	5.67	0.17	0.17	0.30	-3.681E-04	-7.187E-05	-1.521E-04	-1.853E-01	1.675E-01	1.337E-00	-1.495E-03	3.099E-03	-2.166E-00	
30	-325.	6.33	0.17	0.17	0.39	-3.729E-04	-1.005E-04	-1.711E-04	-1.903E-01	1.483E-01	-2.330E-01	-8.428E-04	3.576E-03	4.108E-00	
31	-325.	0.33	0.17	0.17	0.39	-3.727E-04	-1.005E-04	-1.711E-04	-1.891E-01	1.750E-00	1.641E-01	-2.231E-03	3.099E-03	8.199E-03	
32	-325.	1.00	0.20	0.17	0.30	-3.682E-04	-2.605E-04	-2.611E-04	-1.575E-01	-6.259E-20	-1.110E-00	-1.639E-03	3.219E-03	-5.603E-00	
33	-325.	1.67	0.17	0.17	0.30	-3.681E-04	-7.187E-05	-1.521E-04	-1.853E-01	1.675E-01	1.337E-00	-1.495E-03	3.099E-03	-2.166E-00	
34	-325.	2.33	0.17	0.17	0.30	-3.681E-04	-7.187E-05	-1.521E-04	-1.853E-01	1.675E-01	1.337E-00	-1.495E-03	3.099E-03	-2.166E-00	
35	-325.	3.00	0.17	0.17	0.39	-3.729E-04	-1.005E-04	-1.711E-04	-1.903E-01	1.483E-01	-2.330E-01	-8.428E-04	3.576E-03	4.108E-00	
36	-325.	3.67	0.17	0.17	0.39	-3.727E-04	-1.005E-04	-1.711E-04	-1.891E-01	1.750E-00	1.641E-01	-2.231E-03	3.099E-03	8.199E-03	
37	-325.	4.33	0.17	0.17	0.30	-3.682E-04	-2.605E-04	-2.611E-04	-1.575E-01	-6.259E-20	-1.110E-00	-1.639E-03	3.219E-03	-5.603E-00	
38	-325.	5.00	0.17	0.17	0.30	-3.681E-04	-7.187E-05	-1.521E-04	-1.853E-01	1.675E-01	1.337E-00	-1.495E-03	3.099E-03	-2.166E-00	
39	-325.	5.67	0.17	0.17	0.30	-3.681E-04	-7.187E-05	-1.521E-04	-1.853E-01	1.675E-01	1.337E-00	-1.495E-03	3.099E-03	-2.166E-00	
40	-325.	6.33	0.17	0.17	0.39	-3.729E-04	-1.005E-04	-1.711E-04	-1.903E-01	1.483E-01	-2.330E-01	-8.428E-04	3.576E-03	4.108E-00	
41	-325.	0.33	0.17	0.17	0.39	-3.727E-04	-1.005E-04	-1.711E-04	-1.891E-01	1.750E-00	1.641E-01	-2.231E-03	3.099E-03	8.199E-03	
42	-325.	1.00	0.20	0.17	0.30	-3.682E-04	-2.605E-04	-2.611E-04	-1.575E-01	-6.259E-20	-1.110E-00	-1.639E-03	3.219E-03	-5.603E-00	
43	-325.	1.67	0.17	0.17	0.30	-3.681E-04	-7.187E-05	-1.521E-04	-1.853E-01	1.675E-01	1.337E-00	-1.495E-03	3.099E-03	-2.166E-00	
44	-325.	2.33	0.17	0.17	0.30	-3.681E-04	-7.187E-05	-1.521E-04	-1.853E-01	1.675E-01	1.337E-00	-1.495E-03	3.099E-03	-2.166E-00	
45	-325.	3.00	0.17	0.17	0.39	-3.729E-04	-1.005E-04	-1.711E-04	-1.903E-01	1.483E-01	-2.330E-01	-8.428E-04	3.576E-03	4.108E-00	
46	-325.	3.67	0.17	0.17	0.39	-3.727E-04	-1.005E-04	-1.711E-04	-1.891E-01	1.750E-00	1.641E-01	-2.231E-03	3.099E-03	8.199E-03	
47	-325.	4.33	0.17	0.17	0.30	-3.682E-04	-2.605E-04	-2.611E-04	-1.575E-01	-6.259E-20	-1.110E-00	-1.639E-03	3.219E-03	-5.603E-00	
48	-325.	5.00	0.17	0.17	0.30	-3.681E-04	-7.187E-05	-1.521E-04	-1.853E-01	1.675E-01	1.337E-00	-1.495E-03	3.099E-03	-2.166E-00	
49	-325.	5.67	0.17	0.17	0.30	-3.681E-04	-7.187E-05	-1.521E-04	-1.853E-01	1.675E-01	1.337E-00	-1.495E-03	3.099E-03	-2.166E-00	
50	-325.	6.33	0.17	0.17	0.39	-3.729E-04	-1.005E-04	-1.711E-04	-1.903E-01	1.483E-01	-2.330E-01	-8.428E-04	3.576E-03	4.108E-00	
51	-325.	0.33	0.17	0.17	0.39	-3.727E-04	-1.005E-04	-1.711E-04	-1.891E-01	1.750E-00	1.641E-01	-2.231E-03	3.099E-03	8.199E-03	
52	-325.	1.00	0.20	0.17	0.30	-3.682E-04	-2.605E-04	-2.611E-04	-1.575E-01	-6.259E-20	-1.110E-00	-1.639E-03	3.219E-03	-5.603E-00	
53	-325.	1.67	0.17	0.17	0.30	-3.681E-04	-7.187E-05	-1.521E-04	-1.853E-01	1.675E-01	1.337E-00	-1.495E-03	3.099E-03	-2.166E-00	
54	-325.	2.33	0.17	0.17	0.30	-3.681E-04	-7.187E-05	-1.521E-04	-1.853E-01	1.675E-01	1.337E-00	-1.495E-03	3.099E-03	-2.166E-00	
55	-325.	3.00	0.17	0.17	0.39	-3.729E-04	-1.005E-04	-1.711E-04	-1.903E-01	1.483E-01	-2.330E-01	-8.428E-04	3.576E-03	4.108E-00	
56	-325.	3.67	0.17	0.17	0.39	-3.727E-04	-1.005E-04	-1.711E-04	-1.891E-01	1.750E-00	1.641E-01	-2.231E-03	3.099E-03	8.199E-03	
57	-325.	4.33	0.17	0.17	0.30	-3.682E-04	-2.605E-04	-2.611E-04	-1.575E-01	-6.259E-20	-1.110E-00	-1.639E-03	3.219E-03	-5.603E-00	
58	-325.	5.00	0.17	0.17	0.30	-3.681E-04	-7.187E-05	-1.521E-04	-1.853E-01	1.675E-01	1.337E-00	-1.495E-03	3.099E-03	-2.166E-00	
59	-325.	5.67	0.17	0.17	0.30	-3.681E-04	-7.187E-05	-1.521E-04	-1.853E-01	1.675E-01	1.337E-00	-1.495E-03	3.099E-03	-2.166E-00	
60	-325.	6.33	0.17	0.17	0.39	-3.729E-04	-1.005E-04	-1.711E-04	-1.903E-01	1.483E-01	-2.330E-01	-8.428E-04	3.576E-03	4.108E-00	
61	-325.	0.33	0.17	0.17	0.39	-3.727E-04	-1.005E-04	-1.711E-04	-1.8						

