

NASA Contractor Report 187110

# User's Manual for Rocket Combustor Interactive Design (ROCCID) and Analysis Computer Program

## Volume II—Appendixes A-K

J.A. Muss and T.V. Nguyen  
*Gencorp, Aerojet Propulsion Division*  
*Sacramento, California*

and

C.W. Johnson  
*Software and Engineering Associates*  
*Carson City, Nevada*

May 1991

Prepared for  
Lewis Research Center  
Under Contract NAS3-25556



National Aeronautics and  
Space Administration

(NASA-CR-187110) USER'S MANUAL FOR ROCKET  
COMBUSTOR INTERACTIVE DESIGN (ROCCID) AND  
ANALYSIS COMPUTER PROGRAM. VOLUME 2:  
APPENDIXES A-K (Final Report)  
(Aerojet-Gencorp Corp.) 578 p

N91-24231

Unclass

CSCL 21H 63/20 0013055



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**APPENDIX A**  
**INSTRUCTION SUMMARY**

## Instruction Summary

### -- GENERAL --

- \* All input must be in UPPER CASE.
- \* You will be prompted for all input data.
- \* You need only enter data you wish to change.
- \* You can go back and make corrections to input data at any time.

### -- TO ABORT --

- \* CONTROL Y kills everything.
- \* Tilde ( ~ ) saves results already completed.

### -- DATA ENTRY --

- \* General form is NAME[(index)] = data1 [,data2,... ]  
Note : a quantity in brackets indicates optional input  
where : "NAME" is a variable name which is prompted for  
where : "data" can be any of the following:
  - value (a real/integer number)  
i.e., RCHAMB=2.5 ; ICAV=1
  - N\*value (a repeated value)  
i.e., TQW=3\*2000...equivalent to TQW=2000,2000,2000
  - N(value1,value2,...) (a repeated group of values)  
i.e., XP=2(.96,.95)...equivalent to  
XP=.96,.95,.96,.95
  - " (an unchanged array value)  
i.e., XTQW=0,,3.5,...changes the 1st value to 0,  
leaves the 2nd value unchanged, and changes  
the 3rd value to 3.5
  - N\*, (a block of unchanged values)  
i.e., UEO=10,5\*,90...changes the 1st value to 10,  
leaves the 2nd thru 6th values unchanged and  
changes the 7th value to 90 and continues  
filling the array from there.
- \* A line ending with a COMMA signals more input for this array.  
Note: If a COMMA is inadvertently input at the end of a line,  
a <RET> input on the following line will terminate  
input.
- \* <RET> in place of a value signals no change for this variable  
i.e., the current or last input value is retained.
- \* <BACKSPACE> backspaces over characters which have been  
displayed.  
i.e., ==>UEO= is printed on the screen, to start  
changing input at the 9th value, enter a  
<BACKSPACE>, followed by: (9)=400,500,... which  
will appear as ==>UEO(9)=400,500..  
NOTE: <CTRL>H can be used on terminals without a <BACKSPACE>.
- \* @INSERT in place of a value allows the user to INSERT data  
into a table.
- \* @SKIP in place of a value signals the program to skip the  
remaining variables in that application and proceed to  
the next application.
- \* @HELP in place of a value will cause these instructions to be  
repeated.

**-- OVERVIEW --**

- \* The structure of this code is menu driven. Upon picking an application you will be shown the NAMELIST for that application, a description and the VARIABLES in that namelist. You then have the option to proceed, or return.

**-- REPLAY FILES --**

- \* A replay file contains all of the input entered from the terminal. This file can be edited using the EDT editor and used as the input for a subsequent run.
- \* Four special commands are available in the replay input alter mode:
  - @OFF - Stop input from the replay file. (This command can aid in keeping the replay file in-sink when entering a new menu option).
  - @ON - Resume alter input from the replay file (after the next input from the terminal).
  - @GO - Finish processing using the replay file without further keyboard input.
  - @SEARCH 'NAME' - Search through the replay file for 'NAME'. This may be used to get the REPLAY file back in sync. Must be proceeded with the @ON command if @OFF was entered.
- \* Please refer any comments, problems, bugs, etc. to :  
Software and Engineering Associates, Inc.  
Stu Dunn or Curtis Johnson  
(702) 882-1966



**APPENDIX B**  
**ERROR MESSAGES**

## INTERPRETATION OF ERROR MESSAGES

**Message:** ALLEN NOT CONVERGENT IN SUBROUTINE SWIRL, ALLEN=xxx FOR ELEMENT TYPE=xxx

**Remedy:** ALLEN is the "fan" length in SWIRL, the Aerojet swirl coaxial atomization routine. The user should check element inputs for consistency. The iteration counter can be increased in SWIRL, if this problem persists.

**Message:** AX BEYOND THROAT IN SUBROUTINE SHEAR; RUN STOPPED

**Remedy:** AX is the downstream axial position in SHEAR, the Aerojet shear coaxial atomization routine. The message often results when inconsistent element geometry is input. Output contained in the debug file may aid in evaluating the cause of this error.

**Message:** CHAMBR CALLED WITH INVALID MODEL, MCHAM=xxx

**Remedy:** An invalid chamber response model indicator input.

**Message:** COAXIAL ELEMENT FUEL VELOCITY - GAP ITERATION FAILED TO CONVERGE IN SUBROUTINE REDESIGN

**Remedy:** REDESIGN tries to balance the injector design for a directed change operating condition or element design (Section 3.6). This message occurs when the new velocity ratio ( $V_f/V_{o_j}$ ) is less than the minimum (VRATMI), and REDESIGN can not find an annulus gap which satisfies VRATMI. The simplest solution is to reduce VRATMI, but the output and history files should be reviewed first, to determine if the requested design change is reasonable. The iteration counter may also be increased from 20, if it is felt that convergence is slow.

**Message:** COAXIAL ELEMENT LOX FLOWRATE ITERATION CONVERGENCE FAILURE IN SUBROUTINE REDESIGN

**Remedy:** REDESIGN is trying to determine the LOX pressure drop required to meet the engine flowrate with the new fuel annulus design. An iteration counter has been included to prevent ROCCID from getting lost. This condition will result from a slowly converging solution, or more likely from a physically unrealistic fuel annulus size or pressure drop. Check these inputs.

**Message:** COAXIAL ELEMENT LOX FLOWRATE ITERATION CONVERGENCE FAILURE IN SUBROUTINE PDESIGN

**Remedy:** Subroutine PDESIGN calculates the design changes required to satisfy the performance goal. When the new required lox injection velocity has been determined, PDESIGN determines the injection pressure drop required to achieve the new injection velocity, with the existing lox post. This is an iterative process, and a counter has been included to preclude ROCCID from getting into an infinite loop.

**Message:** COMBR CALLED WITH INVALID MODEL, MBURN=xxx

**Remedy:** An invalid burning response model indicator input.

**Message:** COMBUST NONCONVERGENT IN C\*/FLOWRATE CALCULATIONS

**Remedy:** The delivered C\*-C\* efficiency-flowrate iteration in COMBUST utilizes a simple replacement iteration technique. While this technique works well for high performance combustors ( $\text{Eta-C}^* > 0.85$ ) convergence can be slow for lower performing combustors.

**Message:** COMBUST NONCONVERGENT IN TOTAL PRESSURE CALCULATIONS

**Remedy:** The total pressure loss iteration in COMBUST utilizes a simple replacement iteration technique. The maximum number of iterations can be increased if convergence is a persistent problem.

**Message:** CONVERGENCE ERROR IN ANNULAR GAP ITERATION FOR SWIRL COAX ELEMENTS IN SUBROUTINE CORESIZE AFTER xxx ITERATIONS NEL=xxx, GAP=xxx IN

**Remedy:** Subroutine CORESIZE determines the annular gap iteratively for swirl coaxial elements, as discussed in Section 3.2. This message occurs when the specified iteration limit is exceeded. Increasing this limit may resolve this problem, but the iteration history in the debug file (file type .DBG) should be examined.

**Message:** CONVERGENCE ERROR IN ANNULAR GAP ITERATION FOR SHEAR COAX ELEMENTS IN SUBROUTINE CORESIZE AFTER xxx ITERATIONS NEL=xxx, GAP=xxx IN

**Remedy:** Subroutine CORESIZE determines the annular gap iteratively for shear coaxial elements, as discussed in Section 3.2. This message occurs when the specified iteration limit is exceeded. Increasing this limit may resolve this problem, but the iteration history in the debug file (file type .DBG) should be examined.

**Message:** CONVERGENCE ERROR IN ENTROPY ITERATION IN SUBROUTINE  
VGALC  
MANIFOLD P (MPA),T (R), S(J/MOLE-K)=xxx, xxx, xxx  
INJECTED P (MPA),T (R), S(J/MOLE-K)=xxx, xxx, xxx

**Remedy:** Subroutine VGALC iteratively calculates the injection properties of gaseous propellants given the chamber pressure and the manifold pressure and temperature. This message comes from the iteration that determines the injection temperature by matching the injected entropy with the manifold value. The most common cause of this error is that the injected condition lies in the two-phase region of the H-S diagram, i.e. within the dome. This is only a problem because MIPROPS is not capable of determining the fluid quality. The user should check the operating conditions relative to the dome on an H-S diagram. The iteration process can be followed by examining the iteration record contained in the debug file (file type .DBG).

**Message:** CONVERGENCE ERROR IN ENTROPY ITERATION IN SUBROUTINE GASV  
MANIFOLD P (MPA),T (R), S(J/MOLE-K)=xxx, xxx, xxx  
INJECTED P (MPA),T (R), S(J/MOLE-K)=xxx, xxx, xxx

**Remedy:** Subroutine GASV iteratively calculates the injection properties of gaseous propellants given the chamber pressure and manifold temperature. This message comes from the iteration that determines the injection temperature by matching the injected entropy with the manifold value, which was calculated for the guessed manifold pressure. The most common causes of this error are 1) insufficient flow area for the required flow, thereby causing the velocity to become sonic, and 2) the injected condition lies in the two-phase region of the H-S diagram, i.e. within the dome. The latter is a problem because MIPROPS is not capable of determining the fluid quality. The user should check the flow area input. The iteration process can be followed by examining the iteration record contained in the debug file (file type .DBG).

**Message:** CONVERGENCE ERROR IN FLOWRATE ITERATION IN SUBROUTINE  
GASV MANIFOLD P (MPA)=xxx, T (R)= xxx  
INJECTED P (MPA)=xxx, T (R)= xxx  
DESIRED WDOT (LB/S)=xxx, CALCULATED WDOT=xxx

**Remedy:** Subroutine GASV iteratively calculates the injection properties of gaseous propellants given the chamber pressure and manifold temperature. This message comes from the iteration that determines the injected flowrate by adjusting the manifold pressure. The most common causes of this error are 1) insufficient flow area for the required flow, thereby causing the velocity to become sonic, and 2) the injected condition lies in the two-phase region of the H-S diagram, i.e. within the dome. The latter is a problem because MIPROPS is not capable of determining the fluid quality. The user should check the flow area input. The iteration process can be followed by examining the iteration record contained in the debug file (file type .DBG).

**Message:** CONVERGENCE ERROR IN FUEL PRESSURE DROP ITERATION.FOR  
SHEAR COAX ELEMENTS IN SUBROUTINE CORESIZE AFTER xxx  
ITERATIONS

**Remedy:** As discussed in Section 3.2, the core sizing routine, CORESIZE, tries to determine a fuel injection pressure that will increase the fuel injection velocity, thereby satisfying the minimum velocity ratio constraint. Since this process is iterative, a counter has been included to prevent the code from getting caught in a loop. The above message is the result of the number of iterations exceeding the counter, a condition often caused by either a very high oxidizer injection velocity or a large minimum velocity ratio, VRATMIN.

**Message:** CONVERGENCE ERROR IN FUEL PRESSURE DROP ITERATION FOR  
SWIRL COAX ELEMENTS IN SUBROUTINE CORESIZE AFTER xxx  
ITERATIONS

**Remedy:** As discussed in Section 3.2, the core sizing routine, CORESIZE, tries to determine a fuel injection pressure that will increase the fuel injection velocity, thereby satisfying the minimum velocity ratio constraint. Since this process is iterative, a counter has been included to prevent the code from getting caught in a loop. The above message is the result of the number of iterations exceeding the counter, a condition often caused by either a very high oxidizer injection velocity or a large minimum velocity ratio, VRATMIN.

**Message:** CONVERGENCE ERROR IN MACH NUMBER ITERATION IN SUBROUTINE  
RAYLEE: NEW MACH=xxx, OLD MACH=xxx,CALCULATIONS  
PROCEEDING WITH OLD TOTAL PRESSURE LOSS

**Remedy:** Subroutine RAYLEE calculates the change in total pressure due to change in total temperature, area ratio and mass flowrate. It numerically integrates Shapiro's influence coefficient equations from the chamber throat, using the input temperature, mass flow and area profiles. At each spacial location, RAYLEE iterates on the new mach number, using the mach number for the downstream condition. This message occurs when the iteration fails to converge, causing RAYLEE to use the old mach number for the mach number at the current station. There is no fix for this condition, and the message is included for informational purposes.

**Message:** CONVERGENCE FAILURE IN SUBROUTINE FRICTION; F,F1=xxx, xxx

**Remedy:** Subroutine FRICTION iteratively determines the Fanning friction factor from the input nondimensional roughness and Reynold's number using the Coolbrook correlation. A convergence error in FRICTION will cause this message to be printed, where F and F1 are the current and most recent past values of the friction factor.

**Message:** CONVERGENCE ERROR IN SUBROUTINE LEINJ AFTER 50 ITERATIONS  
FREQUENCY=xxx, IEL=xxx, INDX=xxx TRY INCREASING USGF, OGF  
AND/OR DSGF

**Remedy:** USGF, OGF and DSGF control the number of grid points used in the spatial integration in LEINJ. Adjusting any of these parameters in the model control file (file type .CNT) will effect model convergence. IEL refers to the element category, i.e. core (1), baffle (2), barrier (3) or FFC (4), while indx refers to propellant circuit, i.e. fuel (1) or ox (2).

**Message:** CONVERGENCE ERROR IN SUBROUTINE PRESSD FOR INDX=xxx

**Remedy:** Subroutine PRESSD calculates the pressure drop required to achieve the input total flowrate for propellant circuit INDX (INDX=1 for fuel, =2 for ox). Since mixed element patterns can be accomodated, this procedure must be iterative. An iteration counter has been included in PRESSD to prevent the code from running away. The convergence history is contained in the debug file (file type .DBG), and should be examined to diagnose the cause of the convergence failure.

**Message:** CONVERGENCE ERROR IN SUBROUTINE SWIRLPD "a" NOT CONVERGED  
AFTER 1000 ITERATIONS

**Remedy:** SWIRLPD calculates the flowrate, tip Cd, injection velocity and resultant spray cone angle for a swirl coaxial element of prescribed geometry and injection pressure drop. The variable a is used in this calculation procedure, and refers to the same value in the reference by Doumas and Laster. The most common cause of this error is incorrect element geometry input definitions, including NINLETS, CDINLET, DRATIO and AINLET (See Section 4.2 for more description of these input variables).

**Message:** CONVERGENCE ERROR IN SUBROUTINE SWIRLPD,"aPRIME" NOT  
CONVERGED AFTER 1000 ITERATIONS

**Remedy:** SWIRLPD calculates the flowrate, tip Cd, injection velocity and resultant spray cone angle for a swirl coaxial element of prescribed geometry and injection pressure drop. The variable aPRIME (or a') is used in this calculation procedure, and refers to the same value in the reference by Doumas and Laster. The most common cause of this error is incorrect element geometry input definitions, including NINLETS, CDINLET, DRATIO and AINLET (See Section 4.2 for more description of these input variables).

**Message:** CONVERGENCE FAILURE ERROR IN SUBROUTINE EMEST  
ETAMIX REQUIRED=xxx, ETAMIX CALCULATED=xxx, ETAM=xxx

**Remedy:** Subroutine EMEST determines the ETAM (related to Em) required to achieve the input Etamix. This message occurs when the iteration counter exceeds its maximum value. The error message output includes the current value of ETAM and the corresponding value of ETAMIX.

- Message: CONVERGENCE FAILURE IN CHAMBER LENGTH (OR ORIFICE DIAMETER OR INJECTION VELOCITY) ITERATION IN SUBROUTINE PDESIGN
- Remedy: PDESIGN is used to calculate the change in the injector design and operating parameters which will result in the performance goal being met. The change in the parameter is determined by iteratively solving for the effect of the variable change on vaporization efficiency. The above message will result typically unrealistically high or low performance requirements combined with length and/or pressure drop constraints. Review output and history file output along with your inputs (file type .DES) to evaluate if they are 1) consistent, 2) realistic.
- Message: \*\*\* DIST3D FREQUENCY ITERATION NONCONVERGENT AFTER xx ITERATIONS, CONSIDER INCREASING "IDMAX" \*\*\*
- Remedy: IDMAX is the maximum number of successive approximations permitted by DIST3D. IDMAX is contained in the Model Control Variables (file type .CNT). As the message indicates, increasing IDMAX can sometimes permit convergence of the iteration, although large values of IDMAX may result in error accumulation and invalid, negative frequencies.
- Message: ELEMENT PRESSURE DROP CONVERGENCE FAILURE IN SUBROUTINE DPOST
- Remedy: DPOST is trying to size the internal geometry of a shear coaxial element Lox circuit, so as to meet design constraints, e.g., DIVANG, the specified pressure drop and exit diameter, while providing for repeatable, attached flow conditions at the post exit. An iteration counter has been included to preclude getting caught in an infinite loop., This counter can be increased, but the reasonableness of model inputs should first be evaluated by examining the debug output of the iteration convergence process in the Debug file.(DBG).
- Message: ERROR IN CORESPAC, UNKNOWN ELEMENT TYPE=xxx
- Remedy: Subroutine CORESPAC spaces the injection elements radially and circumferentially. This message occurs when the element type is not one of those permitted, i.e. LOL, OFO, FOF, SHD, SHC or SWC.
- Message: ERROR IN CORESPAC, # ELEMENTS NOT DIVISIBLE BY # BAFFLE BLADES; NEL,NBAF=xxx, xxx
- Remedy: Although this error is not likely, the error checking has been included to preclude non-symmetric injector designs.
- Message: \*\*\* ERROR ENCOUNTERED IN NOZZLE GEOMETRY IN SUBROUTINE NOZINI \*\*\*
- Remedy: This message indicates that the input nozzle geometry is inconsistent, e.g. tangency points don't meet. This error is most likely to result when the user has created an input file without the assistance of the IFE, since it checks this condition during input.

Message: \*\*\* ERROR IN HCAVEF FOR F=xxx, PHIF SET TO 0.0

Remedy: HCAVEF is the HIFI subroutine which determines the wave circumferential orientation, relative to the cavity orientation, that result in minimum damping. F is the frequency, in hz, and PHIF is the injector face admittance for the minimally damped orientation. PHIF=0 is equivalent to no cavities being present. This message is mainly information, indicating that the user should consider the results suspect.

Message: \*\*\* ERROR IN HFCS, CHAMBER RESPONSE MINIMUM NOT FOUND FOR MODE M=xxx, N=xxx \*\*\*

Remedy: As discussed in Section 3.4, HFCS centers the chamber response frequency sweep about the calculated resonant frequency of the mode, neglecting any influences of damping devices. When the frequency sweep is complete, HFCS checks that a minimum exists in chamber response curve (not necessarily the fundamental mode, just a minimum). If the minimum is not found, HFCS shifts the center of the frequency sweep to a lower frequency and tries again. If the minimum is not found this time, the error message above is printed. The user should check the chamber response model output in the history file (file type .HIS) to confirm that this is the problem. While there is no remedy for this condition, the last value of the decay coefficient, AL, may be adequate to determine the high frequency stability characteristics for the subject mode, M tangential + N radial.

Message: \*\*\* ERROR IN HFCS, COMBUSTION RESPONSE PEAK NOT FOUND \*\*\*

Remedy: HFCS begins by determining the burning response curve versus frequency, as discussed in Section 3.4. To ensure that a large enough frequency range has been covered, it checks that the burning response magnitude reaches a maximum. If it is not found, HFCS will coarsen the frequency stepsize and look once more for the peak. If it still can't find the peak, this message is printed. The user should examine the burning response model output contained in the history file (file type .HIS), to ensure that this is the case. The user should also check the input model parameters, checking that the frequency range examined includes  $0.5/\text{Tau}$  hz, where  $\text{Tau} = \text{Tausen}$  if the N-Tau model was used, or Tau is the estimated droplet lifetime calculated by CRP.

Message: \*\*\* ERROR IN HFCS, STABILITY CONDITION NOT FOUND IN SUBROUTINE STABC \*\*\* AL=xxx, M=xxx, N=xxx

Remedy: As discussed in Section 3.4, HFCS uses the subroutine STABC to determine the frequency at which the maximum in-phase gain occurs. This message occurs when STABC can not find where the gain function passes between 180 and -180 degrees. The user should verify that the crossing does not exist by examining the STABC output in the history file (file type .HIS). While there is no remedy for this condition, the last value of the decay coefficient, AL, may be adequate to determine the high frequency stability characteristics for the subject mode, M tangential + N radial.

Message: **\*\*\* ERROR IN LFCS, ITERATION COUNTER EXCEEDED BEFORE MARGINAL CONDITION WAS FOUND \*\*\***

Remedy: In an effort to prevent LFCS from getting lost, and thereby pointlessly using excessive amounts of computer time, LFCS is only allowed to adjust the chamber pressure 100 times in its search for the marginal operating pressure (See Section 3.3 for more details). This message occurs when the counter has been exceeded. The user should check that 1) the chamber pressure iteration has not gotten lost or stuck, or 2) the calculations do not indicate that the configuration is either extremely stable or unstable, i.e. throttled to pressures excessively higher or lower than the nominal.

Message: **\*\*\* ERROR IN LFCS, LONGITUDINAL MODE NOT FOUND \*\*\***

Remedy: LFCS begins by determining the chamber response versus frequency, as discussed in Section 3.3. The frequency range is intended to exceed the first longitudinal (1L) resonant frequency. LFCS can confirm that this has occurred by checking for a chamber response minimum. If the minimum is not found, LFCS increases the chamber response frequency stepsize and tries to find the 1L minimum again. If the minimum is not found the second time, this message is printed, and LFCS stops. The user should ensure that the 1L was actually not found by examining the chamber response model output contained in the history file (file type **.HIS**), the user should also examine the frequency stepsize used in both the first and second frequency sweeps.

Message: **\*\*\* ERROR IN LFCS, STABILITY CONDITION NOT FOUND 15 TIMES SUBROUTINE STABC \*\*\***

Remedy: LFCS uses the subroutine STABC to determine the frequency at which the maximum in-phase gain occurs (See Section 3.3). If the system is highly chug stable, a crossing of the gain function between 180 and -180 degrees may not occur. Each time the condition is not found, LFCS will continue to throttle the engine, in an effort to find the marginally stable chamber pressure. LFCS has been designed to accept this error 15 times, before printing the above message and terminating execution. The user should verify that the crossing does not exist by examining the STABC output in the history file (file type **.HIS**). Since this message usually implies extremely large chug stability margins, the user should beware of extremely unstable injection-coupled longitudinal mode stability, i.e. the chamber length is excessively long, so the timelags can not couple with the bulk-flow (low frequency) oscillations.

Message: **\*\*\* ERROR \*\*\* MACH NUMBER GREATER THAN UNITY**

Remedy: This error comes from subroutine MACH, which calculates the mach number for an input area ratio and gamma using the successive approximation technique. The routine is limited to calculations for the subsonic branch, so this message indicates an error in the iteration process.

**Message:** ERROR IN PRELIMD, BOTH PROPELLANTS GASEOUS

**Remedy:** ROCCID is currently not capable of handling gas-gas combustors. This is due to the problems which would arise during the calculation of total timelags

**Message:** ERROR IN REDESIGN, FUEL ORIFICE SIZE CAN NOT BE SPECIFIED FOR SHEAR OR SWIRL ELEMENTS

**Remedy:** Specifying a new fuel orifice diameter for coaxial elements has been avoided in the current code, since this can result in an interminable loop (See Section 3.6). This is message will only occur if the user has modified the code to vary this parameter.

**Message:** ERROR IN SUBROUTINE ALOAD, BAFFLE ELEMENT TYPE xxx NOT RECOGNIZED

**Remedy:** Input baffle element type invalid, i.e. not LOL, OFO, FOF, SHD, SHC or SWC. This error is most common when the user creates input files without the assistance of the IFE. The most likely cause is that the value of the character variable TYPE is not in single quotes (').

**Message:** ERROR IN SUBROUTINE ALOAD, BARRIER ELEMENT TYPE xxx NOT RECOGNIZED

**Remedy:** Input barrier element type invalid, i.e. not LOL, OFO, FOF, SHD, SHC or SWC. This error is most common when the user creates input files without the assistance of the IFE. The most likely cause is that the value of the character variable TYPE is not in single quotes (').

**Message:** ERROR IN SUBROUTINE ALOAD, CORE ELEMENT TYPE xxx NOT RECOGNIZED

**Remedy:** Input core element type invalid, i.e. not LOL, OFO, FOF, SHD, SHC or SWC. This error is most common when the user creates input files without the assistance of the IFE. The most likely cause is that the value of the character variable TYPE is not in single quotes (').

**Message:** ERROR IN SUBROUTINE CORESIZE, COAXIAL ELEMENT SPECIFIED WITH LIQUID FUEL

**Remedy:** As indicated in Sections 2.0 and 5.0, coaxial elements require that the fuel is gaseous, i.e. that the manifold temperature be above the propellant critical temperature. Check the input fuel temperature.

Message: **\*\*\* ERROR IN SUBROUTINE DCAVEF FOR OMEGA=xxx BETAC SET TO 0.0**

Remedy: DCAVEF is the DIST3D subroutine which determines the wave circumferential orientation, relative to the cavity orientation, that result in minimum damping. OMEGA is the frequency, in hz, and BETAC is the effective cavity admittance. BETAC=0 is equivalent to no cavities being present. This message is mainly information, indicating that the user should consider the results suspect.

Message: **\*\*\* ERROR IN SUBROUTINE DINPUT \*\*\* NO DATA AVAILABLE FOR ELEMENT TYPE=xxx**

Remedy: Subroutine DINPUT reads the design definition files at the beginning of a POINTD run (file types .DES and .DEF). This message occurs when the element type is not one of those permitted, i.e. LOL, OFO, FOF, SHD, SHC or SWC.

Message: **\*\*\* ERROR IN SUBROUTINE DLOAD \*\*\* NO DATA AVAILABLE FOR ELEMENT TYPE xxx**

Remedy: Input core element type invalid, i.e. not LOL, OFO, FOF, SHD, SHC or SWC. This error is most common when the user creates input files without the assistance of the IFE. The most likely cause is that the value of the character variable TYPE is not in single quotes (').

Message: **\*\*\* ERROR IN SUBROUTINE NOZADM, CAN NOT GET OUT OF THROAT AFTER xxx ATTEMPTS \*\*\***

Remedy: Subroutine NOZADM numerically calculates the nozzle admittance by integrating from the nozzle throat to the beginning of the constant diameter section of the chamber (nozzle entrance). The first integration step is a pure Newton step (See Appendix K), so a small initial stepsize is desired. Unfortunately, if the stepsize is too small, the integration step occurs within the throat and terms like  $1/(1-M^2)$  become undefined. To avoid this error, NOZADM increases the initial stepsize. If NOZADM can not get out of the throat after several attempts, this message is printed, and the calculations stopped. This error is most common in large diameter nozzles with a large throat entrance radius.

Message: **ERROR IN PRELIMD, CALCULATED CR=xxx <1.0**

Remedy: Although this message is not likely, it has been included to preclude the user from continuing into more detailed analyses if this condition exists. It is most likely to occur with portions of the chamber geometry fixed, and insufficient flowrate or chamber pressure.

**Message:** ERROR IN SUBROUTINE PRELIMD, GASEOUS PRESSURE DROP  
ITERATIONS FOR DPMIN DID NOT CONVERGE AFTER xxx ITERATIONS

**Remedy:** PRELIMD must solve the injection pressure drop required for a specified gaseous propellant injection velocity iteratively. This message occurs when the iteration fails to converge at the minimum pressure drop.

**Message:** ERROR IN SUBROUTINE PRELIMD, GASEOUS PRESSURE DROP  
ITERATIONS FOR DPNOM DID NOT CONVERGE AFTER xxx ITERATIONS

**Remedy:** PRELIMD solves the injection pressure drop required for a specified gaseous propellant injection velocity iteratively. This message occurs when the iteration fails to converge at the minimum pressure drop.

**Message:** ERROR IN SUBROUTINE PRELIMD, PCNOM NOT CONVERGED AFTER xx  
ITERATIONS

**Remedy:** If the user specifies the manifold pressures, PRELIMD solves for the nominal chamber pressure iteratively. The iteration scheme tries to use all the available pressure drop. This message occurs when the iteration fails to converge. Further insight into the convergence iteration can be found in the debug output file (file type .DBG).

**Message:** \*\*\* ERROR IN SUBROUTINE SHEAR, LIQUID FUEL DETECTED \*\*\*

**Remedy:** Adjust fuel temperature and or pressure to adhere to constraints outlined in Section 2.2 or change injector element type

**Message:** ERROR IN SUBROUTINE SHEARPD DUE TO ERROR IN SUBROUTINE  
FRICTION

**Remedy:** Subroutine FRICTION is used by SHEARPD to iteratively determine the Fanning friction factor from the input nondimensional roughness and Reynold's number using the Coolbrook correlation. A convergence error in FRICTION will cause this fatal error in SHEARPD.

**Message:** \*\*\* ERROR IN SUBROUTINE SWIRL, LIQUID FUEL DETECTED \*\*\*

**Remedy:** Adjust fuel temperature and or pressure to adhere to constraints outlined in Section 2.2 or change injector element type

**Message:** ERROR IN SUBROUTINE SPLINT: KHI=KLO - STOPPED

**Remedy:** SPLINT is the cubic spline interpolation routine in Subroutine RAYLEE. This message occurs when the table contains an invalid array of derivative values. It is most likely to occur if this routine is improperly accessed by a user-added analysis model.

**Message:** ERROR IN SUBROUTINE SWIRLSZR, "a" ITERATION DID NOT CONVERGE AFTER 1000 ITERATIONS

**Remedy:** SWIRLSZR determines the geometry of the swirl coaxial post, given element flowrate and injection pressure drop. The variable a is used in this calculation procedure, and refer to the same value in the reference by Doumas and Laster. The most common cause of this error is incorrect element geometry input definitions, including CDINLET and DRATIO (See Section 4.2 for more description of these input variables).

**Message:** ERROR IN SUBROUTINE SWIRLSZR, CONVERGENCE FAILURE IN SWIRL HOLE DIAMETER ITERATION

**Remedy:** SWIRLSZR determines the geometry of the swirl coaxial post, given element flowrate and injection pressure drop. This message occurs when the program is unable to determine a swirl chamber inlet orifice diameter which satisfies other flow constraints (See reference by Doumas and Laster). The user may want to try adjusting the element geometry input definitions CDINLET and DRATIO (See Section 4.2 for more description of these input variables).

**Message:** ERROR IN SUBROUTINE SWILRSZR, aPRIME ITERATION DID NOT CONVERGE AFTER 1000 ITERATIONS

**Remedy:** SWIRLSZR determines the geometry of the swirl coaxial post, given element flowrate and injection pressure drop. The variable aPRIME (or a') is used in this calculation procedure, and refers to the same value in the reference by Doumas and Laster. The most common cause of this error is incorrect element geometry input definitions, including CDINLET and DRATIO (See Section 4.2 for more description of these input variables).

**Message:** ERROR WITH AMINE FLAME CALCULATIONS IN SUBROUTINE VAPRO, RUN STOPPED

**Remedy:** As noted in Section 3.2, COMBUST still contains the dual-flame, monopropellant amine vaporization correction. This message can only occur when an amine fuel is used. It indicates that the calculated generalized length correction term is in some way inconsistent.

Message: ERROR WITH CHAMBER GEOMETRY INPUT, BARREL SECTION LESS OR EQUAL 0  
CALCULATED BARREL LENGTH, IN=xxx  
CALCULATED CONVERGENT SECTION LENGTH, IN=xxx  
CHECK GEOMETRY INPUTS, INCLUDING UNITS:  
RCHAMB, FT=xxx, RTHRT, FT=xxx,  
RNE, FT=xxx, RTE, FT=xxx, ALPHA, DEG=xxx,

Remedy: This error message is generated by subroutine PINPUT. Since ROCCID, in particular the chamber response models can not handle purely conical chambers, this error message has been added. It should be noted that this error may also occur due to inconsistencies in geometry inputs, i.e. barrel lengths less than 0. This error is most likely to result when the user has created an input file without the assistance of the IFE, since it checks this condition during input.

Message: ERROR WITH NOZZLE GEOMETRY AT TANGENT POINT ALPHA=xxx DEG

Remedy: This message, generated by subroutine PINPUT, indicates that the input nozzle geometry is inconsistent, especially where the tangency points should meet. This error is most likely to result when the user has created an input file without the assistance of the IFE, since it checks this condition during input.

Message: ERROR WITH VAPORIZATION INTERPOLATION IN SUBROUTINE  
VAPRO,RUN STOPPED

Remedy: Subroutine VAPRO performs the propellant droplet vaporization using Priem's Generalized Length Correlation (See Section 2.2). This message indicates that the calculated generalized length correction term is in some way inconsistent.

Message: EXPANSION PRESSURE DROP CALCULATION FAILED IN SUBROUTINE  
DPOST

Remedy: DPOST is trying to size the internal geometry of a shear coaxial element Lox circuit, so as to meet design constants, e.g., DIVANG, the specified pressure drop and exit diameter, while providing for repeatable, attached flow conditions at the post exit. An iteration counter has been included to preclude getting caught in an infinite loop., This counter can be increased, but the reasonableness of model inputs should first be evaluated by examining the debug output of the iteration convergence process in the Debug file.(.DBG).

Message: INJR CALLED WITH INVALID MODEL, MINJ=xxx

Remedy: Invalid injector response model index input.

- Message: INPUT ERROR TO FUNCTION AINTP, "X" ARRAY NOT MONOTONICALLY INCREASING X(1)=xxx, xxx, ...
- Remedy: AINTP performs 1-D interpolations, but requires the independent array, X, to be monotonically increasing. This message usually results when the user has input arrays, i.e. XNOZ, XMRA, PCA, without using the IFE, which checks them.
- Message: INPUT ERROR TO FUNCTION GETVAL, ARRAY "XA" NOT MONOTONICALLY DECREASING XA(1)=xxx, xxx, ...
- Remedy: GETVAL performs 1-D power-law interpolations, but requires the independent array, XA, to be monotonically increasing. This message usually results when the user has input array PCA without using the IFE, which checks them.
- Message: INVALID NUMBER OF INPUTS TO SUBROUTINE GETVAL, N=xxx
- Remedy: GETVAL performs 1-D power-law interpolations. The independent and dependent arrays must be contain at least 2 points.
- Message: LESS THAN ONE DROP FORMED OR NEGATIVE RJET IN SUBROUTINE SHEAR
- Remedy: This error is most likely to result from inconsistent element geometry or during application of the model (SHEAR) to unusual element designs (Section 2.1). Check input element dimensions.
- Message: LVAP NOT CONVERGED FOR TRIPLET IN SUBROUTINE TIMELAG, RUN STOPPED
- Remedy: Because of the potential for a dropsizes distribution when triplet injection elements are used, the subroutine TIMELAG must calculate the 20% vaporization length, LVAP, iteratively (See Section 5.1 for more details). This error message occurs when the iteration counter, which is included to prevent the program from getting lost in an infinite loop, is exceeded.
- Message: MACH CALLED WITH ISUB=0, RUN ABORTED
- Remedy: Older versions of subroutine MACH contained a flag ISUB which determined whether the calculation was for the subsonic (ISUB=1) or the supersonic (ISUB=0) branch of the area ratio-mach number calculation. The current routine has retained the flag for compatability, but is only capable of calculating Mach numbers for the subsonic branch.

- Message:** MORE DROPS PRODUCED IN SUBROUTINE SHEAR THAN DIMENSIONED FOR
- Remedy:** This message is often the result of an incorrectly input element geometry, e.g., too small of a fuel annulus gap. If these values are OK, the size of the arrays R, XD, ATLEN and VJL can be increased in subroutine SHEAR from 500.
- Message:** READ ERROR ENCOUNTERED ON UNIT 9 BEFORE THIRD \$END FOUND IN SUBROUTINE VSAVE
- Remedy:** Subroutine VSAVE reads the namelist \$SAVE, contained at the end of the design input file (file type **.DES**). It reads the the current design variables, contained is \$SAVE, which are not contained in the input file, e.g. injection pressure, dropsizes, etc. It must exist, even if there are no values, i.e. PRELIMD hasn't been run yet. The primary cause of the error message is the incorrect manual creation of the **.DES** file, i.e. without the use of the IFE.
- Message:** RJET NOT CONVERGENT IN SUBROUTINE SHEAR
- Remedy:** This error is most likely to result from inconsistent element geometry or during application of the model (SHEAR) to unusual element designs (Section 2.1). Check input element dimensions.
- Message:** SHARP-EDGED ORIFICE CALCULATIONS FAIL IN SUBROUTINE DPOST FOR DMS,DSE,CC0SE=xxx, xxx, xxx
- Remedy:** Subroutine DPOST sizes the lox post internal geometry for shear coaxial elements. This iteration tries to determine the sharp-edged orifice diameter (DSE) with an equivalent contraction coefficient (CC0SE) as the metering section diameter (DMS). The iteration counter is included to prevent the code from entering an infinite loop. Progress of the iteration can be checked in the debug file (file type **.DBG**)
- Message:** \*\*\* SUCCESSION APPROXIMATIONS IN DIST3D LEAD TO NEGATIVE REAL FREQUENCY FOR M=xxx AND N=xxx, RUN STOPPED
- Remedy:** The iteration process in DIST3D can result in an invalid, negative frequency, which will predict negative dissipation. There is currently no solution for this error.
- Message:** TERMINAL ERROR IN SUBROUTINE GASV, TABLES NOT AVAILABLE FOR GAS PHASE PROPERTIES OF xxx
- Remedy:** GASV utilizes the MIPROPS routines, packaged in FLUIDP. If the MIPROPS propellant data doesn't exist, GASV must stop.

**Message:**   **TERMINAL ERROR IN SUBROUTINE VGCALC, TABLES NOT AVAILABLE FOR GAS PHASE PROPERTIES OF xxx**

**Remedy:**    VGCALC utilizes the MIPROPS routines, packaged in FLUIDP. If the MIPROPS propellant data doesn't exist, VGCALC must stop.

**Message:**   **TOO MANY TIME STEPS INPUT TO SUBROUTINE LEINJ, NTINJ=xxx,>LIMIT(50), RETRY WITH SMALLER VALUE**

**Remedy:**    Hardwired matrix sizing in LEINJ limits the number of time integrations per oscillation to 50. Reduce the value in the model control file (file type .CNT) and try again.

**Message:**   **\*\*\* WARNING FROM LFCS, PC CONVERGED WITHOUT MARGINAL CONDITION BEING SATISFIED, FOLLOWING RESULTS SHOULD BE CONSIDERED SUSPECT:**

**Remedy:**    The throttling procedure included in LFCS will occasionally get stuck in the chamber pressure iteration, homing in on a chamber pressure which does not result in a marginally stable system gain (a maximum in-phase magnitude of 1.0). The user should ensure that 1) the converged  $P_c$  is adequate for their minimum needs, and 2) that the system gain amplitude is not very close to 1.0, thereby indicating that the error is only a matter of numerical tolerancing.

## Interactive Front End Error Messages

- ERROR -- Bad integer input --try again  
User input for an integer variable was not an interger.
- ERROR -- Bad menu option chosen --try again  
Menu option chosen was not valid.
- ERROR -- Bad real input --try again  
User input for a real variable was not a real number.
- ERROR -- Both propellants in gaseous state  
For the given operating conditions both propellants are gaseous.
- ERROR -- Cavity geometry not compatible with chamber radius  
Chamber geometry failed the following test:  
IF ICAV=2 THEN  
     $2 * \pi * RCHAMB \geq [NCAV(1) * \sqrt{4 * AC(1) / \pi} + NCAV(2) * \sqrt{4 * AC(2) / \pi} + (NCAV(1) + NCAV(2)) * TPART]$   
OTHERWISE  
     $2 * \pi * RCHAMB \geq [NCAV(1) * (AC(1) / WC(1) + TPART) + NCAV(2) * (AC(2) / WC(2) + TPART)]$
- ERROR -- CCAV array; NCAV entries must be greater than 0.0  
The aperture length for this particular cavity is nonzero,  
but the sound speed (CCAV\*) is zero for the cavity.
- ERROR -- CGAM array; NCAV entries must be greater than 0.0  
The aperture length for this particular cavity is nonzero,  
but the specific heat ratio (CGAM\*) is zero for the cavity.
- ERROR -- Chamber diameter bigger than DCMAX  
RCHAMB was entered at a value larger than DCMAX/2.0.
- ERROR -- Chamber length longer than XLEMAX  
CHAMBL was entered at a value larger than XLEMAX.
- ERROR -- Chamber pressure must be greater than zero (psia)  
Neither chamber pressure nor manifold pressures were entered.
- ERROR -- Contraction ratio less than 1  
Chamber geometry failed the following test:  
RCHAMB  $\geq$  RTHRT
- ERROR -- Could not find correct plot data  
The expected plot data file could not be found. Check for file existence.
- ERROR -- Could not open file  
File could not be opened - either it does not exist, the path to the file is incorrect,  
or the file is locked by another user.

- ERROR -- Data files must be opened before an analysis run can begin.  
The data files have not been opened before starting an analysis run. Enter the SET VARIABLES menu option, then exit that menu. When prompted for saving the data, answer YES.
- ERROR -- DDIF must be less than or equal to XDJ  
Self explanatory.
- ERROR -- DMS must be less than or equal to DIFF  
Self explanatory.
- ERROR -- During a READ  
A read was unsuccessful, most commonly caused by incorrect entries to a namelist.
- ERROR -- Geometry not possible --Chamber radii prevent connecting tangent  
Chamber geometry failed the following test:  
$$RTHRT + RTE * [1 - \cos(\alpha)] \leq RCHAMP - RNE * [1 - \cos(\alpha)]$$
- ERROR -- Geometry not possible --check chamber length  
Input geometry failed the following test:  
$$CHAMBL \leq (RTE + RNE) * \sin(\alpha) + [RCHAMB - RTHRT + (RTE - RNE) * (1 - \cos(\alpha))] / \tan(\alpha)$$
- ERROR -- ICTYP array; NCAV absorbers have not been entered  
More absorbers (NCAV in number) have been requested, but not all have been typed. There must be NCAV entries of absorber types in the ICTYP array.
- ERROR -- Injector geometry is incompatible -check FDJ, TPOST, and XDJ  
The following test failed:  $FDJ - XDJ - 2 * TPOST \leq 0$
- ERROR -- Input out of bounds  
User input was out of the acceptable range for that variable.
- ERROR -- Manifold temperatures must be greater than zero (R)  
Manifold temperatures were not entered, or entered less than -460 degrees F.
- ERROR -- Nominal flow rate is less than minimum  
WDNOM is less than WDMIN.
- ERROR -- Problem reading plot data  
When reading a plot file (\*.PL1, \*.PL2, \*.PL3, etc), the data required to make the plot was not found. Check formatting of the plot file.
- ERROR -- Problem reading plot dimension data  
The program cannot read the actual plot data, check data format.
- ERROR -- RE2 array; NCAV entries must be greater than RS2 and less than the chamber radius  
Either the chamber radius has been set nonzero, and the radius to the outer edge of the absorber segment (RE2) is greater than the chamber radius, or the absorber segment outer radius (RE2) is less than the absorber segment inner radius (RS2).

- ERROR -- Required propellant properties not available**  
Required propellant properties do not exist for this operating condition. Most commonly seen when running fuel through a coaxial injection element.
- ERROR -- RHOAP array; NCAV entries must be greater than 0.0**  
For this particular cavity, the cavity type is either a quarterwave, or user defined (ICTYP\*=1 or 4), the cavity has finite size, but the density (RHOAP\*) is zero.
- ERROR -- RS2 array; NCAV entries must be greater than 0.0 and less than the chamber radius**  
The chamber radius has been set nonzero, and the start of the absorber segment (RS2) has been set greater than the chamber radius or less than zero.
- ERROR -- There are not NCAV(1) 1s in IDCAV**  
There must be NCAV(1) ones in the IDCAV array.
- ERROR -- There are not NCAV(2) 2s in IDCAV**  
There must be NCAV(2) twos in the IDCAV array.
- ERROR -- Unrecognized element type, please try again.**  
The element type entered was not recognized, the type must be: LOL, OFO, FOF, SHD, SHC, SWC.
- ERROR -- Value must be greater than zero**  
A variables was set to zero, when zero is not permitted.
- ERROR -- XMWC array; NCAV entries must be greater than 0.0**  
The aperture length for this particular cavity is nonzero, but the molecular weight (XMWC\*) is zero for the cavity.
- ERROR -- ZCOMB must be less than chamber length**  
The chamber length has been set to a nonzero length, and then combustion plane (ZCOMB) has been set larger than the chamber length.
- ERROR -- ZE1 array; NCAV entries must be greater than ZS1 and less than the chamber length**  
Either the chamber length has been set nonzero, and the distance from the injector face to the end of the absorber segment (ZE1) is greater than the chamber length, or the absorber segment end (ZE1) is less than the absorber segment start (ZS1).
- ERROR -- ZLOW array; NCAV entries must be set (less than backing cavity width - WC)**  
Cavity type is 1, 2, or 3 and ZLOW is either larger than ZUP, or larger than the backing cavity width (WC\*).
- ERROR -- ZS1 array; NCAV entries must be greater than 0.0 and less than the chamber length**  
The chamber length has been set nonzero, and the start of the absorber segment (ZS1) has been set greater than the chamber length or less than zero.

ERROR -- ZUP array; NCAV entries must be set (less than backing cavity width - WC and greater than ZLOW)  
Cavity type is 1, 2, or 3 and ZUP has been set larger than the backing cavity width (WC\*).



**APPENDIX C**  
**NAMelist VARIABLE DEFINITIONS**

# NAMELIST VARIABLE DEFINITIONS

## POINT ANALYSIS

<u>Namelist</u>	<u>Variable</u>	<u>Type</u>	<u>Range</u>	<u>Default</u>	<u>Path</u>	<u>Size</u>	<u>Description</u>
\$MODELS:	MCHAM	INTEGER	1-3	1	1	1	CHAMBER RESPONSE MODEL FLAG; 1-HIFI, 2-DIST3D 3-FDORC
	MBURN	INTEGER	1-2	2	1	1	COMBUSTION RESPONSE MODEL FLAG; 1-CRP, 2-N-TAU
	MINJ	INTEGER	1-2	1	1	1	CHAMBER MODE FLAG; 1-INJ, 2-LEINJ

## \$OPCOND

C-2	FUEL	CHAR*8	N/A		1	1	FUEL NAME, E.G. RP-1, H2, METHANE, PROPANE
	OX	CHAR*8	N/A	LOX	1	1	OXIDIZER NAME, E.G. LOX
	PC	REAL	>0	0.0	1	1	INJECTOR FACE STAGNATION PRESSURE (PSIA)
	XMR	REAL	>0	0.0	1	1	OX/FUEL MIXTURE RATIO
	HGMR*	REAL	>=0	0.0	5	1	OX/FUEL MIXTURE RATIO OF FUEL, IF STAGED COMBUSTION IS USED
	FTMAN	REAL	>-460	-500.0	1	1	FUEL MANIFOLD TEMPERATURE (DEG F)
	XTMAN	REAL	>-460	-500.0	1	1	OXIDIZER MANIFOLD TEMPERATURE (DEG F)
	FBLEED	REAL	0-1	0.0	5	1	FRACTION OF TOTAL FUEL USED FOR UNIFORM FACE BLEED
	EMMAN	REAL	0-1	1.0	1	1	MIXING NONUNIFORMITY DUE TO MANIFOLD MALDISTRIBUTION (1-UNIFORM, NO EFFECT)
	NPERFP	INTEGER	2-30	2	1	1	NUMBER OF POINTS INPUT FOR ISP AND C* VS. MR TABLES
	PMRA	REAL	MONO.	0.0	1	30	ARRAY OF MIXTURE RATIO POINTS FOR PERFORMANCE TABLES (RECOMMEND MR'S FROM COMBUSTION TABLES)
	PISPA	REAL	>=0	0.0	1	30	ARRAY OF ODK-ISP POINTS FOR PERFORMANCE TABLES (SEC)
	PCSA	REAL	>=0	0.0	1	30	ARRAY OF ODK-C* POINTS FOR PERFORMANCE TABLES (FT/SEC)

\*Not used in February 1991 ROCCID version, for future use

POINT ANALYSIS (Continued)

<u>Namelist</u>	<u>Variable</u>	<u>Type</u>	<u>Range</u>	<u>Default</u>	<u>Path</u>	<u>Size</u>	<u>Description</u>
\$GEOM	RCHAMB	REAL	>0	0.0	1	1	CHAMBER RADIUS (FT.)
	RTHRT	REAL	>0	0.0	1	1	THROAT RADIUS (FT.)
	RNE	REAL	>0	0.0	1	1	RADIUS OF CURVATURE AT THE NOZZLE ENTRANCE (FT.)
	RTE	REAL	>0	0.0	1	1	RADIUS OF CURVATURE AT THE THROAT ENTRANCE (FT.)
	ALPHA	REAL	<90	0.0	1	1	CONVERGENCE HALF ANGLE (DEG.)
	CHAMBL	REAL	>0	0.0	1	1	INJECTOR FACE TO THROAT LENGTH (FT.)
	XC	REAL	>0	0.0	4	1	CHAMBER CYLINDRICAL LENGTH (CALCULATED) (FT.)
\$SCORE	TYPE	CHAR*8	N/A		1	1	CORE ELEMENT TYPE: LOL, OFO, FOF, SHD, SHC, SWC
	NEL	INTEGER	>=0	0	1	1	NUMBER OF ELEMENTS
	FDJ	REAL	>=0	0.0	1	1	FUEL ORIFICE OR ANNULUS DIAMETER (IN.)
	FCD	REAL	0-1	1.0	1	1	FUEL ORIFICE Cd
	FIH	REAL	>=0	0.0	1	1	FUEL IMPINGEMENT HEIGHT (IN.)
	FIA	REAL	<45	0.0	1	1	FUEL IMPINGEMENT HALF-ANGLE (DEG.)
	FCANT	REAL	0-45	0.0	1	1	FUEL UNLIKE CANT ANGLE (DEG.)
	FFACET	REAL	>0	0.001	1	1	FUEL FACEPLATE THICKNESS OR ANNULUS LENGTH (IN.)
	XDJ	REAL	>=0	0.0	1	1	OX ORIFICE OR POST DIAMETER (IN.)
	XCD	REAL	0-1	1.0	1	1	OX ORIFICE Cd (IMPINGING ELEMENTS ONLY)
	XIH	REAL	>=0	0.0	1	1	OX IMPINGEMENT HEIGHT (IN.)
	XIA	REAL	<45	0.0	1	1	OX IMPINGEMENT HALF-ANGLE (DEG.)
	XCANT	REAL	0-45	0.0	1	1	OX UNLIKE CANT ANGLE (DEG.)
	XFACET	REAL	>0	0.001	1	1	OX FACEPLATE THICKNESS FOR IMPINGING ELEMENTS OR POST LENGTH SWIRL COAX (IN.)
	TPOST	REAL	>0	0.0	1	1	OX POST WALL THICKNESS FOR COAX ELEMENTS (IN.)
	RECESS	REAL	>=0	0.0	1	1	OX POST RECESS FOR SHEAR COAX ELEMENTS (IN.)
	DMS	REAL	>0	0.0	1	1	OX METERING SECTION DIAMETER FOR SHEAR COAX ELEMENTS (IN.)

POINT ANALYSIS (Continued)

<u>Namelist</u>	<u>Variable</u>	<u>Type</u>	<u>Range</u>	<u>Default</u>	<u>Path</u>	<u>Size</u>	<u>Description</u>
	XMS	REAL	>0	0.0	1	1	OX METERING SECTION LENGTH FOR SHEAR COAX ELEMENTS (IN.)
	DDIF	REAL	>0	0.0	1	1	OX DIFFUSER SECTION DIAMETER FOR SHEAR COAX ELEMENTS (IN.)
	XDL	REAL	>0	0.0	1	1	OX POST DIFFUSER SECTION LENGTH FOR SHEAR COAX ELEMENTS (IN.)
	AINLET	REAL	>0	0.0	1	1	SWIRL COAX SWIRL CHAMBER FEED ORIFICE AREA (PER HOLE) (IN**2)
	CDINLET	REAL	>0	1.0	1	1	SWIRL CHAMBER FEED ORIFICE Cd
	NINLET	INTEGER	>0	0	1	1	NUMBER OF SWIRL CHAMBER FEED ORIFICES (SWIRL COAX ELEMENTS ONLY)
	DRATIO	REAL	>1	1.0	1	1	SWIRL CHAMBER TO ELEMENT EXIT DIAMETER RATIO (DS/XDJ)
	EMUNI	REAL	<1	0.0	4	1	UNIELEMENT RUPE MIXING EFFICIENCY
	CBINT	REAL	0-1	0.0	5	1	INTERACTION INDEX FOR MIXING BETWEEN CORE AND BARRIER

C-4

**\$BAFFLE**

TYPE	CHAR*8	N/A			1	1	BAFFLE INJECTOR ELEMENT TYPE, E.G. LOL, FOF, OFO, SHC, SWC, SHD
NEL	INTEGER	>=0	0		1	1	NUMBER OF ELEMENTS
FDJ	REAL	>=0	0.0		1	1	FUEL ORIFICE OR ANNULUS DIAMETER (IN.)
FCD	REAL	0-1	1.0		1	1	FUEL ORIFICE Cd
FIH	REAL	>=0	0.0		1	1	FUEL IMPINGEMENT HEIGHT (IN.)
FIA	REAL	<45	0.0		1	1	FUEL IMPINGEMENT HALF-ANGLE (DEG.)
FCANT	REAL	0-45	0.0		1	1	FUEL UNLIKE CANT ANGLE (DEG.)
FFACET	REAL	>0	0.001		1	1	FUEL FACEPLATE THICKNESS OR ANNULUS LENGTH (IN.)
FINJ	REAL	>=0	0.0		1	1	FUEL INJECTION POINT RELATIVE TO INJECTOR FACE (IN.)
XDJ	REAL	>=0	0.0		1	1	OX ORIFICE OR POST DIAMETER (IN.)
XCD	REAL	0-1	1.0		1	1	OX ORIFICE Cd (IMPINGING ELEMENTS ONLY)
XIH	REAL	>=0	0.0		1	1	OX IMPINGEMENT HEIGHT (IN.)
XIA	REAL	<45	0.0		1	1	OX IMPINGEMENT HALF-ANGLE (DEG.)
XCANT	REAL	0-45	0.0		1	1	OX UNLIKE CANT ANGLE (DEG.)

POINT ANALYSIS (Continued)

<u>Namelist</u>	<u>Variable</u>	<u>Type</u>	<u>Range</u>	<u>Default</u>	<u>Path</u>	<u>Size</u>	<u>Description</u>
	XFACET	REAL	>0	0.001	1	1	OX FACEPLATE THICKNESS FOR IMPINGING ELEMENTS OR POST LENGTH SWIRL COAX (IN.)
	XINJ	REAL	>=0	0.0	1	1	OX INJECTION POINT RELATIVE TO INJECTOR FACE (IN.)
	TPOST	REAL	>0	0.0	1	1	OX POST WALL THICKNESS FOR COAX ELEMENTS (IN.)
	RECESS	REAL	>=0	0.0	1	1	OX POST RECESS FOR SHEAR COAX ELEMENTS (IN.)
	DMS	REAL	>0	0.0	1	1	OX METERING SECTION DIAMETER FOR SHEAR COAX ELEMENTS (IN.)
	XMS	REAL	>0	0.0	1	1	OX METERING SECTION LENGTH FOR SHEAR COAX ELEMENTS (IN.)
	DDIF	REAL	>0	0.0	1	1	OX DIFFUSER SECTION DIAMETER FOR SHEAR COAX ELEMENTS (IN.)
	XDL	REAL	>0	0.0	1	1	OX POST DIFFUSER SECTION LENGTH FOR SHEAR COAX ELEMENTS (IN.)
	AINLET	REAL	>0	0.0	1	1	SWIRL COAX SWIRL CHAMBER FEED ORIFICE AREA (PER HOLE) (IN**2)
C5	CDINLET	REAL	>0	1.0	1	1	SWIRL CHAMBER FEED ORIFICE Cd
	NINLET	INTEGER	>0	0	1	1	NUMBER OF SWIRL CHAMBER FEED ORIFICES (SWIRL COAX ELEMENTS ONLY)
	DRATIO	REAL	>1	1.0	1	1	SWIRL CHAMBER TO ELEMENT EXIT DIAMETER RATIO (DS/XDJ)
	EMUNI	REAL	<1	0.0	5	1	UNELEMENT RUPE MIXING EFFICIENCY

**\$BARRIER**

TYPE	CHAR*8	N/A	1	1	1	1	BARRIER INJECTOR ELEMENT TYPE, E.G. LOL, FOF, OFO, SHC, SWC, SHD
NEL	INTEGER	>=0	0	1	1	1	NUMBER OF ELEMENTS
FDJ	REAL	>=0	0.0	1	1	1	FUEL ORIFICE OR ANNULUS DIAMETER (IN.)
FCD	REAL	0-1	1.0	1	1	1	FUEL ORIFICE Cd
FIH	REAL	>=0	0.0	1	1	1	FUEL IMPINGEMENT HEIGHT (IN.)
FIA	REAL	<45	0.0	1	1	1	FUEL IMPINGEMENT HALF-ANGLE (DEG.)
FCANT	REAL	0-45	0.0	1	1	1	FUEL UNLIKE CANT ANGLE (DEG.)
FFACET	REAL	>0	0.001	1	1	1	FUEL FACEPLATE THICKNESS OR ANNULUS LENGTH (IN.)
XDJ	REAL	>=0	0.0	1	1	1	OX ORIFICE OR POST DIAMETER (IN.)

POINT ANALYSIS (Continued)

<u>Namelist</u>	<u>Variable</u>	<u>Type</u>	<u>Range</u>	<u>Default</u>	<u>Path</u>	<u>Size</u>	<u>Description</u>
	XCD	REAL	0-1	1.0	1	1	OX ORIFICE Cd (IMPINGING ELEMENTS ONLY)
	XIH	REAL	>=0	0.0	1	1	OX IMPINGEMENT HEIGHT (IN.)
	XIA	REAL	<45	0.0	1	1	OX IMPINGEMENT HALF-ANGLE (DEG.)
	XCANT	REAL	0-45	0.0	1	1	OX UNLIKE CANT ANGLE (DEG.)
	XFACET	REAL	>0	0.001	1	1	OX FACEPLATE THICKNESS FOR IMPINGING ELEMENTS OR POST LENGTH SWIRL COAX (IN.)
	TPOST	REAL	>0	0.0	1	1	OX POST WALL THICKNESS FOR COAX ELEMENTS (IN.)
	RECESS	REAL	>=0	0.0	1	1	OX POST RECESS FOR SHEAR COAX ELEMENTS (IN.)
	DMS	REAL	>0	0.0	1	1	OX METERING SECTION DIAMETER FOR SHEAR COAX ELEMENTS (IN.)
	XMS	REAL	>0	0.0	1	1	OX METERING SECTION LENGTH FOR SHEAR COAX ELEMENTS (IN.)
	DDIF	REAL	>0	0.0	1	1	OX DIFFUSER SECTION DIAMETER FOR SHEAR COAX ELEMENTS (IN.)
	XDL	REAL	>0	0.0	1	1	OX POST DIFFUSER SECTION LENGTH FOR SHEAR COAX ELEMENTS (IN.)
	AINLET	REAL	>0	0.0	1	1	SWIRL COAX SWIRL CHAMBER FEED ORIFICE AREA (PER HOLE) (IN**2)
	CDINLET	REAL	>0	1.0	1	1	SWIRL CHAMBER FEED ORIFICE Cd
	NINLET	INTEGER	>0	0	1	1	NUMBER OF SWIRL CHAMBER FEED ORIFICES (SWIRL COAX ELEMENTS ONLY)
	DRATIO	REAL	>1	1.0	1	1	SWIRL CHAMBER TO ELEMENT EXIT DIAMETER RATIO (DS/XDJ)
	EMUNI	REAL	<1	0.0	5	1	UNELEMENT RUPE MIXING EFFICIENCY
	FFCINT	REAL	0-1	0.0	5	1	INTERACTION INDEX FOR MIXING BETWEEN BARRIER AND FFC

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\$FFC

	TYPE	CHAR*8	N/A		1	1	FFC INJECTOR ELEMENT TYPE, E.G. LOL, SHD
	NEL	INTEGER	>=0	0	1	1	NUMBER OF ELEMENTS
	FDJ	REAL	>=0	0.0	1	1	FUEL ORIFICE OR ANNULUS DIAMETER (IN.)
	FCD	REAL	0-1	1.0	1	1	FUEL ORIFICE Cd
	FIH	REAL	>=0	0.0	1	0	FUEL IMPINGEMENT HEIGHT (IN.)
	FIA	REAL	<45	0.0	1	1	FUEL IMPINGEMENT HALF-ANGLE (DEG.)
	FCANT	REAL	0-45	0.0	1	0	FUEL UNLIKE CANT ANGLE (DEG.)

POINT ANALYSIS (Continued)

<u>Namelist</u>	<u>Variable</u>	<u>Type</u>	<u>Range</u>	<u>Default</u>	<u>Path</u>	<u>Size</u>	<u>Description</u>
	FFACET	REAL	>0	0.001	1	1	FUEL FACEPLATE THICKNESS OR ANNULUS LENGTH (IN.)
	XDJ	REAL	0	0.0	4	1	OX ORIFICE OR POST DIAMETER (IN.)
\$BURN							
	GAMMA	REAL	>0	0.0	4	1	COMBUSTION GAS SPECIFIC HEAT RATIO
	AO	REAL	>0	0.0	4	1	COMBUSTION GAS STAGNATION SOUND SPEED (FT/S)
	GMW	REAL	>0	0.0	4	1	COMBUSTION GAS MOLECULAR WEIGHT (LBm/LB-Mole)
	GPR	REAL	>0	0.0	4	1	COMBUSTION GAS PRANDTL NUMBER
	GK	REAL	>0	0.0	4	1	COMBUSTION GAS THERMAL CONDUCTIVITY (BTU/FT-S-R)
	GMU	REAL	>0	0.0	4	1	COMBUSTION GAS VISCOSITY (LBm/FT-S)
	VGASI	REAL	>0	0.0	4	1	COMBUSTION GAS VELOCITY AT THE INJECTOR FACE (FT/S)
	RML	REAL	>0	0.0	4	1	MASS MEDIAN DROPLET RADIUS (MICRONS)
	VJL	REAL	>0	0.0	4	1	DROPLET INJECTION VELOCITY (FT/S)
	TJL	REAL	>0	0.0	4	1	DROPLET INJECTION TEMPERATURE (DEG. R)
	RHOL	REAL	>0	0.0	4	1	DROPLET DENSITY (LBm/FT**3)
	CPL	REAL	>0	0.0	4	1	DROPLET MEAN HEAT CAPACITY (BTU/LBm-R)
	PCRITL	REAL	>0	0.0	4	1	DROPLET CRITICAL PRESSURE (PSIA)
	TCRITL	REAL	>0	0.0	4	1	DROPLET CRITICAL TEMPERATURE (DEG. R)
	TBOILL	REAL	>0	0.0	4	1	NORMAL BOILING POINT OF DROPLET (DEG. R)
	XMWL	REAL	>0	0.0	4	1	DROPLET MOLECULAR WEIGHT (LBm/LB-Mole)
	HVAPL	REAL	>0	0.0	4	1	DROPLET HEAT OF VAPORIZATION AT NBP (BTU/LBm)
	EN	REAL	>0	0.0	4	1	PRESSURE INTERACTION INDEX
	TAUSEN	REAL	>0	0.0	4	1	SENSITIVE TIME LAG (SEC.)
	ISEN	INTEGER	1-2	0	4	1	INDEX FOR SENSITIVE PROPELLANT CIRCUIT; 1-FUEL, 2-OX

POINT ANALYSIS (Continued)

<u>Namelist</u>	<u>Variable</u>	<u>Type</u>	<u>Range</u>	<u>Default</u>	<u>Path</u>	<u>Size</u>	<u>Description</u>
\$INJ	FMAND	REAL	>0	100.0	1	1	FUEL MANIFOLD CHARACTERISTIC DIAMETER (IN.)
	XMAND	REAL	>0	100.0	1	1	OX MANIFOLD CHARACTERISTIC DIAMETER (IN.)
	FMANL	REAL	>0	1000.0	1	1	FUEL MANIFOLD CHARACTERISTIC LENGTH (IN.)
	XMANL	REAL	>0	1000.0	1	1	OX MANIFOLD CHARACTERISTIC LENGTH (IN.)
	PCA	REAL	>0	0.0	2	3	CHAMBER PRESSURE ARRAY FOR INJECTION VARIABLES, 3 INPUTS IN DECENDING PRESSURE (PSIA)
	FRA	REAL	>0	0.0	3	3	FUEL CIRCUIT RESISTANCE, 3 INPUTS REQUIRED
	FCAPA	REAL	>0	0.0	3	3	FUEL CIRCUIT CAPACITANCE, 3 INPUTS REQUIRED
	NFE	INTEGER	1-4	4	4	1	NUMBER FOR FUEL ELEMENT TYPES
	FTLA	REAL	>0	0.0	3	12	FUEL TOTAL TIME LAG ARRAY, 3*NFE INPUTS REQUIRED (SEC.)
	FINA	REAL	>0	0.0	3	12	FUEL ORIFICE INERTANCE ARRAY, 3*NFE INPUTS REQUIRED (SEC.)
	FFA	REAL	0-1	0.0	4	12	FRACTION OF TOTAL FUEL FLOW FOR EACH ELEMENT TYPE, 3*NFE INPUTS REQUIRED
	NXE	INTEGER	1-4	4	4	1	NUMBER FOR OX ELEMENT TYPES
	XRA	REAL	>0	0.0	3	3	OX CIRCUIT RESISTANCE, 3 INPUTS REQUIRED
	XCAPA	REAL	>0	0.0	3	3	OX CIRCUIT CAPACITANCE, 3 INPUTS REQUIRED
	XTLA	REAL	>0	0.0	3	12	OX TOTAL TIME LAG ARRAY, 3*NXE INPUTS REQUIRED (SEC.)
	XINA	REAL	>0	0.0	3	12	OX ORIFICE INERTANCE ARRAY, 3*NXE INPUTS REQUIRED (SEC.)
	XFA	REAL	0-1	0.0	4	12	FRACTION OF TOTAL OX FLOW FOR EACH ELEMENT TYPE, 3*NXE INPUTS REQUIRED
	XUOR	REAL	>0	0.0	3	8	UPSTREAM SECTION LENGTH (DIMENSIONED 2X4 W/2->PROPELLANT (1-F,2-O) AND 4->ELEMENT TYPE) (IN)
	AUOR	REAL	>0	0.0	3	8	UPSTREAM SECTION FLOW AREA (DIMENSIONED 2X4 W/2->PROPELLANT (1-F,2-O) AND 4->ELEMENT TYPE) (IN**2)
	RUOR	REAL	>0	0.0	3	8	UPSTREAM RADIUS, FOR VISCOUS CALCULATIONS (DIMENSIONED 2X4 W/2->PROPELLANT (1-F,2-O) AND 4->ELEMENT TYPE) (IN)
	XOR	REAL	>0	0.0	3	8	ORIFICE SECTION LENGTH (DIMENSIONED 2X4 W/2->PROPELLANT (1-F,2-O) AND 4->ELEMENT TYPE) (IN)

# POINT ANALYSIS (Continued)

<u>Namelist</u>	<u>Variable</u>	<u>Type</u>	<u>Range</u>	<u>Default</u>	<u>Path</u>	<u>Size</u>	<u>Description</u>
	AOR	REAL	>0	0.0	3	8	ORIFICE SECTION FLOW AREA (DIMENSIONED 2X4 W/2->PROPELLANT (1-F,2-O) AND 4->ELEMENT TYPE) (IN**2)
	ROR	REAL	>0	0.0	3	8	ORIFICE RADIUS, FOR VISCOUS CALCULATIONS (DIMENSIONED 2X4 W/2->PROPELLANT (1-F,2-O) AND 4->ELEMENT TYPE) (IN)
	XDOR	REAL	>0	0.0	3	8	DOWNSTREAM SECTION LENGTH (DIMENSIONED 2X4 W/2->PROPELLANT (1-F,2-O) AND 4->ELEMENT TYPE) (IN)
	ADOR	REAL	>0	0.0	3	8	DOWNSTREAM SECTION FLOW AREA (DIMENSIONED 2X4 W/2->PROPELLANT (1-F,2-O) AND 4->ELEMENT TYPE) (IN**2)
	RDOR	REAL	>0	0.0	3	8	DOWNSTREAM RADIUS, FOR VISCOUS CALCULATIONS (DIMENSIONED 2X4 W/2->PROPELLANT (1-F,2-O) AND 4->ELEMENT TYPE) (IN)
	WDOT	REAL	>0	0.0	4	6	PROPELLANT FLOWRATE ARRAY (DIMENSIONED 2X3 W/2->PROPELLANT (1-F,2-O) AND 3->CHAMBER PRESSURE (PCA)) (LBM/S)

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## \$CHAMBER

NCAV	INTEGER	0-40	0	1	2	ARRAY OF NUMBER OF CAVITIES OF TYPE 1 & 2, TWO INPUTS REQUIRED
XB	REAL	>=0	0.0	3	1	DISTANCE FROM INJECTOR FACE TO CONCENTRATED COMBUSTION PLANE (HIFI) (FT.)
ZS	REAL	>=0	0.0	3	1	DISTANCE FROM INJECTOR FACE TO START OF COMBUSTION DISTRIBUTION (DIST3D) (FT.)
ZE	REAL	>=0	0.0	3	1	DISTANCE FROM INJECTOR FACE TO END OF COMBUSTION DISTRIBUTION (DIST3D) (FT.)
MUB	INTEGER	1-12	1	1	1	NUMBER OF EVENLY SPACED RADIAL BAFFLES
T	REAL	>0	0.001	1	1	BAFFLE BLADE THICKNESS (FT.)
ZB	REAL	>=0	0.0	1	1	BAFFLE LENGTH (FT.)
TPART	REAL	>=0	0.0	1	1	CAVITY PARTITION (SEPARATOR) THICKNESS (WIDTH) (FT.)
NSEC	INTEGER	1-5	1	3	2	ARRAY OF NUMBER OF CAVITIES PROPERTY SECTIONS FOR CAVITY TYPE 1 & 2, TWO INPUTS REQUIRED

POINT ANALYSIS (Continued)

<u>Namelist</u>	<u>Variable</u>	<u>Type</u>	<u>Range</u>	<u>Default</u>	<u>Path</u>	<u>Size</u>	<u>Description</u>
	WC	REAL	>=0	0.0	1	2	1/4 WAVE CAVITY WIDTH ARRAY FOR CAVITY TYPE 1 & 2, TWO INPUTS REQUIRED (FT.)
	AC	REAL	>=0	0.0	1	2	CAVITY SECTION CROSS-SECTIONAL AREA ARRAY FOR CAVITY TYPE 1 & 2 TWO INPUTS REQUIRED (FT**2)
	VCAV	REAL	>=0	0.0	1	2	BACKING CAVITY VOLUME FOR HELMHOLTZ RESONATORS (FT**3)
	ARATIO	REAL	0-1	1.0	1	2	INLET-TO-BACKING CAVITY AREA RATIO FOR HELMHOLTZ RESONATORS
	INLET	INTEGER	0-2	0	1	2	CAVITY INLET DESCRIPTOR ARRAY FOR CAVITY TYPE 1 & 2, TWO INPUTS REQUIRED; 0-SQUARE INLET, 1-ROUNDED INLET, 2-WELL ROUNDED INLET
	IDCAV	INTEGER	0-2	0	1	40	CAVITY RELATIVE LOCATION ARRAY, NCAV(1)+NCAV(2) ENTRIES REQUIRED
	DC1	REAL	>=0	0.0	1	5	1/4 WAVE CAVITY EFFECTIVE LENGTH (DEPTH) OR HELMHOLTZ INLET ORIFICE LENGTH (FT.) ARRAY FOR CAVITY TYPE 1, NSEC(1) INPUTS REQUIRED
	CC1	REAL	>=0	0.0	2	5	CAVITY STAGNATION SOUND SPEED ARRAY FOR CAVITY TYPE 1, NSEC(1) INPUTS REQUIRED (FT/S)
	GAMC1	REAL	>=0	0.0	2	5	CAVITY GAS RATIO OF SPECIFIC HEATS ARRAY FOR CAVITY TYPE 1, NSEC(1) INPUTS REQUIRED
	DC2	REAL	>=0	0.0	1	5	1/4 WAVE CAVITY EFFECTIVE LENGTH (DEPTH) OR HELMHOLTZ INLET ORIFICE LENGTH (FT.) ARRAY FOR CAVITY TYPE 2, NSEC(2) INPUTS REQUIRED (FT)
	CC2	REAL	>=0	0.0	2	5	CAVITY STAGNATION SOUND SPEED ARRAY FOR CAVITY TYPE 2, NSEC(2) INPUTS REQUIRED (FT/S)
	GAMC2	REAL	>=0	0.0	2	5	CAVITY GAS RATIO OF SPECIFIC HEATS ARRAY FOR CAVITY TYPE 2, NSEC(2) INPUTS REQUIRED
	ICAV	INTEGER	0-2	0	1	1	ACOUSTIC CAVITY FLAG; 0 - NO ACOUSTIC CAVITY, 1 - 1/4 WAVE CAVITY, 2 - HELMHOLTZ RESONATOR
\$FDORC	ZCOMB	REAL	>0-XC	0	2	1	AXIAL LOCATION WHERE COMBUSTION IS COMPLETED (FT)
	NZON	INT	1-20	5	2	1	NUMBER OF COMBUSTION ZONES
	FTER	REAL	0.0-1.0	1.0	2	20	FRACTION OF TOTAL ENERGY RELEASE AT END OF EACH COMBUSTION ZONE, NZON INPUTS REQUIRED

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POINT ANALYSIS (Continued)

<u>Namelist</u>	<u>Variable</u>	<u>Type</u>	<u>Range</u>	<u>Default</u>	<u>Path</u>	<u>Size</u>	<u>Description</u>
	NCAV1	INT	0-20	0	1	1	NUMBER OF RADIAL INLET ABSORBERS
	ICTYP1	INT	1-4	0	1	20	ABSORBER TYPE FLAG: 1=1/4 WAVE, 2=HELMHOLTZ, 3=LONG APERTURE, 4=INPUT GEOMETRY AND TEMPERATURE; NCAV1 INPUTS REQUIRED
	ZS1	REAL	0-XC	0.0	1	20	DISTANCE FROM INJECTOR FACE TO START OF ABSORBER SEGMENT, NCAV1 INPUTS REQUIRED (FT)
	ZE1	REAL	ZS1-XC	0.0	1	20	DISTANCE FROM INJECTOR FACE TO END OF ABSORBER SEGMENT, NCAV1 INPUTS REQUIRED (FT)
	AS1	REAL	0.0-360.0	0.0	1	20	ANGLE AT WHICH ABSORBER SEGMENT STARTS, NCAV1 INPUTS REQUIRED (DEG)
	AE1	REAL	0.0-360.0	0.0	1	20	ANGLE AT WHICH ABSORBER SEGMENT ENDS, NCAV1 INPUTS REQUIRED (DEG)
	WC1	REAL	>=0.0	0.0	1	20	BACKING CAVITY WIDTH, =0 FOR 1/4 WAVE CAVITY; NCAV1 INPUTS REQUIRED (FT)
	APL1	REAL	>=0.0	0.0	1	20	INLET APERTURE LENGTH, NCAV1 INPUTS REQUIRED (FT)
	BCL1	REAL	>=0.0	0.0	1	20	BACKING CAVITY LENGTH, =0 FOR 1/4 WAVE CAVITY; NCAV1 INPUTS REQUIRED (FT)
	ZLOW1	REAL	0-WC1	0.0	1	20	DISTANCE FROM CAVITY BOTTOM TO UPPER POINT OF INTERSECTION OF APERTURE WITH BACKING CAVITY, =0 FOR 1/4 WAVE; NCAV1 INPUTS REQUIRED (FT)
	ZUP1	REAL	ZLOW1- WC1	0.0	1	20	DISTANCE FROM CAVITY BOTTOM TO UPPER POINT OF INTERSECTION OF APERTURE WITH BACKING CAVITY, =0 FOR 1/4 WAVE; NCAV1 INPUTS REQUIRED (FT)
	CCAV1	REAL	>0.0	1.0	2	20	BACKING CAVITY SOUND SPEED, NCAV1 INPUTS REQUIRED (FT/S)
	CGAM1	REAL	>0.0	0.0	2	20	BACKING CAVITY RATIO OF SPECIFIC HEATS, NCAV1 INPUTS REQUIRED
	XMWC1	REAL	>0.0	0.0	2	20	BACKING CAVITY GAS MOLECULAR WEIGHT, NCAV1 INPUTS REQUIRED (LBm/LB-MOLE)
	RHOAP1	REAL	>0.0	0.0	2	20	APERTURE DENSITY, NCAV1 INPUTS REQUIRED (LBM/FT**3)
	TSL1	REAL	NONE	0.0	0	20	SLOPE OF MEAN TEMPERATURE PROFILE, NCAV1 INPUTS REQUIRED (R/FT)
	TC1	REAL	>0.0	0.0	0	20	AVERAGE TEMPERATURE IN THE ABSORBER, NCAV1 INPUTS REQUIRED (R)

POINT ANALYSIS (Continued)

<u>Namelist</u>	<u>Variable</u>	<u>Type</u>	<u>Range</u>	<u>Default</u>	<u>Path</u>	<u>Size</u>	<u>Description</u>
	NCAV2	INT	0-20	0	1	1	NUMBER OF AXIAL INLET ABSORBERS ABSORBER TYPE FLAG: 1=1/4 WAVE, 2=HELMHOLTZ, 3=LONG APERTURE, 4=INPUT GEOMETRY AND TEMPERATURE; NCAV2 INPUTS REQUIRED
	ICTYP2	INT	1-4	0	1	20	
	RS2	REAL	0-	0.0	1	20	RADIUS TO INNER EDGE OF ABSORBER SEGMENT, NCAV2 INPUTS REQUIRED (FT)
	RE2	REAL	RCHAMB RS2-	0.0	1	20	RADIUS TO OUTER EDGE OF ABSORBER SEGMENT, NCAV2 INPUTS REQUIRED (FT)
	AS2	REAL	RCHAMB 0.0-360.0	0.0	1	20	ANGLE AT WHICH ABSORBER SEGMENT STARTS, NCAV2 INPUTS REQUIRED (DEG)
	AE2	REAL	0.0-360.0	0.0	1	20	ANGLE AT WHICH ABSORBER SEGMENT ENDS, NCAV2 INPUTS REQUIRED (DEG)
	WC2	REAL	>=0.0	0.0	1	20	BACKING CAVITY WIDTH, =0 FOR 1/4 WAVE CAVITY; NCAV2 INPUTS REQUIRED (FT)
	APL2	REAL	>=0.0	0.0	1	20	INLET APERTURE LENGTH, NCAV2 INPUTS REQUIRED (FT)
	BCL2	REAL	>=0.0	0.0	1	20	BACKING CAVITY LENGTH, =0 FOR 1/4 WAVE; NCAV2 INPUTS REQUIRED (FT)
	ZLOW2	REAL	0-WC2	0.0	1	20	DISTANCE FROM CAVITY BOTTOM TO LOWER POINT OF INTERSECTION OF APERTURE WITH BACKING CAVITY, =0 FOR 1/4 WAVE; NCAV2 INPUTS REQUIRED (FT)
	ZUP2	REAL	ZLOW2- WC2	0.0	1	20	DISTANCE FROM CAVITY BOTTOM TO UPPER POINT OF INTERSECTION OF APERTURE WITH BACKING CAVITY, =0 FOR 1/4 WAVE; NCAV2 INPUTS REQUIRED (FT)
	CCAV2	REAL	>0.0	1.0	2	20	BACKING CAVITY SOUND SPEED, NCAV2 INPUTS REQUIRED (FT/S)
	CGAM2	REAL	>0.0	0.0	2	20	BACKING CAVITY RATIO OF SPECIFIC HEATS, NCAV2 INPUTS REQUIRED
	XMWC2	REAL	>0.0	0.0	2	20	BACKING CAVITY GAS MOLECULAR WEIGHT, NCAV2 INPUTS REQUIRED (LBm/LB-MOLE)
	RHOAP2	REAL	>0.0	0.0	2	20	APERTURE DENSITY, NCAV2 INPUTS REQUIRED (LBm/FT**3)
	TSL2	REAL	NONE	0.0	2	20	SLOPE OF MEAN TEMPERATURE PROFILE, NCAV2 INPUTS REQUIRED (R/FT)
	TC2	REAL	>0.0	0.0	2	20	AVERAGE TEMPERATURE IN THE ABSORBER, NCAV2 INPUTS REQUIRED (R)

POINT ANALYSIS (Continued)

<u>Namelist</u>	<u>Variable</u>	<u>Type</u>	<u>Range</u>	<u>Default</u>	<u>Path</u>	<u>Size</u>	<u>Description</u>
\$MIX	EM	REAL	0-1.0	-1.0,1.0	1	2	OVERALL CORE (EM(1)) AND BARRIER (EM(2)) RUPE MIXING EFFICIENCIES
\$STUFF							
	IBAF	INTEGER	0-1	0	1	1	COMBUSTOR RADIAL BAFFLE BLADES FLAG;0 - NO BAFFLES, 1 - CONTAINS BAFFLES.
\$FINJ							
	IBFE	INTEGER	0-1	0	1	1	BAFFLE ELEMENT FLAG 0 - NO BAFFLE ELEMENTS, 1 - BAFFLE ELEMENTS
	IBRE	INTEGER	0-1	0	1	1	BARRIER ELEMENT FLAG 0 - NO BARRIER ELEMENTS, 1 - BARRIER ELEMENTS
	IFFE	INTEGER	0-1	0	1	1	FUEL FILM AND/OR CAVITY COOLING ELEMENT FLAG 0 - NO FFC ELEMENTS, 1 - FFC ELEMENTS

# **POINT DESIGN VARIABLES**

<u>Namelist</u>	<u>Variable</u>	<u>Type</u>	<u>Range</u>	<u>Default</u>	<u>Path</u>	<u>Size</u>	<u>Description</u>
\$DESIGN	TYPE	CHAR*8	N/A		1	1	CORE ELEMENT TYPE: LOL, OFO, FOF, SHD, SHC, SWC
	FUEL	CHAR*8	N/A		1	1	FUEL NAME, E.G. RP-1, H2, METHANE, PROPANE
	OX	CHAR*8	N/A	LOX	1	1	OXIDIZER NAME, E.G. LOX
	XMR	REAL	>0	0.0	1	1	OX/FUEL MIXTURE RATIO
	HGMR*	REAL	>=0	0.0	5	1	OX/FUEL MIXTURE RATIO OF FUEL, IF STAGED COMBUSTION IS USED
	FTMAN	REAL	>-460	-500.0	1	1	FUEL MANIFOLD TEMPERATURE (DEG F)
	XTMAN	REAL	>-460	-500.0	1	1	OXIDIZER MANIFOLD TEMPERATURE (DEF F)
	PCNOM	REAL	>0	0.0	1	1	INJECTOR FACE STAGNATION PRESSURE (PSIA)
	FPMAN	REAL	>0	0.0	1	1	FUEL MAXIMUM MANIFOLD PRESSURE (PSIA)
	XPMAN	REAL	>0	0.0	1	1	OXIDIZER MAXIMUM MANIFOLD PRESSURE (PSIA)
	WDNOM	REAL	>0	0.0	1	1	FLOW RATE AT NOMINAL CHAMBER PRESSURE (LBM/S)
	WDMIN	REAL	>0	0.0	1	1	FLOW RATE AT THROTTLED OR MINIMUM CHAMBER PRESSURE (LBM/S)
	EREG	REAL	<1	0.0	1	1	ISP-BASED ENERGY RELEASE EFFICIENCY GOAL
	ETACSG	REAL	<1	0.0	1	1	CSTAR EFFICIENCY GOAL
	XLEMAX	REAL	>0	10.0	1	1	MAXIMUM ENGINE LENGTH (FT)
	DCMAX	REAL	>0	10.0	1	1	MAXIMUM COMBUSTION CHAMBER DIAMETER (FT)
	NPERFP	INTEGER	2-30	2	1	1	NUMBER OF POINTS INPUT FOR ISP AND C* VS. MR TABLES
	PMRA	REAL	MONO.	0.0	1	30	ARRAY OF MIXTURE RATIO POINTS FOR PERFORMANCE TABLES (RECOMMEND MR'S FROM COMBUSTION TABLES)
	PISPA	REAL	>=0	0.0	1	30	ARRAY OF ODK-ISP POINTS FOR PERFORMANCE TABLES (SEC)
	PCSA	REAL	>=0	0.0	1	30	ARRAY OF ODK-C* POINTS FOR PERFORMANCE TABLES (FT/SEC)

POINT DESIGN VARIABLES (Continued)

<u>Namelist</u>	<u>Variable</u>	<u>Type</u>	<u>Range</u>	<u>Default</u>	<u>Path</u>	<u>Size</u>	<u>Description</u>
	NNOZ	INTEGER	0-30	0	1	1	NUMBER OF NOZZLE LENGTH POINTS IN PERFORMANCE OPTIMIZATION TABLE
	XNOZ	REAL	MONO.	0.0	1	30	NOZZLE LENGTH ARRAY (NNOZ POINTS REQUIRED) (FT)
	ETANOZ	REAL	0-1	0.0	1	30	NOZZLE EFFICIENCY ARRAY (NNOZ POINTS REQUIRED)
	IBAF	INTEGER	0-1	0	1	1	COMBUSTOR RADIAL BAFFLE BLADES FLAG; 0 - BAFFLES NOT ALLOWED, 1 - BAFFLES ALLOWED.
	ICAVT	INTEGER	-1-2	0	1	1	ACOUSTIC CAVITY FLAG; -1 - 1/4 WAVE AXIAL CAVITY ALLOWED, 0 - CAVITY NOT ALLOWED, 1 - 1/4 WAVE RADIAL CAVITY ALLOWED, 2 - HELMHOLTZ RESONATOR ALLOWED.

**\$FGEOM**

C-15	RCHAMB	REAL	>=0	0.0	1	1	CHAMBER RADIUS (FT.)
	RTHRT	REAL	>=0	0.0	1	1	THROAT RADIUS (FT.)
	RNE	REAL	>=0	0.0	1	1	RADIUS OF CURVATURE AT THE NOZZLE ENTRANCE (FT.)
	RTE	REAL	>=0	0.0	1	1	RADIUS OF CURVATURE AT THE THROAT ENTRANCE (FT.)
	ALPHA	REAL	0-90	0.0	1	1	CONVERGENCE HALF ANGLE (DEG.)
	CHAMBL	REAL	>=0	0.0	1	1	INJECTOR FACE TO THROAT LENGTH (FT.)

# MODEL CONTROL VARIABLES

<u>Namelist</u>	<u>Variable</u>	<u>Type</u>	<u>Range</u>	<u>Default</u>	<u>Path</u>	<u>Size</u>	<u>Description</u>
\$DEBUGC	DEBUG	LOGICAL	T/F	F	1	1	DEBUG OUTPUT GENERATION CONTROL (L1)
\$HIFIC	SHORT	LOGICAL	T/F	F	3	1	TRUE IF SHORT NOZZLE IS ASSUMED, FALSE IF NOT (L1)
	POC	REAL	>=0	0.20	1	2	NORMALIZED CAVITY INLET PRESSURE AMPLITUDE FOR CAVITY TYPE 1 & 2, TWO INPUTS REQUIRED
\$DIST3DC	SHORT	LOGICAL	T/F	F	3	1	TRUE IF SHORT NOZZLE IS ASSUMED, FALSE IF NOT (L1)
	POC	REAL	>=0	0.20	1	2	NORMALIZED CAVITY INLET PRESSURE AMPLITUDE FOR CAVITY TYPE 1 & 2, TWO INPUTS REQUIRED
	PAMP MX	REAL INTEGER	>=0 0-1	0.20 0	1 1	1 1	PEAK-TO-PEAK PRESSURE AMPLITUDE RATIO, P/PC MAIN CHAMBER OSCILLATION INDICATOR; 0- STANDING WAVE, 1-TRAVELING WAVE
	MC	INTEGER	1-20	11	1	1	# OF FOURIER SERIES TERMS TO REPRESENT THE MAIN CHAMBER SOLUTION
	LC	INTEGER	1-20	8	1	1	# OF BESSEL TERMS IN THE MAIN CHAMBER SOLUTION
	MB	INTEGER	1-20	11	1	1	# OF FOURIER SERIES TERMS TO REPRESENT THE SOLUTION IN THE BAFFLE COMP
	LB	INTEGER	1-20	8	1	1	# OF BESSEL TERMS IN THE BAFFLE COMPARTMENT SOLUTION
	IDMAX	INTEGER	1-25	10	1	1	MAXIMUM NUMBER OF ITERATIONS FOR SUCCESSIVE APPROX.

MODEL CONTROL VARIABLES (Continued)

<u>Namelist</u>	<u>Variable</u>	<u>Type</u>	<u>Range</u>	<u>Default</u>	<u>Path</u>	<u>Size</u>	<u>Description</u>
\$CRPC	NDPC	INTEGER	>0	16	1	1	NUMBER OF DROPLETS INJECTED/CYCLE
	NDTFQ	INTEGER	>0	16	1	1	USED TO DETERMINE THE TIME STEP SIZE (DT) WHEN CALCULATING VAPORATION HISTORY OF THE DROPLET (1.0/FREQUENCY/NDTFQ)
	NDTLF	INTEGER	>0	1000	1	1	USED TO DETERMINE THE TIME STEP SIZE (DT) WHEN CALCULATING VAPORATION HISTOR OF THE DROPLET DT (TLIFE/NDTLF) WHERE TLIFE IS THE DROPLET LIFE TIME
	NPRINT	INTEGER	>0	50	1	1	NUMBER OF TIME STEPS BETWEEN OUTPUTS OF THE DROPLET HISTORY FOR FIRST
	NSUMS	INTEGER	1-1E4	3500	1	1	NUMBER OF SUMMATION HISTORIES PER PERIOD
	ISTAND	INTEGER	0-1	0	1	1	MODE TYPE INDICATOR; 0-SPINNING MODE, 1-STANDING MODE
	IVEL	INTEGER	0-1	0	1	1	LINEAR/NON-LINEAR CONTROL; 0-ACOUSTIC VELOCITY EFFECTS NOT INCLUDED, 1-ACOUSTIC VELOCITY EFFECTS (NON-LINEAR) INCLUDED
	PAMPC	REAL	>=0	0.20	1	1	RATIO OF OSCILLATIONS AMPLITUDE TO PC, P/PC
	NRAD	INTEGER	1-20	3	1	1	NUMBER OF RADIAL INJECTION LOCATIONS
	NCIRC	INTEGER	1-50	5	1	1	NUMBER OF CIRCUMFERENTIAL INJECTION LOCATIONS
	PHIF	REAL	>=0	0.0	1	1	PRESSURE OSCILLATION PHASE ANGLE (DEG.)