BOTTOM-POINT PLATE MEMBER STRAINS AND STRESSES FOR ITERATION NO. 3

MEMBERID	COORDINATES X Y	EPS X	EPS Y	FPS XY	STRAINS EPS X	STRAINS EPS Y	FPS XY	SIG X	SIG Y	SIG XY
1	3.3333E-01	1.6666E-01	-	-	-4.2580E-03	-4.2574E-03	1.6214E-08	-5.2813E-00	-5.3125E-01	1.1581E-01
2	1.0000E-00	1.6666E-01	-	-	-4.2597E-03	-4.2569E-03	1.3421E-08	-2.2148E-01	-1.6406E-01	9.5864E-02
3	1.6667E-00	1.6666E-01	-	-	-4.2619E-03	-4.2562E-03	5.1156E-09	-4.3762E-01	-2.8516E-01	3.6540E-02
4	2.3333E-00	1.6666E-01	-	-	-4.2636E-03	-4.2557E-03	4.1084E-09	-6.0832E-01	-2.4609E-01	2.9345E-02
5	3.0000E-00	1.6666E-01	-	-	-4.2645E-03	-4.2554E-03	-9.7114E-10	-7.0004E-01	-2.4609E-01	-6.9367E-03
6	3.6667E-00	1.6666E-01	-	-	-4.2645E-03	-4.2554E-03	2.2789E-09	-6.9676E-01	-2.4609E-01	1.5920E-02
7	4.3333E-00	1.6666E-01	-	-	-4.2635E-03	-4.2557E-03	8.6476E-09	-6.0016E-01	-2.0313E-01	6.1733E-02
8	5.0000E-00	1.6666E-01	-	-	-4.2617E-03	-4.2563E-03	1.1266E-09	-4.2246E-01	-2.4609E-01	8.0474E-02
9	5.6667E-01	1.6666E-01	-	-	-4.2595E-03	-4.2569E-03	2.4391E-09	-2.0192E-01	-1.6406E-01	1.7415E-02
10	6.3333E-00	1.6666E-01	-	-	-4.2577E-03	-4.2575E-03	1.1265E-09	-2.6602E-01	-5.3125E-01	8.1179E-03

PROGRAM LISTING OF INPUT DATA CARDS

~~OPTIONS~~

STATICS PROBLEM

MAXIMUM NO. ITERATIONS = 3

CONVERGENCE PARAMETER = 5.0000E-02

PRIMARY STRUCTURE STRESSES PRESENTED AFTER LAST ITERATION AT PLATE TOP AND BOTTOM SURFACES

TILES OF PRIMARY STRUCTURE

TILE STRESSES PRESENTED AFTER LAST ITERATION

TILE NODE MAP REQUIRED

TILE ELEMENT MAP NOT REQUIRED

TILE NODE COORDINATES NOT REQUIRED

DO NOT PRINT ELEMENT STIFFNESS MATRICES

DO NOT PRINT ASSEMBLED STIFFNESS MATRICES

DO NOT PRINT FILE DEBUGGING INFORMATION

COMPUTE STRESSES FOR ALL TILES

- - - - -
G E O M E T R Y
- - - - -

PLATE	LX =	6.66667E 00	LY =	3.33330E-01	TP =	7.50000E-01
STRINGERS	Y1 =	5.00000E-01	ZS =	0.0	YS =	0.0
	IY* =	0.0	IZ* =	0.0	JX* =	0.0
TILES	NXB =	1	NYB =	1		
	T =	0.0	B1 =	0.0	T2 =	1.16667E 00
	TA =	1.66670E-01	TI =	5.00000E-02	TC =	0.0
BRICK	NB1 =	0	NB2 =	10	ND2 =	1
	NT1 =	0	NT2 =	7		

- - - - -
M A T E R I A L P R O P E R T I E S
- - - - -

PLATE	EP =	1.00000E 07	NU P =	3.00000E-01	GAMMA P =	0.0
STRINGERS	ES =	0.0	NU S =	3.0	GAMMA S =	0.0
ARMESTOR	EX =	6.00000E 04	FY =	6.00000E 04	EZ =	6.00000E 03
UM RSI	NU XY =	5.00000E-01	NU YZ =	1.00000E-01	NU ZX =	1.00000E-02
	GXY =	2.00000E 04	GYZ =	3.20000E 04	GZX =	3.20000E 04
	GAMMA A =	0.0				
	ALPHA X =	0.0	ALPHA Y =	3.0	ALPHA Z =	0.0
ISOLATOR	EI =	9.00000E 01	NU I =	4.90000E-01	GAMMA I =	0.0
RSI	GAMMA R =	0.0	ALPHA FY / ALPHA FX =	0.0	ALPHA RL / ALPHA RK =	1.0

TEMPERATURE DEPENDENT MATERIAL PROPERTIES

		PROPERTY	TEMPERATURE	PROPERTY	TEMPERATURE	PROPERTY	TEMPERATURE
1	ER	ALL	6.000E 04				
1	ER*	ALL	6.000E 03				
1	GR*	ALL	3.200E 04				
1	NU R	ALL	5.000E-C1				
1	NU R*	ALL	1.000E-C2				
1	ALPHA R	ALL	0.0				
1	EC	ALL	0.0				
1	NU C	ALL	0.0				
1	ALPHA C	ALL	0.0				

B O U N D A R Y C O N D I T I O N S

PLATE OUT OF PLANE

PLATE IN PLANE

STRINGERS

EDGE	PLATE OUT OF PLANE	PLATE IN PLANE	STRINGERS
A	PINNED	FREE	FREE
B	PINNED	U HELD, V FREE	FREE
C	FREE	V HELD, U FREE	FREE
D	FREE	FREE	FREE

S T A T I C L O A D I N G

NX = C.0 NY = 0.0 NYX = 0.0 NX = 0.0
PZ = 1.00000E 02 T = C.C M = 0.0 V = 0.0

DEL TEMP S = C.0

R S I - T E M P E R A T U R E S

NO STATIC THERMAL LOADING

UNIFORM TEMPERATURE OPTION

T REFERENCE = 0.0

DEL T U = T - T REF = 0.0

N O D E M A P
S U R F A C E 1

200	202	204	206	208	210	212	214	216	218	220
190	201	203	205	207	209	211	213	215	217	219

N O D E M A P
S U R F A C E 2

178	180	182	184	186	188	190	192	194	196	198
177	179	181	183	185	187	189	191	193	195	197

PRIMARY STRUCTURE

DUDE	X	Y	Z
1	0.0	0.0	0.0
2	3.666670E-01	3.3333300E-01	0.0
3	6.666670E-01	3.3333300E-01	0.0
4	1.333334E-01	3.3333300E-01	0.0
5	-1.333334E-01	3.3333300E-01	0.0
6	2.000001E-01	3.3333300E-01	0.0
7	2.000001E-01	3.3333300E-01	0.0
8	2.666668E-01	3.3333300E-01	0.0
9	2.666668E-01	3.3333300E-01	0.0
10	2.666668E-01	3.3333300E-01	0.0
11	3.333335E-01	3.3333300E-01	0.0
12	3.333335E-01	3.3333300E-01	0.0
13	4.000002E-01	3.3333300E-01	0.0
14	4.000002E-01	3.3333300E-01	0.0
15	4.666669E-01	3.3333300E-01	0.0
16	4.666669E-01	3.3333300E-01	0.0
17	5.333336E-01	3.3333300E-01	0.0
18	5.333336E-01	3.3333300E-01	0.0
19	6.0000C3E-01	3.3333300E-01	0.0
20	6.0000C3E-01	3.3333300E-01	0.0
21	6.666670E-01	3.3333300E-01	0.0
22	6.666670E-01	3.3333300E-01	0.0

DEGREES OF FREEDOM

	DX	DY	DZ	RX	RY	RZ
1	1	0	0	2	3	0
2	4	5	0	6	7	0
3	8	0	9	10	11	0
4	12	13	14	15	16	0
5	17	18	19	20	21	0
6	21	22	23	24	25	0
7	26	27	28	29	29	0
8	30	31	32	33	34	0
9	35	36	37	38	39	0
10	41	42	43	44	45	0
11	45	46	47	48	49	0
12	50	51	52	53	54	0
13	55	56	57	58	59	0
14	60	61	62	63	64	0
15	65	66	67	68	69	0
16	70	71	72	73	74	0
17	75	76	77	78	79	0
18	80	81	82	83	84	0
19	85	86	87	88	89	0
20	90	91	92	93	94	0

PRIMARY STRUCTURE DEFLECTIONS FOR ITERATION NO. 1

NDE	DX			DY			DZ			RX			RY			RZ		
	-1	-2	-3	-4	-5	-6	-7	-8	-9	-10	-11	-12	-13	-14	-15	-16	-17	-
1	0.0	0.0	0.0	0.0	0.0	0.0	0.0	0.0	0.0	0.0	0.0	0.0	0.0	0.0	0.0	0.0	0.0	-
2	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
3	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
4	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
5	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
6	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
7	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
8	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
9	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
10	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
11	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
12	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
13	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
14	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
15	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
16	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
17	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
18	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
19	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
20	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
21	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
22	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-

PRI-MARY STRUCTURE NODES ASSOCIATED WITH TILE NO. 1

1	3	5	7	9	11	13	15	17	19	21
2	4	6	8	10	12	14	16	18	20	22

PRIMARY STRUCTURE DEFLECTIONS FOR ITERATION NO. 2

NODE	CX	DY	DZ	RX			RY			RZ		
				-	-	-	-	-	-	-	-	-
1	-3.901913E-06	7.973046E-08	-	-	-	-	-1.457974E-06	-	-3.361257E-03	-	-	-
2	-3.897959E-06	-	-	-	-	-	-1.453203E-06	-	3.361257E-03	-	-	-
3	-3.676207E-06	-	-	-2.201482E-03	-	-	-2.826525E-05	-	-	-	-	-
4	-3.671434E-06	6.345107E-08	-	-2.201487E-03	-	-	-2.829292E-05	-	3.175853E-03	-	-	-
5	-3.351139E-06	-	-	-4.164275E-03	-	-	-4.925889E-05	-	-	-	-	-
6	-3.346494E-06	4.207058E-08	-	-4.164264E-03	-	-	-4.916536E-05	-	2.661876E-03	-	-	-
7	-2.938193E-06	-	-	-5.699161E-03	-	-	-6.429966E-05	-	-1.907061E-03	-	-	-
8	-2.935175E-06	2.713356E-08	-	-5.699124E-03	-	-	-6.406714E-05	-	1.906994E-03	-	-	-
9	-2.460804E-06	-	-	-6.673034E-03	-	-	-7.339950E-05	-	-	-	-	-
10	-2.459236E-06	1.782072E-08	-	-6.672934E-03	-	-	-7.276677E-05	-	-	-	-	-
11	-1.949114E-06	-	-	-7.006560E-03	-	-	-7.645463E-05	-	-	-	-	-
12	-1.949119E-06	1.465368E-08	-	-7.006425E-03	-	-	-7.562677E-05	-	2.626637E-07	-	-	-
13	-1.437492E-06	-	-	-6.673370E-03	-	-	-7.350260E-05	-	-2.205044E-07	-	-	-
14	-1.439067E-06	1.783919E-08	-	-6.673235E-03	-	-	-7.263929E-05	-	-9.927261E-04	-	-	-
15	-9.603064E-07	-	-	-5.699780E-03	-	-	-6.455116E-05	-	-9.906876E-03	-	-	-
16	-9.6332C4E-07	-	-	-5.699676E-03	-	-	-6.391277E-05	-	-1.906838E-03	-	-	-
17	-5.476637E-07	-	-	-4.164848E-03	-	-	-4.944927E-05	-	-	-	-	-
18	-5.523403F-C7	4.207477E-08	-	-4.164778E-03	-	-	-4.899905E-05	-	-2.662152E-03	-	-	-
19	-2.230393E-07	-	-	-2.201836E-03	-	-	-2.841215E-05	-	-1.76297E-03	-	-	-
20	-2.290736E-C7	6.385153E-08	-	-2.201798E-03	-	-	-2.817533E-05	-	3.176240E-03	-	-	-
21	-	-	-	-	-	-	-1.454991E-06	-	-	-	-	-
22	-	-	-	-7.931C72E-08	-	-	-1.452934E-06	-	-3.361825E-03	-	-	-