# MODEL CONTROL VARIABLES (Continued)

<u>Namelist</u>	<u>Variable</u>	<u>Type</u>	<u>Range</u>	<u>Default</u>	<u>Path</u>	<u>Size</u>	<u>Description</u>
\$LEINJC	IDOMEM	INTEGER	1-2	2	1	1	DOME MODEL FLAG; 1->LUMPED PARAMETER, 2->ACOUSTIC MODE
	PAMPCH	REAL	>=0	0.20	1	1	CHAMBER PEAK-TO-PEAK PRESSURE AMPLITUDE RATIO, P'/PC
	NTINJ	INTEGER	0-50	18	1	1	NUMBER OF TIME STEPS PER CYCLE
	USGF	REAL	0-1	0.020	1	1	UPSTREAM SECTION GRID SIZE FACTOR (NUOR EQUALS XUOR/USGF/(VSOUND/FREQ))
	OGF	REAL	0-1	0.020	1	1	ORIFICE SECTION GRID SIZE FACTOR (NOR EQUALS XOR/OGF/(VSOUND/FREQ))
	DSGF	REAL	0-1	0.020	1	1	DOWNSTREAM SECTION GRID SIZE FACTOR (NDOR EQUALS XDOR/DSGF/(VSOUND /FREQ))
	FCDO	REAL	>0	1.0	1	4	FUEL ORIFICE CD ARRAY (4 ENTRIES - CORE, BAFFLE, BARRIER AND FFC)
	XCDO	REAL	>0	1.0	1	4	OX ORIFICE CD ARRAY (4 ENTRIES - CORE, BAFFLE, BARRIER AND FFC)
\$COMBUSTC	FALVM	REAL	>=0	1.0	1	4	ARRAY OF MULTIPLIERS FOR FUEL ATOMIZATION LENGTH USED IN VAPORIZATION CALCULATIONS (4 ENTRIES - CORE, BAFFLE, BARRIER AND FFC)
	XALVM	REAL	>=0	1.0	1	4	ARRAY OF MULTIPLIERS FOR OX ATOMIZATION LENGTH USED IN VAPORIZATION CALCULATIONS (4 ENTRIES - CORE, BAFFLE, BARRIER AND FFC)
	FALTM	REAL	>=0	1.0	1	4	ARRAY OF MULTIPLIERS FOR FUEL ATOMIZATION LENGTH USED IN TOTAL TIMELAG CALCULATIONS (4 ENTRIES - CORE, BAFFLE, BARRIER AND FFC)
	XALTM	REAL	>=0	1.0	1	4	ARRAY OF MULTIPLIERS FOR OX ATOMIZATION LENGTH USED IN TOTAL TIMELAG CALCULATIONS (4 ENTRIES - CORE, BAFFLE, BARRIER AND FFC)
	FRMM	REAL	>=0	1.0	1	4	ARRAY OF MULTIPLIERS FOR FUEL DROPSIZES (4 ENTRIES - CORE, BAFFLE, BARRIER AND FFC)

C-18

<u>Namelist</u>	<u>Variable</u>	<u>Type</u>	<u>Range</u>	<u>MODEL CONTROL VARIABLES (Continued)</u>			<u>Description</u>
				<u>Default</u>	<u>Path</u>	<u>Size</u>	
	XRMM	REAL	>0	1.0	1	4	ARRAY OF MULTIPLIERS FOR OX DROPSIZES (4 ENTRIES - CORE, BAFLE, BARRIER AND FFC)
	EMMUL	REAL	>0	1.0	1	2	ARRAY OF MULTIPLIERS ON CORE (1) AND BARRIER (2) OVERALL MIXING EFFICIENCIES (EM)
	AOMUL	REAL	>0	1.0	1	1	MULTIPLIER ON CALCULATED VALUE OF MEAN CHAMBER SOUND SPEED (AO)
	ENMUL	REAL	>0	1.0	1	1	MULTIPLIER ON CALCULATED VALUE OF PRESSURE INTERACTION INDEX (EN)
	TAUMUL	REAL	>0	1.0	1	1	MULTIPLIER ON CALCULATED VALUES OF SENSITIVE TIME LAG (TAUSEN)
	CCMUL	REAL	>0	1.0	1	1	MULTIPLIER ON CALCULATED VALUES OF CAVITY SOUND SPEEDS (CC1, CC2)
	COMBXB	REAL	0-1.0	0.50	1	1	FRACTION OF C* EFFICIENCY AT WHICH THE COMBUSTION PLANE (XB) IS PLACED.
\$FDORCC	SHORT	LOGICAL		T/F	F	1	0 TRUE IF SHORT NOZZLE IS ASSUMED, FALSE IF REAL NOZZLE IS USED
	EPSIL	REAL	>0	20	1	1	PEAK-TO-PEAK PRESSURE AMPLITUDE (PERCENT)
	ERROR	REAL	>0	0.1	1	1	MAXIMUM % ERROR BETWEEN SUCCESSIVE ITERATIONS
	LTS	INT	1-10	5	1	1	# OF RADIAL SERIES TERMS
	MTS	INT	1-9	5	1	1	# OF TANGENTIAL SERIES TERMS
	NTS	INT	1-20	10	1	1	# OF LONGITUDINAL SERIES TERMS
	ITMAX	INT	>0	100	1	1	MAXIMUM # OF SUCCESSIVE ITERATIONS
	RELX	REAL	0<RELX<1.0	0.65	1	1	SOLUTION RELAXATION FACTOR
	MTYPE	INT	0-1	0	1	1	TANGENTIAL WAVE TYPE, 0-STANDING, 1- TRAVELING
	NEET	INT	1-10	3	1	1	# OF EIGENFUNCTION EXPANSION TERMS
	NXFST	INT	1-10	3	1	1	# OF FOURIER SERIES TERMS IN THE CAVITY X- DIRECTION
	NYFST	INT	1-10	3	1	1	# OF FOURIER SERIES TERMS IN THE CAVITY Y- DIRECTION
	NAFST	INT	1-10	3	1	1	# OF FOURIER SERIES TERMS IN THE CAVITY APERTURE
	NHAT	INT	>=0	0	1	1	LONGITUDINAL MODE NUMBER
	MORE	LOGICAL		T/F	F	1	0 TRUE FOR ADDITIONAL FDORC OUTPUT TO HISTORY FILE

MORE LOGICAL T/F F 1 0 TRUE FOR ADDITIONAL FDORC OUTPUT TO HISTORY FILE

# DEFAULT DESIGN CONTROL VARIABLES

<u>Namelist</u>	<u>Variable</u>	<u>Type</u>	<u>Range</u>	<u>Default</u>	<u>Path</u>	<u>Size</u>	<u>Description</u>
\$CONTROL	DPPCS	REAL	<0-1	0.15	1	1	RATIO OF INJECTION PRESSURE DROP TO PC AT THROTTLED PC FOR LOW FREQ STABILITY
	RNR	REAL	>0	1.0	1	1	NONDIMENSIONAL NOZZLE ENTRANCE RADIUS OF CURVATURE (RNE/RTHRT)
	RTR	REAL	>0	1.0	1	1	NONDIMENSIONAL THROAT ENTRANCE RADIUS OF CURVATURE (RTE/RTHRT)
	CONVA	REAL	0-90	25.0	1	1	THROAT CONVERGENCE HALF-ANGLE (DEG)
	XMSTAB	REAL	>0	0.20	5	1	MINIMUM MASS FRACTION FOR ELEMENT TO EFFECT STABILITY
\$LOLC	FCD	REAL	<1	0.78	1	1	FUEL DISCHARGE COEFFICIENT
	XCD	REAL	<1	0.78	1	1	OX DISCHARGE COEFFICIENT
	FIA	REAL	<45	30.0	1	1	FUEL IMPINGEMENT ANGLE (DEG)
	XIA	REAL	<45	30.0	1	1	OX IMPINGEMENT ANGLE (DEG)
	FIHOD	REAL	>0	2.0	1	1	FUEL IMPINGEMENT HEIGHT TO ORIFICE DIAMETER RATIO
	XIHOD	REAL	>0	2.0	1	1	OX IMPINGEMENT HEIGHT TO ORIFICE DIAMETER RATIO
	FCANT	REAL	0-45	16.0	1	1	FUEL CANT ANGLE (DEG)
	XCANT	REAL	0-45	16.0	1	1	OX CANT ANGLE (DEG)
	FLOD	REAL	>0	5.0	1	1	FUEL ORIFICE LENGTH TO DIAMETER RATIO
	XL0D	REAL	>0	5.0	1	1	OX ORIFICE LENGTH TO DIAMETER RATIO
	EMUNI	REAL	<1	0.65	1	1	UNIELEMENT RUPE MIXING EFFICIENCY

# DEFAULT DESIGN CONTROL VARIABLES (Continued)

<u>Namelist</u>	<u>Variable</u>	<u>Type</u>	<u>Range</u>	<u>Default</u>	<u>Path</u>	<u>Size</u>	<u>Description</u>
\$TRIPC	FCD	REAL	<1	0.78	1	1	FUEL DISCHARGE COEFFICIENT
	XCD	REAL	<1	0.78	1	1	OX DISCHARGE COEFFICIENT
	ULIA	REAL	<45	35.0	1	1	UNLIKE IMPINGEMENT HALF-ANGLE (DEG)
	UILOD	REAL	>0	2.0	1	1	TRIPLET IMPINGEMENT HEIGHT TO ORIFICE DIAMETER RATIO; OX L/D FOR OFO'S AND FUEL L/D FOR FOF
	FLOD	REAL	>0	5.0	1	1	FUEL ORIFICE LENGTH TO DIAMETER RATIO
	XL0D	REAL	>0	5.0	1	1	OX ORIFICE LENGTH TO DIAMETER RATIO
	DODF	REAL	>0	1.0	1	1	TRIPLET ORIFICE OX DIAMETER TO FUEL DIAMETER RATIO
	EMUNI	REAL	<1	0.77	1	1	UNELEMENT RUPE MIXING EFFICIENCY
\$SHDC	FCD	REAL	<1	0.78	1	1	FUEL DISCHARGE COEFFICIENT
	XCD	REAL	<1	0.78	1	1	OX DISCHARGE COEFFICIENT
	FLOD	REAL	>0	5.0	1	1	FUEL ORIFICE LENGTH TO DIAMETER RATIO
	XL0D	REAL	>0	5.0	1	1	OX ORIFICE LENGTH TO DIAMETER RATIO
C-22	EMUNI	REAL	<1	0.50	1	1	UNELEMENT RUPE MIXING EFFICIENCY
\$SHEARC	FCD	REAL	<1	0.93	1	1	FUEL DISCHARGE COEFFICIENT
	VRATMI	REAL	>0	10.0	1	1	MINIMUM GAS TO LIQUID VELOCITY RATIO
	GAPM	REAL	>0	0.010	1	1	MINIMUM FUEL ANNULUS GAP WIDTH
	TPOST	REAL	>0	0.020	1	1	OX POST TIP THICKNESS (IN)
	FLOD	REAL	>0	5.0	1	1	FUEL ANNULUS LENGTH TO ANNULAR GAP RATIO
	DIVANG	REAL	0-7	0.0	1	1	OX POST EXIT DIVERGENCE ANGLE (DEG)
	EMUNI	REAL	<1	0.80	5	1	UNELEMENT RUPE MIXING EFFICIENCY

DEFAULT DESIGN CONTROL VARIABLES (Continued)

<u>Namelist</u>	<u>Variable</u>	<u>Type</u>	<u>Range</u>	<u>Default</u>	<u>Path</u>	<u>Size</u>	<u>Description</u>
\$SWIRLC	FCD	REAL	<1	0.93	1	1	FUEL DISCHARGE COEFFICIENT
	VRATMI	REAL	>0	10.0	1	1	MINIMUM GAS TO LIQUID VELOCITY RATIO
	GAPM	REAL	>0	0.010	1	1	MINIMUM FUEL ANNULUS GAP WIDTH (IN)
	TPOST	REAL	>0	0.020	1	1	OX POST TIP THICKNESS (IN)
	FLOD	REAL	>0	5.0	1	1	FUEL ANNULUS LENGTH TO ANNULAR GAP RATIO
	XCA	REAL	<45	24.0	1	1	SWIRL CONE HALF-ANGLE (DEG)
	NINLET	INTEGER	>0	0	1	1	THIS VARIABLE HAS BEEN REMOVED
	DRATIO	REAL	>0	1.20	1	1	SWIRL CHAMBER TO OX POST EXIT DIAMETER RATIO
	CDINLET	REAL	<1	0.90	1	1	SWIRL CHAMBER INLET ORIFICE CD
	EMUNI	REAL	<1	0.80	1	1	UNELEMENT RUPE MIXING EFFICIENCY
\$SAID	TBAF	REAL	>0	0.033	1	1	BAFFLE THICKNESS (FT)
	TPART	REAL	>=0	0.0	1	1	CAVITY PARTITION THICKNESS (FT)
	INLET	INTEGER	0-2	0	1	1	1/4 WAVE/HELMHOLTZ CAVITY INLET DESCRIPTOR; 0 - SHARP EDGE, 1 - ROUNDED, 2 - WELL ROUNDED



**APPENDIX D**  
**CREATING COMBUSTION GAS TABLES**

## Creating Combustion Gas Tables

Some of the analysis and design codes of the ROCCID package require combustion gas properties. To keep the user from entering these properties manually, the program interpolates the data, from tables of gas properties versus mixture ratio, temperature and pressure for a given set of propellants. Tables of equilibrium gas temperature as a function of mixture ratio and pressure are also required. The code already provides tables for the following propellant combinations:

LOX/HYDROGEN  
LOX/METHANE  
LOX/PROPANE  
LOX/RP-1

It is likely at some point, that propellant combinations other than those listed above will be needed, or tables with better resolution for the above propellants will be required. For these reasons the ROCCID program has the ability to automatically create new tables using an ODE (one dimensional equilibrium) module.

To make new tables using the ODE module, first enter the utilities menu off the ROCCID base menu. Next enter option 1, the ODE module. The program will now prompt you to enter the mixture ratio, pressure, and temperature arrays and the number of mixture ratio entries. Enter these data using standard IFE syntax. The array sizes are limited to 30, 6, and 10 for mixture ratio, pressure and temperature respectively. Selecting temperature below 600 R is not recommended since ODE may have trouble calculating the gas properties below this temperature. Next the module will ask you if you want to select reactants. Answer Y for yes. The program will then list its propellant library and ask you to pick propellants. For each propellant chosen, the program will display the ODE propellant card. You then have the option of changing the reference temperature, enthalpy, and percent composition of that propellant. Finally before the ODE run begins, you are prompted for omitting species. Simply press RETURN and the ODE will begin.

ODE first uses the TP option to create full tables of viscosity, molecular weight, Prandtl number, thermal conductivity, specific heat ratio, enthalpy, and entropy versus mixture ratio, pressure and temperature. Then ODE uses a very similar run deck using the HP option to create a table of equilibrium temperatures versus pressure and mixture ratio. The input decks are saved for the above to run in files COMTBL.INP and COMTB1.INP respectively.

After ODE has completed the tables, the program asks the user to input a fuel name and an oxidizer name (up to 8 characters each). The module will then create a file based on the propellant names as follows:

**<oxidizer name>\_<fuel name>.DAT**

For example, if the oxidizer is LOX and the fuel is RP-1 the filename created will be LOX\_RP-1.DAT.

In order for the ROCCID program to use the generated file, the user must manually enter the filename and a one line propellant descriptor into PROP.FIL, contained in the propellant property library directory. It is important that the filename is appended first, then the one line descriptor. Correct additions to this file will automatically add the new propellant options to the Propellant Menu in both the Point Analysis and Point Design sections of the ROCCID code.

**APPENDIX E**  
**FILES NAMING CONVENTIONS**

## Files Naming Conventions

<u>FILE TYPE</u>	<u>UNIT #</u>	<u>DESCRIPTION</u>
*.DAT	3	Replay files
*.INP	7	Point analysis input
*.HIS	8	History/submodel output
*.DES	9	Point design input
*.OUT	10	Summarized output
*.CNT	12	Analysis model control parameters
*.DEF	13	Point design control parameters
*.DBG	14	Debug output
*.TDK	15	TDK input data
PROPELLANT <sup>(1)</sup> .DAT	20	Combustion gas data
*.PL1	21	Steady state performance plot data
*.PL2	22	Chug stability plot data
*.PL3	23	High frequency stability plot data
*.COF	25	MIPROPS propellant data

(1) Propellant combinaions, i.e., LOX\_H2.DAT OR LOX\_METHANE.DAT, See Appendix D

**APPENDIX F**  
**ROCCID PROGRAM FLOW CHARTS**

**Part A**  
**Point Analysis Module Flow Charts**

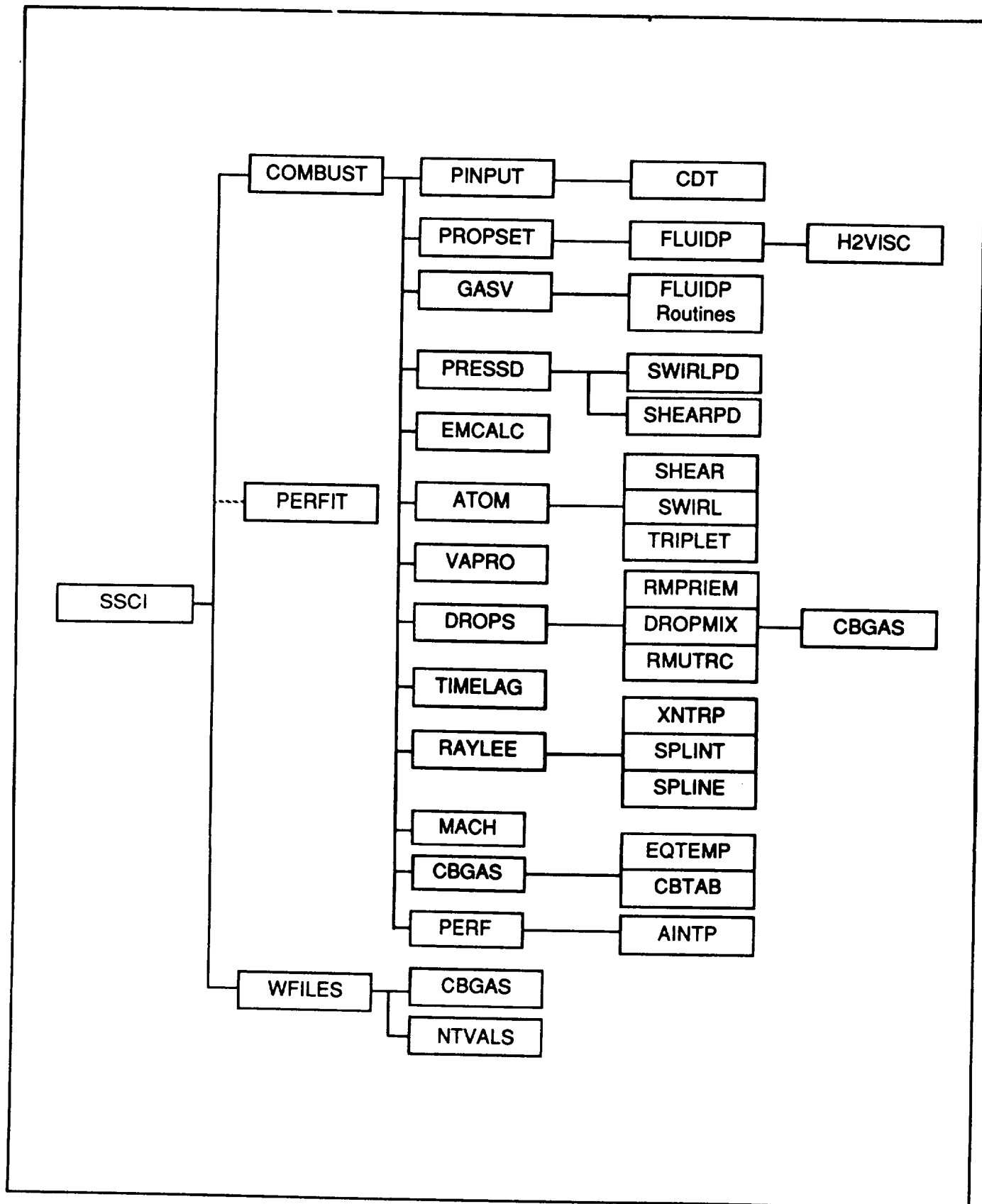
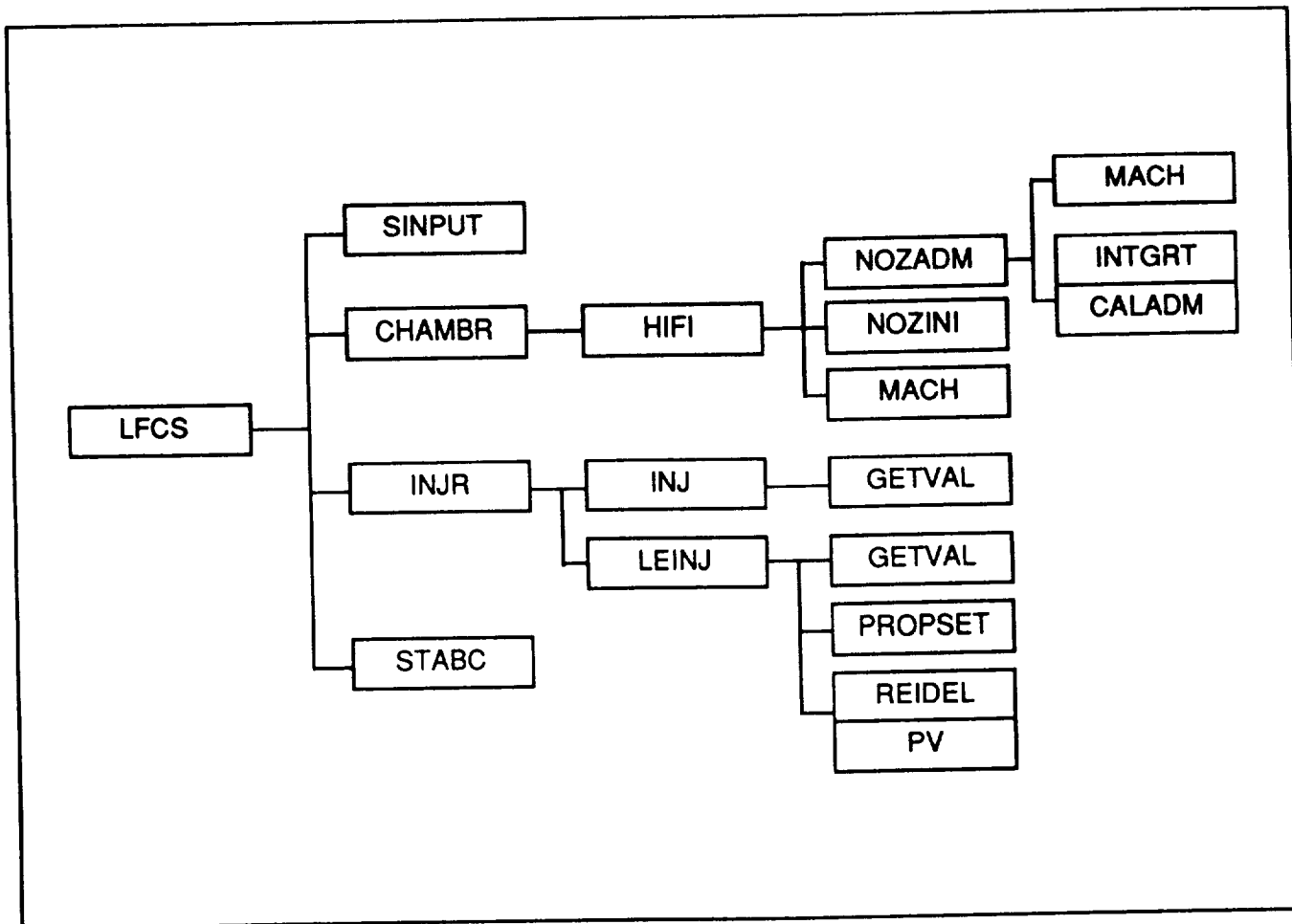
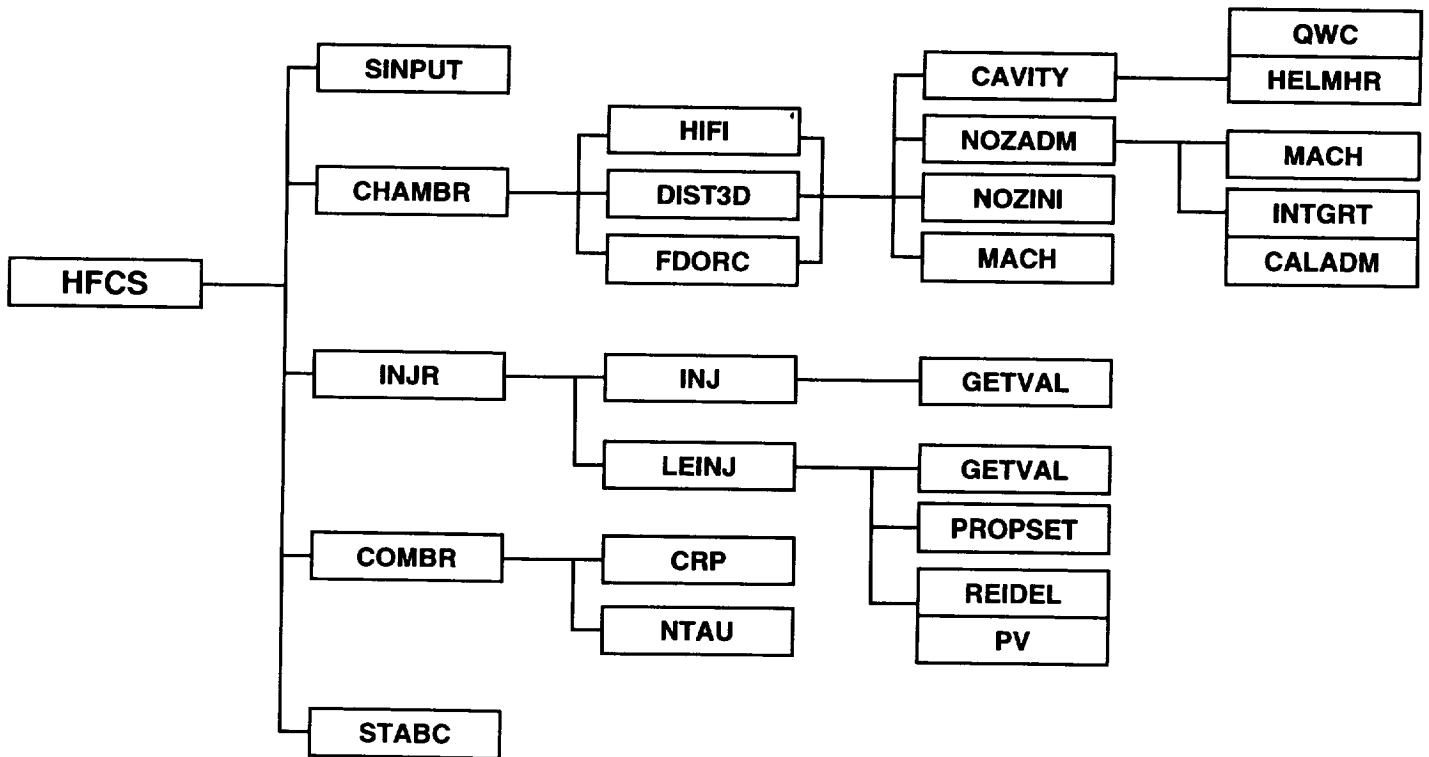


Figure F-1. Flowchart of Module SSCI

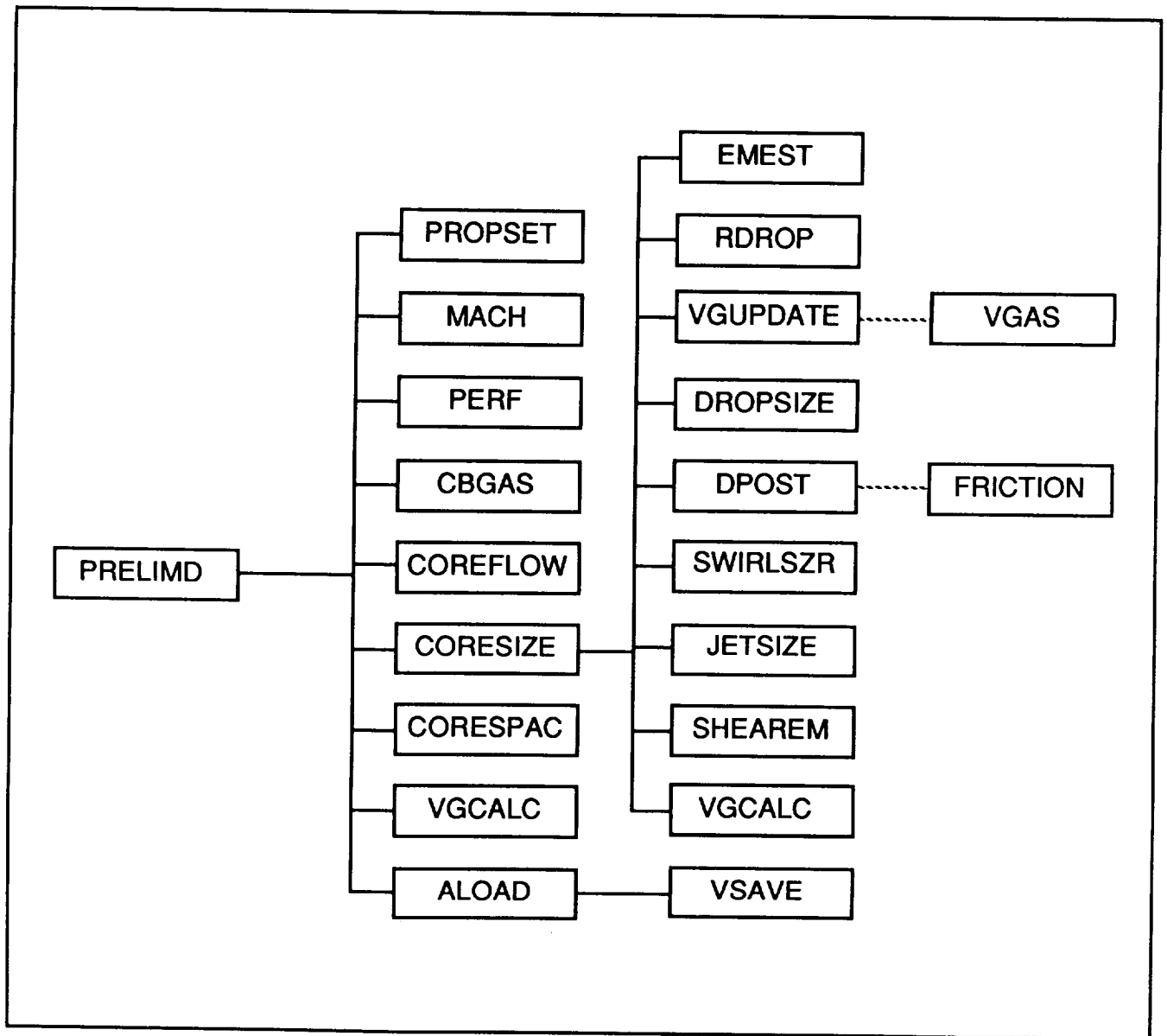


**Figure F-2. Flowchart of Module LFCS**

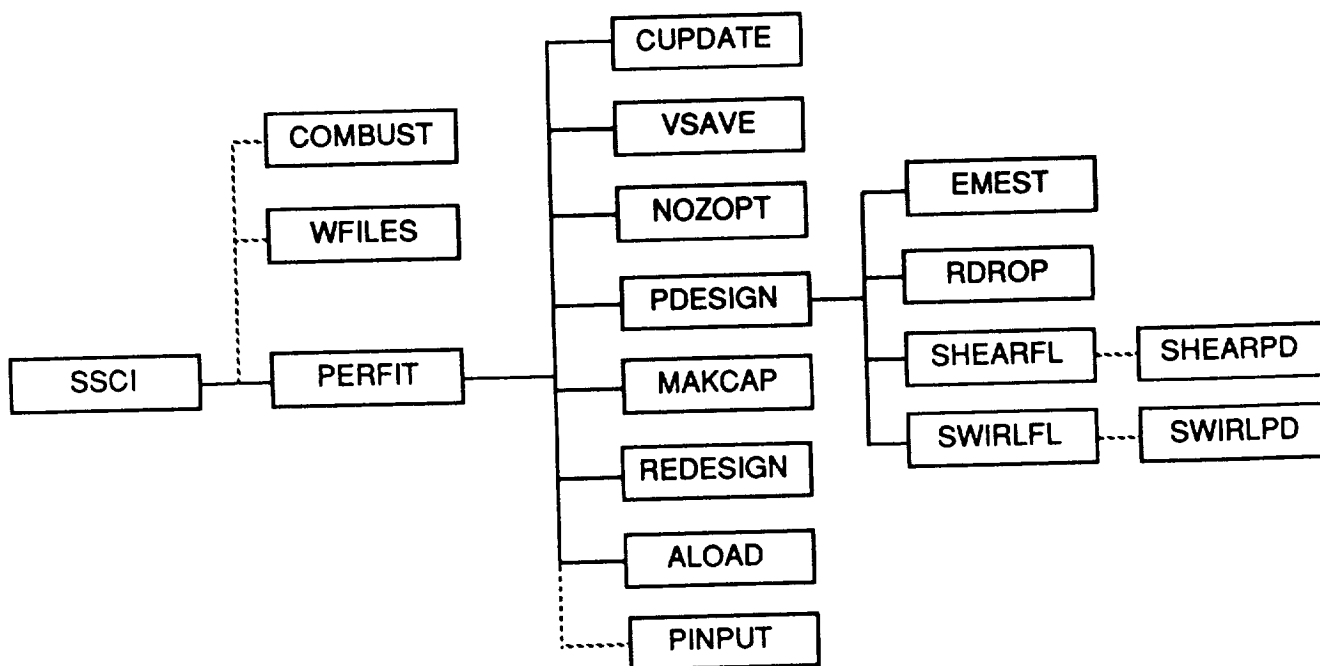


**Figure F-3. Flowchart of Module HFCS**

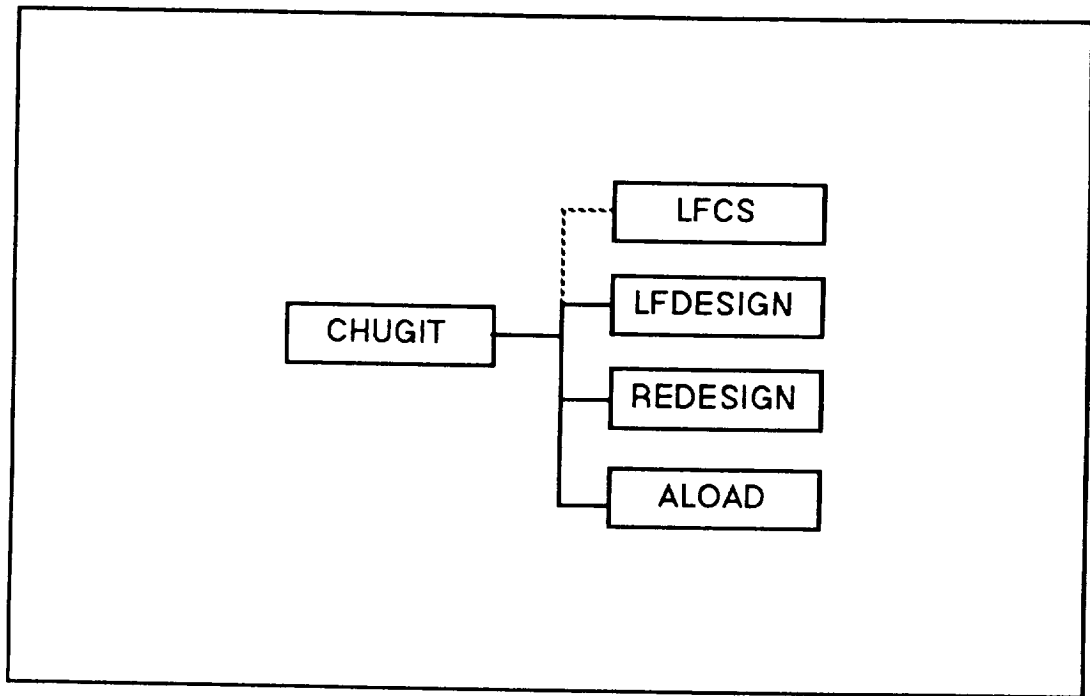
**Part B**  
**Point Design Module Flow Charts**



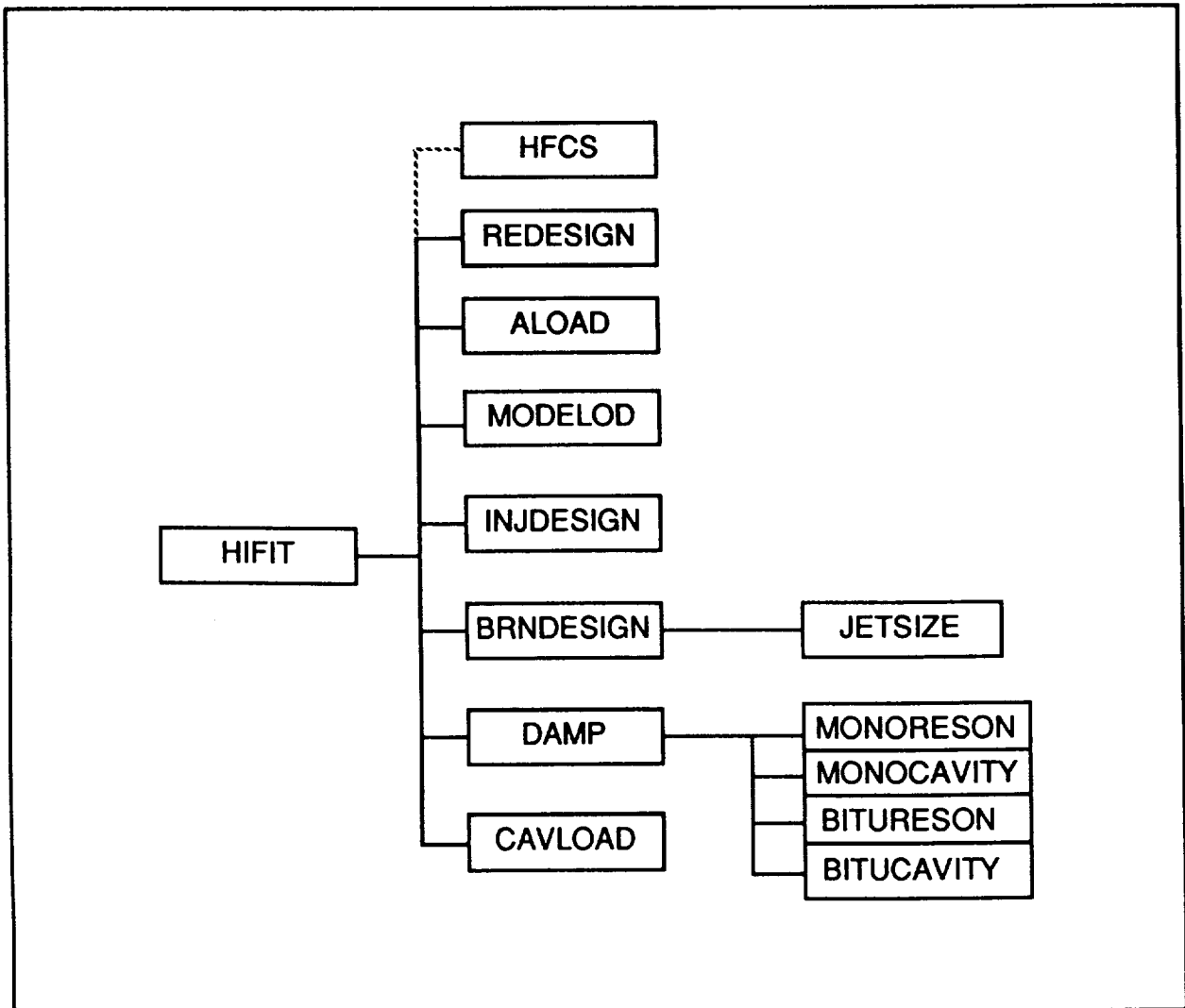
**Figure F-4. Flowchart for Module PRELIMD**



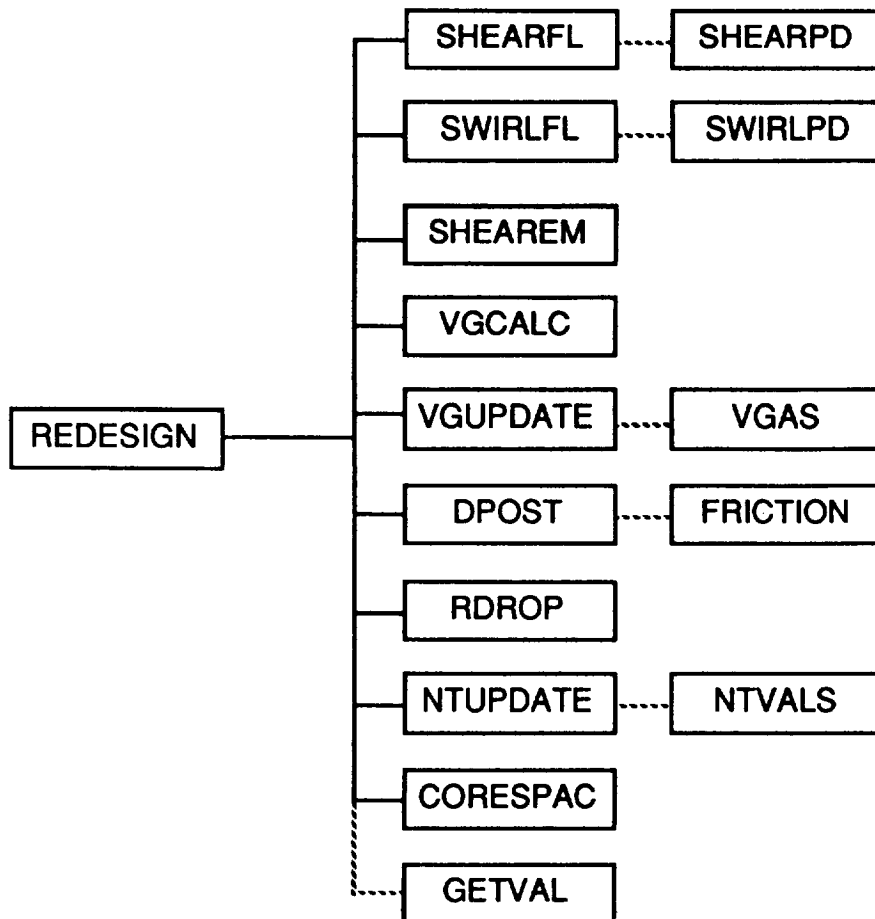
**Figure F-5. Flowchart for Module PERFIT**



**Figure F-6. Flowchart for Module CHUGIT**



**Figure F-7. Flowchart for Module HIFIT**



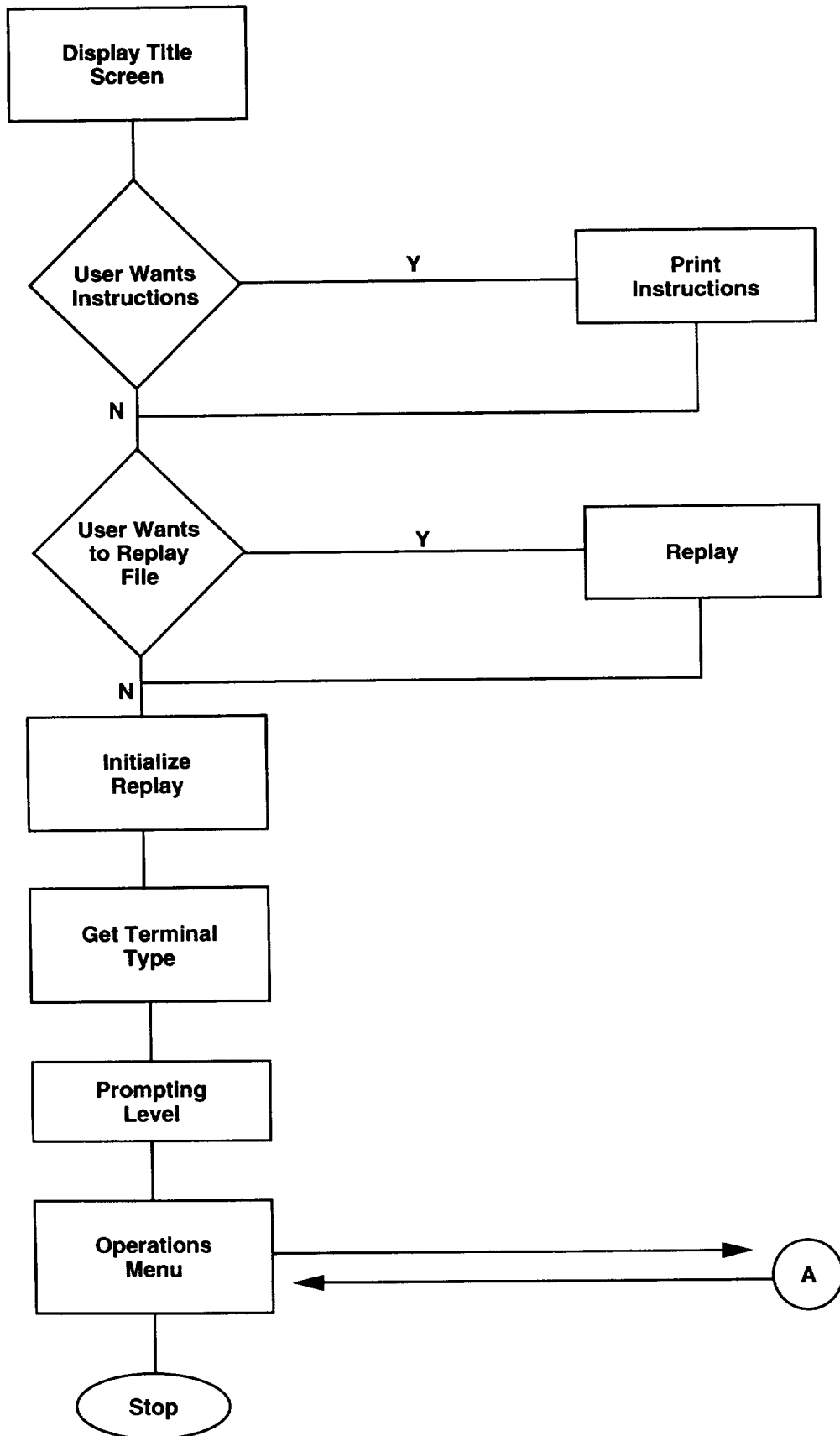
**Figure F-8. Flowchart for Module REDESIGN**