MAXIMUM DEFLECTION = -3.36177E-C3 FOR DOF 93

MAXIMUM DEFLECTION DIFFERENCE = -2.15776E-C4 FOR DOF 50

MAXIMUM CONVERGENCE PARAMETER = 6.41852E-C2

SOLUTION HAS NOT CONVERGED

PRIMARY STRUCTURE DEFLECTIONS FOR ITERATION NO. 3

NUDE	DX	DY	DZ	RX	RY	RZ
-1	-3.838843E-06	-	-	-	-	-
-2	-3.83492E-06	7.958687E-08	-	-	-	3.364080E-03
-3	-3.610429E-06	-	-	-2.203349E-03	1.455494E-06	3.364078E-03
-4	-3.294500E-06	6.373557E-08	-	-2.828105E-05	-	-
-5	-3.29982E-06	-	-	-2.829883E-05	3.178601E-03	3.178587E-03
-6	-	4.221585E-08	-	-2.920547E-05	-	-
-7	-2.898414E-06	-	-	-4.167896E-03	4.928574E-05	2.664342E-03
-8	-2.895486E-06	2.840901E-08	-	-4.167885E-03	-4.920547E-05	-
-9	-2.419676E-06	-	-	-5.704246E-03	-6.433637E-05	-1.908939E-03
-10	-2.418157E-06	1.937944E-08	-	-5.704213E-03	-6.412840E-05	1.908872E-03
-11	-1.917612E-06	-	-	-	-	-
-12	-1.917616E-06	1.631173E-08	-	-	-	-
-13	-1.0415613E-06	-	-	-	-	-
-14	-1.417140E-06	1.939791E-08	-	-6.679095E-03	7.343816E-05	3.942634E-04
-15	-9.470772E-07	-	-	-6.679010E-03	-7.286872E-05	9.941910E-04
-16	-9.500015E-07	2.845171E-08	-	-7.012971E-03	7.650600E-05	2.646018E-07
-17	-5.412911E-07	-	-	-7.012848E-03	-7.574550E-05	2.219912E-07
-18	-5.458402E-07	4.29206CE-08	-	-6.679438E-03	-	-
-19	-2.211529E-07	-	-	-5.704869E-03	6.458355E-05	-1.908754E-03
-20	-2.260555E-07	6.412620E-08	-	-5.704772E-03	-6.399328E-05	-1.908718E-03
-21	-	-	-	-	-	-
-22	-	7.917902E-08	-	-	-1.455612E-06	-3.364645E-03
		-	-	-	1.453408E-06	-3.364590E-03

MAXIMUM DEFLECTION = -3.364590E-03 FOR DOF 93

MAXIMUM DEFLECTION DIFFERENCE = 4.4224CE-06 FOR DOF 50

MAXIMUM CONVERGENCE PARAMETER = 1.00MP2E-C3

SOLUTION HAS CONVERGED

TPS DISPLACEMENTS FOR TILE NO. 1 AND ITERATION NO. 3

NOCE	X COMPONENT(U)	Y COMPONENT(V)	Z COMPONENT(W)
1	5.1257691E-03	5.4560314E-07	0.0
2	1.2576941E-03	4.6622353E-07	-2.2033486E-03
3	1.1883602E-03	-1.0603594E-05	-1.0675798E-05
4	1.1863595E-03	1.0675798E-05	-4.1678958E-03
5	5.9584647E-04	-1.848210E-05	-1.678846E-03
6	1.8494960E-04	1.8494960E-05	-5.7042465E-03
7	7.1296352E-04	-2.4126130E-05	-5.7042465E-03
8	7.1294140E-04	2.4076551E-05	-5.7042465E-03
9	3.7042890E-04	-2.7539313E-05	-6.6790953E-03
10	3.7003296E-04	2.7345130E-05	-6.6790953E-03
11	-1.18343690E-06	-2.8689741E-05	-7.0129707E-03
12	-1.18343690E-06	2.8420851E-05	-7.0129707E-03
13	-3.7406676E-04	-2.7586095E-05	-6.6794381E-03
14	-3.7406676E-04	2.7289876E-05	-6.6793077E-03
15	-7.672956E-04	-2.4218825E-05	-5.7048686E-03
16	-7.1671531E-04	2.4025925E-05	-5.7047717E-03
17	-6.9977315E-04	-1.8548948E-05	-4.1684695E-03
18	-5.9975588E-04	1.8437131E-05	-4.1684061E-03
19	-1.0623651E-03	-1.0657704E-05	-2.2036997E-03
20	-1.11923490E-03	1.0634339E-05	-2.2036657E-03
21	-1.02617419E-03	5.4585439E-07	0.0
22	-1.2617209E-03	-4.584898E-07	0.0
23	-1.19288568E-03	-3.6905985E-04	-3.7478809E-03
24	-1.19288899E-03	3.6939513E-04	-3.7476560E-03
25	-1.17317153E-03	-3.40217583E-04	-3.6907944E-03
26	-1.173176229E-03	3.4074089E-04	-3.6905597E-03
27	-1.14811209E-03	-3.08027276E-04	-7.4508972E-03
28	-1.148117549E-03	3.0878210E-04	-7.0866355E-03
29	-1.08823159E-03	-2.7959980E-04	-8.8766739E-03
30	-1.08834010E-03	2.8061983E-04	-8.8774636E-03
31	-5.7616225E-04	-2.6274496E-04	-9.7944960E-03
32	-5.7614408E-04	2.6372981E-04	-9.7941980E-03
33	-3.2821521E-07	-2.5797938E-04	-1.011035E-02
34	-1.1892517E-07	2.5797938E-04	-1.0110315E-02
35	5.7595822E-04	-2.6276428E-04	-9.7948052E-03
36	5.7593700E-04	2.6369118E-04	-8.8774811E-03
37	1.0882223E-03	-2.7955766E-04	-8.8774636E-03
38	1.0881955E-03	2.8062542E-04	-8.8771619E-03
39	1.4811466E-03	-3.0784123E-04	-7.4514151E-03
40	1.4810288E-03	3.0894089E-04	-7.4511021E-03
41	1.7317953E-03	-3.3979258E-04	-5.6911409E-03
42	1.7316577E-03	3.4115440E-04	-5.6909099E-03
43	1.9283283E-03	-3.6838348E-04	-3.7480515E-03
44	1.9284511E-03	3.7009967E-04	-3.748020E-03
45	-1.4385517F-03	-1.6514414E-04	-6.8438686E-03
46	-1.4386310E-03	1.6520663E-04	-6.8435557E-03
47	-1.3081105E-03	-1.4886131E-04	-8.5816644E-03
48	-1.3081587E-03	1.4912240E-04	-8.581347AE-03
49	-1.1065399E-03	-1.2930391F-04	-1.0192931E-02
50	-1.1065831E-03	1.2975477E-04	-1.0192588E-02

STRESSES FOR ISOLATOR AND ARRESTOR FOR TILE NO. 1 AND ITERATION NO. 3

LOCAL COORDINATES		Z	X	Y	XX	YY	ZZ	STRESSES	XY	YZ	ZX	
5.2578E-01	2.6289E-01	3.9434E-02	-1.0207E	02	-1.0199E	02	-1.0637E	02	-5.3123E-04	1.1811E-01	-1.8883E-00	
5.2578E-01	7.0441E-02	3.9434E-02	-1.0208E	02	-1.0199E	02	-1.0637E	02	-5.3118E-04	-1.1792E-01	-1.8883E-00	
1.4088E-01	2.6289E-01	3.9434E-02	-1.0642E	02	-1.089E	02	-1.2610E-04	-1.3092E-04	-1.9813E-00	-1.9813E-00	-1.9813E-00	
1.4088E-01	7.0441E-02	3.9434E-02	-1.0642E	02	-1.0633E	02	-1.1090E	02	-5.3209E-04	-1.0933E-05	-1.8852E-00	
5.2578E-01	2.6289E-01	1.0566E-02	-1.0416E	02	-1.0413E	02	-1.0844E	02	-7.1102E-05	-1.1810E-01	-1.8851E-00	
5.2578E-01	7.0441E-02	1.0566E-02	-1.0417E	02	-1.0414E	02	-1.0845E	02	-7.1112E-05	-1.1809E-01	-1.8851E-00	
1.4088E-01	2.6289E-01	1.0566E-02	-1.0863E	02	-1.0860E	02	-1.1309E	02	-1.2609E-05	-1.1242E-01	-1.9881E-00	
1.4088E-01	7.0441E-02	1.0566E-02	-1.0863E	02	-1.0860E	02	-1.1309E	02	-1.2590E-05	-1.0886E-01	-1.9881E-00	
ELEMENT NUMBER		1	ELEMENT NUMBER		ELEMENT NUMBER		ELEMENT NUMBER		ELEMENT NUMBER		ELEMENT NUMBER	
1.1925E-00	2.6289E-01	3.9434E-02	-9.5839E	01	-9.5761E	01	-9.9871E	01	-6.1842E-04	-1.0427E-01	-1.6344E-00	
1.1925E-00	7.0441E-02	3.9434E-02	-9.5766E	01	-9.875E	01	-9.1226E	01	-1.1226E-04	-1.6344E-00	-1.6344E-00	
8.3755E-01	2.6289E-01	3.9434E-02	-9.9211E	01	-9.9123E	01	-1.0339E	02	-6.1797E-04	-1.7889E-01	-1.7889E-00	
8.0755E-01	7.0441E-02	3.9434E-02	-9.9215E	01	-9.9133E	01	-1.0339E	02	-6.1828E-04	-1.1916E-01	-1.7889E-00	
1.1925E-00	2.6289E-01	1.0566E-02	-9.7959E	01	-9.7921E	01	-1.0197E	02	-1.3852E-05	-1.0486E-01	-1.6398E-00	
1.1925E-00	7.0441E-02	1.0566E-02	-9.7963E	01	-9.7925E	01	-1.0197E	02	-1.3852E-05	-1.0486E-01	-1.6398E-00	
8.3755E-01	2.6289E-01	1.0566E-02	-1.0145E	02	-1.0141E	02	-1.0560E	02	-1.6456E-05	-1.1224E-01	-1.7943E-00	
8.0755E-01	7.0441E-02	1.0566E-02	-1.0145E	02	-1.0142E	02	-1.0561E	02	-1.6426E-05	-1.1192E-01	-1.7943E-00	
ELEMENT NUMBER		2	ELEMENT NUMBER		ELEMENT NUMBER		ELEMENT NUMBER		ELEMENT NUMBER		ELEMENT NUMBER	
1.1925E-00	2.6289E-01	3.9434E-02	-9.1963E	01	-9.1902E	01	-9.5846E	01	-5.5080E-04	-2.0995E-02	-1.2399E-00	
1.1925E-00	7.0441E-02	3.9434E-02	-9.3739E	01	-9.1906E	01	-9.5850E	01	-5.5449E-04	-9.1458E-02	-1.2399E-00	
1.4742E-00	2.6289E-01	3.9434E-02	-9.3739E	01	-9.3674E	01	-9.7699E	01	-5.5190E-04	-8.8869E-02	-1.4756E-00	
1.4742E-00	7.0441E-02	3.9434E-02	-9.3743E	01	-9.3678E	01	-9.7704E	01	-5.5373E-04	-8.8341E-02	-1.4756E-00	
1.3591E-00	2.6289E-01	1.0566E-02	-9.4213E	01	-9.4172E	01	-9.8061E	01	-4.2444E-05	-9.2670E-02	-1.2428E-00	
1.3591E-00	7.0441E-02	1.0566E-02	-9.4217E	01	-9.4176E	01	-9.8065E	01	-4.2877E-05	-9.1499E-02	-1.2428E-00	
1.4742E-00	2.6289E-01	1.0566E-02	-9.6090E	01	-9.6094E	01	-1.0000E	02	-1.4478E-05	-9.8656E-02	-1.4778E-00	
1.4742E-00	7.0441E-02	1.0566E-02	-9.6094E	01	-9.6052E	01	-1.0000E	02	-1.4902E-05	-9.8330E-02	-1.4778E-00	
ELEMENT NUMBER		3	ELEMENT NUMBER		ELEMENT NUMBER		ELEMENT NUMBER		ELEMENT NUMBER		ELEMENT NUMBER	
1.0591E-00	2.6289E-01	3.9434E-02	-9.1963E	01	-9.1902E	01	-9.5846E	01	-5.5080E-04	-2.0995E-02	-1.2399E-00	
1.0591E-00	7.0441E-02	3.9434E-02	-9.3739E	01	-9.3674E	01	-9.7699E	01	-5.5190E-04	-9.1458E-02	-1.2399E-00	
1.4742E-00	2.6289E-01	3.9434E-02	-9.3739E	01	-9.3678E	01	-9.7704E	01	-5.5373E-04	-8.8869E-02	-1.4756E-00	
1.4742E-00	7.0441E-02	3.9434E-02	-9.3743E	01	-9.3682E	01	-9.7704E	01	-5.5373E-04	-8.8341E-02	-1.4756E-00	
1.3591E-00	2.6289E-01	1.0566E-02	-9.4213E	01	-9.4172E	01	-9.8061E	01	-4.2444E-05	-9.2670E-02	-1.2428E-00	
1.3591E-00	7.0441E-02	1.0566E-02	-9.4217E	01	-9.4176E	01	-9.8065E	01	-4.2877E-05	-9.1499E-02	-1.2428E-00	
1.4742E-00	2.6289E-01	1.0566E-02	-9.6090E	01	-9.6094E	01	-1.0000E	02	-1.4478E-05	-9.8656E-02	-1.4778E-00	
1.4742E-00	7.0441E-02	1.0566E-02	-9.6094E	01	-9.6052E	01	-1.0000E	02	-1.4902E-05	-9.8330E-02	-1.4778E-00	
ELEMENT NUMBER		4	ELEMENT NUMBER		ELEMENT NUMBER		ELEMENT NUMBER		ELEMENT NUMBER		ELEMENT NUMBER	
2.5258E-00	2.6289E-01	3.9434E-02	-3.9434E-02	-8.9875E	01	-8.9827E	01	-9.3683E	01	-3.3072E-04	-6.4091E-02	-7.2296E-01
2.5258E-00	7.0441E-02	3.9434E-02	-3.9434E-02	-9.0789E	01	-8.9831E	01	-9.3687E	01	-3.2842E-04	-6.3348E-02	-7.2297E-01
2.1409E-00	2.6289E-01	3.9434E-02	-3.9434E-02	-9.0793E	01	-9.0738E	01	-9.4637E	01	-3.2452E-04	-6.8140E-02	-7.2100E-01
2.1409E-00	7.0441E-02	3.9434E-02	-3.9434E-02	-9.2246E	01	-9.2204E	01	-9.4642E	01	-3.2670E-04	-6.8743E-02	-7.2100E-01
2.5258E-00	2.6289E-01	1.0566E-02	-1.0566E-02	-9.2250E	01	-9.2207E	01	-9.6009E	01	-2.8702E-05	-6.4050E-02	-7.2446E-01
2.5258E-00	7.0441E-02	1.0566E-02	-1.0566E-02	-9.3224E	01	-9.3177E	01	-9.9976E	01	-1.9976E-05	-6.8335E-02	-7.2447E-01
2.1409E-00	2.6289E-01	1.0566E-02	-1.0566E-02	-9.3224E	01	-9.3181E	01	-9.7023E	01	-2.9040E-05	-6.8192E-02	-7.0225E-00
ELEMENT NUMBER		5	ELEMENT NUMBER		ELEMENT NUMBER		ELEMENT NUMBER		ELEMENT NUMBER		ELEMENT NUMBER	
3.0124E-00	2.6289E-01	3.9434E-02	-3.9434E-02	-9.055E	01	-8.9055E	01	-9.913E	01	-9.2836E	01	-1.0732E-02
3.0124E-00	7.0441E-02	3.9434E-02	-3.9434E-02	-9.056E	01	-8.9239E	01	-9.9313E	01	-9.2120E-02	-7.9974E-02	-1.0515E-01
2.4775E-00	2.6289E-01	3.9434E-02	-3.9434E-02	-9.056E	01	-8.9239E	01	-9.9313E	01	-9.2120E-02	-7.9974E-02	-1.0515E-01
2.4775E-00	7.0441E-02	3.9434E-02	-3.9434E-02	-9.150E	01	-9.150E	01	-9.5239E	01	-1.0252E-02	-8.1349E-02	-1.0521E-01
2.1924E-00	2.6289E-01	1.0566E-02	-1.0566E-02	-9.150E	01	-9.1458E	01	-9.5236E	01	-1.0252E-02	-8.1349E-02	-1.0521E-01
2.1924E-00	7.0441E-02	1.0566E-02	-1.0566E-02	-9.150E	01	-9.1461E	01	-9.5236E	01	-1.0252E-02	-8.1349E-02	-1.0521E-01
2.1924E-00	2.6289E-01	1.0566E-02	-1.0566E-02	-9.150E	01	-9.1461E	01	-9.5236E	01	-1.0252E-02	-8.1349E-02	-1.0521E-01
2.1924E-00	7.0441E-02	1.0566E-02	-1.0566E-02	-9.150E	01	-9.1461E	01	-9.5236E	01	-1.0252E-02	-8.1349E-02	-1.0521E-01