**Part C**  
**Main IFE Flow Charts**

### ANALYSIS REQUEST MENU (BMENU)

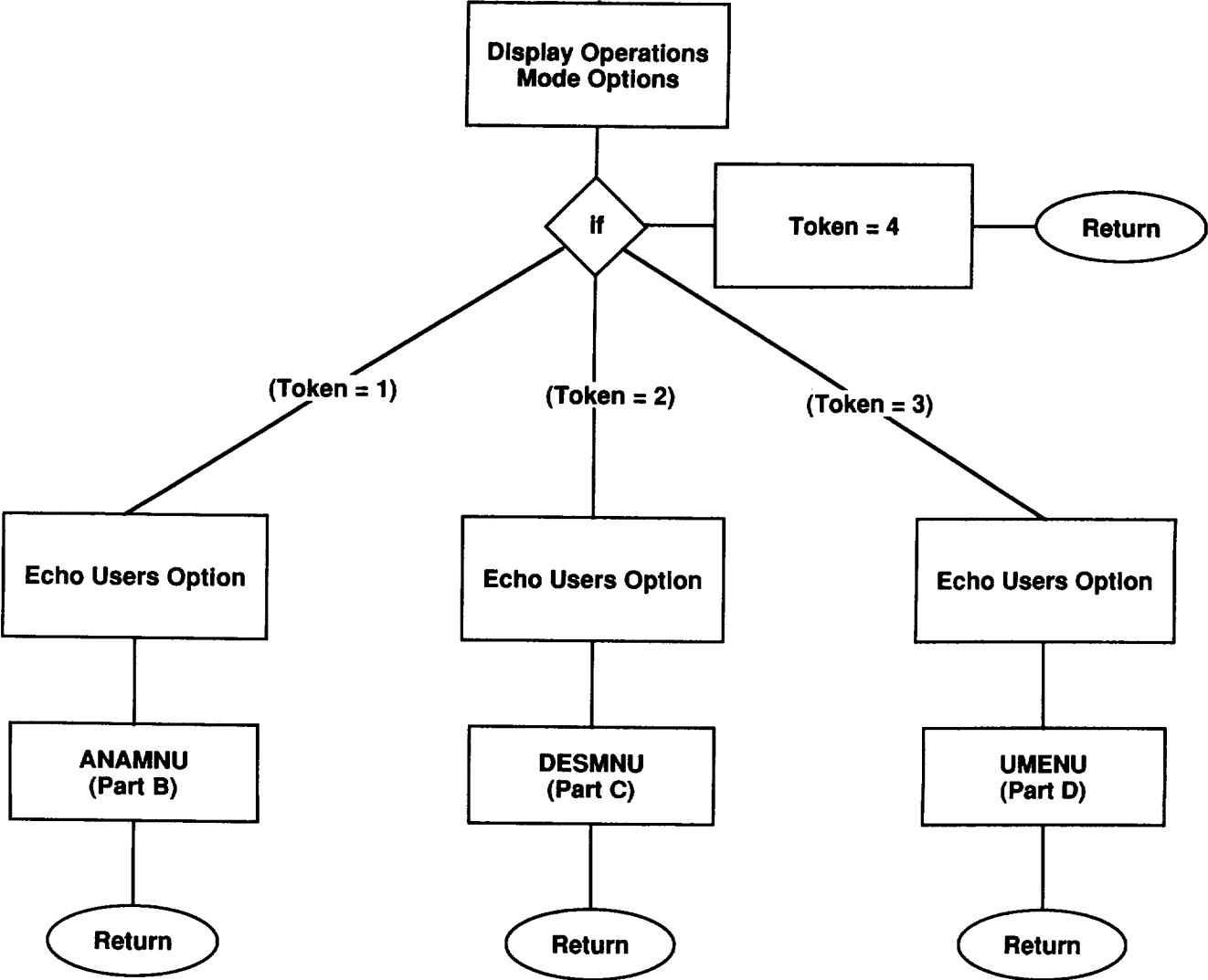
1. Point Analysis
2. Point Design
3. Utility Routines
4. Stop

# ROCCID



# B MENU

A



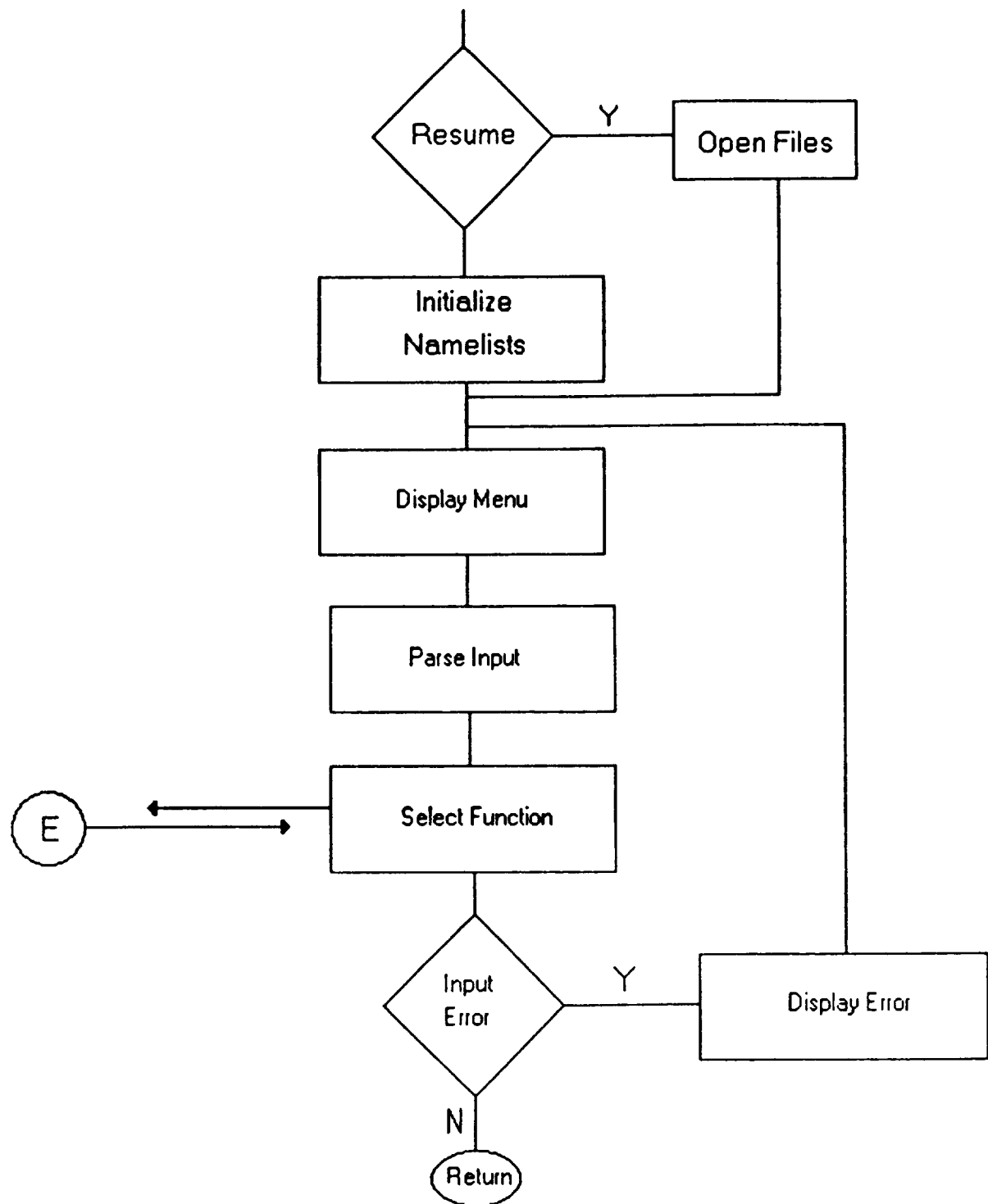
**Part D**  
**IFE Point Analysis Section Flow Charts**

## POINT ANALYSIS MENU OPTIONS (ANAMNU)

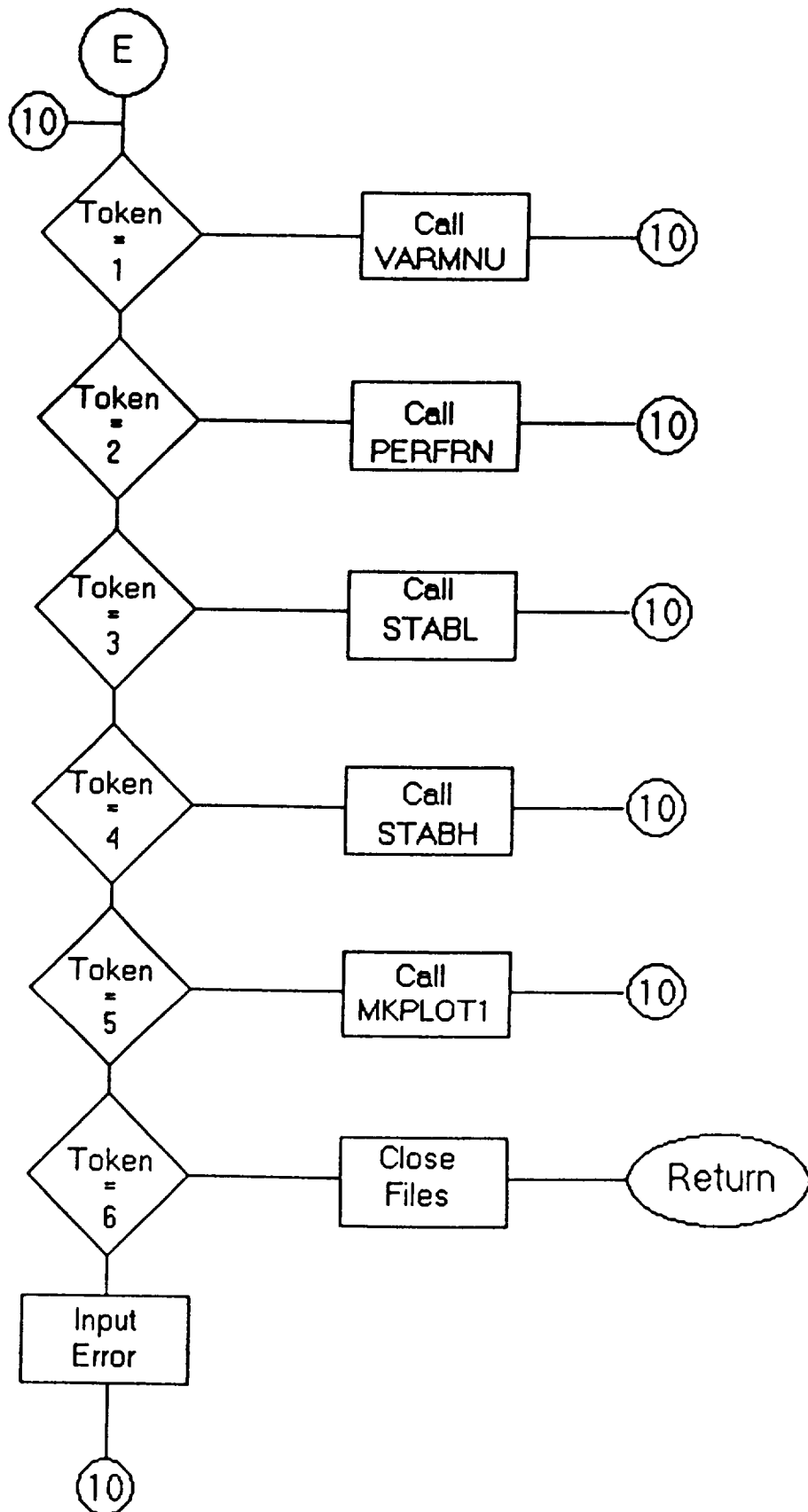
1. Set Variables
2. Run Steady State Performance
3. Run Low Frequency Stability
4. Run High Frequency Stability
5. Plot Current Data
6. Previous Menu

# ANALYSIS MENU

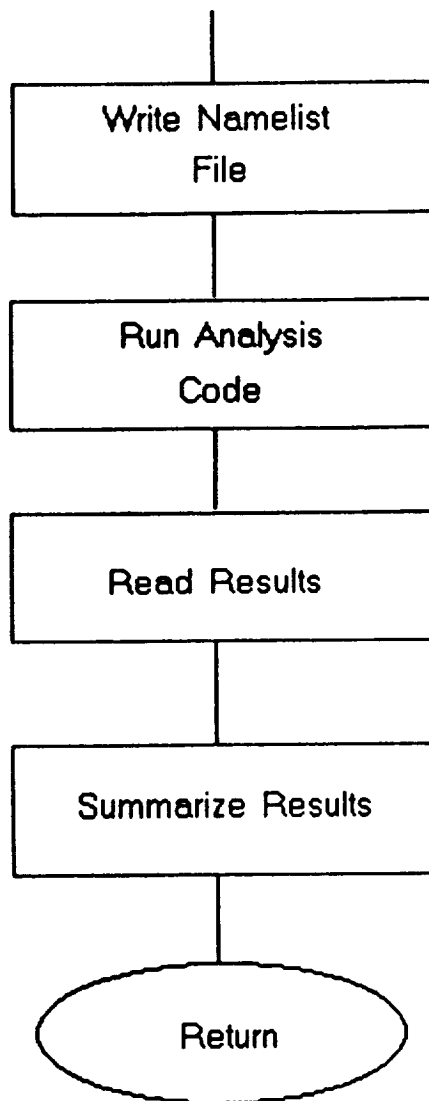
## (ANAMNU)



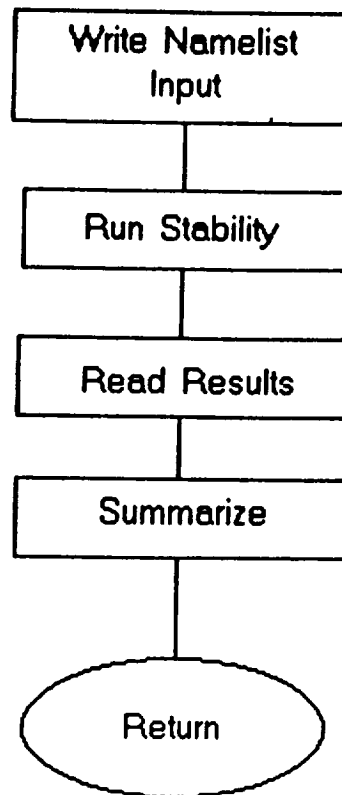
# ANAMNU.



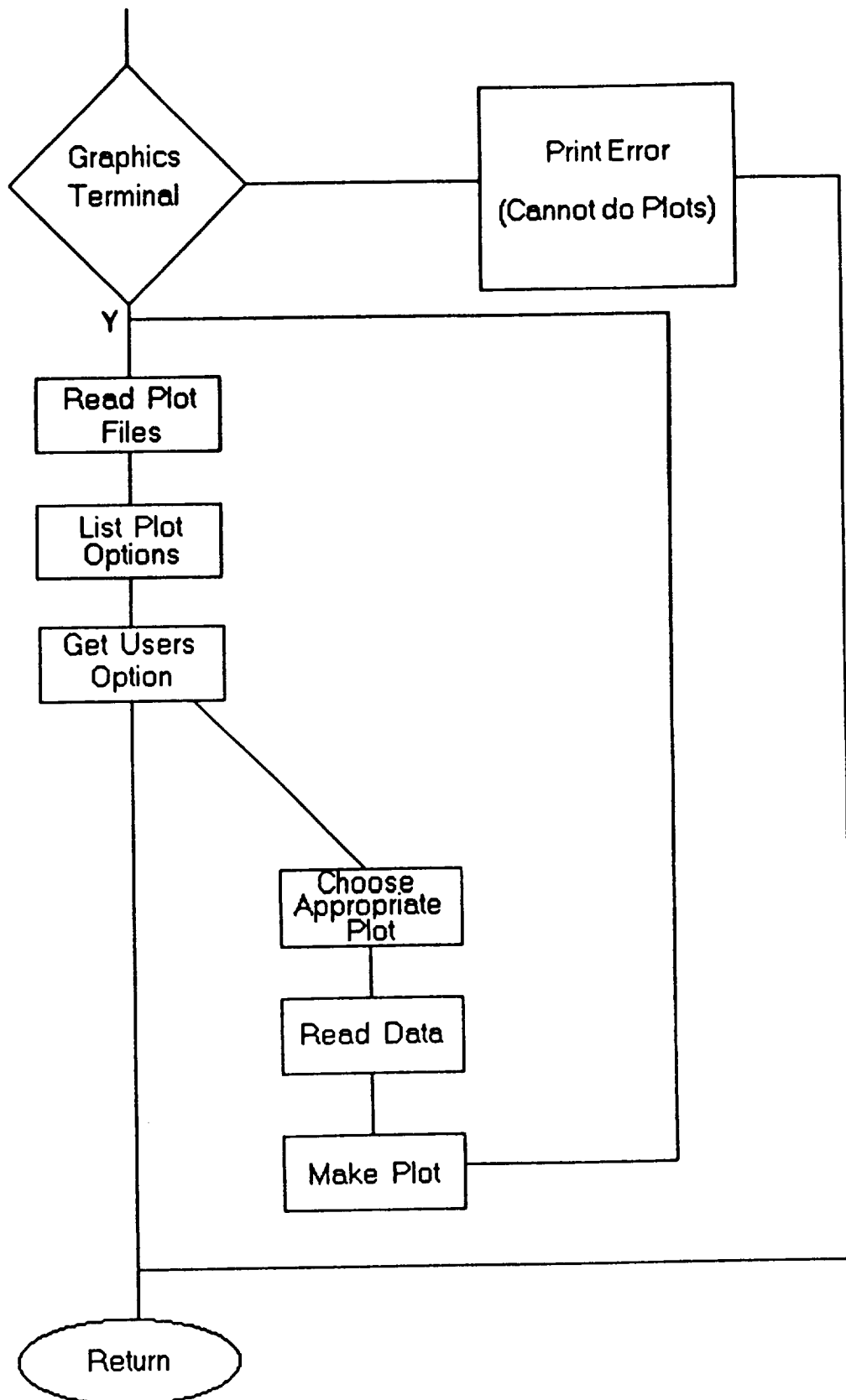
# SUBROUTINE PERFRN



# SUBROUTINE STAB



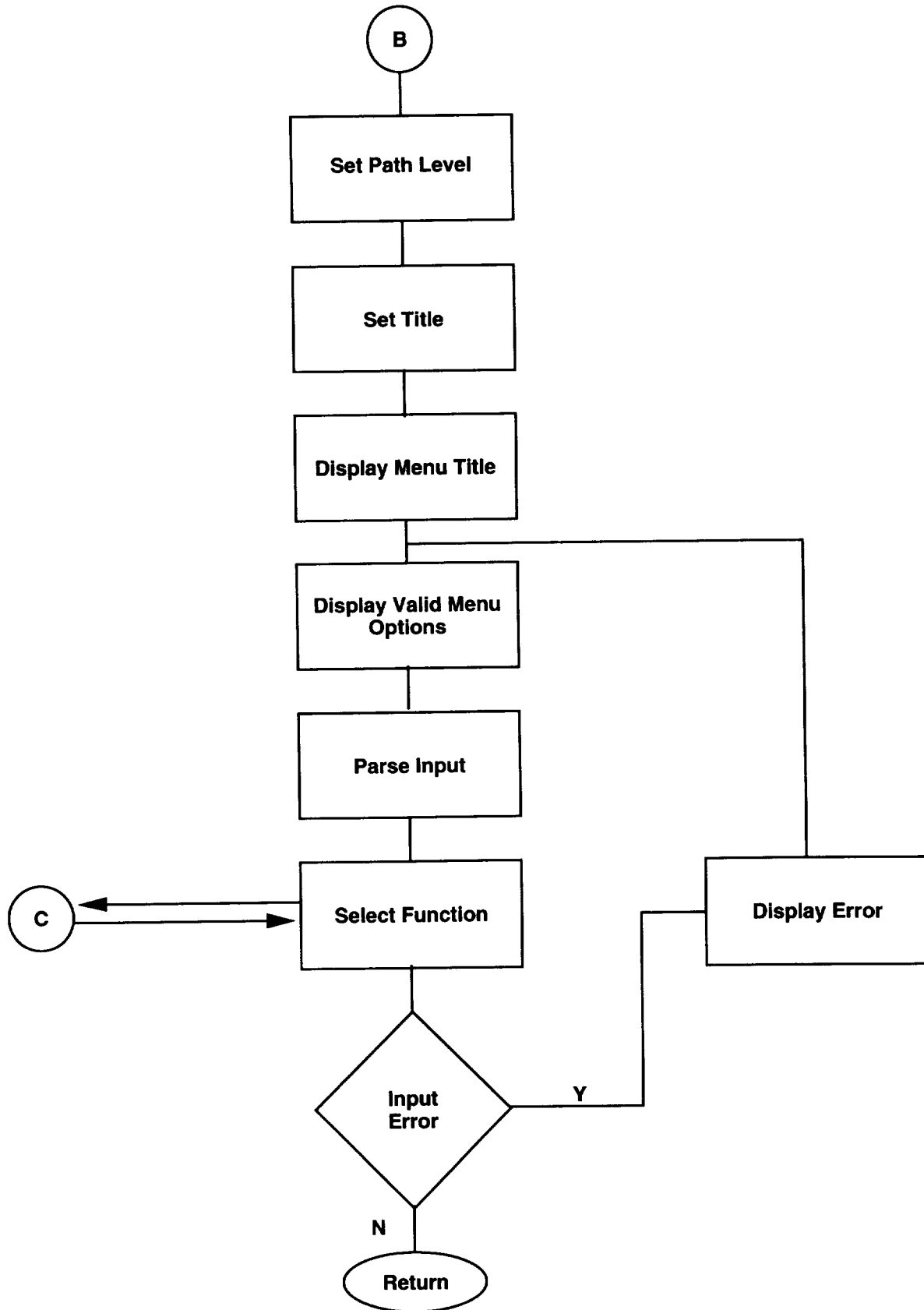
# SUBROUTINE DISPLAY



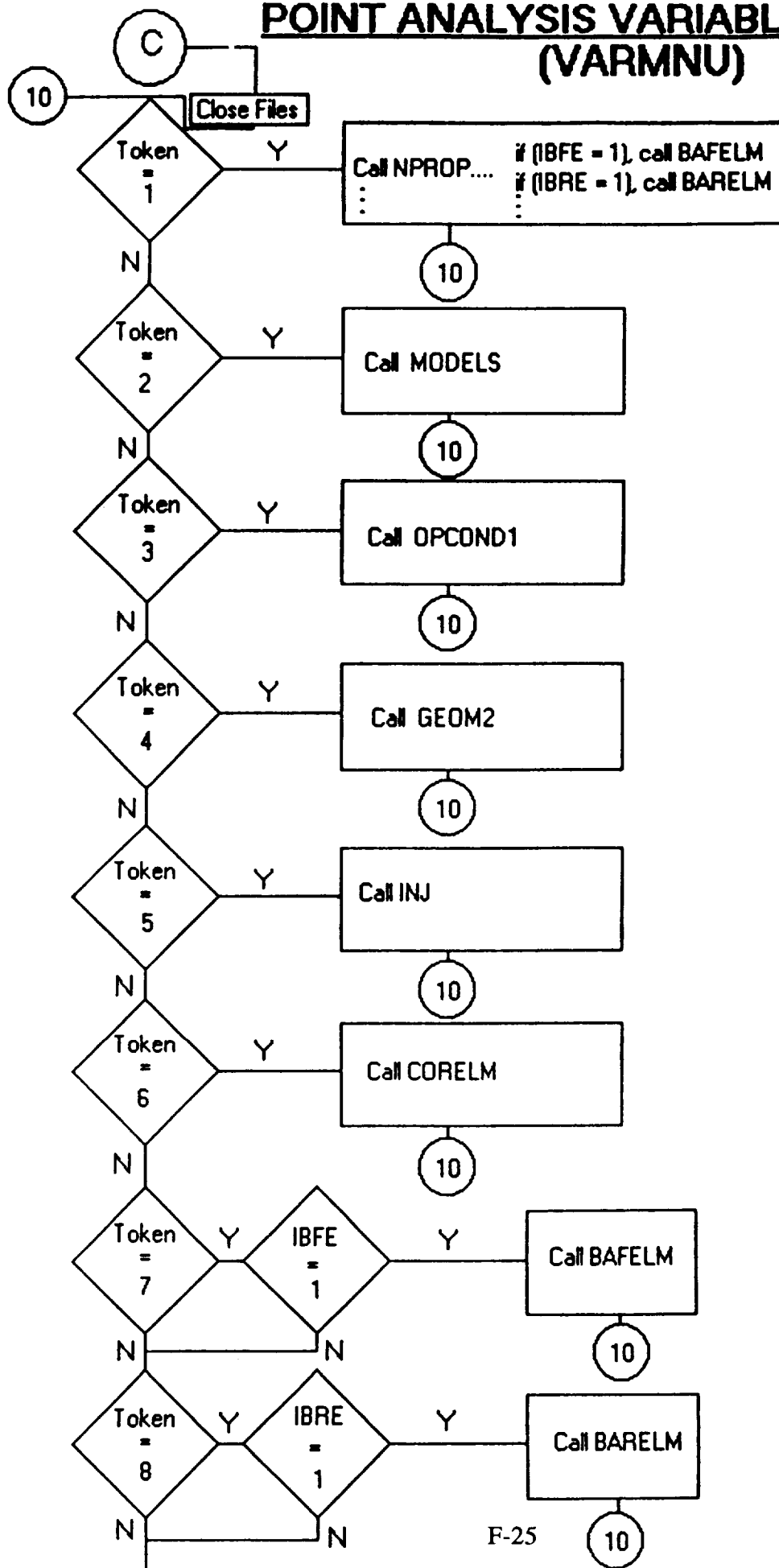
## POINT ANALYSIS VARIABLES MENU OPTIONS

1. Complete Setup
2. Set Model Selection
3. Set Operating Conditions
4. Set Geometry
5. Set Injector Element Types
6. Set Core Element
7. Set Baffle Element (if IBFE =1)
8. Set Barrier Element (if IBRE =1)
9. Set Fuel Film/Cavity Cooling Element (if IFFE =1)
10. Set Stability Aid Type (if MCHAM=1 or 2)
11. Set Manifold Description
12. Set Baffle Configuration (if IBAF =1)
13. Set 1/4 Wave Cavity Configuration (if ICAV =1 and MCHAM=1 or 2)
14. Set Helmholtz Resonator Configuration (if ICAV =2 and MCHAM=1 pr 2)
15. Set FDORC Variables (if MCHAM=3)
16. Set Model Control Variables
17. Previous Menu

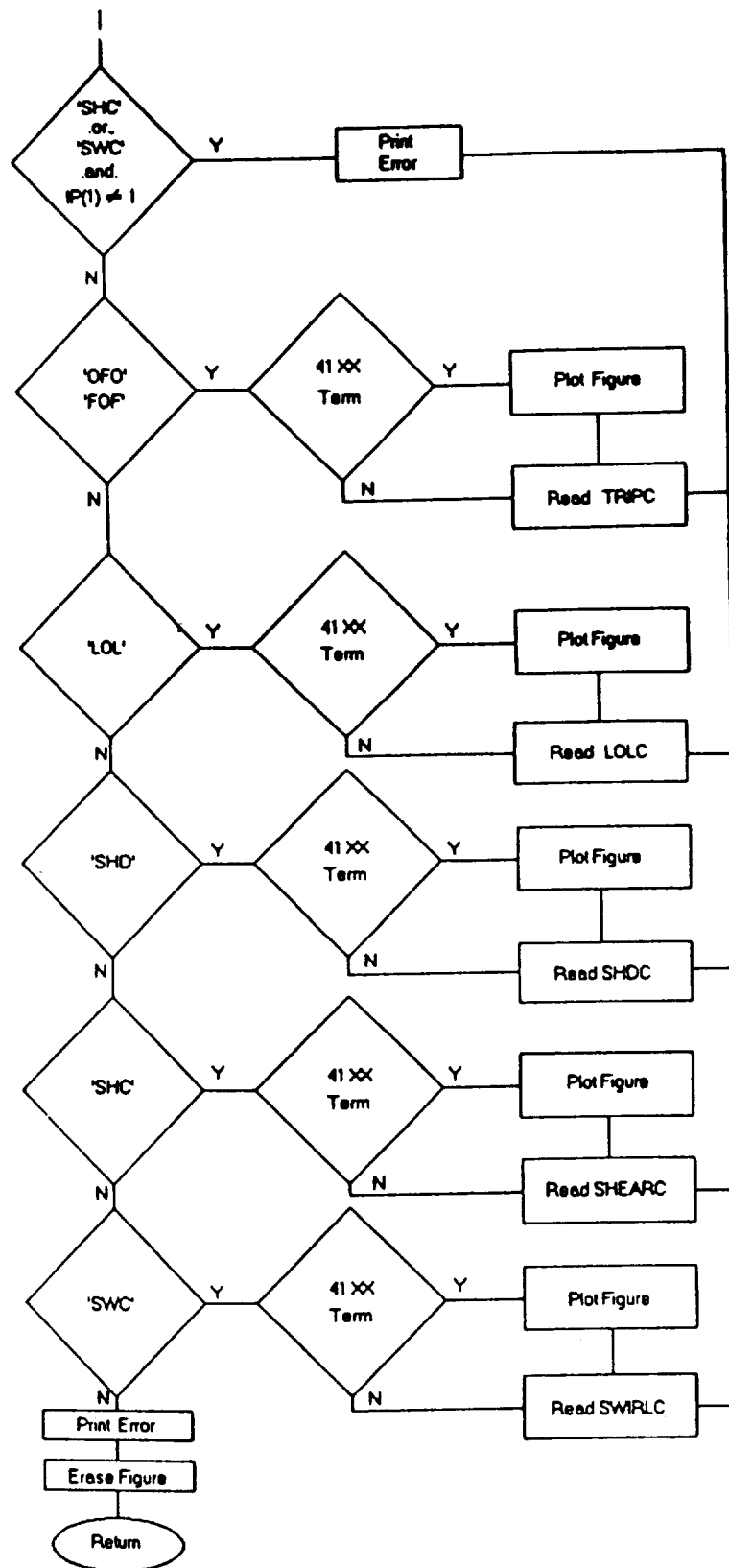
# VARMNU



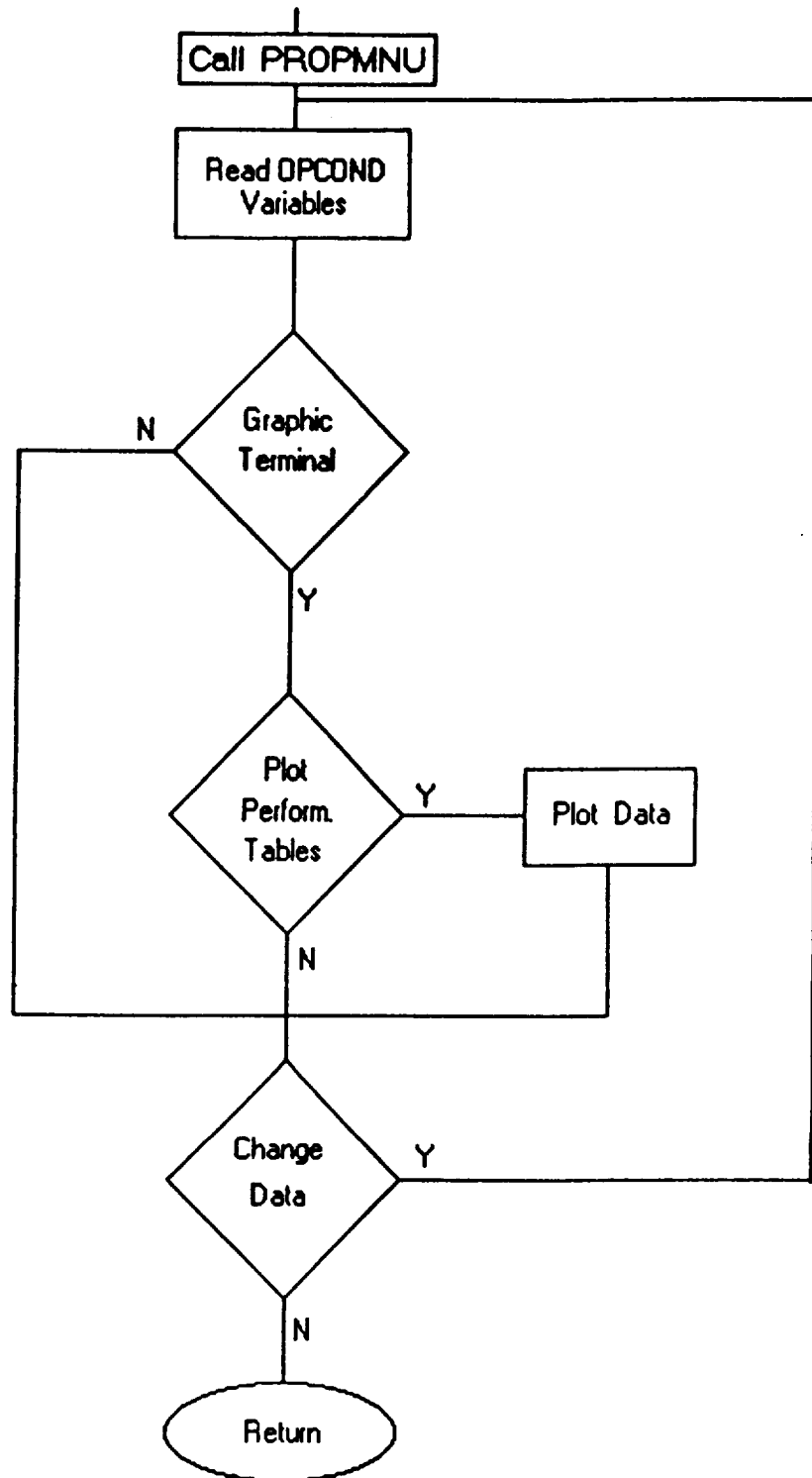
# POINT ANALYSIS VARIABLE FUNCTIONS (VARMNU)