MEM	TE4P	LOCAL COORDINATES	Z	STRAINS			STRESSES			STRESSES			STRESSES			STRESSES			
				X	XX	YY	Z	XX	YY	ZZ	XY	YZ	ZX	XY	YZ	ZX	XY	YZ	
1	2	3	4	0.02	9.585E-05	1.080E-03	-7.235E-02	-1.053E	0.02	-1.053E	0.02	-1.097E	0.02	2.936E	-0.07	1.191E	-0.04	-1.938E	0.00
0.02	0.02	0.02	0.02	0.02	4.355E-05	1.010E-03	-6.705E-02	-0.862E	0.01	-0.862E	0.01	-1.077E	0.01	-1.714E	-0.04	-1.359E	0.00		
0.17	0.17	0.17	0.17	0.17	8.233E-05	9.467E-04	-6.455E-02	-0.940E	0.01	-0.940E	0.01	-1.068E	0.01	-1.578E	-0.07	-2.086E	-0.07		
0.33	0.33	0.33	0.33	0.33	1.152E-05	1.528E-04	-6.213E-02	-0.904E	0.01	-0.904E	0.01	-1.056E	0.01	-1.723E	-0.04	-2.106E	-0.04		
0.67	0.67	0.67	0.67	0.67	1.567E-05	8.652E-04	-6.021E-02	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
1.00	1.00	1.00	1.00	1.00	2.000E-05	1.529E-04	-5.865E-02	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
1.33	1.33	1.33	1.33	1.33	2.447E-05	8.923E-04	-5.702E-02	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
1.67	1.67	1.67	1.67	1.67	2.894E-05	9.467E-04	-5.545E-02	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
2.00	2.00	2.00	2.00	2.00	3.331E-05	8.240E-05	-5.388E-02	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
2.33	2.33	2.33	2.33	2.33	3.768E-05	4.353E-05	-5.231E-02	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
2.67	2.67	2.67	2.67	2.67	4.205E-05	1.017E-05	-5.074E-02	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
3.00	3.00	3.00	3.00	3.00	4.642E-05	5.067E-05	-4.917E-02	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
3.33	3.33	3.33	3.33	3.33	5.079E-05	9.582E-05	-4.760E-02	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
3.67	3.67	3.67	3.67	3.67	5.516E-05	1.017E-04	-4.603E-02	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
4.00	4.00	4.00	4.00	4.00	5.953E-05	5.067E-05	-4.446E-02	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
4.33	4.33	4.33	4.33	4.33	6.391E-05	1.017E-04	-4.289E-02	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
4.67	4.67	4.67	4.67	4.67	7.022E-05	1.529E-04	-4.132E-02	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
5.00	5.00	5.00	5.00	5.00	7.457E-05	8.652E-04	-3.975E-02	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
5.33	5.33	5.33	5.33	5.33	7.894E-05	1.017E-04	-3.818E-02	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
5.67	5.67	5.67	5.67	5.67	8.331E-05	5.067E-05	-3.661E-02	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
6.00	6.00	6.00	6.00	6.00	8.768E-05	1.017E-04	-3.504E-02	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
6.33	6.33	6.33	6.33	6.33	9.205E-05	5.067E-05	-3.347E-02	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
6.67	6.67	6.67	6.67	6.67	9.642E-05	1.017E-04	-3.190E-02	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
7.00	7.00	7.00	7.00	7.00	1.007E-04	5.067E-05	-3.033E-02	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
7.33	7.33	7.33	7.33	7.33	1.054E-04	1.017E-04	-2.876E-02	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
7.67	7.67	7.67	7.67	7.67	1.101E-04	5.067E-05	-2.719E-02	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
8.00	8.00	8.00	8.00	8.00	1.148E-04	1.017E-04	-2.562E-02	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
8.33	8.33	8.33	8.33	8.33	1.195E-04	5.067E-05	-2.405E-02	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
8.67	8.67	8.67	8.67	8.67	1.242E-04	1.017E-04	-2.248E-02	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
9.00	9.00	9.00	9.00	9.00	1.289E-04	5.067E-05	-2.091E-02	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
9.33	9.33	9.33	9.33	9.33	1.336E-04	1.017E-04	-1.934E-02	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
9.67	9.67	9.67	9.67	9.67	1.383E-04	5.067E-05	-1.776E-02	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
1.00	1.00	1.00	1.00	1.00	1.430E-04	1.017E-04	-1.619E-02	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
1.13	1.13	1.13	1.13	1.13	1.477E-04	5.067E-05	-1.462E-02	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
1.27	1.27	1.27	1.27	1.27	1.524E-04	1.017E-04	-1.305E-02	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
1.41	1.41	1.41	1.41	1.41	1.571E-04	5.067E-05	-1.148E-02	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
1.55	1.55	1.55	1.55	1.55	1.618E-04	1.017E-04	-9.919E-03	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
1.69	1.69	1.69	1.69	1.69	1.665E-04	5.067E-05	-8.391E-03	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
1.83	1.83	1.83	1.83	1.83	1.712E-04	1.017E-04	-6.863E-03	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
1.97	1.97	1.97	1.97	1.97	1.759E-04	5.067E-05	-5.335E-03	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
2.11	2.11	2.11	2.11	2.11	1.806E-04	1.017E-04	-3.808E-03	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
2.25	2.25	2.25	2.25	2.25	1.853E-04	5.067E-05	-2.301E-03	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
2.39	2.39	2.39	2.39	2.39	1.900E-04	1.017E-04	-8.943E-04	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
2.53	2.53	2.53	2.53	2.53	1.947E-04	5.067E-05	-4.466E-04	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
2.67	2.67	2.67	2.67	2.67	1.994E-04	1.017E-04	-1.987E-04	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
2.81	2.81	2.81	2.81	2.81	2.041E-04	5.067E-05	-6.419E-05	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
2.95	2.95	2.95	2.95	2.95	2.088E-04	1.017E-04	-1.172E-05	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
3.09	3.09	3.09	3.09	3.09	2.135E-04	5.067E-05	-1.705E-05	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
3.23	3.23	3.23	3.23	3.23	2.182E-04	1.017E-04	-2.237E-05	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
3.37	3.37	3.37	3.37	3.37	2.229E-04	5.067E-05	-2.769E-05	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
3.51	3.51	3.51	3.51	3.51	2.276E-04	1.017E-04	-3.301E-05	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
3.65	3.65	3.65	3.65	3.65	2.323E-04	5.067E-05	-3.833E-05	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
3.79	3.79	3.79	3.79	3.79	2.370E-04	1.017E-04	-4.365E-05	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
3.93	3.93	3.93	3.93	3.93	2.417E-04	5.067E-05	-4.897E-05	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
4.07	4.07	4.07	4.07	4.07	2.464E-04	1.017E-04	-5.429E-05	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
4.21	4.21	4.21	4.21	4.21	2.511E-04	5.067E-05	-5.961E-05	-0.904E	0.01	-0.904E	0.01	-1.066E	0.01	-1.707E	-0.04	-2.106E	-0.04		
4.35	4.35	4.35	4.35	4.35	2.558E-04	1.017E-04	-												

TOP-POINT PLATE MEMBER STRAINS AND STRESSES FOR ITERATION NO. 3

MEMBER	COORDINATES		STRAINS		STRESSES		SIG XY
	X	Y	EPS X	EPS Y	EPS XY	SIG X	
1	-3.3333E-01	-1.66666E-01	-1.0400E-04	-3.0404E-05	-1.7294E-09	-1.0426E-03	-8.7436E-00
2	1.0000E+00	1.66666E-01	-2.8878E-04	8.7388E-05	-2.8538E-08	-2.8853E-03	-8.3048E-00
3	1.66667E+00	1.66666E-01	-4.24333E-04	1.2777E-04	-4.6300E-08	-4.2418E-03	-5.1730E-00
4	2.3333E+00	1.66666E-01	-5.1380E-04	1.5463E-04	-9.0042E-08	-5.1364E-03	-5.3959E-00
5	3.0000E+00	1.66666E-01	-5.5616E-04	1.6799E-04	-5.9231E-08	-5.5820E-03	-5.3321E-00
6	3.66667E+00	1.66666E-01	-5.8836E-04	1.6798E-04	-2.0045E-08	-5.5821E-03	-5.1957E-00
7	4.3333E+00	1.66666E-01	-5.1396E-04	1.5468E-04	-4.8764E-08	-5.1382E-03	-5.3562E-00
8	5.0000E+00	1.66666E-01	-4.2456E-04	1.2785E-04	-4.8311E-08	-4.2440E-03	-5.2660E-00
9	5.66667E+00	1.66666E-01	-2.8889E-04	8.7419E-05	-5.5284E-08	-2.8864E-03	-8.2666E-00
10	6.3333E+00	1.66666E-01	-1.0406E-04	3.0421E-05	-6.6454E-08	-1.0432E-03	-8.7649E-00

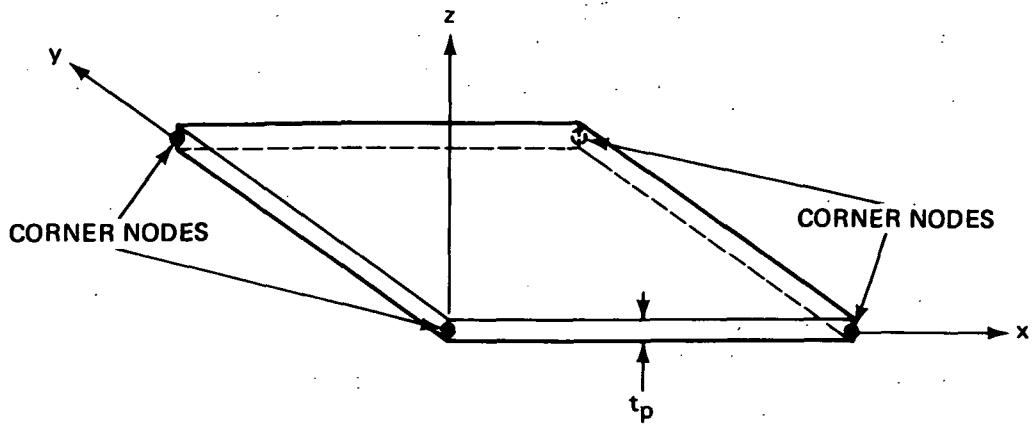
BOTTOM-POINT PLATE MEMBER STRAINS AND STRESSES FOR ITERATION NO. 3

MEMBER	COORDINATES		STRAINS		STRESSES		SIG XY	
	X	Y	EPS X	EPS Y	EPS XY	SIG X		
1	3.3333E-01	-1.6666E-01	-1.0467E-04	-2.9974E-05	-2.5926E-09	-1.0514E-03	-1.5686E-01	1.8519E-C2
2	1.0000E+00	1.6666E-01	-2.8974E-04	-8.7068E-05	-2.6673E-08	-2.8969E-03	-1.6196E+00	1.9052E-01
3	1.6667E+00	-1.6666E-01	4.2555E-04	-1.2756E-04	4.6589E-08	-4.2559E-03	-1.1875E+00	3.3278E-01
4	2.3333E+00	1.6666E-01	-5.1521E-04	-1.5449E-04	8.9941E-08	-5.1523E-03	-8.0867E-01	6.4243E-01
5	3.0000E+00	1.6666E-01	5.5987E-04	-1.6789E-04	5.9203E-08	-5.5989E-03	-8.0227E-01	4.2288E-01
6	3.6667E+00	1.6666E-01	5.5987E-04	-1.6787E-04	2.0064E-08	-5.5989E-03	-9.3916E-01	1.4331E-01
7	4.3333E+00	1.6666E-01	5.1539E-04	-1.5454E-04	-4.8651E-08	-5.1541E-03	-8.4817E-01	-3.4751E-01
8	5.0000E+00	1.6666E-01	4.2578E-04	-1.2763E-04	-4.867CE-08	-4.2581E-03	-1.0976E+00	-3.4764E-01
9	5.6667E+00	1.6666E-01	2.8985E-04	-8.7097E-05	-5.3538E-08	-2.8980E-03	-1.5727E+00	-3.8296E-C1
10	6.3333E+00	1.6666E-01	1.0473E-04	-2.9991E-05	-6.2518E-08	-1.0520E-03	-1.5701E+01	-4.4656E-C1