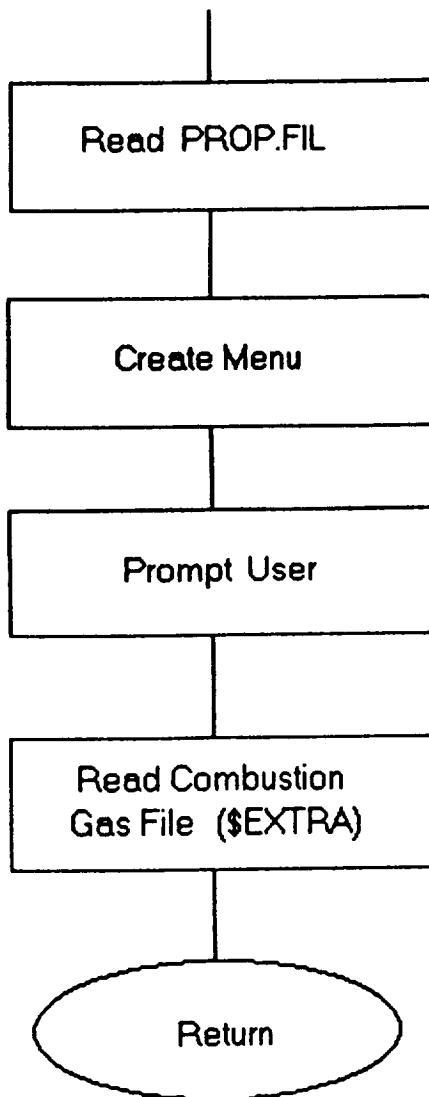
# SUBROUTINE ELMDEF



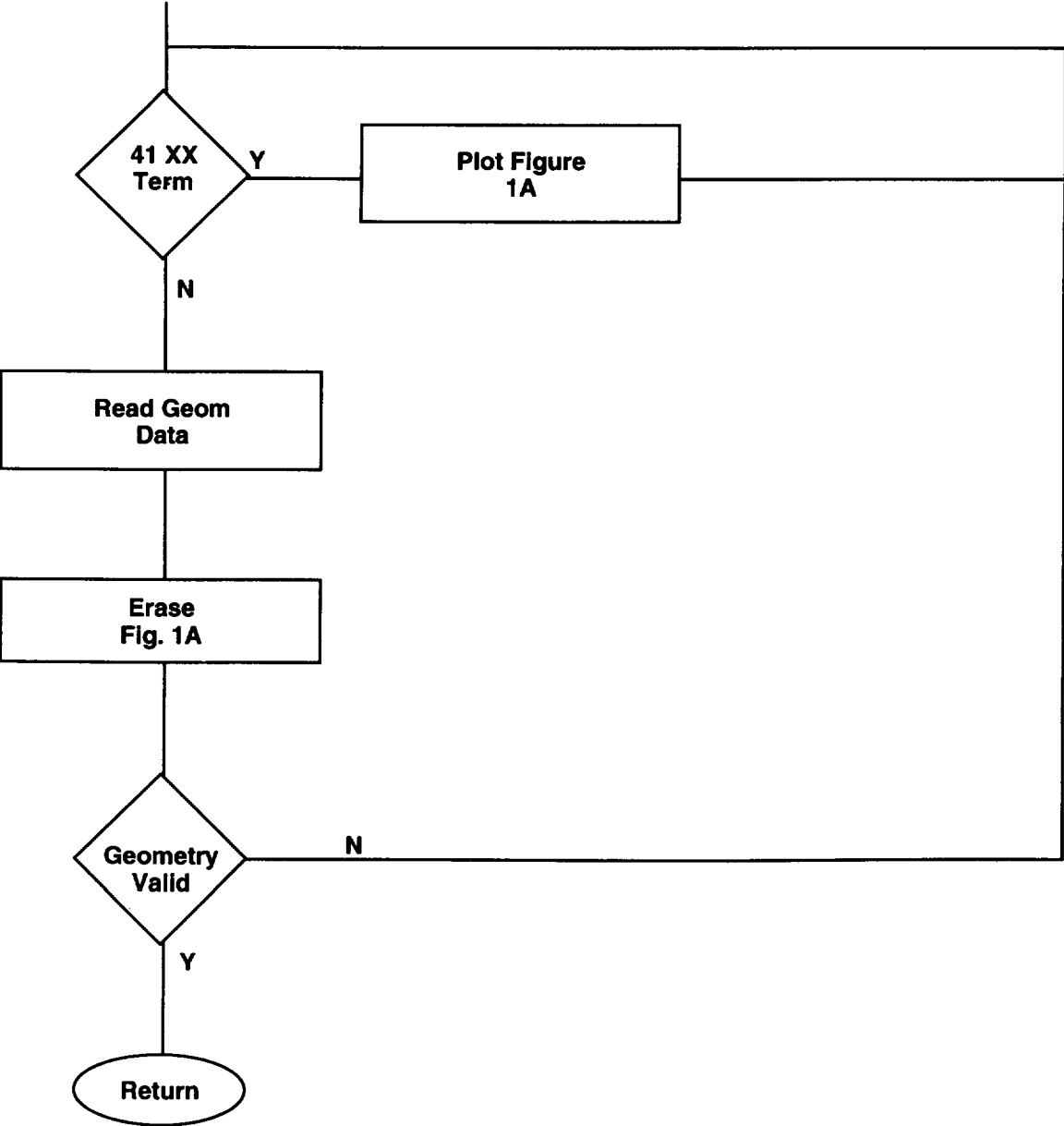
# SUBROUTINE OPCOND1



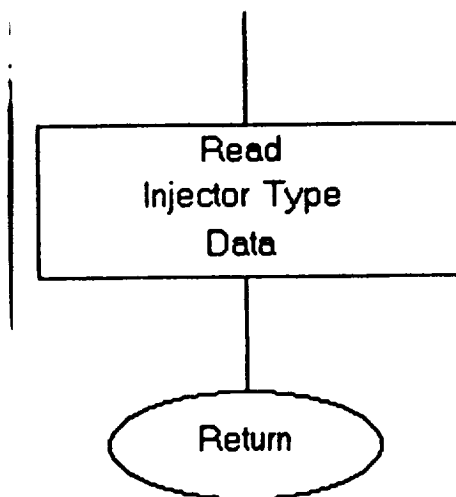
# SUBROUTINE PROPMNU



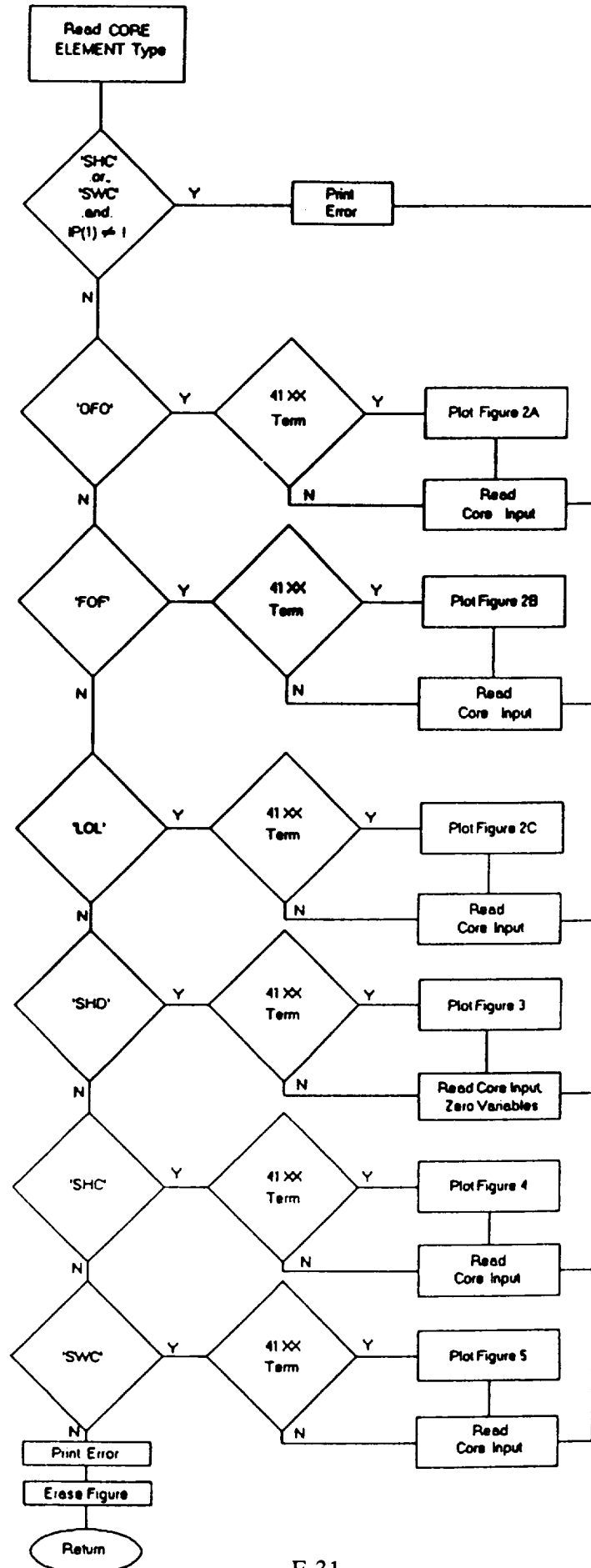
# SUBROUTINE GEOM2



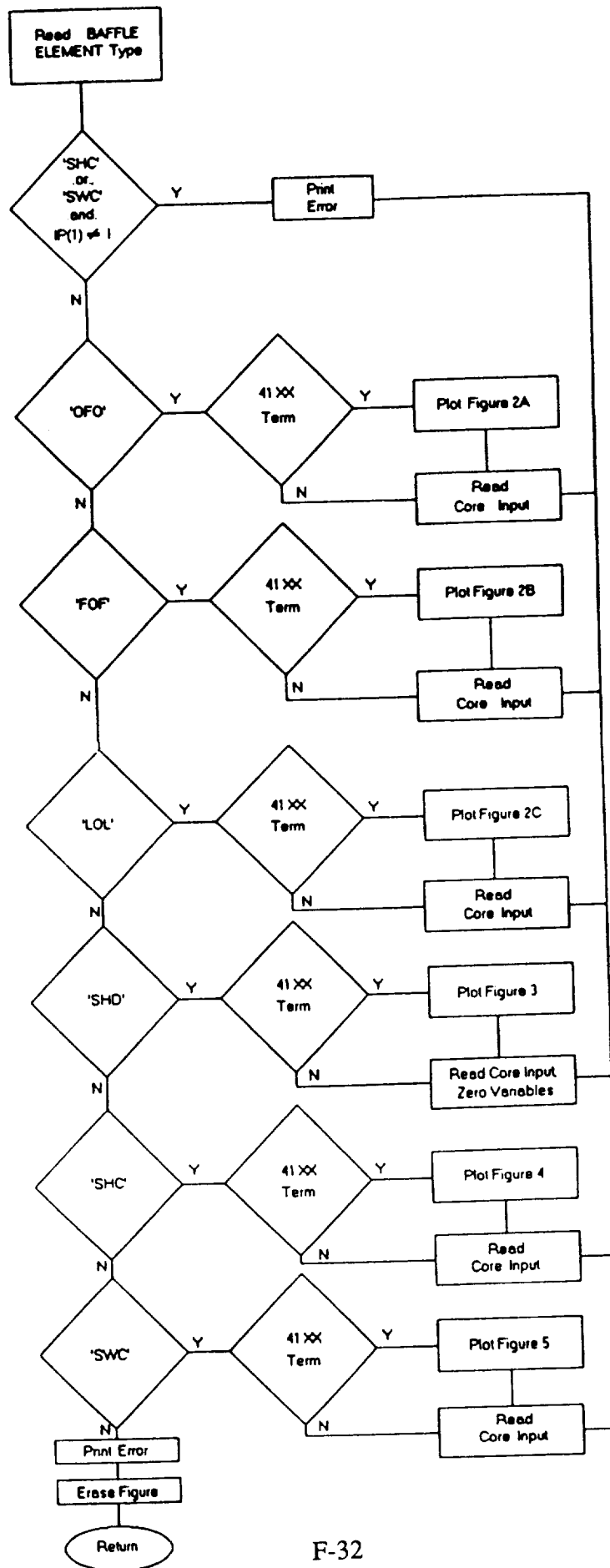
# SUBROUTINE INJ



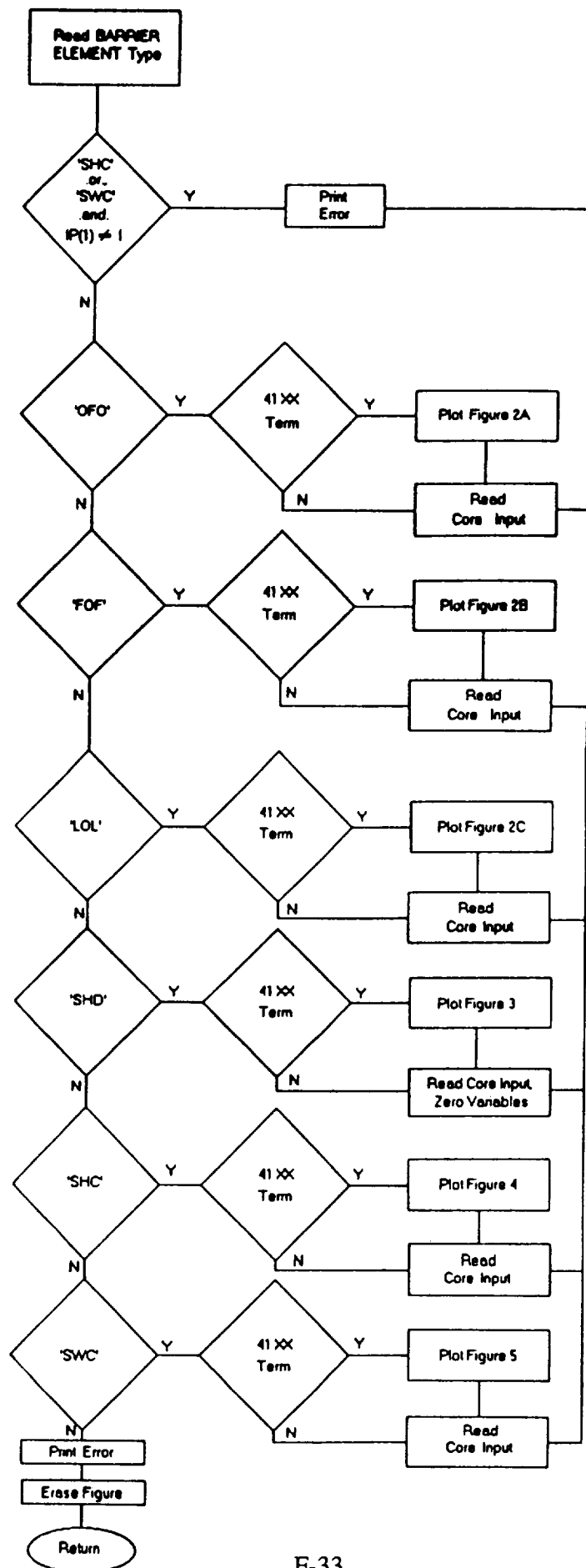
# SUBROUTINE CORELM



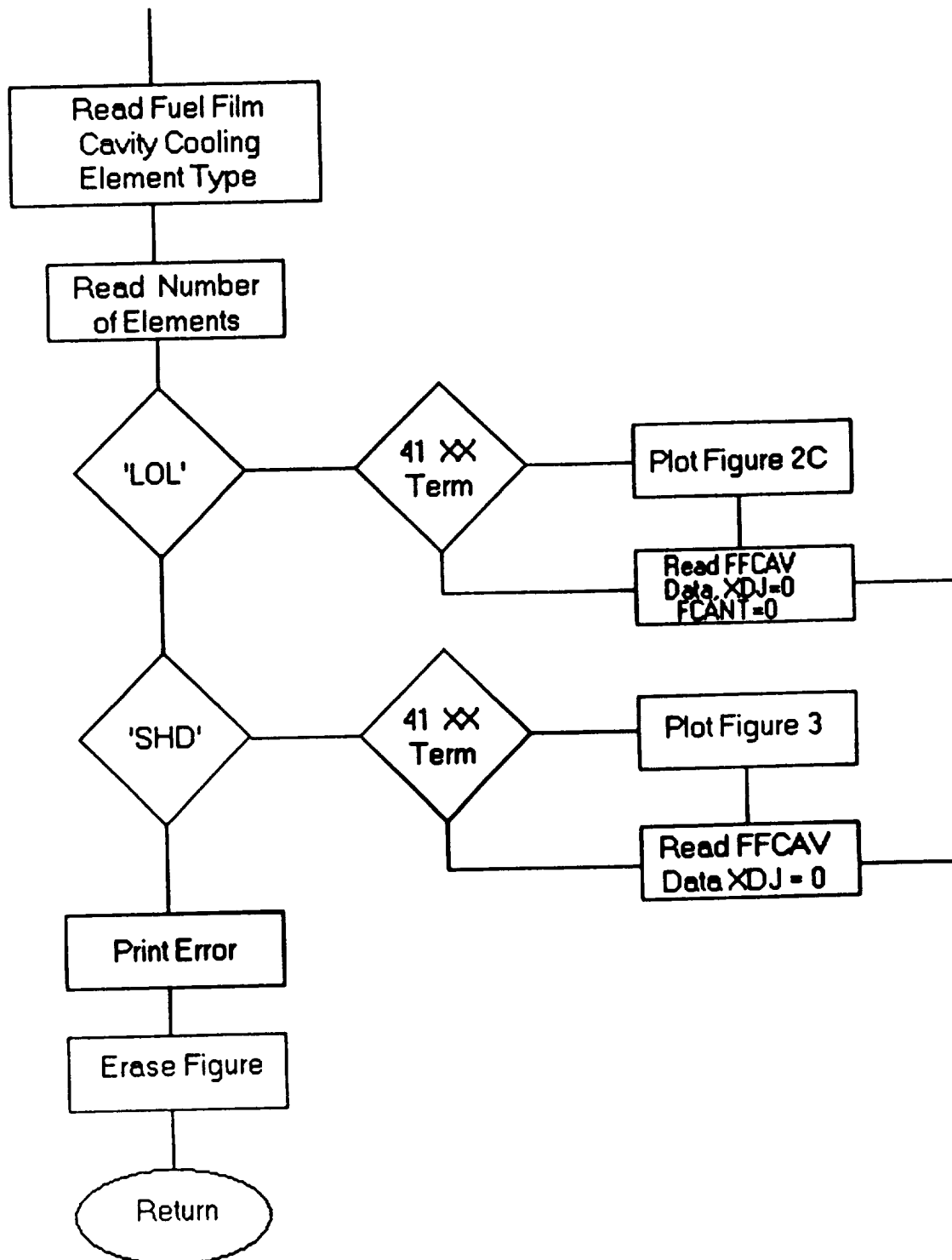
# SUBROUTINE BAFELM



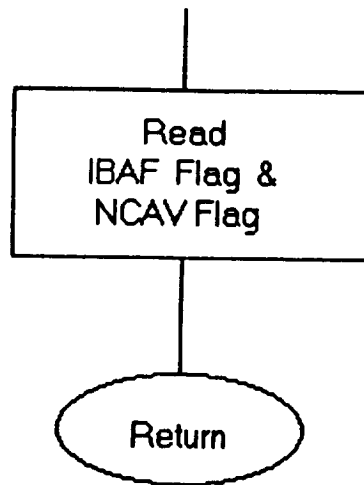
# SUBROUTINE BARELM



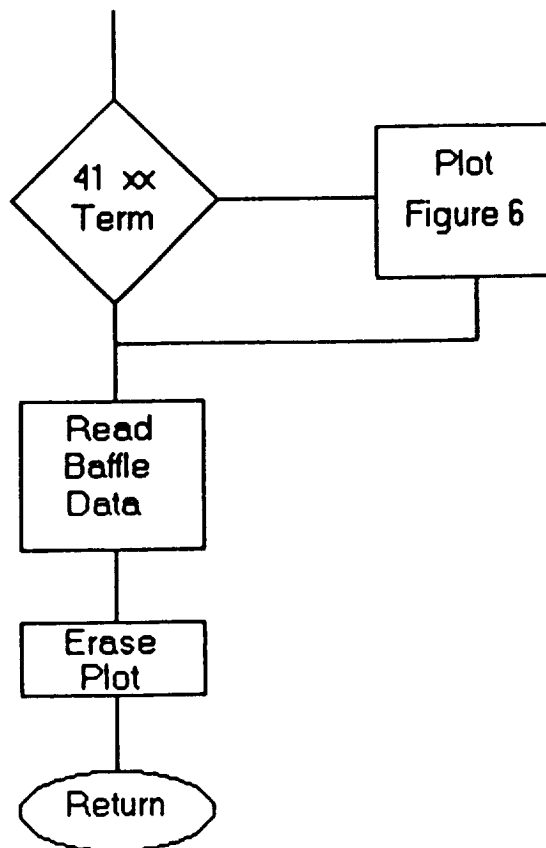
# SUBROUTINE FFCAVE



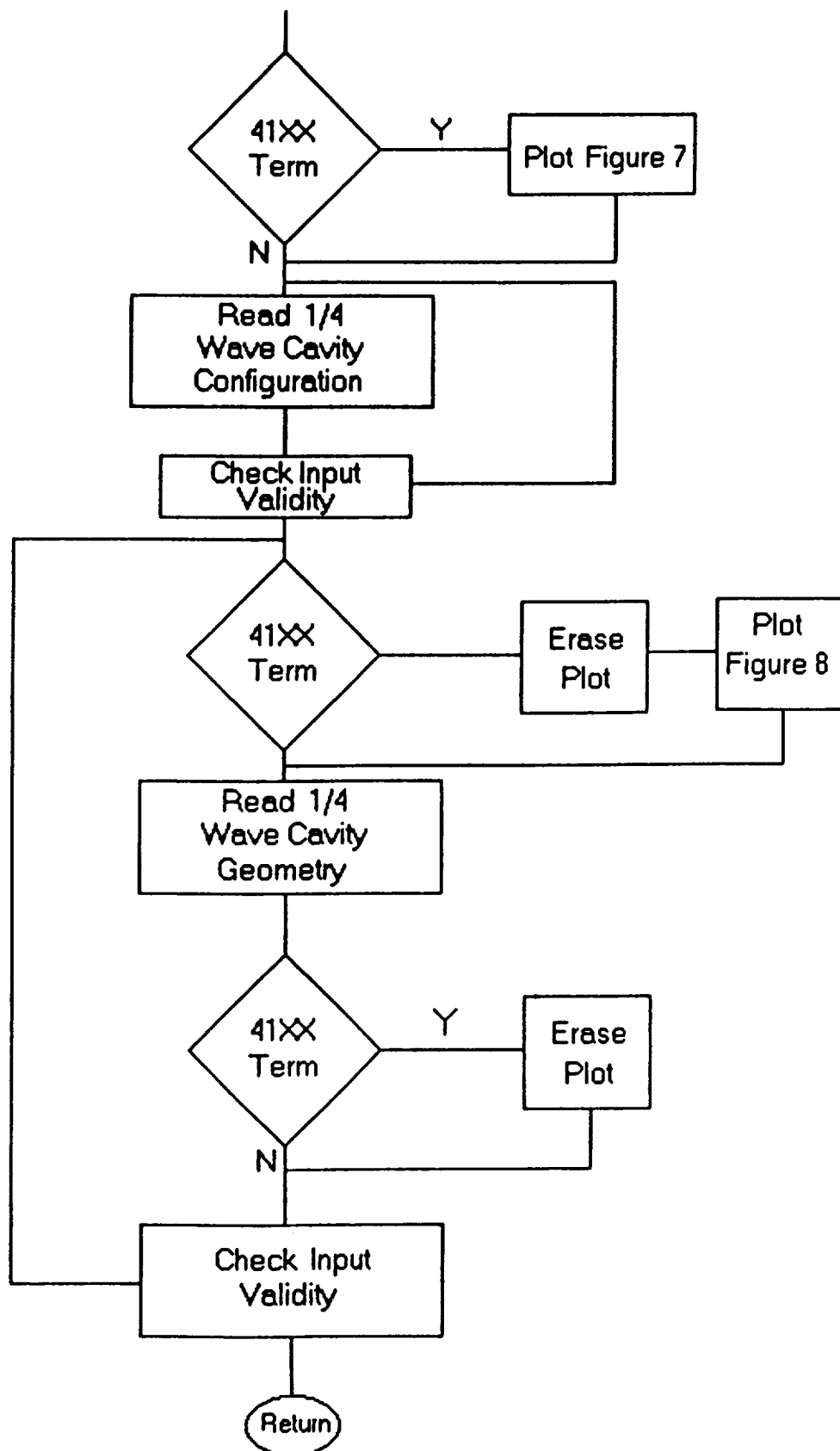
# SUBROUTINE STBAID



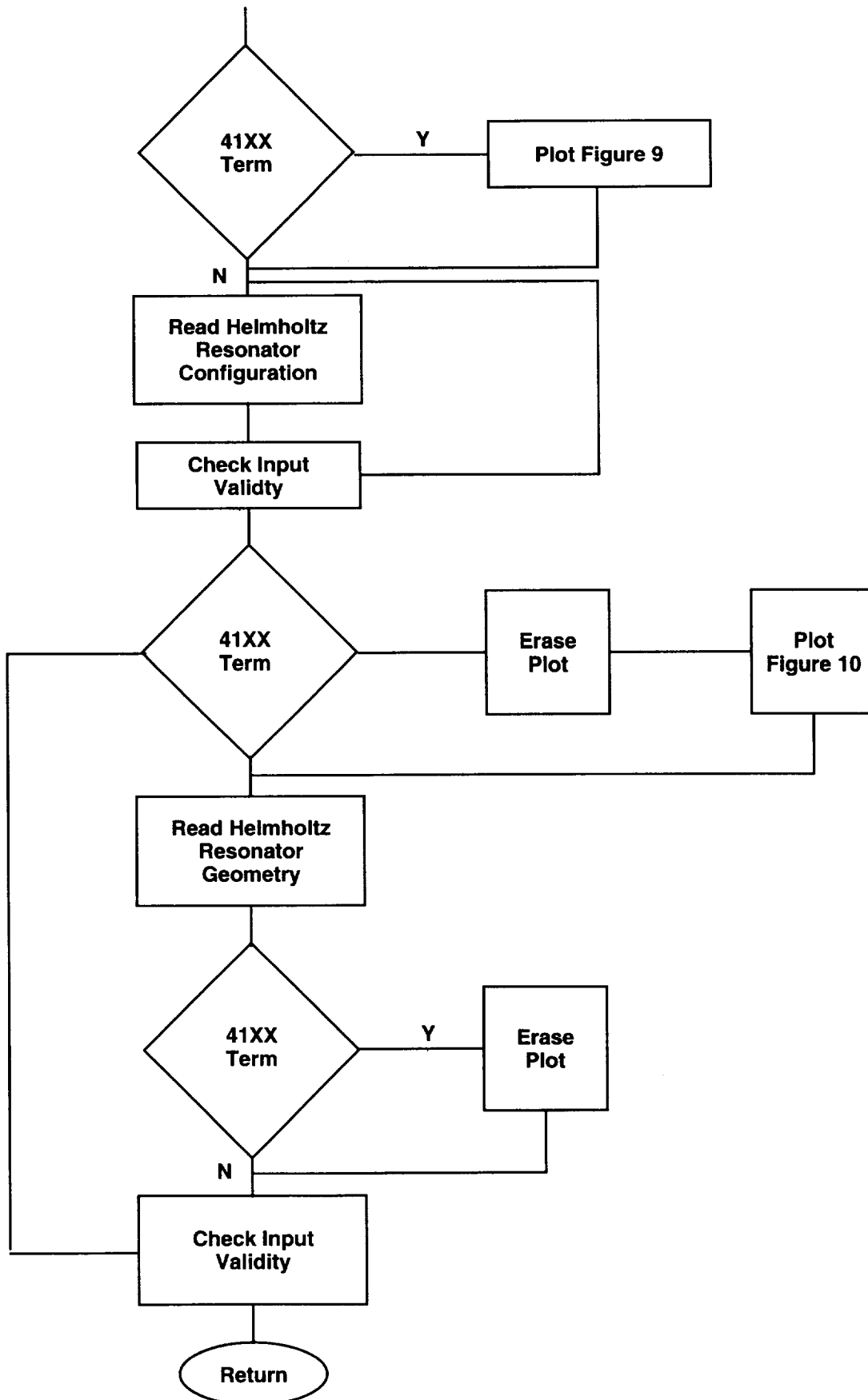
# SUBROUTINE RADBAF



# SUBROUTINE CAV



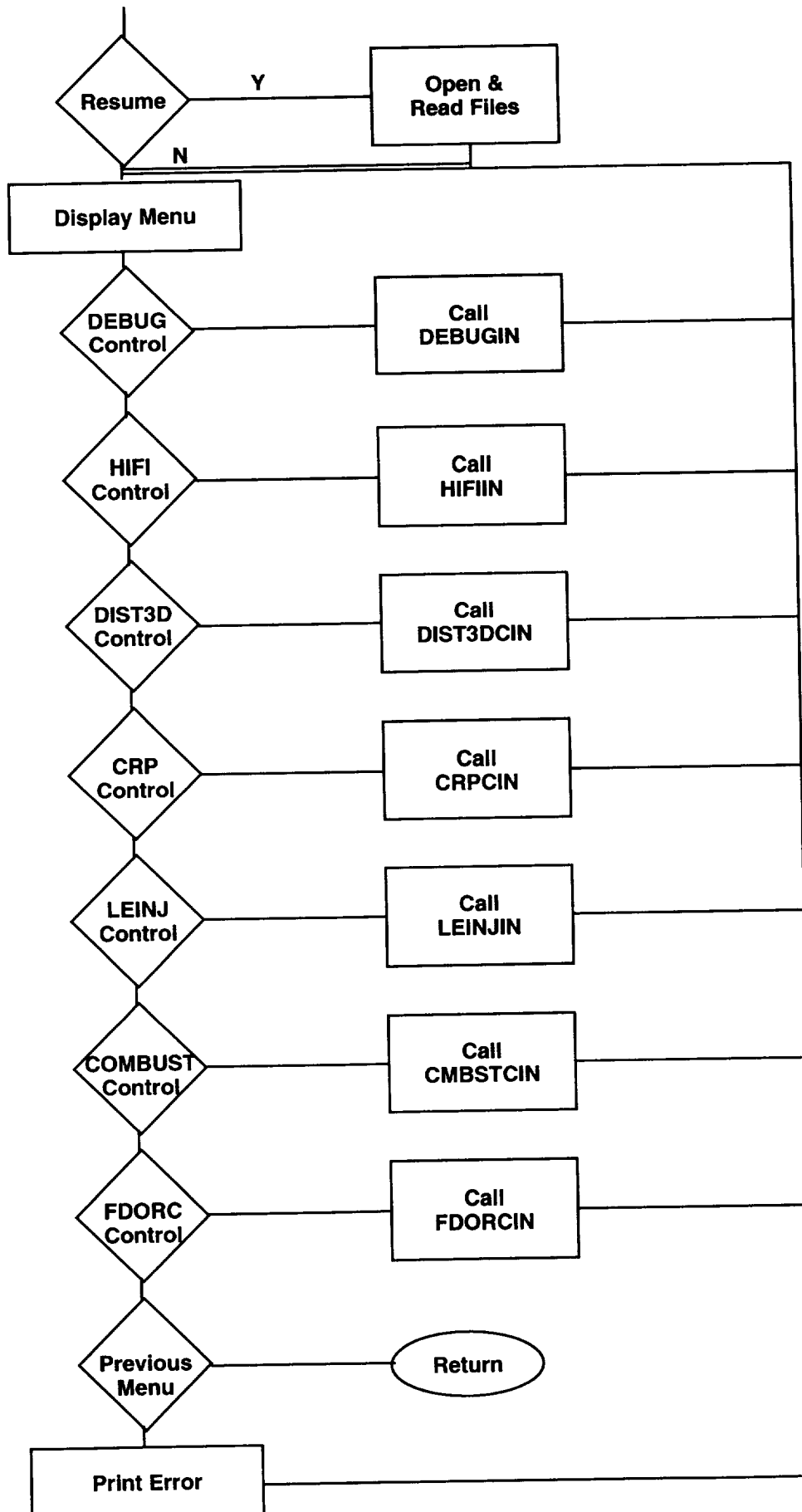
# SUBROUTINE HELMRS



## CONTROL VARIABLE OPTIONS (AUXMNU)

1. Set DEBUG Control
2. Set HIFI Control
3. Set DIST3D Control
4. set CRP Control
5. Set LEINJ Control
6. Set COMBUST Control
7. Set FDORC Control
8. Previous Menu

# AUXMNU



**Part E**  
**IFE Point Design Section Flow Charts**

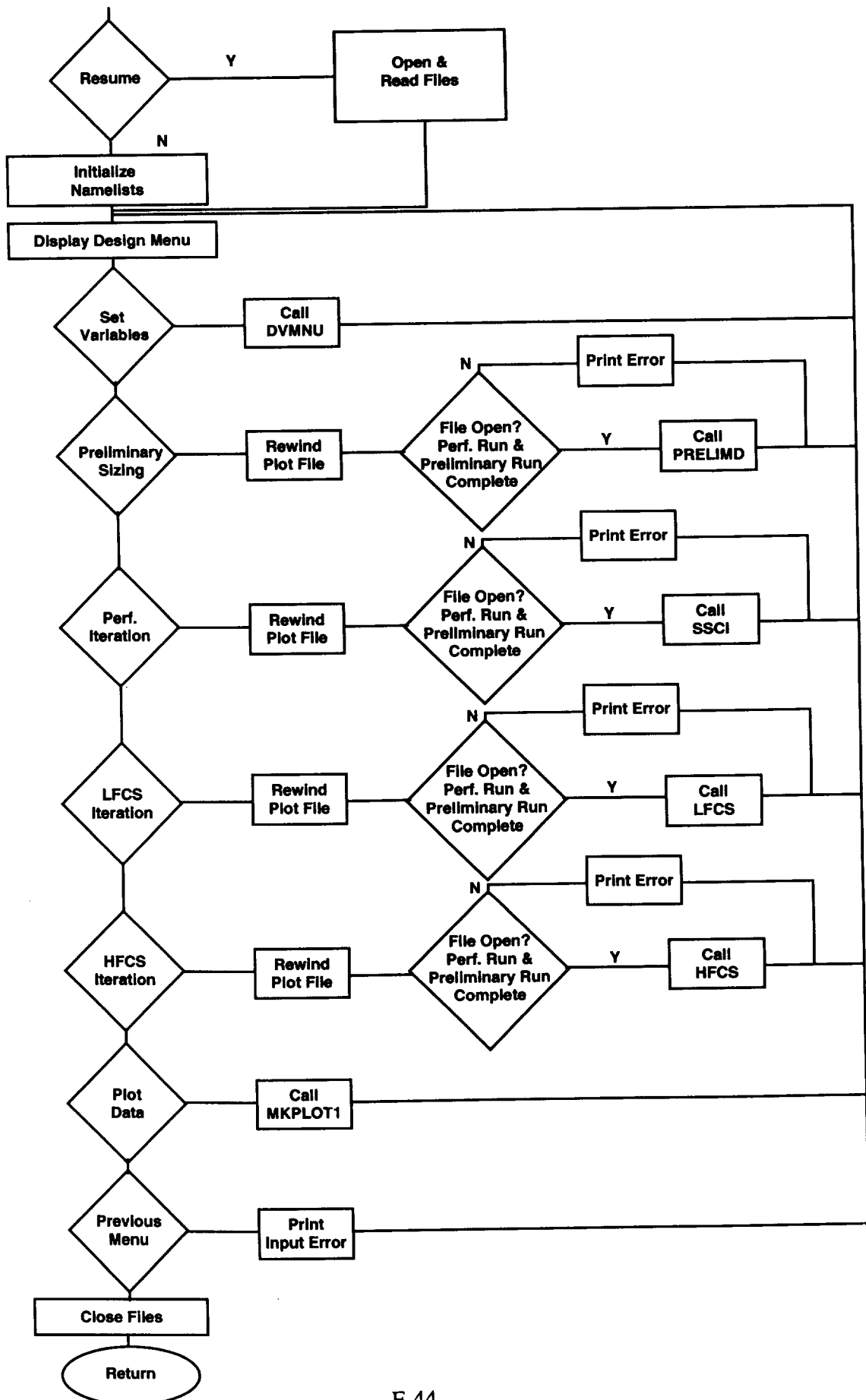
## POINT DESIGN MENU OPTIONS (DESMNU)

- 1 Set Variables
2. Preliminary Sizing
3. Steady State Performance Iteration
4. Chug Stability Iteration
5. High Frequency Stability Iteration
6. Plot Output
7. Previous Menu

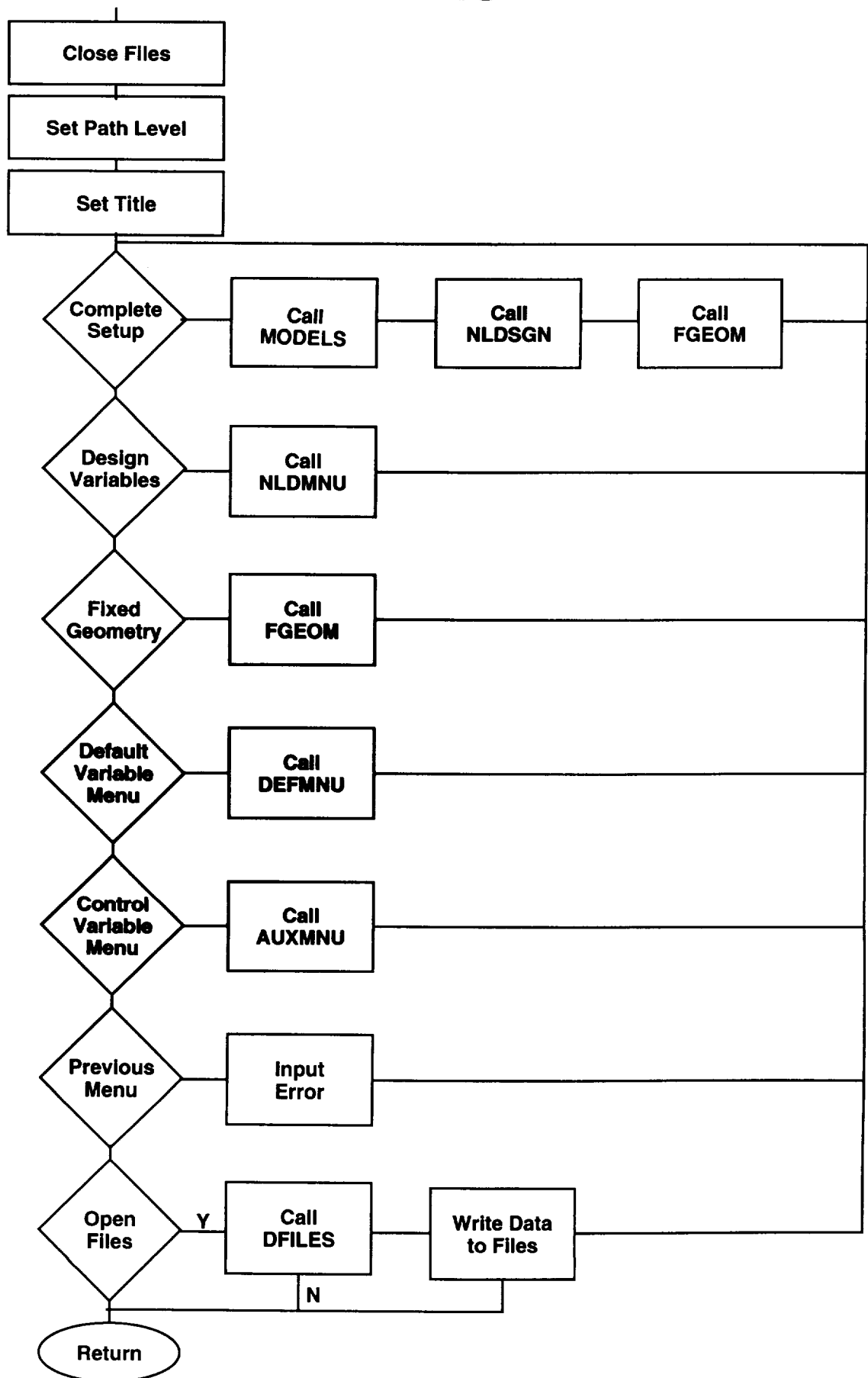
## POINT DESIGN VARIABLE MENU OPTIONS

1. Complete Setup
2. Set Models
3. Set Design Variables
4. Set Fixed Geometry
5. Set Default Variables
6. Set Control Variables
7. Previous Menu

# DESMNU MENU



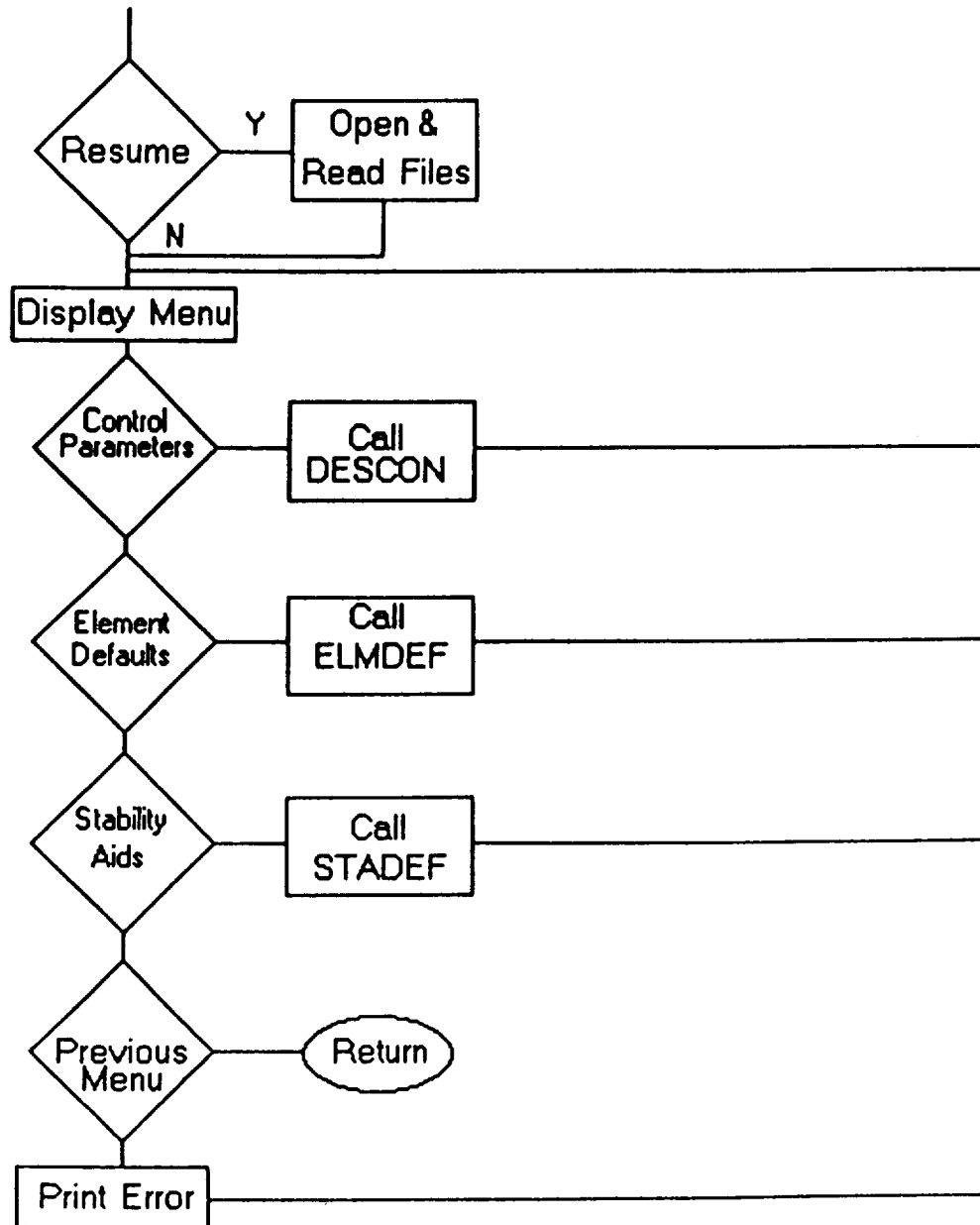
# DVMNU



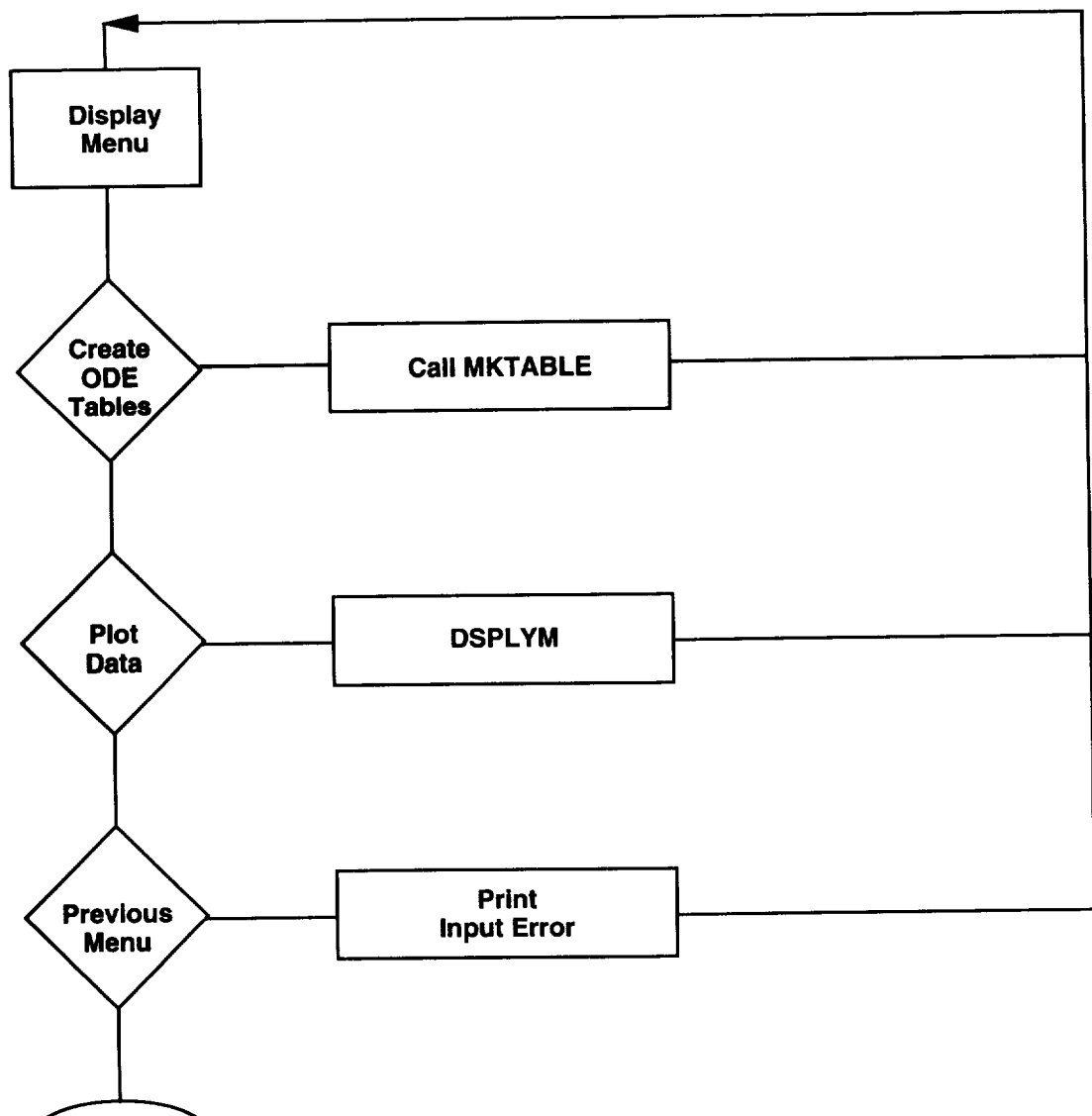
## DESIGN DEFAULT MENU OPTIONS (DEFMNU)

1. Set Design Control Inputs
2. Set Element Parameters
3. Set Stability Aid Parameters
4. Return to Previous Menu

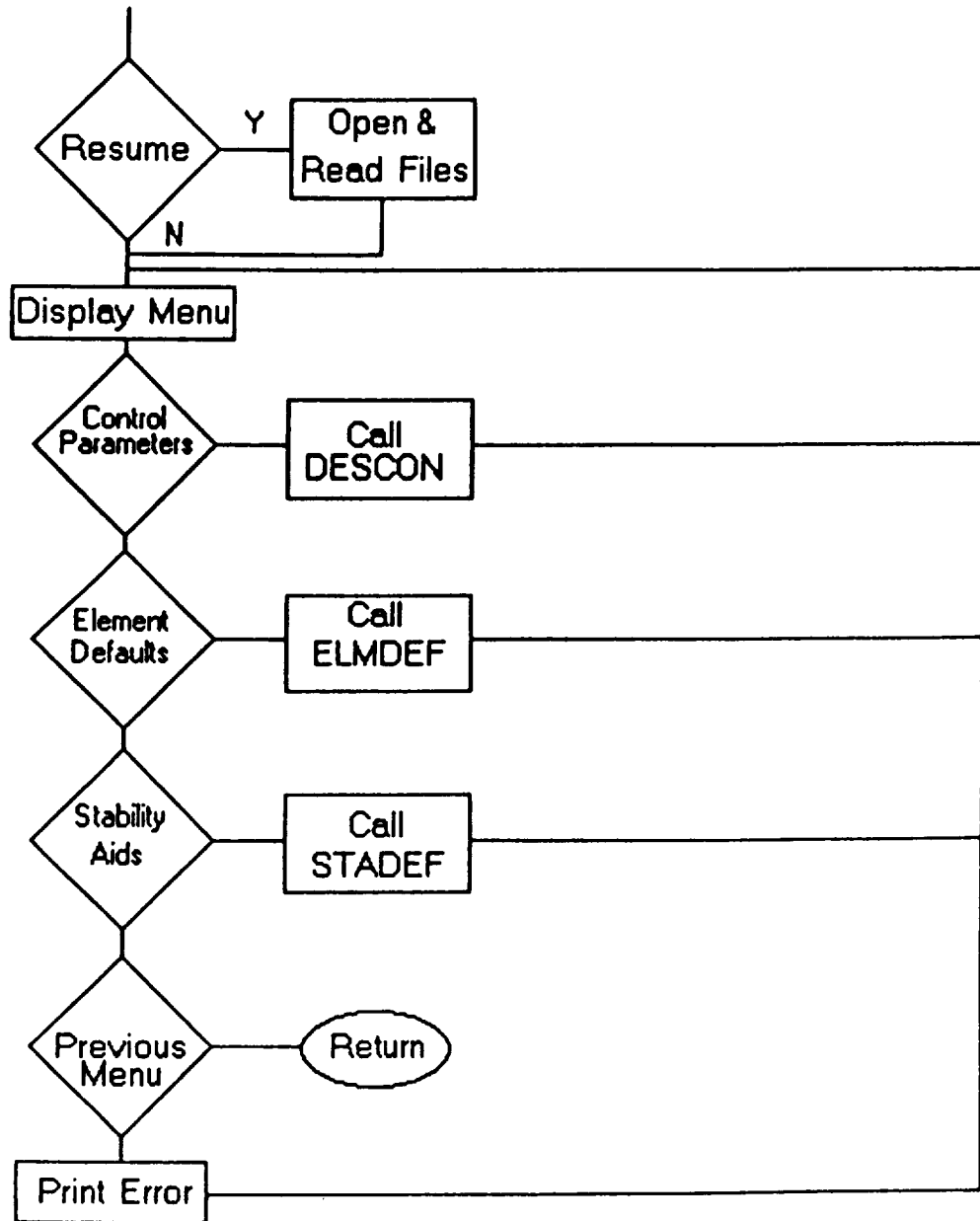
# DEFMNU



# UMENU



# DEF MNU

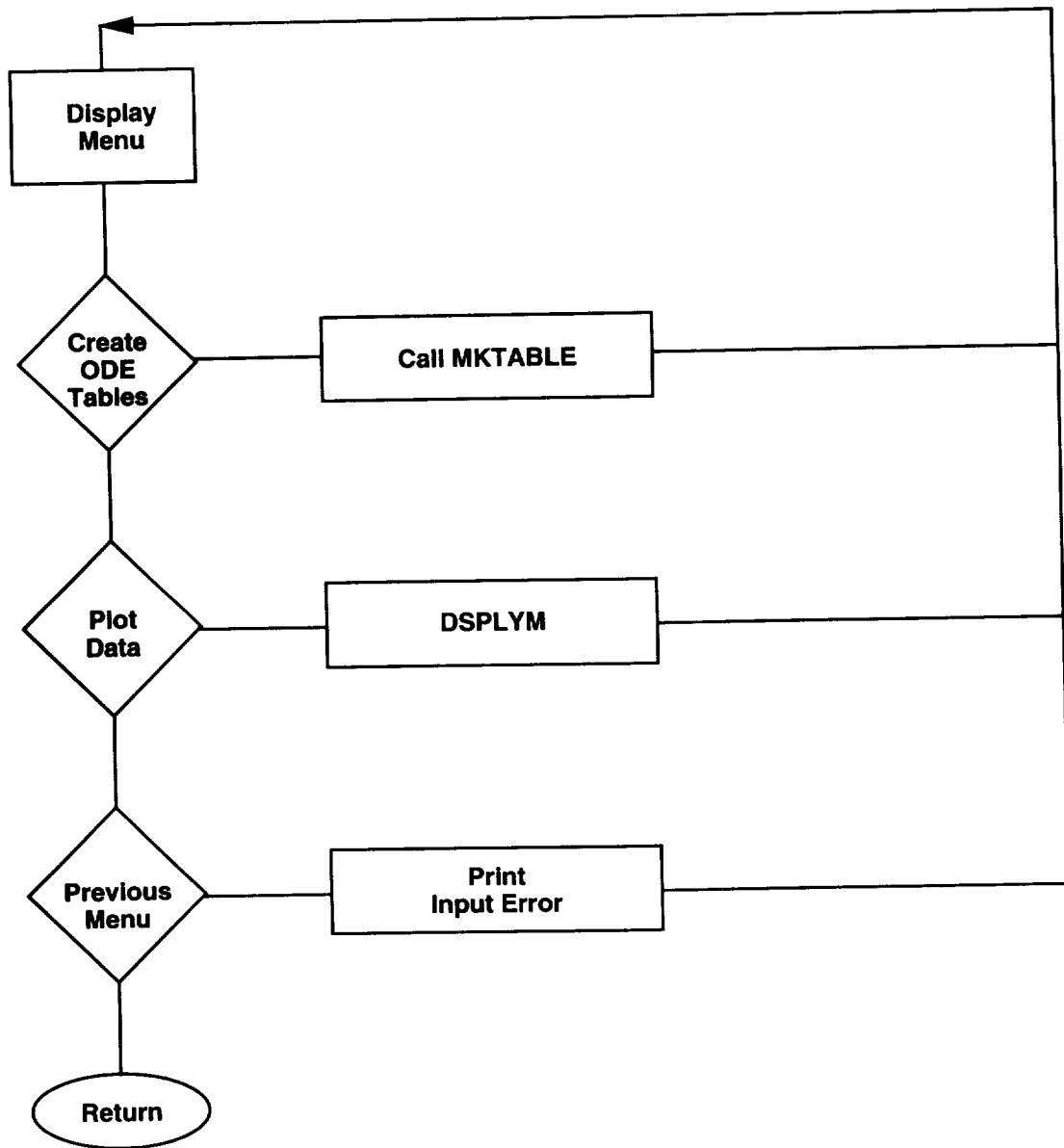


**Part F**  
**IFE Utility Programs**

### UTILITIES MENU (UMENU)

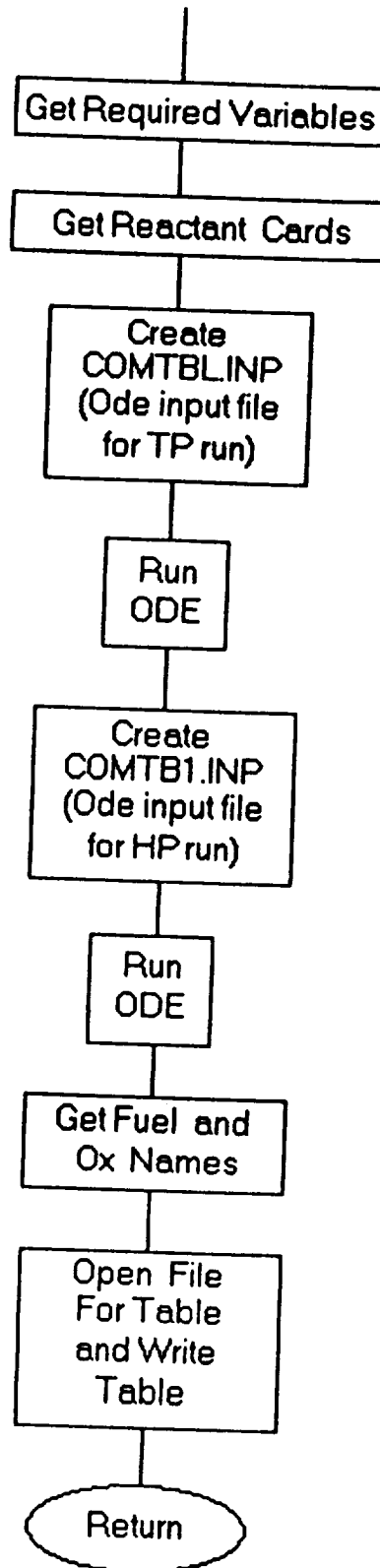
1. Create ODE Combustion Gas Tables
2. Display Results
3. Return to Previous Menu