APPENDIX C

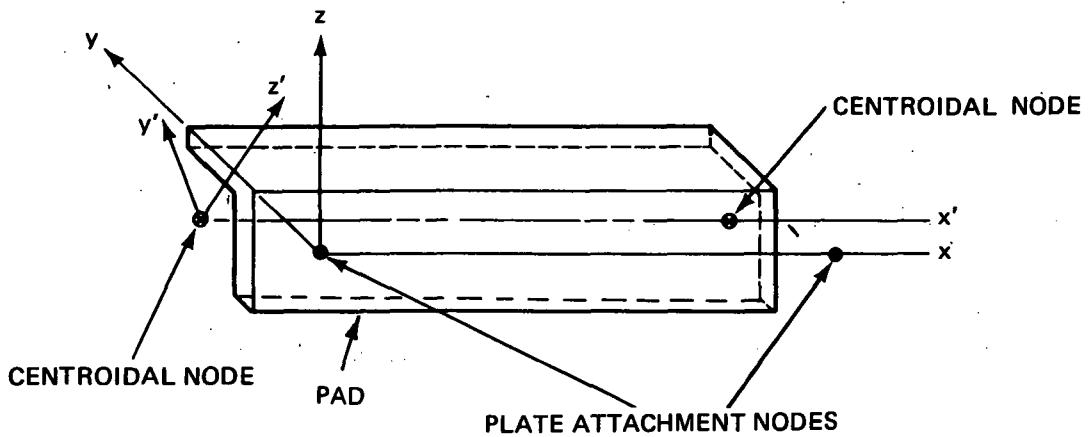
FINITE ELEMENT TYPES USED IN RE*S*I*ST

The active degrees-of-freedom (dof) for an element are defined as the dof for which there is at least one nonzero term in the corresponding row of the element stiffness matrix. The nodes and active dof per node, for each type element included in the RESIST computer program, are defined in (Figures C-1 through C-4).



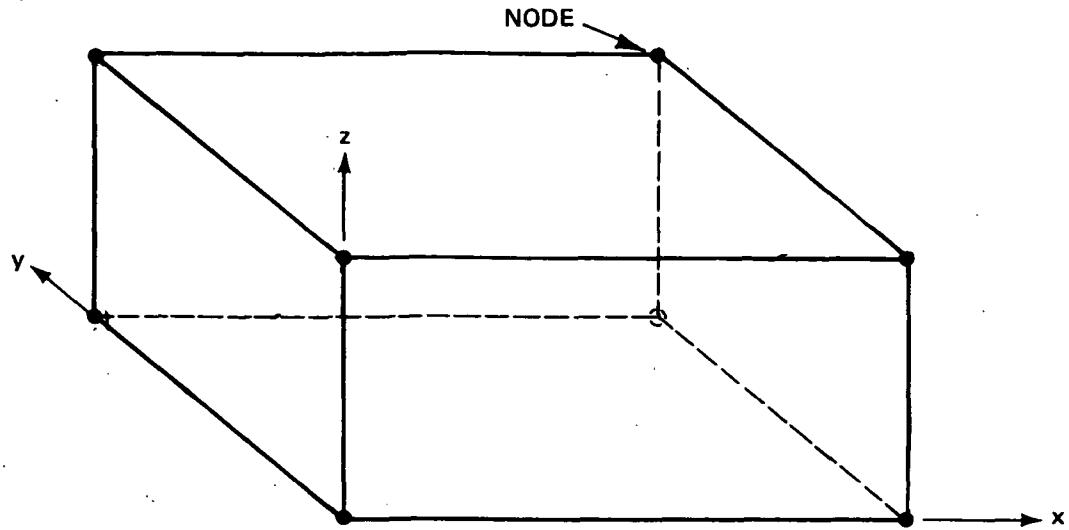
- BENDING AND MEMBRANE STIFFNESS
- FOUR NODES AT PLATE MIDDLE SURFACE
- ACTIVE DOF PER NODE ARE x , y AND z DEFLECTIONS AND x AND y ROTATIONS
- LUMPED MASSES AT NODES

Figure C-1 Primary Structure Isotropic Plate Element



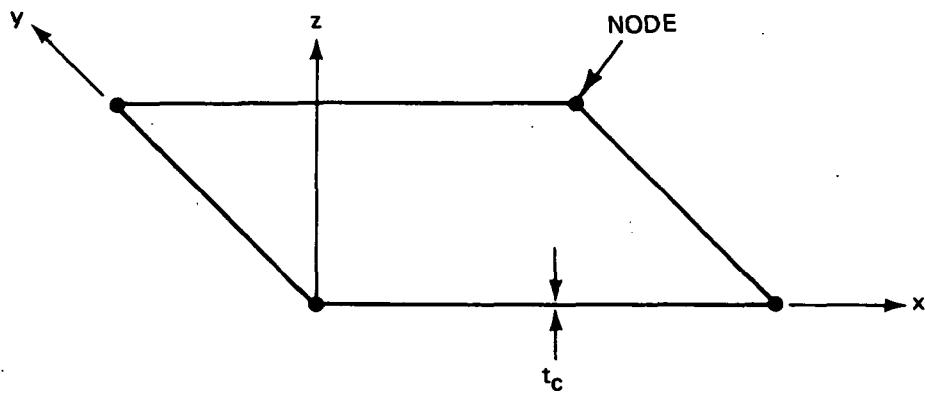
- ARBITRARY CROSS SECTION BENDING, TWISTING AND AXIAL STIFFNESS
- TWO CENTROIDAL NODES AND TWO ATTACHMENT NODES
- ACTIVE DOF ARE x , y AND z DEFLECTIONS AND ROTATIONS AT ATTACHMENT NODES ONLY. NO ACTIVE DOF AT CENTROIDAL NODES
- LUMPED MASSES AT CENTROIDAL NODES

Figure C-2 Primary Structure Uniform Stiffener Element



- THREE DIMENSIONAL STIFFNESS ELEMENT
- EIGHT CORNER NODES
- ACTIVE DOF PER NODE ARE x , y AND z DEFLECTIONS
- LUMPED MASSES AT CORNER NODES

Figure C-3 Three Dimensional Orthotropic TPS Element



- TWO DIMENSIONAL MEMBRANE ELEMENT
- FOUR CORNER NODES
- ACTIVE DOF PER NODE ARE x AND y DEFLECTIONS ONLY
- INERTIA EFFECTS NEGLECTED FOR THESE ELEMENTS

Figure C-4 TPS Thin Membrane Coating Element

GRUMMAN AEROSPACE CORPORATION

BETHPAGE, NEW YORK 11714