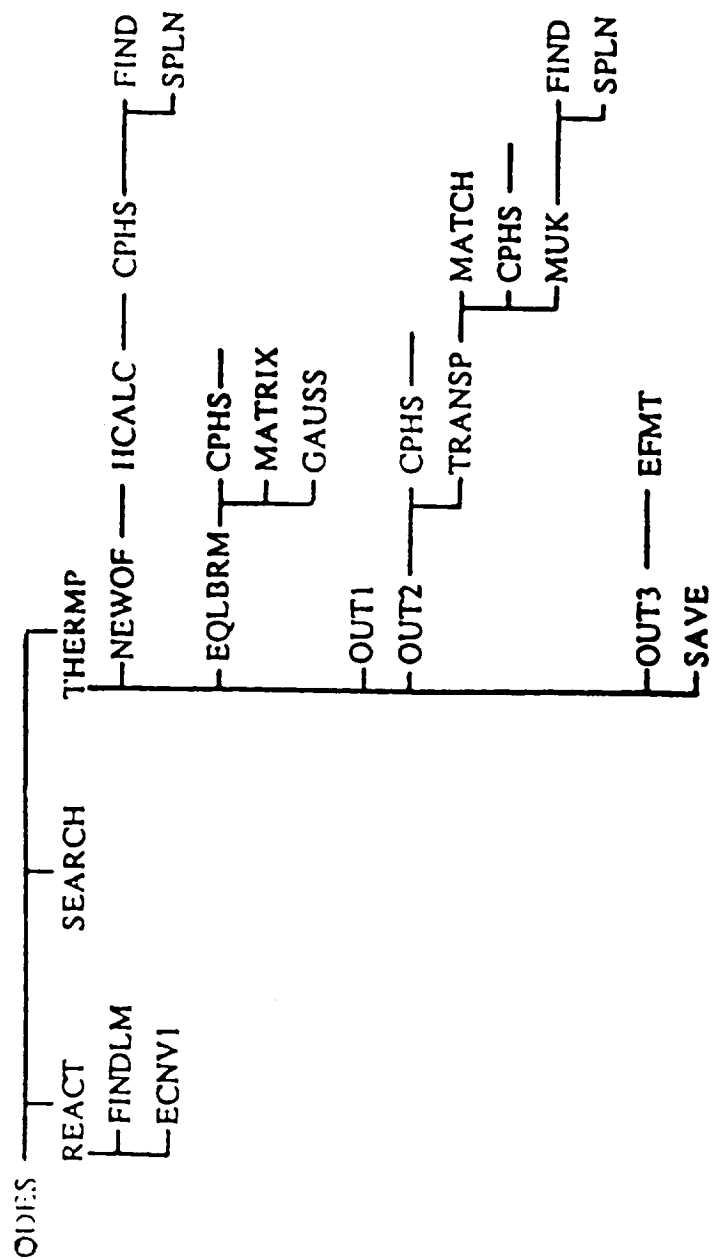
# UMENU



# SUBROUTINE MKTABL

(Makes Combustion Gas Tables)





ROCCID ODE Module Flowchart

**Appendix G**  
**Subroutine Description**

**Part A**  
**POINTA Routines**

## Point Analysis Modules

<u>Routine Name</u>	<u>Source Name</u>	<u>Purpose</u>
COMBUSTD	COMBUST	POINTA BLOCK DATA
HFCS	HFCS	HIGH FREQUENCY STABILITY CONTROL PROGRAM
INJR	INJR	INJECTION RESPONSE MODEL SELECTION
LFCS	LFCS	LOW FREQUENCY STABILITY CONTROL PROGRAM
PROPSET	PROPSET	SETS PROPELLANT PROPERTIES
SINPUT	SINPUT	STABILITY MODEL INPUT READ
SSCI	SSCI	STEADY STATE COMBUSTION ITERATION CONTROL
STABC	STABC	DETERMINES STABILITY FROM COMPONENT TRANSFER FUNCTIONS
WFILES	WFILES	WRITES DATA FROM SSCI FOR USE BY STABILITY MODELS AND TDK
CBGAS	CBGAS	DETERMINES COMBUSTION GAS PROPERTIES
CHAMBR	CHAMBR	CHAMBER RESPONSE MODEL SELECTION ROUTINE
COMBR	COMBR	COMBUSTION RESPONSE SELECTION MODULE
COMBUST	COMBUST	STEADY STATE COMBUSTION MAIN PROGRAM
CRP	CRP	COMBUSTION RESPONSE PREDICTION MODEL
DIST3D	DIST3D	DISTRIBUTED COMBUSTION CHAMBER RESPONSE WITH BAFFLES
DROPS	DROPS	PROPELLANT DROPSIZE MODEL SELECTION ROUTINE
EMCALC	EMCALC	STREAMTUBE MASS FRACTION AND MIXTURE RATIO CALCULATION
EQTEMP	EQTEMP	GETS EQUILIBRIUM COMBUSTION GAS TEMPERATURE FROM TABLES
FLUIDP	FLUIDP	MAIN ROUTINE FOR MIPROPS
GASV	GASV	GASEOUS PROPELLANT INJECTION PRESSURE AND VELOCITY CALCULATION
H2VISC	H2VISC	H2 VISCOSITY DATA
HIFI	HIFI	HIGH FREQUENCY CHAMBER RESPONSE MODEL
INJ	INJ	LUMPED PARAMETER INJECTION RESPONSE MODE
LEINJ	LEINJ	NASA/LERC NON-LINEAR INJECTION RESPONSE MODEL
LIMITS	FLUIDP	CHECKS DATA TABLE LIMITS

# Point Analysis Modules (Continued)

<u>Routine Name</u>	<u>Source Name</u>	<u>Purpose</u>
MACH	MACH	CALCULATES MACH NUMBER FROM GAMMA AND AREA RATIO
MAKCAP	MAKCAP	CONVERTS STRINGS TO CAPITAL LETTERS
NOZADM	NOZADM	MAIN ROUTINE FOR NOZZLE ADMITTANCE MODEL
NOZINI	NOZADM	INITIALIZES PARAMETERS FOR NOZADM
NTAU	NTAU	CONVERTS N-TAU DATA INTO COMBUSTION RESPONSE
NTAU2	DIST3D	CALCULATES BURNING ADMITTANCE
NTVALS	NTVALS	SMITH-READON N-TAU CORRELATION DATA
PERF1	PERF1	INTERPOLATES PERFORMANCE DATA
PINPUT	PINPUT	STEADY STATE COMBUSTION MODEL INPUT
PMELT	FLUIDP	MIPROPS CALCULATION SUBROUTINE
PRESSD	PRESSD	LIQUID PRESSURE DROP CALCULATION ROUTINE
PV	CRP	VAPOR PRESSURE CALCULATION CORRELATION
RAYLEE	RAYLEE	QUASI 1-D GAS DYNAMICS MODEL FOR TOTAL PRESSURE LOSS CALCULATIONS
REIDEL	CRP	DETERMINES PARAMETERS FOR PV CORRELATION
RMPRIEM	DROPS	PREIM DROPSIZE CORRELATION
RMUTRC	DROPS	UTRC DROPSIZE CORRELATION
ROOT	DIST3D	DETERMINES FUNCTION ROOTS
SHEARPD	PRESSD	DETERMINES FLOW CONDITION AND PRESSURE DROP IN SHEAR COAX ELEMENTS
SL	CRP	HEAT OF VAPORIZATION CORRELATION
SOUND	FLUIDP	DETERMINES SOUND SPEED OF PROPELLANT
SPLINE	RAYLEE	SETUP ROUTINE FOR CUBIC SPLINE INTERPOLATION
SPLINT	RAYLEE	CUBIC SPLINE INTERPOLATION ROUTINE
STOIC	ATOM	SETS STOICHIOMETRIC MR FOR DIFFERENT PROPELLANT COMBINTAIONS
SUM	CRP	SUMS OSCILLATORY VAPORIZATION TO DETERMINE RESPONSE
SWIRLPD	PRESSD	SWIRL COAX ELEMENT PRESSURE DROP/FLOWRATE CALCULATION
TIMELAG	TIMELAG	DETERMINES TIMELAGS FROM COMBUST DATA
TMELT	FLUIDP	MIPROPS CALCULATION SUBROUTINE
VAPRO	VAPRO	GENERALIZED LENGTH VAPORIZATION ROUTINE
VDISP	DIST3D	VISCOUS DISIPATION CALCULATIONS
VISC	FLUIDP	VISCOSITY CALCULATION ROUTINE

# Point Analysis Modules (Continued)

<u>Routine Name</u>	<u>Source Name</u>	<u>Purpose</u>
XNTRP ACOUSTIC	RAYLEE HIFI	LINEAR INTERPOLATION SUBROUTINE CALCULATES ACOUSTIC BEHAVIOR WITHIN A SECTION
AINTP ATOM BESSCAL BI1 BI2 BI3 BJ CALADM CAVITY CBTAB CCOSH CDT CHI1 CHI2 CHI3 CP	AINTP ATOM DIST3D DIST3D DIST3D DIST3D DIST3D NOZADM CAVITY CBTAB DIST3D CDT DIST3D DIST3D DIST3D FLUIDP	LINEAR INTERPOLATION FUNCTION MAIN ATOMIZATION ROUTINE CALCULATES BESSAL FUNCTIONS NUMERICAL INTEGRATION FUNCTION NUMERICAL INTEGRATION FUNCTION NUMERICAL INTEGRATION FUNCTION NUMERICAL INTEGRATION FUNCTION COMPUTES NOZZLE ADMITTANCE FROM "CHI" ACOUSTIC CAVITY CONTROL MODEL GETS COMBUSTION GAS PROPERTIES FROM TABLES HYPERBOLIC COSINE THROAT CD CORRELATION NUMERICAL INTEGRATION FUNCTION NUMERICAL INTEGRATION FUNCTION NUMERICAL INTEGRATION FUNCTION CONSTANT PRESSURE SPECIFIC HEAT CALCULATION
CRPDBG CV DCAVEF DILV DROPMIX ENTHAL ENTROP EXCESV FDATA FDCV FINDD FRICTION FWDOT GETVAL HCAVEF HELMHR INTGRT	CRP FLUIDP DIST3D FLUIDP DROPS FLUIDP FLUIDP FLUIDP FLUIDP FLUIDP FLUIDP PRESSD CRP GETVAL HIFI CAVITY NOZADM	CRP DEBUG OUTPUT GENERATION CONSTANT VOLUME SPECIFIC HEAT CALCULATION DIST3D CAVITY ORIENTATION EFFECTIVENESS MIPROPS CALCULATION SUBROUTINE DROPMIX DROPSIZE CORRELATIONS ENTHALPY CALCULATION ENTROPY CALCULATION EXCESS VISCOSITY ROUTINE LOADS PROPELLANT DATA MIPROPS CALCULATION SUBROUTINE DENSITY CALCULATION MOODY FRICTION FACTOR CHART OSCILLATORY FLOWRATE CALCULATION POWER-LAW INTERPOLATION FUNCTION HIFI CAVITY ORIENTATION EFFECTIVENESS HELMHOLTZ RESONATOR MODEL NUMERICAL INTEGRATION ROUTINE

## Point Analysis Modules (Continued)

<u>Routine Name</u>	<u>Source Name</u>	<u>Purpose</u>
MAXMIN	RAYLEE	FINDS MAX AND MIN OF DATA SET
PROPS	FLUIDP	MIPROPS CALCULATION SUBROUTINE
QWC	CAVITY	1/4 WAVE CAVITY MODEL
REGULA	FLUIDP	MIPROPS CALCULATION SUBROUTINE
SATL	FLUIDP	MIPROPS CALCULATION SUBROUTINE
SATV	FLUIDP	MIPROPS CALCULATION SUBROUTINE
SHEAR	ATOM	AEROJET SHEAR COAX ATOMIZATION MODEL
SSATL	FLUIDP	MIPROPS CALCULATION SUBROUTINE
SSATV	FLUIDP	MIPROPS CALCULATION SUBROUTINE
SWIRL	ATOM	AEROJET SWIRL COAX ATOMIZATION MODEL
TRIPLET	ATOM	AEROJET TRIPLET ATOMIZATION MODEL
VISCE	FLUIDP	MIPROPS CALCULATION SUBROUTINE
VPN	FLUIDP	MIPROPS CALCULATION SUBROUTINE
VSCITY0	FLUIDP	MIPROPS CALCULATION SUBROUTINE
BJ0	CRP	BESSEL FUNCTION CALCULATION ROUTINE
BJ1	CRP	BESSEL FUNCTION CALCULATION ROUTINE
CPI	FLUIDP	MIPROPS CALCULATION SUBROUTINE
FINDM	FLUIDP	MIPROPS CALCULATION SUBROUTINE
SHI	FLUIDP	MIPROPS CALCULATION SUBROUTINE
HIF12FD	HIF12FD	TRANSFER/HIF1/DIST3D CAVITY DATA TO \$FDORC
FDORC	FDORC	MAIN ROUTINE FOR FDORC
CAV	FDORC	FDORC ROUTINE FOR VARIABLE TEMPERATURE CAVITIES
FNTAU	FDORC	FDORC ROUTINE TO CALCULATE N AND TAU

**Part B**  
**POINTD Routines**

## Point Design Modules

<u>Routine Name</u>	<u>Source Name</u>	<u>Purpose</u>
CHUGIT CLOAD	CHUGIT CLOAD	CHUG STABILITY ITERATION CONTROL LOADS DESIGN COMMONS WITH POINTA DATA FOR STABILITY MODELS
DINPUT HIFIT	DINPUT HIFIT	READS POINT DESIGN INPUT HIGH FREQUENCY STABILITY ITERATION CONTROL
LFDESIGN MODEL0D PERFIT	LFDESIGN MODEL0D PERFIT	LOW FREQUENCY REDESIGN MODULE LOADS H.F. UNSTABLE MODE TABLE STEADY STATE PERFORMANCE ITERATION CONTROL
PRELIMD REDESIGN	PRELIMD REDESIGN	MAIN PRELIMINARY DESIGN ROUTINE RESIZES INJECTOR/CHAMBER FOR SELECTED DESIGN CHANGE
VGALC VGUPDATE	VGALC VGUPDATE	GASEOUS INJECTION VELOCITY ROUTINE UPDATES GASEOUS VELOCITY AND PRESSURE DROP IN POINTD USING GASV
VSAVE	VSAVE	SAVES "MEMORY" VARIABLES IN \$\$SAVE SO RUNS CAN BE CONTINUED
ALOAD	ALOAD	WRITES DESIGN DATA IN ANALYSIS INPUT FILE FORMAT
BITUCAVITY BITURES0N BRNDESIGN	BITUCAVITY BITURES0N BRNDESIGN	BITUNE 1/4 WAVE CAVITY DESIGN MODULE BITUNE HELMHOLTZ RESONATOR DESIGN MODULE BURNING-COUPLED H.F. STABILITY REDESIGN MODULE
CAVLOAD COREFLOW CORESIZE	CAVLOAD COREFLOW CORESIZE	LOADS ACOUSTIC CAVITY DESIGN VARIABLES DETERMINES CORE ELEMENT FLOWRATE AND MR SIZES CORE ELEMENTS DURING PRELIMINARY DESIGN
CORESPAC	CORESPAC	CHECKS SPACING AND LA YOUT FOR CORE ELEMENTS
CUPDATE	CUPDATE	UPDATES DESIGN COMMONS WITH POINTA DATA FROM SSC1
DAMP DPOST	DAMP DPOST	DAMPING DEVICE DESIGN MODULE SHEAR COAX LOX POST DESIGN ROUTINE

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## Point Design Modules (Continued)

<u>Routine Name</u>	<u>Source Name</u>	<u>Purpose</u>
DROPSIZE EMEST	DROPSIZE EMEST	DETERMINES DROPSIZE TO MEET INPUT RESPONSE ESTIMATES OVERALL EM REQUIRED FOR INPUT ETA-MIX
INJDESIGN	INJDESIGN	INJECTION-COUPLED H.F. STABILITY REDESIGN MODULE
JETSIZE	JETSIZE	DETERMINES ORIFICE SIZE TO GENERATE INPUT DROPSIZE
MONOCAVITY MONORESON	MONOCAVITY MONORESON	MONOTUNE 1/4 WAVE CAVITY DESIGN MODULE MONOTUNE HELMHOLTZ RESONATOR DESIGN MODULE
NOZOPT	NOZOPT	NOZZLE-COMBUSTOR LENGTH OPTIMIZATION ROUTINE
NTUPDATE	NTUPDATE	ESTIMATES CHANGES TO N & TAU FROM DESIGN CHANGES W/ NTVALS
PDESIGN RDROP	PDESIGN RDROP	PERFORMANCE ITERATION REDESIGN MODULE DETERMINES DROPSIZE FROM INPUT ORIFICE DIAMETER
SHEAREM	SHEAREM	CALCULATES SHEAR COAX UNI-ELEMENT EM FROM CORRELATION
SHEARFL	SHEARFL	DETERMINES FLOWRATE OF FIXED SHEAR COAX POST W/ SHAERPD
SWIRLFL	SWIRLFL	DETERMINES FLOWRATE OF FIXED SWIRL COAX POST W/ SWIRLPD
SWIRLSZR	SWIRLSZR	SWIRL COAX POST DESIGN ROUTINE

**Part C**  
**IFE Routines**

## IFE Modules

<u>Routine Name</u>	<u>Purpose</u>
Abort	Closes files and stops program
Afiler	Opens files for Point Analysis (asks opening questions)
Afiles	Opens and closes files for Point Analysis
Alfas	Sets alphanumeric letters size on TEK terminal
Allfigs	Draws figures
Anamnu	Point analysis main menu
Ansi	Sets TEK terminal in ANSI mode
Arrow	Draws arrows for figures
Askint	Asks user and gets an Integer
Auxlib	Plotting library
Auxmnu	Control variable menu
Backup	Backs up replay file N lines
Bafelm	Main routine for baffle element data
Bafflein	Computer written code to get baffle element data
Barelm	Main routine for barrier element data
Barrierin	Computer written code to get barrier element data
Bivar	bi-variable interpolation routine
Blanks	Deletes leading blanks
Blink	Starts blinking output on VT100 terminal
Bmenu	Base menu
Bottom	Reads to the bottom of a file
Burnin	Computer written code to get burn data
Caps	Changes everything to capitals
Cav	Main routine for cavity data
Cfdorcin	Reads \$FDORC variables
Chambin	Computer written routine to chamber data
Checki	Check limits on an integer
Clear	Clear screen
Clears	Clear graphics screen
Cmbstcin	Reads and checks COMBUSTC variables
Cntrl	Reads and writes title
Controlin	Computer written code to get control variables
Corein	Computer written code to get core element variables
Corelm	Main routine for core element data
Crack	Reads data IFE table
Crpcin	Computer written code to get crp control info.

## IFE Modules (Continued)

<u>Routine Name</u>	<u>Purpose</u>
Curmv	Moves cursor on a VT100 terminal
Cycle	Changes revision number manually (Needed for SUN)
Dasorc	Determine source of nominal namelist data
Debugin	Computer written code to set debug flag
Decodr	Decode character strings to variables
Defmnu	Default menu for Point Design variables
Descon	Gets design control variables
Designin	Computer written code to get design control variables
Desmnu	Point Design main menu
Dfiler	Opens files for Point design (asks opening questions)
Dfiles	Opens and closes files for Point Design
Dgicomb	Design side routine to get combustion data
Dialog	Asks initial questions
Dispcs	Asks for C* and ISP on the design side
Dist3dcin	Computer written code to get DIST3D control variables
Dnozeff	Prompts user for nozzle efficiency curve
Dopmnu	Design Operating conditions menu
Dpropmnu	Design Propellant selection menu
Dsplay	Displays present variable value
Dsplym	Plotting utility
Dsplymp	Online plotting option
Dvmnu	Design variable menu
Effmnu	Design side efficiency menu
Elmdef	Reads element data for design side
Encode	4107 primitive device driver
Erase	erases screen
Errsns	Dummy routine to comply with VAX used in error message
Fcolon	Finds colon when reading plot data files
Fdorcin	Reads \$FDORC variables
Ffcave	Main routine for reading Fuel Film/Cavity data
Ffcin	Computer written code to get FFC data
Fgeom1	Reads fixed geometry
Fgeomin	Computer written code to get fixed geometry
Figs	Plots figures
Filler	Interprets input (from replay file)
Find	Finds variable in list
Findsp	Finds space
Finjin	Computer written code to get fixed injector data

## IFE Modules (Continued)

<u>Routine Name</u>	<u>Purpose</u>
Geom2	Gets Point Analysis geometry
Geomin	Computer written code to get fixed injector data
Getchr	Gets a character
Getcom	Get a command file
Getdat	Read data
Getfil	Opens a file
Getint	Reads an integer
Getit	Get a variable
Getset1	Reads a set of plot data from a plot file
Getvar	Reads a data input line
Gfdorc	Prompts user and checks \$FDORC variables
Gtcomb	Read gas combustion data defaults
Gtcomb1	Read gas combustion data defaults and tables
Gtdata	Searches plot data file for correct data set
Gtdsna	Prompts user for Point Design C*/ISP data
Gtdsne	Prompts user for Point Desing efficiency data
Gting	Get an integer
Gtline	Read a Character line
Gtopca	Prompts user for Point Analysis C*/ISP data
Gtreal	Get a real variable
Helmrs	Get Helmholtz resonator data
Hifiin	Computer written code to get HIFI control data
Ifed	Block Data, contained in ROCCID for
Inject1	Routine which reads Injector data
Injin	Computer written code to get injector data
Inperr	Informs of input error
Instr	Displays instructions
Ioerr	Gives VAX status of IO error
Iter	Secant search iteration scheme
Leinjin	Computer written code to get Lewis injector data
Load	Loads variable into array
Lolcin	Computer written code to get LOL data for Design
Ltcphs	Part of ODE
Makfig	Loads the figures at the beginning of a session
Manif	Gets manifold data
Mkplot	Reads plot data 1 file
Mkplot1	Reads plot data 3 files
Mktabl	ODE utility

## IFE Modules (Continued)

<u>Routine Name</u>	<u>Purpose</u>
Model	Gets model control variables
Modelsin	Computer written code to read model control data
Nchars	Finds number of non-blank characters (space)
Newfil	Read old replay files
Newfrm	Start new plotting frame
Nldmnu	\$Design variable menu
Nldsgn	\$Design variable prompts
Not	Yes/No function (reads yes of no and acts)
Nozeff	Nozzle efficiency vs length routine
Nrmvid	Puts VT100 terminal in normal mode
Nulfld	Advances index past null fields
Odecst	Sets up and runs ODE to calculate C* and Isp arrays
Odein	Computer written code to read ODE input data
Odemain	Main calling routine for ODE
Opcoin	Computer written code to read operating conditions
Opcond1	Main routine to get operating conditions
Outcm	Writes character variable out
Outrm	Writes real variable out
Parse	Parse data line
Perfm	Driver for steady state routines
Perr	Print error
Plotm	Plotting routine
Pltset	Reads plot attributes (size, titles, etc.) from data file
Propmnu	Point Analysis side propellant menu
Putburn	Writes \$BURN
Putcham	Writes \$CHAMBER
Putgeom	Writes \$GEOM
Putinj	Writes \$INJ
Putnls	Writes \$BURN, \$CHAMBER, \$GEOM, and \$INJ
Radbf	Routine which gets radial baffle data
Rcheck	Check variable range
Reactc	Prepare reactant cards for an ODE run
Readit	Reads PROCESS.BIN
Resum	Opens and reads one resume file
Retq	Pauses and waits for a return
Retry	Determine if the user wants to RE-TRY resuming data
Rkam	Adams - Moulant integration scheme
Roccid	Main program

## IFE Modules (Continued)

<u>Routine Name</u>	<u>Purpose</u>
Roccid_Trace.msg	Vax traceback generators for Roccid (message file)
Rplcat	Replicate the first column of an array into all columns
Saidin	Computer written code to read stability aid data
Search	Search through the replay file to find a "name"
Setit	Set a namelist value
Shdcin	Computer written code to read showerhead control data
Shearcin	Computer written code to read swirl coax control data
Stabh	Driving routine for Point Analysis high freq. stability
Stabl	Driving routine for Point Analysis low freq. stability
Stadef	Gets \$\$AID data
Stbaid	Gets stability aid data (Point Analysis)
String	Read a character string
Stuffin	Computer written code to read baffle flags
Swirlcin	Computer written code to read swirl coax flags
Tekdfl	Sets TEK screen dialog area
Tekdia	Turns TEK dialog on
Tekdof	Turns TEK dialog off
Tekmod	Puts TEK terminal in graphics mode
Title	Gets run title
Tripcin	Computer written code to read triplet control
Trivar	Tri-variable interpolation routine
Trmtyp	Ask user for terminal type
Umenu	Utility menu
Varmnu	Point Analysis variable menu
Vax1_auxlib	SEA modified Calcomp library
Vax1_clean	File deletion routine
Vax1_getfil	Reads replay file entries
Vax1_newfil	Adds entries to replay file list
Vax1_seaplib	SEA vax-tektronics plotting drivers

**Part D**  
**ODE Routines**

## ODE Routines

The utilities module of ROCCID contains an ODE option which allows calculation of thermodynamic and transport gas properties for various rocket engine propellants. Complete documentation of the ODE subroutines have been published in "Two Dimensional Kinetics (TDK) Nozzle Performance Computer Program, Volumes I - III" by Nickerson, Dang, and Coats of Software and Engineering Associates, under NASA contract NAS 8-36863 (March 1989). The following describes only subroutines that are used directly with ROCCID.

<u>Routine Name</u>	<u>Purpose</u>
ODEMAIN	Drives the ODE routines used to calculate ISP and C*
ODES	Reads input for calculating ISP and C* and runs the rocket chemical equilibrium analysis
OUT1	ODE output routine (used only in the ODE utility)
RKTOUT	Stores ISP and C* arrays
ROCKET	Calculates ISP and C* using equilibrium gas properties
THERMP	Calculates thermochemical properties of equilibrium gases



**APPENDIX H**  
**ROCCID INSTALLATION INSTRUCTIONS**

## ROCCID Installation Instructions for VAX/VMS Computer Systems

ROCCID is currently configured with a multiple directory structure (Fig. H.1). The main directory contains the executable code (ROCCID.EXE), a linker map (ROCCID.MAP), a command file to link the code (LINK\_ROCCID.COM), object code for the main program (ROCCID.OBJ), the IFE binary control data file (PROCESS.BIN) and four object libraries, (POINTA.OLB, POINTD.OLB, IFE.OLB and ODE.OLB) and the date stamp (VDATE.BIN). It also has 6 subdirectories. The POINTA, POINTD, IFE and ODE subdirectories contain the source code modules for those respective segments of the code, and the object libraries contain the compiled code in the respective subdirectories. The PROPDAT subdirectory contains propellant property and combustion gas data used by all modules of the program. The SAMPLES subdirectory contains sample case input and output files, and the UPDATE directory contains utility code used to generate/update the defaults in the IFE control file, PROCESS.BIN.

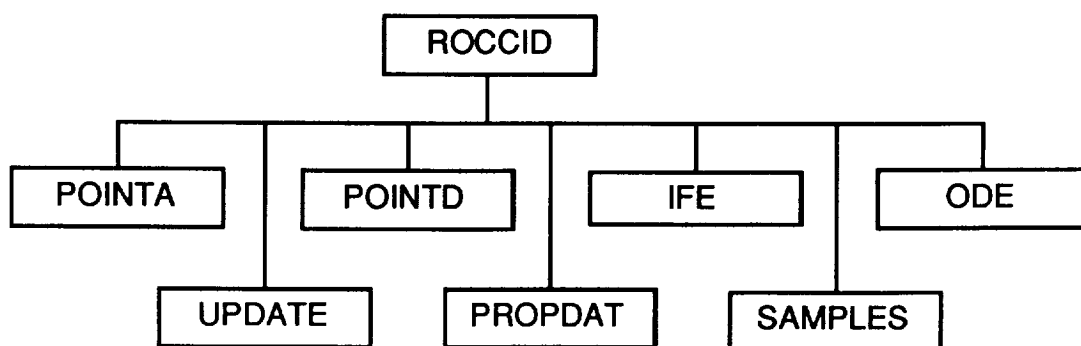
Several modifications must be made to the files to ensure that the code will look in the correct place for files. Table H.1 contains a list of routines and files which must be changed. Data files only need to be changed, but FORTRAN files (file type **.FOR**) must be changed, compiled and replaced in the appropriate libraries prior to running the command file to link ROCCID, LINK\_ROCCID.COM. In order to minimize potential problems during the initial installation, it is recommended that the contents of the all object libraries be deleted and replaced with object code generated under your machine's operating system. The object code for main calling program, ROCCID, which is contained in the IFE subdirectory, should be copied to the main directory and must not be included in the IFE object library (IFE.OLB).

The VAX/VMS message file ROCCID\_TRACE.MSG is contained in the IFE subdirectory. It has been included to facilitate error handling, and it generates a traceback message for controlled aborts in ROCCID. To compile it, use the VAX/VMS message compiler:

Message ROCCID\_Traceback.msg

Then insert the object code (**.OBJ**) into the IFE object library. If you are not running on a VAX/VMS System, you must either A) write a similar routine for your operating system or B) comment - out all references to ROCCID-TRACEB in ABORT.FOR.

Accompanying the ROCCID code in another program called FILGEN (located in subdirectory **[.UPDATE]**), which creates a binary file of variable definitions for the interactive front end. This binary file, named PROCESS.BIN must be read by ROCCID before any analyses can be executed, therefore FILGEN must be run whenever ROCCID is transported to a new computer.



**Figure H-1. ROCCID Directory Structure**

FILGEN reads an ASCII file of variable definitions (**VARIABLES.DAT**) which provides the information contained in **PROCESS.BIN**. This input file can be modified to add new variable names or definitions, or to modify the attributes of current variables. The number of variables and the order of the variables must not change. The list below describes the parameters used to define each variable in **VARIABLES.DAT**:

**01.NAME:** Entry for the variable name (upper case only)

**02.DESCR:** A brief 1 to 2 line variable description. For each additional line use the end-of-line character @, as the line continuation character. The description may not contain the equal symbol (=).

**03.IND-VAR:** The independent variable entry. Indicate if the variable is an independent variable (enter **FLAG**, **COUNTER**, or dependent variable name) or is not (enter **NO**). Typically, this entry is used to describe variable arrays used to enter tables of data, e.g. gas property tables, wall table data, initial condition profiles, etc.

**FLAG** indicates that the variable is a flag

**COUNTER** indicates that this variable is a **COUNTER** for arrays (tables) input.

**YES** indicates that other variables are required inputs in conjunction with this variable.

**NO** indicates that there are no other variables required in addition to this variable.

**04.TYPE:** Defines the variable type; **INTEGER**, **REAL**, OR **LOGICAL**.

**05.DIM:** The length of array variables or **NONE** for scalar quantities.

**06.RANGE:** This sets the variable limits of applicability. General form of this expression is (lower limit:upper limit). The upper and lower values must be integers. Examples of acceptable forms include **(1:2)**, **(T:F)**, **>0**, **<0**, **MONOTONIC INCREASING**, and **MONOTONIC DECREASING**.

- 07.APPLIC:** Describes the general application of variable. Application must match one of the applications listed in Table 1.
- 08.DESTIN:** Name of \$NAMELIST to which this variable belongs.
- 09.DEFAULT:** Default value for this variable. Use .FALSE. or .TRUE. to indicate set defaults for logical type variables.
- 10.SUBSET:** The name of variable(s) which makes this variable a required input. If this condition does not exist enter NONE. See SUPERSET for more details on usage.
- 11.SUPERSET:** Sets condition and variable(s) which are required. Multiple conditions and/or options are specified using the ( ) to enclose condition and required variable(s); and commas to separate multiple sets of conditions and variable requirements. Conditional statements are separated by colons and variable lists are separated by commas. If these conditions do not exist, enter NONE. Typically, the SUPERSET/SUBSET combination is used to describe a variable which flags a variety of options, which require different sets of variable value and corresponding variables to be used in conjunction with the option value; the SUBSET entry describes the backward tract to the option variable. Two examples are given below.

1) A variable array, ID, is read in and is only used if the number of array elements, NID, is specified. For the variable NID, its SUPERSET is the variable ID. For the variable ID, its SUPERSET and SUBSET entries are NONE.

2) An option to specify different wall contours to a nozzle performance program is set by the variable IWALL. IWALL has possible values of 1,3, and 4. The associated variables for IWALL=1 are RWTD, THETA, EPS, RZNORM. For IWALL=3, the required variables to be used are RWTD, THETA, RMAX...

SUPERSET:(IWALL=1:RWTD, THETA, EPS, ZNORM),

(IWALL=3:RWTD, THETA, RMAX, ZMAX,RZNORM),

(IWALL=4:RWTD, THETA, THE, RS, ZS, NWS, RZNORN)

**12.LEVEL:** Defines the usage or path level. Enter 1 for commonly used, 2 for sometimes used, and 3 for rarely usage only.

The FILGEN executable is created by ,compile and link all the FORTRAN files (file type **.FOR**) in the **[.UPDATE]** subdirectory. If necessary, modify the **VARIABLES.DAT** file to reflect any changes to the variable names or definitions. The successful execution of FILGEN requires a driver file **MASTER.DAT**, which must reside in the same subdirectory as FILGEN. After running FILGEN, copy the new **PROCESS.BIN** file into the **[ROCCID]** directory, where ROCCID will read it.

If you have any questions about this installation procedure, please contact Jeff Muss at (916) 355-3663.

**TABLE H.1**  
**ROCCID Files Which Contain Path Statements**

**File Name:** [ROCCID.IFE]AFILER.FOR

```
OPEN(UNIT=12,FILE='DISK$3:[UTILITY.ROCCID]DEFAULT.CNT',
9100 WRITE(6,*) "'DISK$3:[UTILITY.ROCCID]DEFAULT.CNT" NOT FOUND, ',
```

**File Name:** [ROCCID.IFE]DASORC.FOR

```
DATA LIBNAM /DISK$3:[UTILITY.ROCCID.PROPDAT]REACTLIB.DAT/
```

**File Name:** [ROCCID.IFE]DFILER.FOR

```
OPEN(UNIT=13,FILE='DISK$3:[UTILITY.ROCCID]DEFAULT.DEF',
30 OPEN(UNIT=12,FILE='DISK$3:[UTILITY.ROCCID]DEFAULT.CNT',
9100 WRITE(6,*) "'DISK$3:[UTILITY.ROCCID]DEFAULT.CNT" NOT FOUND, ',
9200 Write (6,*)' "DISK$3:[UTILITY.ROCCID]DEFAULT.DEF" NOT FOUND, ',
```

**File Name:** [ROCCID.IFE]DFILES.FOR

```
30 OPEN(UNIT=12,FILE='DISK$3:[UTILITY.ROCCID]DEFAULT.CNT',
9100 WRITE(6,*) "'DISK$3:[UTILITY.ROCCID]DEFAULT.CNT" NOT FOUND, ',
```

**File Name:** [ROCCID.IFE]DPROPMNU.FOR

```
OPEN(UNIT=45, FILE='DISK$3:[UTILITY.ROCCID.PROPDAT]PROP.FIL',
```

**File Name:** [ROCCID.IFE]GETCOM.FOR

```
DATA DIRECT /DISK$3:[UTILITY.ROCCID]COMFIL.DAT/
```

**File Name:** [ROCCID.IFE]PROPMNU.FOR

```
OPEN(UNIT=45, FILE='DISK$3:[UTILITY.ROCCID.PROPDAT]PROP.FIL',
```

**File Name:** [ROCCID.IFE]REACTC.FOR

```
DATA FILNAM /DISK$3:[UTILITY.ROCCID.PROPDAT]REACTLIB.DAT/
```

**File Name:** [ROCCID.IFE]READIT.FOR

```
DATA FILNAM /DISK$3:[UTILITY.ROCCID]PROCESS.BIN/
```

**TABLE H.1 (continued)**

**ROCCID Files Which Contain Path Statements**

**File Name:** [ROCCID.IFE]TITLE.FOR

OPEN(UNIT=91, FILE='DISK\$3:[UTILITY.ROCCID]VDATE.BIN',

**File Name:** [ROCCID.ODE]ODEMAIN.FOR

OPEN(UNIT = 25,FILE='DISK\$3:[UTILITY.ROCCID.PROPDAT]JANNAF.DAT',

**File Name:** [ROCCID.POINTA]FLUIDP.FOR

3 OPEN(UNIT=25,FILE='DISK\$3:[UTILITY.ROCCID.PROPDAT]PH2.COF',  
4 OPEN(UNIT=25,FILE='DISK\$3:[UTILITY.ROCCID.PROPDAT]METH.COF',  
7 OPEN(UNIT=25,FILE='DISK\$3:[UTILITY.ROCCID.PROPDAT]O2.COF',  
9 OPEN(UNIT=25,FILE='DISK\$3:[UTILITY.ROCCID.PROPDAT]C3H8.COF',

**File Name:** [ROCCID.POINTA]PINPUT.FOR

OPEN(UNIT=57,FILE='DISK\$3:[UTILITY.ROCCID]DEFAULT.CNT',

**File Name:** [ROCCID.POINTA]PROPSET.FOR

PROFFILE='DISK\$3:[UTILITY.ROCCID.PROPDAT]//OX(:IOEND)//' \_'//

**File Name:** [ROCCID.PROPDAT]PROP.FIL

DISK\$3:[UTILITY.ROCCID.PROPDAT]LOX\_RP-1.DAT  
DISK\$3:[UTILITY.ROCCID.PROPDAT]LOX\_H2.DAT  
DISK\$3:[UTILITY.ROCCID.PROPDAT]LOX\_METHANE.DAT  
DISK\$3:[UTILITY.ROCCID.PROPDAT]LOX\_PROPANE.DAT

## ROCCID UNIX TRANSPORTING INSTRUCTIONS

### Transporting ROCCID to a UNIX Computer

ROCCID is available for the VAX computer only. If you wish to use ROCCID on any other computer you must transport the code. Although transporting codes typically is not difficult, it requires a strong understanding of both FORTRAN and the FORTRAN compiler on your computer in order to do it efficiently. A short discussion of FORTRAN statements which are likely to cause problems follows.

#### OPEN Statements

The VAX/VMS compiler's OPEN statement arguments are often incompatible with UNIX OPENs. For example, to make a VAX/VMS OPEN statement so the file can only be read (to prevent overwriting the file) the command is

```
OPEN(...STATUS='OLD',READONLY....)
```

The same statement on a SUN/UNIX is

```
OPEN(...STATUS='READONLY'....).
```

Another likely syntax problem on the OPEN statements is the VAX/VMS CARRIAGE CONTROL statement which will probably have to be deleted.

To ensure the OPEN statements are compatible with your compiler, first consult your FORTRAN manual to find the correct formatting, then "grep" through the ROCCID source code and check each OPEN statement in the source.

Along with the OPEN statements the file paths (sub-directory tree) must also be changed to a UNIX format. An improper path is the most likely cause if the code cannot open a file.

## REAL\*16 Declarations

One subroutine in ROCCID uses two REAL\*16 variables, which will not be supported on most UNIX computers. The existing SUN version of ROCCID runs fine with the variables declared as REAL\*8, so that fix will probably work on other UNIX machines. These variables are found in the function CPI in the FLUIDP.FOR file.

## DATA Statements

Another common transporting problem is the setting of hollerith variables in DATA statements. For example, a VAX/VMS DATA statement of

```
DATA  BS/'08'X/
```

reverses to

```
DATA  BS/X'08'/
```

on the SUN computer.

## READ and WRITE Statements

READ and WRITE statements can also require modification to work on different computers. For example, a SGI computer requires a NAMELIST read to be formatted

```
READ(10,NML=CRPC)
```

ROCCID currently is written as

```
READ(10, CRPC).
```

## GRAPHICS

Another problem in transporting the ROCCID code is the graphics. Since transporting the graphics to different computers is a BIG problem, it is best to comment out all graphics subroutines and just run the code without graphics. The following is a list of subroutines

which should be commented out:

```
PLOTM
TEKMOD
TEKDIA
FIG*
ERASE
ANSI
```

The interactive front end of ROCCID was developed on a SUN computer, therefore, most of the source should transport to most UNIX computers relatively easily. An expert programmer should be able to transport the code within one man-week.



**APPENDIX I**  
**SAMPLE CASE OUTPUT**

**Part A**  
**POINTA Sample Case ACASE1**

ROCCID  
Rocket Combuster Interactive Design Methodology  
Version 23-FEB-81

DIRECT INPUT ECHO FROM SUBROUTINE DINPUT.

ROCCID POINT DESIGN TEST CASE 1  
LOX/RP-1 LIKE DOUBLET PAIR WITH FIXED PC  
APPROXIMATES -0100 SUBSCALE DOUBLET

\$MODELS

1  
2  
1  
\$END

\$DESIGN

PISA= 8.9800E+01, 1.0940E+02, 1.3940E+02, 1.5880E+02, 1.7780E+02,  
2.0870E+02, 2.3550E+02, 2.5470E+02, 2.6700E+02, 2.7040E+02,  
2.7180E+02, 2.7180E+02, 2.7130E+02, 2.8980E+02, 2.8820E+02,  
2.8880E+02, 2.6500E+02, 2.8210E+02, 2.5920E+02, 2.5880E+02,  
2.8980E+02, 2.3940E+02, 2.2760E+02, 2.1100E+02, 2.0140E+02,  
1.5560E+02, 1.0870E+02, 8.0000E+01,  
PCSA= 1.7780E+03, 2.4580E+03, 3.4670E+03, 4.0740E+03, 4.5800E+03,  
5.0700E+03, 5.4540E+03, 5.7340E+03, 5.9080E+03, 5.9810E+03,  
5.9880E+03, 5.9540E+03, 5.9340E+03, 5.9090E+03, 5.8820E+03,  
5.8550E+03, 5.8280E+03, 5.7590E+03, 5.7000E+03, 5.6430E+03,  
5.5220E+03, 5.2780E+03, 5.0470E+03, 4.8760E+03, 4.4200E+03,  
3.4570E+03, 2.4940E+03, 1.4650E+03,  
ETANOZ= 8.8200E-01, 8.3000E-01, 8.0900E-01, 7.5000E-01, 8.1200E-01,  
8.8300E-01, 8.8000E-01, 9.2200E-01, 9.3900E-01, 9.5200E-01,  
9.8200E-01, 9.7000E-01, 9.7700E-01, 9.8100E-01,

'LOL'

'RP-1'

'LOX'

'LOX'

XMR= 2.8800E+00

FTMAN= 7.1000E+01

XTMAN= -2.7800E+02

PCNOM= 2.1180E+03

WONOM= 1.7838E+02

WDMIN= 1.2894E+02

ETACSG= 9.8880E-01

XLEMAX= 4.0000E+00

DOMAX= 7.5000E-01

28

NPERFP=

PMRA=

1.0000E-01, 9.0000E-01, 9.0000E-01, 8.0000E-01, 1.0000E+00,  
1.2600E+00, 1.5000E+00, 1.7600E+00, 2.0000E+00, 2.2000E+00,  
2.4000E+00, 2.5000E+00, 2.6000E+00, 2.7000E+00, 2.8000E+00,  
2.8000E+00, 3.0000E+00, 3.2000E+00, 3.4000E+00, 3.6000E+00,  
4.0000E+00, 5.0000E+00, 6.0000E+00, 8.0000E+00, 1.0000E+01,  
1.5000E+01, 2.0000E+01, 5.0000E+01,

14

NNOZ=

XNOZ=

9.6300E-01, 7.2700E-01, 1.4630E+00, 2.1800E+00, 2.9070E+00,  
9.8300E+00, 4.3600E+00, 5.0870E+00, 5.8130E+00, 6.8770E+00,  
7.2670E+00, 7.9830E+00, 8.7170E+00, 9.0630E+00,

\$END

\$FGEOM

RNE=

RTE=

ALPHA=

1.1100E-01

1.1100E-01

9.0000E+01

```

$END
$FDORG
NZON=
FTER= 20*1.0000E+00
$END
$CONTROL
DPPCS= 2.3000E-01
RNR= 1.0000E+00
RTR= 1.0000E+00
CONVA= 2.6000E+01
XMSTAB= 2.0000E-01
$END
$LOLC
FCD= 9.1000E-01
XCD= 9.4000E-01
FIA= 3.0000E+01
XIA= 3.0000E+01
FIHOD= 2.0000E+00
XIHOD= 2.0000E+00
FCANT= 1.6000E+01
XCANT= 1.6000E+01
FLOD= 5.0000E+00
XLOD= 5.0000E+00
EMJN1= 6.5000E-01
$END
$SAID
TBAF= 3.3000E-02
$END

```

END OF INPUT ECHO

# POINTD DESIGN MODEL INPUTS

## RUN DESCRIPTORS

ROCCID POINT DESIGN TEST CASE 1  
 LOX/RP-1 LIKE DOUBLET PAIR WITH FIXED PC  
 APPROXIMATES .0100 SUBSCALE DOUBLET

## SELECTED STABILITY MODELS

BURNING MODEL=N-TAU INJECTION MODEL=INJ CHAMBER MODEL=HIFI  
 DEBUG OUTPUT=F

## PROPELLENT DESCRIPTION

FUEL=RP-1 Tman., F= 71.00  
 OX=LOX Tman., F=279.00

## OPERATING CONSTRAINTS

CORE ELEMENT TYPE=LOL INJECTED MR= 2.8500  
 NOMINAL PC=2118.00 PSIA  
 NOMINAL FLOWRATE= 1.793E+02 Lbm/s THROTTLED FLOWRATE= 1.299E+02 Lbm/s  
 MAXIMUM CHAMBER DIAMETER= 0.7500 FT MAXIMUM ENGINE LENGTH= 4.0000 FT  
 C\* EFFICIENCY GOAL= 95.88 %

## STABILITY AID PREFERENCE

NO BAFFLES NO CAVITIES

## FIXED CHAMBER GEOMETRY

NOZZLE ENTRANCE RADIUS OF CURVATURE = 0.1110 FT  
 THROAT ENTRANCE RADIUS OF CURVATURE = 0.1110 FT  
 CONVERGENCE HALF-ANGLE =30.0000 DEGREES

## DESIGN CONTROL PARAMETERS

DELTA-P/PC AT PCMIN= 0.2300  
 NON-DIMENSIONAL NOZZLE ENTRANCE RADIUS OF CURVATURE (RNE/RTHRT)= 1.0000  
 NON-DIMENSIONAL THROAT ENTRANCE RADIUS OF CURVATURE (RTE/RTHRT)= 1.0000  
 NOZZLE CONVERGENCE HALF-ANGLE=25.0000 DEGREES

## LIKE DOUBLET PAIR DESIGN VARIABLES

FUEL Cd= 0.9100 OX Cd= 0.9400 UNIELEMENT Em= 0.6500  
 FUEL IMPINGEMENT HALF-ANGLE=30.0000 DEG  
 OX IMPINGEMENT HALF-ANGLE=30.0000 DEG  
 FUEL IMPINGEMENT POINT HEIGHT TO ORIFICE DIAMETER RATIO= 2.0000  
 OX IMPINGEMENT POINT HEIGHT TO ORIFICE DIAMETER RATIO= 2.0000  
 FUEL UNLIKE CANT ANGLE=18.0000 DEG  
 OX UNLIKE CANT ANGLE=18.0000 DEG  
 FUEL ORIFICE LENGTH TO ORIFICE DIAMETER RATIO= 5.0000

OX ORIFICE LENGTH TO ORIFICE DIAMETER RATIO= 5.0000

STABILITY AID DESIGN VARIABLES

BAFFLE BLADE THICKNESS= 0.0330 IN  
1/4 WAVE CAVITY PARTITION THICKNESS= 0.0000 FT  
ACOUSTIC CAVITY TYPE 1 INLET=SHARP EDGED  
ACOUSTIC CAVITY TYPE 2 INLET=SHARP EDGED

# PRELIMINARY DESIGN CHAMBER SIZING RESULTS

NOMINAL CHAMBER PRESSURE = 2.110E+03 PSIA  
 THROTTLED CHAMBER PRESSURE = 1.886E+03 PSIA  
 FUEL MANIFOLD PRESSURE = 2.780E+03 PSIA  
 OXIDIZER MANIFOLD PRESSURE = 2.780E+03 PSIA  
 CHAMBER RADIUS = 2.786E-01 FT  
 THROAT RADIUS = 1.862E-01 FT  
 NOZZLE ENTRANCE RADIUS OF CURVATURE = 1.110E-01 FT  
 THROAT ENTRANCE RADIUS OF CURVATURE = 1.110E-01 FT  
 NOZZLE CONVERGENCE HALF-ANGLE = 3.000E+01 DEG  
 INJECTOR-TO-THROAT CHAMBER LENGTH = 1.183E+00 FT  
 BARREL SECTION LENGTH = 9.601E-01 FT

## IMPINGING ELEMENT SIZING RESULTS

ELEMENT TYPE = LOL  
 NO. OF ELEMENTS = 93  
 FUEL 20% VAPORIZATION LENGTH = 6.360E-01 IN  
 OX 20% VAPORIZATION LENGTH = 5.363E-01 IN  
 FUEL ORIFICE DIAMETER = 5.328E-02 IN  
 OX ORIFICE DIAMETER = 8.171E-02 IN

## CORE ELEMENT SPACING RESULTS

ELEMENT TYPE = LOL  
 NUMBER OF ELEMENTS = 93  
 FUEL ORIFICE/ANNULUS DIAMETER = 5.328E-02 IN  
 OXIDIZER ORIFICE DIAMETER = 8.171E-02 IN  
 FUEL INJECTION VELOCITY = 3.633E+02 FT/S  
 OXIDIZER INJECTION VELOCITY = 2.978E+02 FT/S

ROW	# ELEMENTS	MID-ROW RADIUS (IN)
1	6	8.316E-01
2	12	1.382E+00
3	18	1.953E+00
4	24	2.513E+00
5	30	3.074E+00

ROCKET COMBUSTOR INTERACTIVE DESIGN METHODOLOGY  
Version 23-FEB-91

I-8

XIH=	1.4182E-01	
XIA=	3.0000E+01	
XCANT=	1.6000E+01	
XFACET=	3.6380E-01	
EMUNI=	8.5000E-01	
SEND		
\$BAFFLE		
SEND		
\$BARRIER		
SEND		
\$FFC		
SEND		
\$BURN		
SEND		
\$INJ		
FMAND=	6.7076E+00	
XMAND=	6.7076E+00	
FMANL=	3.3636E+00	
XMANL=	3.3636E+00	
PCA=	2.1180E+03, 1.5347E+03, 7.8734E+02	
SEND		
\$CHAMBER		
MJB=	3	
T=	3.3000E-02	
SEND		
\$FDORG		
NZON=	5	
FTER=	20*1.0000E+00	
SEND		
\$MIX		
EM(1)=	9.180E-01, 1.000E+00	
SEND		
\$DEBUGG		
DEBUG=	F	
SEND		
\$HIFIG		
SHORT=	F	
POG=	2.0000E-01, 2.0000E-01	
SEND		
\$DISTSDC		
SHORT=	F	
POG=	2.0000E-01, 2.0000E-01	
PAMP=	2.0000E-01	
MC=	11	
LC=	8	
MB=	11	
LB=	8	
IDMAX=	10	
SEND		
\$CRPG		
NOPO=	16	
NOTFQ=	16	
NOTLF=	1000	
NPRINT=	50	
NSLMS=	3500	
PAMPO=	2.0000E-01	
NRIAD=	3	
NCIRO=	5	
SEND		
\$LEINJC		
IDOMEN=	2	
PAMPCH=	2.0000E-01	
NTINJ=	16	
USGF=	2.0000E-02	

OGF=	2.000E-02				
DGGF=	2.000E-02				
FGDO=	1.000E+00,	1.000E+00,	1.000E+00,	1.000E+00,	1.000E+00
XCDO=	1.000E+00,	1.000E+00,	1.000E+00,	1.000E+00,	1.000E+00
\$END					
\$COMBUSTC					
FALVM=	1.000E+00,	1.000E+00,	1.000E+00,	1.000E+00,	1.000E+00
XALVM=	1.000E+00,	1.000E+00,	1.000E+00,	1.000E+00,	1.000E+00
FALTM=	1.000E+00,	1.000E+00,	1.000E+00,	1.000E+00,	1.000E+00
XALTM=	1.000E+00,	1.000E+00,	1.000E+00,	1.000E+00,	1.000E+00
FNAM=	1.000E+00,	1.000E+00,	1.000E+00,	1.000E+00,	1.000E+00
XNAM=	1.000E+00,	1.000E+00,	1.000E+00,	1.000E+00,	1.000E+00
ENAMULT=	1.000E+00,	1.000E+00,	1.000E+00,	1.000E+00,	1.000E+00
COMBXB=	5.000E-01				
ACMULT=	1.000E+00				
COMULT=	1.000E+00				
ENAMULT=	1.000E+00				
TAUMULT=	1.000E+00				
\$END					
\$FORCC					
SHORT=	F				
EPSIL=	2.000E+01				
ERROR=	1.000E-01				
LTS=	5				
MTS=	5				
NTS=	10				
ITMAX=	100				
RELX=	6.500E-01				
NEET=	3				
NXFST=	3				
MYFST=	3				
NAFST=	3				
MORE=	F				
\$END					

END OF INPUT ECHO

# STEADY STATE COMBUSTION ANALYSIS PROGRAM

## RUN DESCRIPTORS

ROCCID POINT DESIGN TEST CASE 1  
LOX/RP-1 LIKE DOUBLET PAIR WITH FIXED PC  
APPROXIMATES .0100 SUBSCALE DOUBLET

## PROPELLANT DESCRIPTION

FUEL=RP-1 Tman., F= 71.00  
OX=LOX Tman., F=-279.00

## CHAMBER GEOMETRY

CHAMBER RADIUS = 3.3639 IN.  
CYLINDRICAL SECTION =11.5206 IN.  
NOZZLE ENTRANCE RADIUS OF CURVATURE = 1.3320 IN.  
CONVERGENCE HALF-ANGLE =30.0000 DEG.  
THROAT RADIUS = 2.2228 IN.  
CONVERGENT SECTION LENGTH = 2.6790 IN.  
THROAT ENTRANCE RADIUS OF CURVATURE = 1.3320 IN.  
CONTRACTION RATIO = 2.28

## INJECTOR DATA

INJECTOR CORE CONTAINS 93 LOL ELEMENTS

FUEL SIDE: Orifice Diam. =5.329E-02 In.  
Impingement Half-angle =30.00 Deg.  
OX SIDE: Orifice Diam. =8.170E-02 In.  
Impingement Half-angle =30.00 Deg.  
Cd =0.9100  
Unlike Cant Angle =16.00 Deg.  
Cd =0.9400  
Unlike Cant Angle =16.00 Deg.  
Impingement Height =0.092 In.  
Faceplate Thickness = 0.2307 In.  
Impingement Height =0.142 In.  
Faceplate Thickness = 0.3638 In.

## MIXING EFFICIENCIES

CORE MIXING EFFICIENCY=0.9180 BARRIER MIXING EFFICIENCY=1.0000

## COMBUST CONTROL PARAMETERS

MULTIPLIERS:  
FUEL ATOMIZATION LENGTH FOR VAPORIZATION: CORE 1.000 BAFFLE 1.000 BARRIER 1.000 FFC 1.000  
OX ATOMIZATION LENGTH FOR VAPORIZATION: 1.000 1.000 1.000 1.000  
FUEL ATOMIZATION LENGTH FOR TIMELAGS: 1.000 1.000 1.000 1.000  
OX ATOMIZATION LENGTH FOR TIMELAGS: 1.000 1.000 1.000 1.000  
FUEL DROPSIZE: 1.000 1.000 1.000 1.000  
OX DROPSIZE: 1.000 1.000 1.000 1.000  
MIXING (Em): 1.000 1.000 1.000 1.000  
AO-Multiplier=1.000 CO-Multiplier=1.000 N-Multiplier=1.000 Tau-Multiplier=1.000  
Eta-C\* for XB=0.500

BEGIN STEADY STATE COMBUSTION ANALYSIS  
PC=2118.00 PSIA

PROPELLANT PROPERTIES

FUEL=RP-1	Phase=Liquid	Tman.. F= 71.00	
	Injected Density= 49.69 Lbm/Cu. Ft	Viscosity=1.360E-03 Lbm/Ft-s	Surface Tension=1.667E-03 Lbf/Ft
OX=LOX	Phase=Liquid	Tman.. F=-279.00	
	Injected Density= 70.18 Lbm/Cu. Ft	Viscosity=1.166E-04 Lbm/Ft-s	Surface Tension=7.326E-04 Lbf/Ft

OPERATING CONDITIONS

PC FACE=2118.00 PSIA PC THROAT=2034.15 MIXTURE RATIO= 2.660  
 FUEL INJECTION PRESSURE DROP= 633.67 Psia FUEL INJECTION VELOCITY= 343.06 Ft/s  
 OX INJECTION PRESSURE DROP= 633.67 Psia OX INJECTION VELOCITY= 289.25 Ft/s

FUEL FLOWRATE= 44.864 OX FLOWRATE= 129.296

ATOMIZATION OUTPUT

DROPSIZE MODEL=AEROJET

ELEMENT TYPE 1 IS LOL			
FUEL:	ATOMIZATION LENGTH, In.=0.64239	ATOMIZATION LENGTH FOR VAPORIZATION, In.=0.64239	DROPLET RADIUS, Microns= 61.99
OX:	ATOMIZATION LENGTH, In.=1.64611	ATOMIZATION LENGTH FOR VAPORIZATION, In.=1.64611	DROPLET RADIUS, Microns= 68.69

# VAPORIZATION CALCULATIONS

X (In.)	CORE-LOL		BAFFLE-		BARRIER-		FFG-	
	%FUEL VAP	%OX VAP	%FUEL VAP	%OX VAP	%FUEL VAP	%OX VAP	%FUEL VAP	%OX VAP
0.0000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
0.2839	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
0.5677	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
0.8516	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.1355	10.421	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.4194	22.348	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.7032	31.303	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.9871	38.802	0.000	0.000	0.000	0.000	0.000	0.000	0.000
2.2710	44.860	31.536	0.000	0.000	0.000	0.000	0.000	0.000
2.5548	49.928	48.782	0.000	0.000	0.000	0.000	0.000	0.000
2.8387	54.297	60.033	0.000	0.000	0.000	0.000	0.000	0.000
3.1226	58.099	67.871	0.000	0.000	0.000	0.000	0.000	0.000
3.4065	61.462	73.310	0.000	0.000	0.000	0.000	0.000	0.000
3.6903	64.386	77.993	0.000	0.000	0.000	0.000	0.000	0.000
3.9742	67.004	81.238	0.000	0.000	0.000	0.000	0.000	0.000
4.2581	69.246	83.986	0.000	0.000	0.000	0.000	0.000	0.000
4.5420	71.372	86.236	0.000	0.000	0.000	0.000	0.000	0.000
4.8258	73.322	88.311	0.000	0.000	0.000	0.000	0.000	0.000
5.1097	75.154	89.847	0.000	0.000	0.000	0.000	0.000	0.000
5.3936	76.762	91.126	0.000	0.000	0.000	0.000	0.000	0.000
5.6774	78.256	92.243	0.000	0.000	0.000	0.000	0.000	0.000
5.9613	79.475	93.184	0.000	0.000	0.000	0.000	0.000	0.000
6.2452	80.681	94.018	0.000	0.000	0.000	0.000	0.000	0.000
6.5291	81.881	94.887	0.000	0.000	0.000	0.000	0.000	0.000
6.8129	82.956	95.356	0.000	0.000	0.000	0.000	0.000	0.000
7.0968	83.785	95.918	0.000	0.000	0.000	0.000	0.000	0.000
7.3807	84.613	96.414	0.000	0.000	0.000	0.000	0.000	0.000
7.6645	85.442	96.816	0.000	0.000	0.000	0.000	0.000	0.000
7.9484	86.270	97.211	0.000	0.000	0.000	0.000	0.000	0.000
8.2323	87.099	97.478	0.000	0.000	0.000	0.000	0.000	0.000
8.5162	87.919	97.745	0.000	0.000	0.000	0.000	0.000	0.000
8.8000	88.478	98.012	0.000	0.000	0.000	0.000	0.000	0.000
9.0839	88.040	98.280	0.000	0.000	0.000	0.000	0.000	0.000
9.3678	88.600	98.460	0.000	0.000	0.000	0.000	0.000	0.000
9.6516	90.160	98.626	0.000	0.000	0.000	0.000	0.000	0.000
9.9355	90.608	98.781	0.000	0.000	0.000	0.000	0.000	0.000
10.2194	91.046	98.924	0.000	0.000	0.000	0.000	0.000	0.000
10.5033	91.466	99.066	0.000	0.000	0.000	0.000	0.000	0.000
10.7871	91.824	99.148	0.000	0.000	0.000	0.000	0.000	0.000
11.0710	92.288	99.242	0.000	0.000	0.000	0.000	0.000	0.000
11.3548	92.593	99.336	0.000	0.000	0.000	0.000	0.000	0.000
11.6388	92.934	99.419	0.000	0.000	0.000	0.000	0.000	0.000
11.9226	93.298	99.478	0.000	0.000	0.000	0.000	0.000	0.000
12.2065	93.681	99.538	0.000	0.000	0.000	0.000	0.000	0.000
12.4904	94.018	99.599	0.000	0.000	0.000	0.000	0.000	0.000
12.7742	94.291	99.658	0.000	0.000	0.000	0.000	0.000	0.000
13.0581	94.683	99.677	0.000	0.000	0.000	0.000	0.000	0.000
13.3420	94.836	99.716	0.000	0.000	0.000	0.000	0.000	0.000
13.6259	95.109	99.747	0.000	0.000	0.000	0.000	0.000	0.000
13.9097	95.381	99.774	0.000	0.000	0.000	0.000	0.000	0.000
14.1936	95.623	99.801	0.000	0.000	0.000	0.000	0.000	0.000

OVERALL VAPORIZATION EFFICIENCIES  
 FUEL= 96.62% OX= 99.80%

MASS DISTRIBUTION PROFILE						
X (IN)	CORE (lbm/s)		BARRIER (lbm/s)		LOCAL VAPOR MIXTURE RATIO	ETA-C*
	FUEL	OX	FUEL	OX		
0.0000	0.000	0.000	0.000	0.000	0.00	0.0000
0.2039	0.000	0.000	0.000	0.000	0.00	0.0000
0.5077	0.000	0.000	0.000	0.000	0.00	0.0000
0.8510	0.000	0.000	0.000	0.000	0.00	0.0000
1.1355	4.878	0.000	0.000	0.000	0.00	0.0086
1.4184	10.033	0.000	0.000	0.000	0.00	0.0141
1.7032	14.053	0.000	0.000	0.000	0.00	0.0198
1.9871	17.465	0.000	0.000	0.000	0.00	0.0246
2.2710	20.138	40.775	0.000	0.000	2.02	0.3497
2.5548	22.415	63.048	0.000	0.000	2.81	0.4901
2.8387	24.377	77.620	0.000	0.000	3.18	0.5750
3.1226	26.083	87.386	0.000	0.000	3.35	0.6344
3.4065	27.593	94.786	0.000	0.000	3.44	0.6814
3.6903	28.906	100.765	0.000	0.000	3.49	0.7200
3.9742	30.081	105.036	0.000	0.000	3.48	0.7500
4.2581	31.066	108.563	0.000	0.000	3.48	0.7752
4.5420	32.042	111.500	0.000	0.000	3.48	0.7973
4.8258	32.917	114.182	0.000	0.000	3.47	0.8176
5.1097	33.740	116.189	0.000	0.000	3.44	0.8343
5.3936	34.462	117.623	0.000	0.000	3.42	0.8488
5.6774	35.133	119.287	0.000	0.000	3.39	0.8616
5.9613	35.680	120.419	0.000	0.000	3.37	0.8718
6.2452	36.221	121.562	0.000	0.000	3.36	0.8821
6.5291	36.747	122.426	0.000	0.000	3.33	0.8909
6.8129	37.243	123.290	0.000	0.000	3.31	0.8995
7.0968	37.615	124.016	0.000	0.000	3.30	0.9062
7.3807	37.987	124.680	0.000	0.000	3.28	0.9126
7.6645	38.359	125.178	0.000	0.000	3.26	0.9184
7.9484	38.731	125.889	0.000	0.000	3.25	0.9242
8.2323	39.103	126.036	0.000	0.000	3.22	0.9292
8.5162	39.471	126.380	0.000	0.000	3.20	0.9342
8.8000	39.722	126.726	0.000	0.000	3.19	0.9391
9.0839	39.974	127.046	0.000	0.000	3.18	0.9418
9.3678	40.225	127.305	0.000	0.000	3.16	0.9453
9.6516	40.477	127.518	0.000	0.000	3.15	0.9486
9.9355	40.678	127.719	0.000	0.000	3.14	0.9514
10.2194	40.878	127.905	0.000	0.000	3.13	0.9540
10.5033	41.072	128.075	0.000	0.000	3.12	0.9566
10.7871	41.268	128.185	0.000	0.000	3.11	0.9590
11.0710	41.423	128.316	0.000	0.000	3.10	0.9609
11.3548	41.589	128.437	0.000	0.000	3.09	0.9628
11.6386	41.722	128.545	0.000	0.000	3.08	0.9647
11.9226	41.895	128.623	0.000	0.000	3.07	0.9665
12.2065	42.048	128.700	0.000	0.000	3.06	0.9684
12.4904	42.208	128.777	0.000	0.000	3.06	0.9702
12.7742	42.381	128.828	0.000	0.000	3.04	0.9715
13.0581	42.464	128.878	0.000	0.000	3.04	0.9728
13.3420	42.576	128.928	0.000	0.000	3.03	0.9741
13.6259	42.698	128.988	0.000	0.000	3.02	0.9754
13.9097	42.821	129.004	0.000	0.000	3.01	0.9767
14.1936	42.928	129.036	0.000	0.000	3.01	0.9778

# AXIAL PRESSURE PROFILE

X (in)	MACH #	Ptotal (psia)	Pstatic (psia)	Ttotal (R)	Tstatic (R)	Wdot (Lbm/s)	Local Radius (in)
0.91	0.000	2118.74	2119.89	1685.15	1689.71	0.58	3.354
1.21	0.004	2118.72	2119.85	1691.89	1688.48	8.00	3.354
1.51	0.009	2118.66	2119.51	1770.23	1784.55	12.10	3.354
1.81	0.009	2118.65	2119.51	1691.89	1688.48	12.49	3.354
2.11	0.036	2117.17	2118.54	3559.19	3547.55	34.80	3.354
2.42	0.113	2109.63	2089.35	6661.31	6635.43	76.31	3.354
2.72	0.142	2095.15	2072.25	6617.73	6780.31	95.34	3.354
3.02	0.165	2087.07	2055.88	6661.31	6832.01	109.80	3.354
3.32	0.181	2080.97	2043.48	6635.66	6805.42	119.88	3.354
3.62	0.194	2075.45	2032.23	6629.70	6788.10	128.14	3.354
3.93	0.205	2071.07	2023.28	6628.95	6784.41	134.28	3.354
4.23	0.213	2067.37	2015.71	6626.97	6783.63	139.20	3.354
4.53	0.220	2064.09	2008.99	6628.61	6784.74	143.38	3.354
4.83	0.226	2061.03	2002.70	6630.60	6785.83	147.17	3.354
5.13	0.232	2058.54	1997.58	6634.78	6789.38	150.12	3.354
5.44	0.236	2056.39	1993.15	6638.55	6802.69	152.82	3.354
5.74	0.240	2054.49	1989.23	6641.91	6805.59	154.76	3.354
6.04	0.243	2052.90	1985.95	6643.83	6807.14	156.57	3.354
6.34	0.246	2051.36	1982.78	6646.05	6808.98	158.28	3.354
6.64	0.249	2050.02	1980.00	6648.64	6811.25	159.73	3.354
6.95	0.251	2048.78	1977.44	6650.48	6812.79	161.07	3.354
7.25	0.253	2047.75	1975.29	6651.91	6813.97	162.18	3.354
7.55	0.255	2046.80	1973.33	6653.77	6815.60	163.18	3.354
7.85	0.257	2045.89	1971.45	6655.70	6817.30	164.13	3.354
8.15	0.258	2045.10	1969.81	6657.87	6819.27	164.94	3.354
8.46	0.260	2044.35	1968.25	6660.35	6821.55	165.71	3.354
8.76	0.261	2043.72	1966.84	6660.90	6821.95	166.38	3.354
9.06	0.262	2043.14	1965.73	6660.90	6821.81	166.97	3.354
9.36	0.263	2042.61	1964.64	6660.94	6821.73	167.52	3.354
9.66	0.264	2042.13	1963.65	6660.97	6821.65	168.01	3.354
9.97	0.265	2041.72	1962.79	6661.00	6821.58	168.44	3.354
10.27	0.266	2041.39	1961.97	6661.03	6821.52	168.84	3.354
10.57	0.266	2040.98	1961.20	6661.08	6821.48	169.23	3.354
10.87	0.267	2040.64	1960.54	6661.08	6821.41	169.55	3.354
11.17	0.268	2040.36	1959.93	6661.10	6821.35	169.84	3.354
11.46	0.268	2040.08	1959.38	6661.12	6821.32	170.12	3.354
11.76	0.273	2039.81	1958.27	6661.15	6820.20	170.39	3.330
12.06	0.292	2039.54	1943.78	6661.17	6818.11	170.64	3.228
12.36	0.329	2039.20	1918.54	6661.20	6804.67	170.90	3.064
12.66	0.378	2038.86	1881.29	6661.22	6788.68	171.11	2.888
12.96	0.440	2038.45	1827.89	6661.24	6765.25	171.29	2.714
13.29	0.528	2037.88	1744.80	6661.26	6727.87	171.47	2.540
13.59	0.657	2037.05	1599.18	6661.28	6667.81	171.65	2.368
13.89	0.821	2035.80	1399.45	6661.30	6582.51	171.81	2.257
14.19	0.998	2034.15	1174.85	6661.31	6419.78	171.97	2.223

# PERFORMANCE SUMMARY

## C\* EFFICIENCY CALCULATIONS (ODK)

INJECTED MR= 2.8800 CSTAR=5880.40 CORE Em=9.9168 BARRIER Em=1.0000 MASS FRACTION= 1.0000  
 CORE: OVERALL MR= 2.8800 VAPOR MR= 9.9999 CSTAR-MIX=5804.37 MASS FRACTION= 0.0000  
 BARRIER: OVERALL MR= 0.0000 VAPOR MR=99.9999 CSTAR-MIX= 0.00  
 ENGINE: OVERALL MR= 2.8800 VAPOR MR= 9.9999 CSTAR-DEL=5730.31  
 C\* EFFICIENCY = 9.776E-01

## ISP EFFICIENCY CALCULATIONS

ISP-ODK, INJ = 2.888E+02 SEC.  
 ISP-ODK, M.Z. INJ = 2.881E+02 SEC.  
 VAPORIZATION EFFICIENCY = 9.819E-01  
 ENERGY RELEASE EFFICIENCY = 9.786E-01  
 ISP-ODK, M.Z. VAPOR = 2.848E+02 SEC.  
 MIXING EFFICIENCY = 9.988E-01

NOTE: ISP-DEL = ISP-ODK, INJ. \* ERE \* ETADIV \* DELISP-BL

# TIME-LAG CALCULATIONS, Milliseconds

Cohem, In.=0.104E-03      FUEL Cohem, In.=2.293E+02      OX Cohem, In.=0.104E+01

## ELEMENT 1 IS TYPE-LOL

FUEL:      Cinj, In.=1.100E-02      Lvap, In.= 0.422      ATOMIZATION LENGTH USED, In.= 0.424E-01  
             Timp=2.500E-02      Tatom=2.040E-01      Tvap=1.025E-01      Total=3.330E-01

OX:      Cinj, In.=1.100E-02      Lvap, In.= 0.154      ATOMIZATION LENGTH USED, In.= 1.040E+00  
             Timp=4.700E-02      Tatom=5.310E-01      Tvap=4.433E-02      Total=6.233E-01

## EFFECTIVE TIMELAGS, Milliseconds

FUEL:      Cinj, In.=1.100E-02      Lvap, In.= 0.422      Total=3.330E-01  
             Timp=2.500E-02      Tatom=2.040E-01      Tvap=1.025E-01

OX:      Cinj, In.=1.100E-02      Lvap, In.= 0.154      Total=6.233E-01  
             Timp=4.700E-02      Tatom=5.310E-01      Tvap=4.433E-02

# CHAMBER-NOZZLE OPTIMIZATION RESULTS

CHAMBER LENGTH (FEET)	ETA-C*	ETA-NOZ	OVERALL EFFICIENCY
0.0000	0.0000	0.8807	0.0000
0.1667	0.0393	0.8726	0.0343
0.3333	0.7523	0.8646	0.8505
0.5000	0.8732	0.8537	0.7454
0.6667	0.9251	0.8419	0.7789
0.8333	0.9520	0.8302	0.7804
1.0000	0.9670	0.8185	0.7915
1.1667	0.9770	0.8067	0.7872
1.3333	0.9850	0.7915	0.7796
1.5000	0.9929	0.7773	0.7718
1.6667	1.0000	0.7631	0.7631
1.8333	1.0000	0.7485	0.7485
2.0000	1.0000	0.7299	0.7299
2.1667	1.0000	0.7114	0.7114
2.3333	1.0000	0.6926	0.6926
2.5000	1.0000	0.6742	0.6742
2.6667	1.0000	0.6461	0.6461
2.8333	1.0000	0.6142	0.6142
3.0000	1.0000	0.5823	0.5823
3.1667	1.0000	0.5504	0.5504
3.3333	1.0000	0.5071	0.5071
3.5000	1.0000	0.4439	0.4439
3.6667	1.0000	0.3808	0.3808
3.8333	1.0000	0.3176	0.3176
4.0000	1.0000	0.2544	0.2544

OPTIMUM CHAMBER LENGTH= 1.0000 FT  
 MAXIMUM OVERALL EFFICIENCY= 0.7915

# REDESIGNED CHAMBER RESULTS

NOMINAL CHAMBER PRESSURE = 2.118E+03 PSIA  
 THROTTLED CHAMBER PRESSURE = 1.535E+03 PSIA  
 FUEL INJECTION PRESSURE DROP = 6.337E+02 PSI  
 OX INJECTION PRESSURE DROP = 6.337E+02 PSI  
 CHAMBER RADIUS = 2.785E-01 FT  
 THROAT RADIUS = 1.852E-01 FT  
 NOZZLE ENTRANCE RADIUS OF CURVATURE = 1.110E-01 FT  
 THROAT ENTRANCE RADIUS OF CURVATURE = 1.110E-01 FT  
 NOZZLE CONVERGENCE HALF-ANGLE = 3.000E+01 DEG  
 INJECTOR-TO-THROAT CHAMBER LENGTH = 1.183E+00 FT  
 BARREL SECTION LENGTH = 9.801E-01 FT

## IMPINGING ELEMENT SIZING RESULTS

ELEMENT TYPE = LOL  
 NO. OF ELEMENTS = 60  
 FUEL ORIFICE DIAMETER = 6.634E-02 IN  
 OX ORIFICE DIAMETER = 1.017E-01 IN

## CORE ELEMENT SPACING RESULTS

ELEMENT TYPE = LOL  
 NUMBER OF ELEMENTS = 60  
 FUEL ORIFICE/ANNULUS DIAMETER = 6.634E-02 IN  
 OXIDIZER ORIFICE DIAMETER = 1.017E-01 IN  
 FUEL INJECTION VELOCITY = 3.430E+02 FT/S  
 OXIDIZER INJECTION VELOCITY = 2.892E+02 FT/S

ROW	# ELEMENTS	MID-ROW RADIUS (IN)
1	6	9.115E-01
2	12	1.609E+00
3	18	2.307E+00
4	24	3.005E+00

ROCKET COMBUSTOR INTERACTIVE DESIGN METHODOLOGY  
Version 23-FEB-91

I-20

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XIH= 1.7819E-01
XIA= 3.0000E+01
XCANT= 1.9000E+01
XFACET= 4.4047E-01
EMUNI= 8.5000E-01
$END
$BAFFLE
$END
$BARRIER
$END
$FFC
$END
$BURN
$END
$INJ
FMAND= 6.7076E+00
XMAND= 6.7076E+00
FMANL= 3.3538E+00
XMANL= 3.3538E+00
PCA= 2.1190E+03, 1.5347E+03, 7.8734E+02
$END
$CHAMBER
MJB= 3
T= 3.3000E-02
$END
$FDORC
NZON= 5
FTER= 20*1.0000E+00
$END
$MIX
EM(1)= 8.930E-01, 1.000E+00
$END
$DEBUGC
DEBUG= F
$END
$SHIFC
SHORT= F
POC= 2.0000E-01, 2.0000E-01
$END
$D18T8DC
SHORT= F
POC= 2.0000E-01, 2.0000E-01
PAMP= 2.0000E-01
MC= 11
LG= 8
MB= 11
LB= 8
IDMAX= 10
$END
$CRPC
NDPC= 18
NDTFQ= 18
NDTLF= 1000
NPRINT= 50
NSUMS= 3500
PAMPC= 2.0000E-01
NRAD= 3
NCIRC= 5
$END
$LEINJC
IDMEN= 2
PAMPCH= 2.0000E-01
NTINJ= 18
USGF= 2.0000E-02

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OBF= 2.0000E-02
DSGF= 2.0000E-02
FCDO= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00
XCDO= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00
$END
$COMBUSTC
FALVM= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00
XALVM= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00
FALTM= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00
XALTM= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00
FRMM= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00
XRMM= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00
ENMULT= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00
COMBXB= 5.0000E-01
ACMULT= 1.0000E+00
CCMULT= 1.0000E+00
ENMULT= 1.0000E+00
TAMULT= 1.0000E+00
$END
$FDORCC
SHORT= F
EPSIL= 2.0000E+01
ERROR= 1.0000E-01
LTS= 5
MTS= 5
NTS= 10
ITMAX= 100
RELX= 6.5000E-01
NEET= 3
NXFST= 3
NYFST= 3
NAFST= 3
MORE= F
$END

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END OF INPUT ECHO

# STEADY STATE COMBUSTION ANALYSIS PROGRAM

## RUN DESCRIPTORS

ROCCID POINT DESIGN TEST CASE 1  
LOX/RP-1 LIKE DOUBLET PAIR WITH FIXED PC  
APPROXIMATES .0100 SUBSCALE DOUBLET

## PROPELLANT DESCRIPTION

FUEL=RP-1 Tman., F= 71.00  
OX=LOX Tman., F=-279.00

## CHAMBER GEOMETRY

CHAMBER RADIUS = 3.3539 IN.  
CYLINDRICAL SECTION =11.5206 IN.  
NOZZLE ENTRANCE RADIUS OF CURVATURE = 1.3320 IN.  
CONVERGENCE HALF-ANGLE =30.0000 DEG.  
THROAT RADIUS = 2.2228 IN.  
CONVERGENT SECTION LENGTH = 2.6730 IN.  
THROAT ENTRANCE RADIUS OF CURVATURE = 1.3320 IN.  
CONTRACTION RATIO = 2.28

## INJECTOR DATA

INJECTOR CORE CONTAINS 80 LOL ELEMENTS

FUEL SIDE: Orifice Diam. =6.634E-02 in.  
Impingement Half-angle =30.00 Deg.  
OX SIDE: Orifice Diam. =1.017E-01 in.  
Impingement Half-angle =30.00 Deg.  
Cd =0.9100  
Unlike Cant Angle =16.00 Deg.  
Cd =0.9400  
Unlike Cant Angle =16.00 Deg.  
Impingement Height =0.116 in.  
Faceplate Thickness = 0.2873 in.  
Impingement Height =0.176 in.  
Faceplate Thickness = 0.4406 in.

## MIXING EFFICIENCIES

CORE MIXING EFFICIENCY=0.8930 BARRIER MIXING EFFICIENCY=1.0000

## COMBUST CONTROL PARAMETERS

MULTIPLIERS:  
FUEL ATOMIZATION LENGTH FOR VAPORIZATION: CORE 1.000 Baffle 1.000 FFC 1.000  
OX ATOMIZATION LENGTH FOR VAPORIZATION: 1.000 1.000 1.000  
FUEL ATOMIZATION LENGTH FOR TIMELAGS: 1.000 1.000 1.000  
OX ATOMIZATION LENGTH FOR TIMELAGS: 1.000 1.000 1.000  
FUEL DROPSIZE: 1.000 1.000 1.000  
OX DROPSIZE: 1.000 1.000 1.000  
MIXING (Em): 1.000 1.000 1.000  
AO-Multiplier=1.000 CC-Multiplier=1.000 N-Multiplier=1.000 Tau-Multiplier=1.000  
Eta-C\* for XB=0.500

BEGIN STEADY STATE COMBUSTION ANALYSIS  
PC=2118.00 PSIA

PROPELLANT PROPERTIES

FUEL-RP-1	Phase=Liquid	Tman., F= 71.00	
	Injected Density= 49.69 Lbm/Cu. Ft	Viscosity=1.380E-03 Lbm/Ft.-S	Surface Tension=1.857E-03 Lbf/Ft
OX=LOX	Phase=Liquid	Tman., F=-279.00	
	Injected Density= 70.16 Lbm/Cu. Ft	Viscosity=1.186E-04 Lbm/Ft.-S	Surface Tension=7.326E-04 Lbf/Ft

OPERATING CONDITIONS

PC FACE=2118.00 PSIA PC THROAT=2030.94 MIXTURE RATIO= 2.680  
 FUEL INJECTION PRESSURE DROP= 656.62 Psia FUEL INJECTION VELOCITY= 349.73 Ft/S  
 OX INJECTION PRESSURE DROP= 656.62 Psia OX INJECTION VELOCITY= 294.66 Ft/S  
 FUEL FLOWRATE= 45.732 OX FLOWRATE= 131.707

ATOMIZATION OUTPUT

DROPSIZE MODEL=AEROJET

ELEMENT TYPE 1 IS LOL  
 FUEL: ATOMIZATION LENGTH, In.=1.07694 ATOMIZATION LENGTH FOR VAPORIZATION, In.=1.07694 DROPLET RADIUS, Microns= 75.16  
 OX: ATOMIZATION LENGTH, In.=2.36005 ATOMIZATION LENGTH FOR VAPORIZATION, In.=2.36005 DROPLET RADIUS, Microns= 83.28

VAPORIZATION CALCULATIONS

X (in.)	CORE=LOL		BAFFLE=		BARRIER=		FFC=	
	%FUEL VAP	%OX VAP	%FUEL VAP	%OX VAP	%FUEL VAP	%OX VAP	%FUEL VAP	%OX VAP
0.0000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
0.2539	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
0.5077	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
0.7616	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.0155	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.2694	6.599	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.5232	16.464	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.7771	26.294	0.000	0.000	0.000	0.000	0.000	0.000	0.000
2.0310	32.325	0.000	0.000	0.000	0.000	0.000	0.000	0.000
2.2848	37.965	1.798	0.000	0.000	0.000	0.000	0.000	0.000
2.5387	42.919	27.016	0.000	0.000	0.000	0.000	0.000	0.000
2.7925	46.687	42.304	0.000	0.000	0.000	0.000	0.000	0.000
3.0464	50.475	52.613	0.000	0.000	0.000	0.000	0.000	0.000
3.3003	53.740	60.293	0.000	0.000	0.000	0.000	0.000	0.000
3.5542	56.574	66.206	0.000	0.000	0.000	0.000	0.000	0.000
3.8081	59.407	70.763	0.000	0.000	0.000	0.000	0.000	0.000
4.0620	61.724	74.721	0.000	0.000	0.000	0.000	0.000	0.000
4.3159	63.904	77.981	0.000	0.000	0.000	0.000	0.000	0.000
4.5698	66.064	80.486	0.000	0.000	0.000	0.000	0.000	0.000
4.8237	67.735	82.841	0.000	0.000	0.000	0.000	0.000	0.000
5.0776	69.406	84.634	0.000	0.000	0.000	0.000	0.000	0.000
5.3315	71.015	86.227	0.000	0.000	0.000	0.000	0.000	0.000
5.5854	72.468	87.914	0.000	0.000	0.000	0.000	0.000	0.000
5.8393	73.921	89.059	0.000	0.000	0.000	0.000	0.000	0.000
6.0932	75.239	90.203	0.000	0.000	0.000	0.000	0.000	0.000
6.3471	76.436	91.100	0.000	0.000	0.000	0.000	0.000	0.000
6.6010	77.637	91.996	0.000	0.000	0.000	0.000	0.000	0.000
6.8549	78.609	92.861	0.000	0.000	0.000	0.000	0.000	0.000
7.1088	79.517	93.325	0.000	0.000	0.000	0.000	0.000	0.000
7.3627	80.424	93.958	0.000	0.000	0.000	0.000	0.000	0.000
7.6166	81.296	94.490	0.000	0.000	0.000	0.000	0.000	0.000
7.8705	82.168	94.966	0.000	0.000	0.000	0.000	0.000	0.000
8.1244	82.870	95.485	0.000	0.000	0.000	0.000	0.000	0.000
8.3783	83.587	95.867	0.000	0.000	0.000	0.000	0.000	0.000
8.6322	84.205	96.255	0.000	0.000	0.000	0.000	0.000	0.000
8.8861	84.823	96.558	0.000	0.000	0.000	0.000	0.000	0.000
9.1400	85.440	96.866	0.000	0.000	0.000	0.000	0.000	0.000
9.3939	86.058	97.185	0.000	0.000	0.000	0.000	0.000	0.000
9.6478	86.675	97.508	0.000	0.000	0.000	0.000	0.000	0.000
9.9017	87.293	97.598	0.000	0.000	0.000	0.000	0.000	0.000
10.1556	87.907	97.758	0.000	0.000	0.000	0.000	0.000	0.000
10.4095	88.348	97.997	0.000	0.000	0.000	0.000	0.000	0.000
10.6634	88.813	98.215	0.000	0.000	0.000	0.000	0.000	0.000
10.9173	89.280	98.382	0.000	0.000	0.000	0.000	0.000	0.000
11.1712	89.748	98.538	0.000	0.000	0.000	0.000	0.000	0.000
11.4251	90.212	98.688	0.000	0.000	0.000	0.000	0.000	0.000
11.6790	90.577	98.798	0.000	0.000	0.000	0.000	0.000	0.000
11.9329	90.943	98.917	0.000	0.000	0.000	0.000	0.000	0.000
12.1868	91.308	99.028	0.000	0.000	0.000	0.000	0.000	0.000
12.4407	91.675	99.113	0.000	0.000	0.000	0.000	0.000	0.000
12.6946	92.030	99.191	0.000	0.000	0.000	0.000	0.000	0.000

OVERALL VAPORIZATION EFFICIENCIES  
FUEL= 92.03% OX= 96.16%

# MASS DISTRIBUTION PROFILE

X (IN)	CORE (lbm/s)		BARRIER (lbm/s)		LOCAL VAPOR MIXTURE RATIO		ETA-C*
	FUEL	OX	FUEL	OX			
0.0000	0.000	0.000	0.000	0.000	0.00	0.0000	0.0000
0.2839	0.000	0.000	0.000	0.000	0.00	0.0000	0.0000
0.5677	0.000	0.000	0.000	0.000	0.00	0.0000	0.0000
0.8516	0.000	0.000	0.000	0.000	0.00	0.0000	0.0000
1.1355	0.000	0.000	0.000	0.000	0.00	0.0000	0.0000
1.4194	3.933	0.000	0.000	0.000	0.00	0.0084	0.0084
1.7032	8.453	0.000	0.000	0.000	0.00	0.0117	0.0117
1.9871	12.026	0.000	0.000	0.000	0.00	0.0166	0.0166
2.2710	14.763	0.000	0.000	0.000	0.00	0.0204	0.0204
2.5548	17.362	2.369	0.000	0.000	0.14	0.0362	0.0362
2.8387	19.628	36.562	0.000	0.000	1.81	0.3041	0.3041
3.1226	21.355	55.835	0.000	0.000	2.81	0.4361	0.4361
3.4065	23.083	69.296	0.000	0.000	3.00	0.5149	0.5149
3.6903	24.576	79.410	0.000	0.000	3.23	0.5734	0.5734
3.9742	25.872	87.196	0.000	0.000	3.37	0.6193	0.6193
4.2581	27.166	93.199	0.000	0.000	3.43	0.6574	0.6574
4.5420	28.227	96.412	0.000	0.000	3.49	0.6897	0.6897
4.8258	29.224	102.707	0.000	0.000	3.51	0.7176	0.7176
5.1097	30.212	105.962	0.000	0.000	3.51	0.7410	0.7410
5.3936	30.976	109.106	0.000	0.000	3.52	0.7616	0.7616
5.6774	31.741	111.338	0.000	0.000	3.51	0.7765	0.7765
5.9613	32.476	113.557	0.000	0.000	3.50	0.7950	0.7950
6.2452	33.141	115.769	0.000	0.000	3.49	0.8109	0.8109
6.5291	33.805	117.927	0.000	0.000	3.47	0.8237	0.8237
6.8129	34.408	119.985	0.000	0.000	3.45	0.8358	0.8358
7.0968	34.956	121.165	0.000	0.000	3.43	0.8461	0.8461
7.3807	35.504	122.041	0.000	0.000	3.41	0.8593	0.8593
7.6645	36.049	122.916	0.000	0.000	3.39	0.8643	0.8643
7.9484	36.364	123.790	0.000	0.000	3.37	0.8719	0.8719
8.2323	36.779	124.449	0.000	0.000	3.35	0.8796	0.8796
8.5162	37.178	125.105	0.000	0.000	3.33	0.8828	0.8828
8.8000	37.577	125.761	0.000	0.000	3.31	0.8981	0.8981
9.0839	37.943	126.290	0.000	0.000	3.30	0.9040	0.9040
9.3678	38.226	126.814	0.000	0.000	3.29	0.9089	0.9089
9.6516	38.508	127.213	0.000	0.000	3.28	0.9133	0.9133
9.9355	38.791	127.606	0.000	0.000	3.27	0.9176	0.9176
10.2194	39.073	127.900	0.000	0.000	3.25	0.9219	0.9219
10.5033	39.356	128.000	0.000	0.000	3.24	0.9257	0.9257
10.7871	39.638	128.299	0.000	0.000	3.22	0.9294	0.9294
11.0710	39.920	128.531	0.000	0.000	3.20	0.9332	0.9332
11.3549	40.201	128.793	0.000	0.000	3.19	0.9362	0.9362
11.6388	40.402	129.088	0.000	0.000	3.18	0.9394	0.9394
11.9226	40.616	129.356	0.000	0.000	3.17	0.9423	0.9423
12.2065	40.829	129.578	0.000	0.000	3.16	0.9451	0.9451
12.4904	41.043	129.761	0.000	0.000	3.15	0.9478	0.9478
12.7742	41.265	129.953	0.000	0.000	3.14	0.9501	0.9501
13.0581	41.422	130.124	0.000	0.000	3.13	0.9523	0.9523
13.3420	41.590	130.290	0.000	0.000	3.12	0.9544	0.9544
13.6258	41.757	130.427	0.000	0.000	3.11	0.9564	0.9564
13.9097	41.924	130.556	0.000	0.000	3.10	0.9583	0.9583
14.1936	42.087	130.641	0.000	0.000			

# AXIAL PRESSURE PROFILE

X (in)	MACH $\phi$	Ptotal (psia)	Pstatic (psia)	Ttotal (R)	Tstatic (R)	Wdot (Lbm/s)	Local Radius (in)
1.21	0.000	2118.88	2118.88	1881.88	1885.35	0.88	3.354
1.51	0.004	2118.88	2118.88	1888.04	1888.48	5.48	3.354
1.81	0.007	2118.84	2118.86	1881.88	1885.34	9.82	3.354
2.11	0.010	2118.88	2118.88	1788.84	1782.88	18.81	3.354
2.42	0.010	2118.87	2118.42	1881.88	1885.34	14.42	3.354
2.72	0.042	2118.84	2118.36	1878.22	1880.80	38.98	3.354
3.02	0.103	2108.14	2084.48	1881.88	1832.81	71.07	3.354
3.32	0.130	2088.70	2078.48	1888.88	1827.80	88.15	3.354
3.62	0.151	2082.03	2088.78	1888.78	1826.88	101.88	3.354
3.93	0.167	2086.38	2084.88	1888.01	1812.13	111.87	3.354
4.23	0.178	2081.82	2044.87	1887.01	1802.17	118.88	3.354
4.53	0.190	2077.14	2038.78	1888.22	1782.50	128.38	3.354
4.83	0.198	2073.25	2027.81	1888.88	1787.18	132.03	3.354
5.13	0.207	2088.87	2021.18	1884.88	1787.33	138.84	3.354
5.44	0.214	2088.88	2014.81	1882.82	1784.84	140.88	3.354
5.74	0.218	2084.81	2008.88	1888.38	1788.81	143.88	3.354
6.04	0.224	2081.88	2004.88	1888.33	1787.31	148.88	3.354
6.34	0.229	2088.84	1988.81	1887.88	1788.28	148.72	3.354
6.64	0.233	2087.73	1988.88	1881.88	1781.88	151.87	3.354
6.95	0.237	2088.88	1982.31	1884.28	1783.88	154.08	3.354
7.25	0.240	2084.34	1988.88	1887.82	1788.81	158.88	3.354
7.55	0.243	2082.81	1988.04	1888.80	1788.34	157.48	3.354
7.85	0.245	2081.71	1988.88	1884.48	1800.88	158.88	3.354
8.15	0.248	2080.44	1988.84	1888.84	1802.88	160.23	3.354
8.45	0.250	2048.38	1978.78	1885.71	1803.88	181.41	3.354
8.75	0.252	2048.38	1978.88	1887.87	1808.38	182.82	3.354
9.05	0.254	2047.30	1974.43	1888.30	1808.78	183.88	3.354
9.35	0.256	2048.48	1972.73	1888.48	1807.73	184.88	3.354
9.65	0.257	2045.87	1971.08	1881.88	1808.87	188.35	3.354
9.97	0.258	2044.88	1988.88	1888.88	1808.88	188.07	3.354
10.27	0.260	2044.28	1988.18	1884.48	1811.22	188.80	3.354
10.57	0.261	2043.81	1988.77	1888.88	1812.84	187.48	3.354
10.87	0.262	2043.04	1988.80	1887.74	1814.18	188.07	3.354
11.17	0.264	2042.48	1984.37	1888.84	1818.78	188.88	3.354
11.48	0.264	2041.92	1983.27	1888.78	1818.80	188.20	3.354
11.78	0.270	2041.41	1958.73	1888.88	1818.88	188.72	3.354
12.08	0.288	2040.88	1948.88	1888.81	1818.80	170.22	3.228
12.38	0.328	2040.81	1921.88	1888.84	1800.43	170.87	3.084
12.68	0.378	2038.80	1884.28	1888.87	1784.78	171.08	2.888
12.98	0.437	2038.78	1888.88	1888.88	1781.81	171.48	2.714
13.28	0.523	2037.70	1747.48	1888.88	1724.48	172.15	2.640
13.58	0.664	2038.14	1602.15	1881.03	1685.42	172.15	2.388
13.88	0.817	2033.88	1402.87	1881.08	1551.88	172.45	2.257
14.18	0.888	2030.84	1172.78	1881.08	1418.81	172.73	2.223

# PERFORMANCE SUMMARY

## C\* EFFICIENCY CALCULATIONS (OOK)

INJECTED MR= 2.800 CSTAR=5880.40 CORE Em=0.8930 BARRIER Em=1.0000  
 CORE: OVERALL MR= 2.800 VAPOR MR= 3.1041 CSTAR-MIX=5769.16 MASS FRACTION= 1.0000  
 BARRIER: OVERALL MR= 0.0000 VAPOR MR=58.9000 CSTAR-MIX= 0.00 MASS FRACTION= 0.0000  
 ENGINE: OVERALL MR= 2.800 VAPOR MR= 3.1041 CSTAR-DEL=5815.00  
 C\* EFFICIENCY = 9.563E-01

## ISP EFFICIENCY CALCULATIONS

ISP-ODK, INJ = 2.668E+02 SEC.  
 ISP-ODK, M.Z. INJ = 2.652E+02 SEC. ISP-ODK, M.Z. VAPOR = 2.628E+02 SEC.  
 VAPORIZATION EFFICIENCY = 9.647E-01 MIXING EFFICIENCY = 9.935E-01  
 ENERGY RELEASE EFFICIENCY = 9.584E-01

NOTE: ISP-DEL = ISP-ODK, INJ. \* ERE \* ETADIV - DELISP-BL

Cochem, In.=0.164E-03	FUEL Cochem, In.=2.293E+02	OX Cochem, In.=8.194E+01
0.000164	229.3	81.94

**ELEMENT 1 IS TYPE=LOL**

FUEL:	CLIN, IN.=	1.564E-02	LVAP, IN.=	0.596	ATOMIZATION LENGTH USED, IN.=	1.077E+00
FUEL:	TEMP=	3.162E-02	TATON=	2.566E-01	TVAP=	1.348E-01
					Total=	4.230E-01

```

OX:  CInj, In.=1.590E-02  Lvap, In.= 0.206  ATOMIZATION LENGTH USED, In.= 2.360E+00
      TImp=6.749E-02      Tatom=6.669E-01      Tvp=5.833E-02      Total=7.826E-01

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FUEL:	Clinj, In.=1.554E-02	Lvap, In.= 0.555	
	Temp=5.162E-02	Tatom=2.555E-01	Tvap=1.345E-01
			Total=4.230E-01
OX:	Clinj, In.=1.555E-02	Lvap, In.= 0.205	
	Temp=5.749E-02	Tatom=5.555E-01	Tvap=5.533E-02
			Total=7.525E-01

# CHAMBER-NOZZLE OPTIMIZATION RESULTS

CHAMBER LENGTH (FEET)	ETA-C*	ETA-NOZ	OVERALL EFFICIENCY
0.0000	0.0000	0.8807	0.0000
0.1667	0.0168	0.8728	0.0148
0.3333	0.6227	0.8848	0.6384
0.5000	0.7872	0.8837	0.8805
0.6667	0.8733	0.8419	0.7353
0.8333	0.9142	0.8302	0.7590
1.0000	0.9402	0.8166	0.7886
1.1667	0.9670	0.8057	0.7711
1.3333	0.9704	0.7915	0.7681
1.5000	0.9838	0.7773	0.7647
1.6667	0.9872	0.7631	0.7610
1.8333	1.0000	0.7465	0.7485
2.0000	1.0000	0.7299	0.7299
2.1667	1.0000	0.7114	0.7114
2.3333	1.0000	0.6928	0.6928
2.5000	1.0000	0.6742	0.6742
2.6667	1.0000	0.6461	0.6461
2.8333	1.0000	0.6142	0.6142
3.0000	1.0000	0.5823	0.5823
3.1667	1.0000	0.5504	0.5504
3.3333	1.0000	0.5071	0.5071
3.5000	1.0000	0.4439	0.4439
3.6667	1.0000	0.3808	0.3808
3.8333	1.0000	0.3176	0.3176
4.0000	1.0000	0.2544	0.2544

OPTIMUM CHAMBER LENGTH= 1.1667 FT  
 MAXIMUM OVERALL EFFICIENCY= 0.7711

# REDESIGNED CHAMBER RESULTS

NOMINAL CHAMBER PRESSURE	=	2.110E+03	PSIA
THROTTLED CHAMBER PRESSURE	=	1.936E+03	PSIA
FUEL INJECTION PRESSURE DROP	=	6.586E+02	PSI
OX INJECTION PRESSURE DROP	=	6.586E+02	PSI
CHAMBER RADIUS	=	2.786E-01	FT
THROAT RADIUS	=	1.852E-01	FT
NOZZLE ENTRANCE RADIUS OF CURVATURE	=	1.110E-01	FT
THROAT ENTRANCE RADIUS OF CURVATURE	=	1.110E-01	FT
NOZZLE CONVERGENCE HALF-ANGLE	=	3.000E+01	DEG
INJECTOR-TO-THROAT CHAMBER LENGTH	=	1.188E+00	FT
BARREL SECTION LENGTH	=	9.646E-01	FT

## IMPINGING ELEMENT SIZING RESULTS

ELEMENT TYPE	=	LOL
NO. OF ELEMENTS	=	60
FUEL ORIFICE DIAMETER	=	6.694E-02 IN
OX ORIFICE DIAMETER	=	1.017E-01 IN

DIRECT INPUT ECHO FROM SUBROUTINE PINPUT

ROCCID  
Rocket Combustor Interactive Design Methodology  
Version 23-FEB-91

ROCCID POINT DESIGN TEST CASE 1  
LOX/RP-1 LIKE DOUBLET PAIR WITH FIXED PC  
APPROXIMATES .0100 SUBSCALE DOUBLET

\$MODELS									
1									
2									
1									
\$MCHAM=									
\$MBURN=									
\$MINJ=									
\$END									
\$OPCOND									
\$PSPA=	8.8800E+01,	1.0940E+02,	1.3940E+02,	1.5880E+02,	1.7780E+02,				
	2.0970E+02,	2.3550E+02,	2.5470E+02,	2.6700E+02,	2.7040E+02,				
	2.7180E+02,	2.7190E+02,	2.7130E+02,	2.8980E+02,	2.8620E+02,				
	2.8860E+02,	2.8500E+02,	2.8210E+02,	2.5920E+02,	2.5630E+02,				
	2.5080E+02,	2.3840E+02,	2.2750E+02,	2.1100E+02,	2.0140E+02,				
	1.5560E+02,	1.0970E+02,	8.0000E+01,						
	1.7780E+03,	2.4560E+03,	3.4870E+03,	4.0740E+03,	4.5600E+03,				
	5.0700E+03,	5.4540E+03,	5.7340E+03,	5.9080E+03,	5.9810E+03,				
	5.9690E+03,	5.9640E+03,	6.9340E+03,	5.9080E+03,	5.8620E+03,				
	5.8550E+03,	5.8260E+03,	5.7590E+03,	5.7000E+03,	5.6430E+03,				
	5.5320E+03,	5.2750E+03,	5.0470E+03,	4.6780E+03,	4.4200E+03,				
	3.4570E+03,	2.4940E+03,	1.4650E+03,						
\$FUEL=									
\$LOX									
\$PC=	2.1180E+03								
\$XMR=	2.8800E+00								
\$FTMAN=	7.1000E+01								
\$XTMAN=	-2.7900E+02								
\$EMMAN=	1.0000E+00								
\$NPERFP=	28								
\$PMRA=	1.0000E-01,	3.0000E-01,	8.0000E-01,	8.0000E-01,	1.0000E+00,				
	1.2500E+00,	1.5000E+00,	1.7500E+00,	2.0000E+00,	2.2000E+00,				
	2.4000E+00,	2.6000E+00,	2.8000E+00,	3.0000E+00,	3.2000E+00,				
	2.8000E+00,	3.0000E+00,	3.2000E+00,	3.4000E+00,	3.6000E+00,				
	4.0000E+00,	5.0000E+00,	6.0000E+00,	8.0000E+00,	1.0000E+01,				
	1.6000E+01,	2.0000E+01,	5.0000E+01,						
\$END									
\$GEOM									
\$CHAMB=	2.7949E-01								
\$RTHRT=	1.8523E-01								
\$RNE=	1.1100E-01								
\$RTE=	1.1100E-01								
\$ALPHA=	3.0000E+01								
\$CHAMBL=	1.1878E+00								
\$XC=	9.6477E-01								
\$END									
\$SCORE									
\$TYPE=	'LOL'								
\$NEL=	60								
\$FDJ=	8.6345E-02								
\$FCD=	9.1000E-01								
\$FIH=	1.1491E-01								
\$FIA=	3.0000E+01								
\$FGANT=	1.6000E+01								
\$FFACET=	2.8728E-01								
\$XDJ=	1.0172E-01								
\$XCD=	9.4000E-01								

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XIH= 1.7618E-01
XIA= 3.0000E+01
XCANT= 1.6000E+01
XFACET= 4.4047E-01
EMJN1= 6.5000E-01
$END
$BAFFLE
$END
$BARRIER
$END
$FFC
$END
$BURN
$END
$INJ
FMAND= 6.7076E+00
XMAND= 6.7076E+00
FMANL= 3.3538E+00
XMANL= 3.3538E+00
PCA= 2.1160E+03, 1.5947E+03, 7.6734E+02
$END
$CHAMBER
MJB= 3
T= 3.3000E-02
$END
$FDORC
NZON= 5
PTER= 20*1.0000E+00
$END
$MIX
EM(1)= 8.930E-01, 1.000E+00
$END
$DEBUGC
DEBUG= F
$END
$HIFIG
SHORT= F
POC= 2.0000E-01, 2.0000E-01
$END
$DIST3DC
SHORT= F
POC= 2.0000E-01, 2.0000E-01
PAMP= 2.0000E-01
MC= 11
LC= 8
MB= 11
LB= 8
IDMAX= 10
$END
$CRPC
NDPC= 16
NDTFQ= 16
NDTLF= 1000
NPRINT= 50
NBUMS= 3500
PAMPC= 2.0000E-01
NRAD= 3
NCIRC= 5
$END
$LEINJC
IDMEM= 2
PAMPC= 2.0000E-01
NTINJ= 16
USGF= 2.0000E-02

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OGF=	2.0000E-02				
D8GF=	2.0000E-02				
FCDO=	1.0000E+00,	1.0000E+00,	1.0000E+00,	1.0000E+00,	1.0000E+00
XCOO=	1.0000E+00,	1.0000E+00,	1.0000E+00,	1.0000E+00,	1.0000E+00
\$END					
\$COMBUSTC					
FALVM=	1.0000E+00,	1.0000E+00,	1.0000E+00,	1.0000E+00,	1.0000E+00
XALVM=	1.0000E+00,	1.0000E+00,	1.0000E+00,	1.0000E+00,	1.0000E+00
FALTM=	1.0000E+00,	1.0000E+00,	1.0000E+00,	1.0000E+00,	1.0000E+00
XALTM=	1.0000E+00,	1.0000E+00,	1.0000E+00,	1.0000E+00,	1.0000E+00
FRMM=	1.0000E+00,	1.0000E+00,	1.0000E+00,	1.0000E+00,	1.0000E+00
XFRMM=	1.0000E+00,	1.0000E+00,	1.0000E+00,	1.0000E+00,	1.0000E+00
ENMULT=	1.0000E+00,	1.0000E+00,	1.0000E+00,	1.0000E+00,	1.0000E+00
COMBXS=	5.0000E-01				
ACMULT=	1.0000E+00				
CCMULT=	1.0000E+00				
ENMULT=	1.0000E+00				
TAUMULT=	1.0000E+00				
\$END					
\$FDORCC					
SHORT=	F				
EPSIL=	2.0000E+01				
ERROR=	1.0000E-01				
LT8=	5				
MT8=	5				
NT8=	10				
ITMAX=	100				
RELX=	6.5000E-01				
NEET=	3				
NXFST=	3				
NYFST=	3				
NAFST=	3				
MORE=	F				
\$END					

END OF INPUT ECHO

# STEADY STATE COMBUSTION ANALYSIS PROGRAM

## RUN DESCRIPTORS

ROGID POINT DESIGN TEST CASE 1  
LOX/RP-1 LIKE DOUBLET PAIR WITH FIXED PC  
APPROXIMATES .0100 SUBSCALE DOUBLET

## PROPELLANT DESCRIPTION

FUEL=RP-1 Tman., F= 71.00  
OX=LOX Tman., F=-279.00

## CHAMBER GEOMETRY

CHAMBER RADIUS = 3.3530 IN.  
CYLINDRICAL SECTION =11.5770 IN.  
NOZZLE ENTRANCE RADIUS OF CURVATURE = 1.3320 IN.  
CONVERGENCE HALF-ANGLE =30.0000 DEG.  
THROAT RADIUS = 2.2228 IN.  
CONVERGENT SECTION LENGTH = 2.8730 IN.  
THROAT ENTRANCE RADIUS OF CURVATURE = 1.3320 IN.  
CONTRACTION RATIO = 2.28

## INJECTOR DATA

INJECTOR CORE CONTAINS 80 LOL ELEMENTS

FUEL SIDE: Orifice Diam. =6.624E-02 in.  
Impingement Half-angle =30.00 Deg.  
Cd =0.9100  
OX SIDE: Orifice Diam. =1.017E-01 in.  
Impingement Half-angle =30.00 Deg.  
Cd =0.9400  
Unlike Cant Angle =16.00 Deg.  
Unlike Cant Angle =16.00 Deg.  
Impingement Height =0.115 in.  
Faceplate Thickness = 0.2673 in.  
Impingement Height =0.176 in.  
Faceplate Thickness = 0.4405 in.

## MIXING EFFICIENCIES

CORE MIXING EFFICIENCY=0.8880 BARRIER MIXING EFFICIENCY=1.0000

## COMBUST CONTROL PARAMETERS

MULTIPLIERS:  
FUEL ATOMIZATION LENGTH FOR VAPORIZATION: 1.000  
OX ATOMIZATION LENGTH FOR VAPORIZATION: 1.000  
FUEL ATOMIZATION LENGTH FOR TIMELAGS: 1.000  
OX ATOMIZATION LENGTH FOR TIMELAGS: 1.000  
FUEL DROPSIZE: 1.000  
OX DROPSIZE: 1.000  
MIXING (Em): 1.000  
BAFFLE 1.000  
BARRIER 1.000  
FFC 1.000

AC-Multiplier=1.000 CC-Multiplier=1.000 N-Multiplier=1.000  
Eta-C\* for XB=0.500 Tau-Multiplier=1.000

BEGIN STEADY STATE COMBUSTION ANALYSIS  
PC=2119.00 PSIA

PROPELLANT PROPERTIES

FUEL=RP-1	Phase=Liquid	Tman.. F= 71.00	
	Injected Density= 49.89 Lbm/Cu. Ft	Viscosity=1.360E-03 Lbm/Ft-s	Surface Tension=1.657E-03 Lbf/Ft
OX=LOX	Phase=Liquid	Tman.. F=-279.00	
	Injected Density= 70.18 Lbm/Cu. Ft	Viscosity=1.166E-04 Lbm/Ft-s	Surface Tension=7.326E-04 Lbf/Ft

OPERATING CONDITIONS

PC FACE=2119.00 PSIA PC THROAT=2031.07 MIXTURE RATIO= 2.860  
FUEL INJECTION PRESSURE DROP= 658.47 Psia FUEL INJECTION VELOCITY= 349.70 Ft/s  
OX INJECTION PRESSURE DROP= 658.47 Psia OX INJECTION VELOCITY= 294.85 Ft/s

FUEL FLOWRATE= 46.721 OX FLOWRATE= 131.677

ATOMIZATION OUTPUT

DROPSIZE MODEL=AEROJET

ELEMENT TYPE 1 IS LOL			
FUEL:	ATOMIZATION LENGTH, In.=1.07692	ATOMIZATION LENGTH FOR VAPORIZATION, In.=1.07692	DROPLET RADIUS, Microns= 75.16
OX:	ATOMIZATION LENGTH, In.=2.36002	ATOMIZATION LENGTH FOR VAPORIZATION, In.=2.36002	DROPLET RADIUS, Microns= 83.28

VAPORIZATION CALCULATIONS

X (in.)	CORE=LOL		BAFFLE=		BARRIER=		FFC=	
	%FUEL VAP	%OX VAP	%FUEL VAP	%OX VAP	%FUEL VAP	%OX VAP	%FUEL VAP	%OX VAP
0.0000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
0.2500	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
0.5000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
0.7500	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.0000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.2500	6.848	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.5000	16.674	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.7500	28.464	0.000	0.000	0.000	0.000	0.000	0.000	0.000
2.0000	32.518	0.000	0.000	0.000	0.000	0.000	0.000	0.000
2.2500	38.181	3.124	0.000	0.000	0.000	0.000	0.000	0.000
2.5000	43.071	27.675	0.000	0.000	0.000	0.000	0.000	0.000
2.7500	46.864	42.849	0.000	0.000	0.000	0.000	0.000	0.000
3.0000	50.858	52.986	0.000	0.000	0.000	0.000	0.000	0.000
3.2500	53.868	60.604	0.000	0.000	0.000	0.000	0.000	0.000
3.5000	56.733	66.463	0.000	0.000	0.000	0.000	0.000	0.000
3.7500	59.578	71.002	0.000	0.000	0.000	0.000	0.000	0.000
4.0000	61.864	74.931	0.000	0.000	0.000	0.000	0.000	0.000
4.2500	64.053	78.151	0.000	0.000	0.000	0.000	0.000	0.000
4.5000	66.165	80.641	0.000	0.000	0.000	0.000	0.000	0.000
4.7500	67.863	82.970	0.000	0.000	0.000	0.000	0.000	0.000
5.0000	69.541	84.670	0.000	0.000	0.000	0.000	0.000	0.000
5.2500	71.138	86.370	0.000	0.000	0.000	0.000	0.000	0.000
5.5000	72.597	88.015	0.000	0.000	0.000	0.000	0.000	0.000
5.7500	74.056	89.165	0.000	0.000	0.000	0.000	0.000	0.000
6.0000	76.355	90.290	0.000	0.000	0.000	0.000	0.000	0.000
6.2500	78.559	91.190	0.000	0.000	0.000	0.000	0.000	0.000
6.5000	77.762	92.086	0.000	0.000	0.000	0.000	0.000	0.000
6.7500	76.708	92.733	0.000	0.000	0.000	0.000	0.000	0.000
7.0000	78.619	93.400	0.000	0.000	0.000	0.000	0.000	0.000
7.2500	80.626	94.060	0.000	0.000	0.000	0.000	0.000	0.000
7.5000	81.401	94.650	0.000	0.000	0.000	0.000	0.000	0.000
7.7500	82.277	95.050	0.000	0.000	0.000	0.000	0.000	0.000
8.0000	83.050	95.540	0.000	0.000	0.000	0.000	0.000	0.000
8.2500	83.670	95.940	0.000	0.000	0.000	0.000	0.000	0.000
8.5000	84.290	96.330	0.000	0.000	0.000	0.000	0.000	0.000
8.7500	84.910	96.630	0.000	0.000	0.000	0.000	0.000	0.000
9.0000	85.530	96.930	0.000	0.000	0.000	0.000	0.000	0.000
9.2500	86.150	97.220	0.000	0.000	0.000	0.000	0.000	0.000
9.5000	86.770	97.420	0.000	0.000	0.000	0.000	0.000	0.000
9.7500	87.390	97.620	0.000	0.000	0.000	0.000	0.000	0.000
10.0000	87.975	97.820	0.000	0.000	0.000	0.000	0.000	0.000
10.2500	88.413	98.029	0.000	0.000	0.000	0.000	0.000	0.000
10.5000	88.882	98.239	0.000	0.000	0.000	0.000	0.000	0.000
10.7500	89.351	98.407	0.000	0.000	0.000	0.000	0.000	0.000
11.0000	89.820	98.558	0.000	0.000	0.000	0.000	0.000	0.000
11.2500	90.270	98.699	0.000	0.000	0.000	0.000	0.000	0.000
11.5000	90.637	98.819	0.000	0.000	0.000	0.000	0.000	0.000
11.7500	91.006	98.935	0.000	0.000	0.000	0.000	0.000	0.000
12.0000	91.372	99.047	0.000	0.000	0.000	0.000	0.000	0.000
12.2500	91.738	99.126	0.000	0.000	0.000	0.000	0.000	0.000
12.5000	92.079	99.205	0.000	0.000	0.000	0.000	0.000	0.000

OVERALL VAPORIZATION EFFICIENCIES  
FUEL= 92.06% OX= 99.20%

MASS DISTRIBUTION PROFILE

X (IN)	CORE (lbm/s)		BARRIER (lbm/s)		LOCAL VAPOR MIXTURE RATIO	ETA-C*
	FUEL	OX	FUEL	OX		
0.0000	0.000	0.000	0.000	0.000	0.00	0.0000
0.2500	0.000	0.000	0.000	0.000	0.00	0.0000
0.5000	0.000	0.000	0.000	0.000	0.00	0.0000
0.6550	0.000	0.000	0.000	0.000	0.00	0.0000
1.1400	0.000	0.000	0.000	0.000	0.00	0.0000
1.4250	4.045	0.000	0.000	0.000	0.00	0.0058
1.7100	6.538	0.000	0.000	0.000	0.00	0.0118
1.9950	12.099	0.000	0.000	0.000	0.00	0.0167
2.2800	14.868	0.000	0.000	0.000	0.00	0.0205
2.5650	17.448	4.114	0.000	0.000	0.24	0.0468
2.8500	19.992	36.442	0.000	0.000	1.85	0.3105
3.1350	21.427	56.422	0.000	0.000	2.63	0.4398
3.4200	23.161	69.770	0.000	0.000	3.01	0.5178
3.7050	24.938	79.602	0.000	0.000	3.24	0.5758
3.9900	26.939	87.618	0.000	0.000	3.37	0.6214
4.2750	27.240	93.493	0.000	0.000	3.43	0.6595
4.5600	28.285	98.667	0.000	0.000	3.49	0.6915
4.8450	29.286	102.907	0.000	0.000	3.51	0.7192
5.1300	30.260	108.185	0.000	0.000	3.51	0.7426
5.4150	31.028	109.253	0.000	0.000	3.52	0.7629
5.7000	31.785	111.491	0.000	0.000	3.51	0.7798
5.9850	32.525	113.729	0.000	0.000	3.50	0.7964
6.2700	33.192	116.895	0.000	0.000	3.49	0.8120
6.5550	33.859	117.409	0.000	0.000	3.47	0.8248
6.8400	34.453	118.691	0.000	0.000	3.45	0.8368
7.1250	35.003	120.076	0.000	0.000	3.43	0.8471
7.4100	35.554	121.230	0.000	0.000	3.41	0.8573
7.6950	35.986	122.108	0.000	0.000	3.39	0.8651
7.9800	36.403	122.986	0.000	0.000	3.38	0.8728
8.2650	36.817	123.642	0.000	0.000	3.36	0.8804
8.5500	37.218	124.600	0.000	0.000	3.35	0.8870
8.8350	37.616	125.158	0.000	0.000	3.33	0.8938
9.1200	37.971	125.803	0.000	0.000	3.31	0.8997
9.4050	38.255	126.330	0.000	0.000	3.30	0.9046
9.6900	38.538	126.844	0.000	0.000	3.29	0.9085
9.9750	38.822	127.239	0.000	0.000	3.28	0.9138
10.2600	39.105	127.634	0.000	0.000	3.26	0.9182
10.5450	39.389	128.016	0.000	0.000	3.25	0.9228
10.8300	39.672	128.279	0.000	0.000	3.23	0.9263
11.1150	39.956	128.542	0.000	0.000	3.22	0.9300
11.4000	40.228	128.808	0.000	0.000	3.20	0.9336
11.6850	40.423	129.081	0.000	0.000	3.19	0.9387
11.9700	40.638	129.358	0.000	0.000	3.18	0.9398
12.2550	40.882	129.579	0.000	0.000	3.17	0.9427
12.5400	41.087	129.778	0.000	0.000	3.16	0.9458
12.8250	41.272	129.950	0.000	0.000	3.15	0.9482
13.1100	41.440	130.122	0.000	0.000	3.14	0.9504
13.3950	41.608	130.275	0.000	0.000	3.13	0.9528
13.6800	41.776	130.422	0.000	0.000	3.12	0.9548
13.9650	41.944	130.528	0.000	0.000	3.11	0.9567
14.2500	42.089	130.629	0.000	0.000	3.10	0.9586

# AXIAL PRESSURE PROFILE

X (in)	MACH $\phi$	Ptotal (psia)	Pstatic (psia)	Total (R)	Tstatic (R)	Wdot (Lbm/s)	Local Radius (in)
1.20	0.000	2118.88	2119.85	1891.89	1885.41	0.59	3.354
1.30	0.001	2118.88	2119.84	1891.89	1885.41	1.99	3.354
1.61	0.005	2118.88	2119.59	1898.27	1888.78	7.02	3.354
1.91	0.008	2118.82	2119.51	1891.89	1885.41	11.01	3.354
2.21	0.010	2118.56	2119.38	1747.29	1740.60	14.53	3.354
2.51	0.013	2118.49	2119.28	1891.89	1885.40	18.01	3.354
2.81	0.065	2119.63	2109.53	5194.34	5173.41	51.27	3.354
3.11	0.112	2103.85	2088.87	8788.26	8768.52	76.48	3.354
3.41	0.137	2096.60	2076.26	8861.19	8829.62	92.65	3.354
3.71	0.156	2080.31	2062.52	8858.63	8823.79	104.72	3.354
4.01	0.171	2084.92	2051.58	8842.59	8808.98	114.13	3.354
4.32	0.183	2080.22	2042.04	8834.93	8800.08	121.68	3.354
4.62	0.193	2075.99	2033.41	8828.31	8790.82	128.07	3.354
4.92	0.202	2072.30	2025.87	8823.87	8787.40	138.34	3.354
5.22	0.208	2069.10	2019.33	8824.00	8786.64	137.69	3.354
5.52	0.216	2066.26	2013.51	8823.08	8785.30	141.45	3.354
5.82	0.220	2063.64	2008.54	8825.78	8787.45	144.52	3.354
6.12	0.228	2081.31	2003.35	8828.49	8787.59	147.68	3.354
6.42	0.230	2059.13	1998.68	8828.92	8788.92	150.30	3.354
6.72	0.234	2057.25	1994.98	8832.30	8782.44	152.52	3.354
7.02	0.238	2055.55	1991.50	8836.13	8794.88	154.48	3.354
7.39	0.241	2063.94	1988.18	8838.65	8797.90	156.31	3.354
7.69	0.244	2052.61	1985.43	8841.10	8800.13	157.80	3.354
7.99	0.246	2051.40	1982.94	8842.78	8801.50	159.15	3.354
8.23	0.249	2050.17	1980.38	8844.28	8802.73	160.51	3.354
8.59	0.251	2049.12	1978.22	8846.20	8804.39	161.64	3.354
8.83	0.253	2048.09	1976.07	8848.12	8806.08	162.76	3.354
9.13	0.255	2047.10	1974.04	8849.59	8807.28	163.81	3.354
9.43	0.256	2046.30	1972.38	8850.72	8808.22	164.66	3.354
9.73	0.258	2045.52	1970.75	8851.92	8809.22	165.48	3.354
10.04	0.259	2044.64	1969.34	8853.40	8810.53	166.20	3.354
10.34	0.260	2044.14	1967.69	8854.62	8811.78	166.92	3.354
10.64	0.261	2043.49	1966.54	8855.37	8813.16	167.59	3.354
10.94	0.263	2042.83	1965.38	8858.13	8814.77	168.18	3.354
11.24	0.264	2042.36	1964.15	8859.89	8816.37	168.74	3.354
11.64	0.265	2041.84	1963.11	8860.86	8817.23	169.28	3.354
11.84	0.270	2041.34	1959.41	8860.68	8816.00	169.78	3.328
12.14	0.290	2040.82	1946.44	8860.91	8810.89	170.27	3.225
12.44	0.327	2040.25	1920.98	8860.94	8800.44	170.71	3.061
12.74	0.378	2039.58	1883.50	8860.97	8784.71	171.12	2.886
13.05	0.438	2038.73	1829.62	8860.99	8761.49	171.49	2.712
13.35	0.524	2037.68	1746.33	8861.02	8724.23	171.83	2.539
13.65	0.655	2036.12	1600.73	8861.04	8684.99	172.18	2.387
13.95	0.818	2035.99	1402.09	8861.06	8651.43	172.48	2.257
14.25	0.999	2031.07	1172.87	8861.08	8419.51	172.73	2.223

# PERFORMANCE SUMMARY

## C\* EFFICIENCY CALCULATIONS (OOK)

INJECTED MR= 2.8500 CSTAR=5550.40 CORE Em=0.8930 BARRIER Em=1.0000  
 CORE: OVERALL MR= 2.8500 VAPOR MR= 3.1029 CSTAR-MIX=5769.49 MASS FRACTION= 1.0000  
 BARRIER: OVERALL MR= 0.0000 VAPOR MR=99.9000 CSTAR-MIX= 0.00 MASS FRACTION= 0.0000  
 ENGINE: OVERALL MR= 2.8500 VAPOR MR= 3.1029 CSTAR-DEL=5617.64  
 C\* EFFICIENCY = 9.566E-01

## ISP EFFICIENCY CALCULATIONS

ISP-ODK, INJ = 2.669E+02 SEC.  
 ISP-ODK, M.Z. INJ = 2.652E+02 SEC.  
 VAPORIZATION EFFICIENCY = 9.650E-01  
 ENERGY RELEASE EFFICIENCY = 9.567E-01  
 ISP-ODK, M.Z. VAPOR = 2.628E+02 SEC.  
 MIXING EFFICIENCY = 9.935E-01

NOTE: ISP-DEL = ISP-ODK, INJ. \* ERE \* ETADIV - DELISP.BL

TIME-LAG CALCULATIONS, Milliseconds

Cohem, In.=0.104E-03      FUEL Cohem, In.=2.293E+02      OX Cohem, In.=0.104E+01

ELEMENT 1 IS TYPE=LOL

FUEL:      CInj, In.=1.593E-02      Lvap, In.= 0.566      ATOMIZATION LENGTH USED, In.= 1.077E+00  
              TImp=3.102E-02      Tatom=2.566E-01      Tvap=1.346E-01      Total=4.231E-01

OX:      CInj, In.=1.596E-02      Lvap, In.= 0.206      ATOMIZATION LENGTH USED, In.= 2.360E+00  
              TImp=6.750E-02      Tatom=6.670E-01      Tvap=6.633E-02      Total=7.629E-01

EFFECTIVE TIMELAGS, Milliseconds

FUEL:      CInj, In.=1.593E-02      Lvap, In.= 0.566      Total=4.231E-01  
              TImp=3.102E-02      Tatom=2.566E-01      Tvap=1.346E-01

OX:      CInj, In.=1.596E-02      Lvap, In.= 0.206      Total=7.629E-01  
              TImp=6.750E-02      Tatom=6.670E-01      Tvap=6.633E-02

# CHAMBER-NOZZLE OPTIMIZATION RESULTS

CHAMBER LENGTH (FEET)	ETA-C*	ETA-NOZ	OVERALL EFFICIENCY
0.0000	0.0000	0.8807	0.0000
0.0167	0.0166	0.8726	0.0148
0.0333	0.6228	0.8846	0.6364
0.5000	0.7972	0.8837	0.6806
0.6667	0.8738	0.8419	0.7368
0.8333	0.9143	0.8302	0.7590
1.0000	0.8401	0.8186	0.7695
1.1667	0.9570	0.8087	0.7710
1.3333	0.9899	0.7915	0.7677
1.5000	0.9828	0.7773	0.7639
1.6667	0.9957	0.7631	0.7588
1.8333	1.0000	0.7485	0.7485
2.0000	1.0000	0.7299	0.7299
2.1667	1.0000	0.7114	0.7114
2.3333	1.0000	0.6926	0.6926
2.5000	1.0000	0.6742	0.6742
2.6667	1.0000	0.6481	0.6481
2.8333	1.0000	0.6142	0.6142
3.0000	1.0000	0.5823	0.5823
3.1667	1.0000	0.5504	0.5504
3.3333	1.0000	0.5071	0.5071
3.5000	1.0000	0.4439	0.4439
3.6667	1.0000	0.3606	0.3606
3.8333	1.0000	0.3176	0.3176
4.0000	1.0000	0.2544	0.2544

OPTIMUM CHAMBER LENGTH= 1.1667 FT  
 MAXIMUM OVERALL EFFICIENCY= 0.7710

# REDESIGNED CHAMBER RESULTS

NOMINAL CHAMBER PRESSURE	=	2.141E+03	PSIA
THROTTLED CHAMBER PRESSURE	=	1.636E+03	PSIA
FUEL INJECTION PRESSURE DROP	=	6.729E+02	PSI
OX INJECTION PRESSURE DROP	=	6.729E+02	PSI
CHAMBER RADIUS	=	2.786E-01	FT
THROAT RADIUS	=	1.852E-01	FT
NOZZLE ENTRANCE RADIUS OF CURVATURE	=	1.110E-01	FT
THROAT ENTRANCE RADIUS OF CURVATURE	=	1.110E-01	FT
NOZZLE CONVERGENCE HALF-ANGLE	=	3.000E+01	DEG
INJECTOR-TO-THROAT CHAMBER LENGTH	=	1.166E+00	FT
BARREL SECTION LENGTH	=	9.646E-01	FT

## IMPINGING ELEMENT SIZING RESULTS

ELEMENT TYPE	=	LOL
NO. OF ELEMENTS	=	60
FUEL ORIFICE DIAMETER	=	6.634E-02 IN
OX ORIFICE DIAMETER	=	1.017E-01 IN

ROCCID  
Rocket Combustor Interactive Design Methodology  
Version 23-FEB-91

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XIH= 1.7619E-01
XIA= 3.0000E+01
XCANT= 1.6000E+01
XFACET= 4.4047E-01
EMUNI= 6.5000E-01
$END
$BAFFLE
$END
$BARRIER
$END
$FFC
$END
$BURN
$END
$INJ
FMAND= 6.7076E+00
XMAND= 6.7076E+00
FMANL= 3.3538E+00
XMANL= 3.3538E+00
PCA= 2.1411E+03, 1.5514E+03, 7.6734E+02
$END
$CHAMBER
MUB= 3
T= 3.3000E-02
$END
$FDORC
NZON= 5
FTER= 20*1.0000E+00
$END
$MIX
EM(1)= 8.930E-01, 1.000E+00
$END
$DEBUG
DEBUG= F
$END
$HIFIC
SHORT= F
POC= 2.0000E-01, 2.0000E-01
$END
$DIST3DC
SHORT= F
POC= 2.0000E-01, 2.0000E-01
PAMP= 2.0000E-01
MC= 11
LC= 8
MB= 11
LB= 8
IDMAX= 10
$END
$CRPC
NDFC= 16
NOTFQ= 16
NDTLF= 1000
NPRINT= 50
NBLMS= 3500
PAMP= 2.0000E-01
NRAD= 3
NCIRC= 5
$END
$LEINJC
IDOMEM= 2
PAMPCH= 2.0000E-01
NTINJ= 16
USGF= 2.0000E-02

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CBF= 2.000E-02
DBGF= 2.000E-02
FCDO= 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00
XCDO= 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00
$END
$COMBUSTC
FALVM= 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00
XALVM= 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00
FALTM= 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00
XALTM= 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00
FRMM= 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00
XRMM= 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00
EMMUL= 1.000E+00, 1.000E+00, 1.000E+00
COMBXS= 5.000E-01
ACMUL= 1.000E+00
CCMUL= 1.000E+00
ENMUL= 1.000E+00
TAMUL= 1.000E+00
$END
$FDORCG
SHORT= F
EPSIL= 2.000E+01
ERROR= 1.000E-01
LTS= 5
MTS= 5
NTS= 10
ITMAX= 100
RELX= 6.500E-01
NEET= 3
NXFST= 3
NYFST= 3
NAFST= 3
MORE= F
$END

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END OF INPUT ECHO

# STEADY STATE COMBUSTION ANALYSIS PROGRAM

## RUN DESCRIPTORS

ROCCID POINT DESIGN TEST CASE 1  
LOX/RP-1 LIKE DOUBLET PAIR WITH FIXED PC  
APPROXIMATES -0100 SUBSCALE DOUBLET

## PROPELLANT DESCRIPTION

FUEL=RP-1 Tman., F= 71.00  
OX=LOX Tman., F=-279.00

## CHAMBER GEOMETRY

CHAMBER RADIUS = 3.3539 IN.  
CYLINDRICAL SECTION =11.5770 IN.  
NOZZLE ENTRANCE RADIUS OF CURVATURE = 1.3320 IN.  
CONVERGENCE HALF-ANGLE =30.0000 DEG.  
THROAT RADIUS = 2.2226 IN.  
CONVERGENT SECTION LENGTH = 2.6730 IN.  
THROAT ENTRANCE RADIUS OF CURVATURE = 1.3320 IN.  
CONTRACTION RATIO = 2.28

## INJECTOR DATA

INJECTOR CORE CONTAINS 60 LOL ELEMENTS

FUEL SIDE: Orifice Diam. =6.94E-02 in.  
Impingement Half-angle =30.00 Deg.  
OX SIDE: Orifice Diam. =1.017E-01 in.  
Impingement Half-angle =30.00 Deg.  
Cd =0.9100  
Unlike Cant Angle =16.00 Deg.  
Cd =0.9400  
Unlike Cant Angle =16.00 Deg.  
Impingement Height =0.115 in.  
Faceplate Thickness = 0.2673 in.  
Impingement Height =0.176 in.  
Faceplate Thickness = 0.4405 in.

## MIXING EFFICIENCIES

CORE MIXING EFFICIENCY=0.9930 BARRIER MIXING EFFICIENCY=1.0000

## COMBUST CONTROL PARAMETERS

MULTIPLIERS:  
FUEL ATOMIZATION LENGTH FOR VAPORIZATION: CORE 1.000  
OX ATOMIZATION LENGTH FOR VAPORIZATION: 1.000  
FUEL ATOMIZATION LENGTH FOR TIMELAGS: 1.000  
OX ATOMIZATION LENGTH FOR TIMELAGS: 1.000  
FUEL DROP SIZE: 1.000  
OX DROP SIZE: 1.000  
MIXING (Em): 1.000  
AO-Multiplier=1.000 CC-Multiplier=1.000 N-Multiplier=1.000 Tau-Multiplier=1.000  
Eta-C\* for XB=0.500

BEGIN STEADY STATE COMBUSTION ANALYSIS  
PC=2141.10 PSIA

PROPELLANT PROPERTIES

FUEL=RP-1	Phase=Liquid	T <sub>man..</sub> F= 71.00	
	Injected Density= 49.69 Lbm/Cu. Ft	Viscosity=1.380E-03 Lbm/Ft-S	Surface Tension=1.657E-03 Lbf/Ft
OX=LOX	Phase=Liquid	T <sub>man..</sub> F=-279.00	
	Injected Density= 70.20 Lbm/Cu. Ft	Viscosity=1.167E-04 Lbm/Ft-S	Surface Tension=7.326E-04 Lbf/Ft

OPERATING CONDITIONS

PC FACE=2141.10 PSIA PC THROAT=2069.29 MIXTURE RATIO= 2.860  
FUEL INJECTION PRESSURE DROP= 673.16 Psia FUEL INJECTION VELOCITY= 363.66 Ft/s  
OX INJECTION PRESSURE DROP= 673.16 Psia OX INJECTION VELOCITY= 296.06 Ft/s  
FUEL FLOWRATE= 46.216 OX FLOWRATE= 163.109

ATOMIZATION OUTPUT

DROPSIZE MODEL=AEROJET

ELEMENT TYPE 1 IS LOL			
FUEL:	ATOMIZATION LENGTH, In.=1.07825	ATOMIZATION LENGTH FOR VAPORIZATION, In.=1.07825	DROPLET RADIUS, Microns= 75.07
OX:	ATOMIZATION LENGTH, In.=2.36256	ATOMIZATION LENGTH FOR VAPORIZATION, In.=2.36256	DROPLET RADIUS, Microns= 63.19

# VAPORIZATION CALCULATIONS

X (In.)	CORE-LOL		BAFFLE-		BARRIER-		FFC-	
	%FUEL VAP	%OX VAP	%FUEL VAP	%OX VAP	%FUEL VAP	%OX VAP	%FUEL VAP	%OX VAP
0.0000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
0.2500	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
0.5000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
0.7500	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.0000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.2500	8.786	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.5000	16.647	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.7500	26.447	0.000	0.000	0.000	0.000	0.000	0.000	0.000
2.0000	32.505	0.000	0.000	0.000	0.000	0.000	0.000	0.000
2.2500	36.163	2.803	0.000	0.000	0.000	0.000	0.000	0.000
2.5000	43.068	27.537	0.000	0.000	0.000	0.000	0.000	0.000
2.7500	46.864	42.768	0.000	0.000	0.000	0.000	0.000	0.000
3.0000	50.659	52.930	0.000	0.000	0.000	0.000	0.000	0.000
3.2500	53.892	60.664	0.000	0.000	0.000	0.000	0.000	0.000
3.5000	56.736	68.436	0.000	0.000	0.000	0.000	0.000	0.000
3.7500	59.585	70.860	0.000	0.000	0.000	0.000	0.000	0.000
4.0000	61.671	74.915	0.000	0.000	0.000	0.000	0.000	0.000
4.2500	64.061	78.140	0.000	0.000	0.000	0.000	0.000	0.000
4.5000	66.192	80.932	0.000	0.000	0.000	0.000	0.000	0.000
4.7500	67.871	82.988	0.000	0.000	0.000	0.000	0.000	0.000
5.0000	69.550	84.988	0.000	0.000	0.000	0.000	0.000	0.000
5.2500	71.147	86.366	0.000	0.000	0.000	0.000	0.000	0.000
5.5000	72.807	88.013	0.000	0.000	0.000	0.000	0.000	0.000
5.7500	74.087	89.164	0.000	0.000	0.000	0.000	0.000	0.000
6.0000	75.365	90.269	0.000	0.000	0.000	0.000	0.000	0.000
6.2500	76.589	91.190	0.000	0.000	0.000	0.000	0.000	0.000
6.5000	77.774	92.067	0.000	0.000	0.000	0.000	0.000	0.000
6.7500	78.717	92.734	0.000	0.000	0.000	0.000	0.000	0.000
7.0000	79.629	93.401	0.000	0.000	0.000	0.000	0.000	0.000
7.2500	80.838	94.061	0.000	0.000	0.000	0.000	0.000	0.000
7.5000	81.412	94.561	0.000	0.000	0.000	0.000	0.000	0.000
7.7500	82.268	95.061	0.000	0.000	0.000	0.000	0.000	0.000
8.0000	83.068	95.641	0.000	0.000	0.000	0.000	0.000	0.000
8.2500	83.878	95.841	0.000	0.000	0.000	0.000	0.000	0.000
8.5000	84.289	96.331	0.000	0.000	0.000	0.000	0.000	0.000
8.7500	84.918	96.831	0.000	0.000	0.000	0.000	0.000	0.000
9.0000	85.540	96.832	0.000	0.000	0.000	0.000	0.000	0.000
9.2500	86.160	97.221	0.000	0.000	0.000	0.000	0.000	0.000
9.5000	86.781	97.421	0.000	0.000	0.000	0.000	0.000	0.000
9.7500	87.401	97.621	0.000	0.000	0.000	0.000	0.000	0.000
10.0000	87.982	97.821	0.000	0.000	0.000	0.000	0.000	0.000
10.2500	88.421	98.031	0.000	0.000	0.000	0.000	0.000	0.000
10.5000	88.891	98.241	0.000	0.000	0.000	0.000	0.000	0.000
10.7500	89.360	98.408	0.000	0.000	0.000	0.000	0.000	0.000
11.0000	89.830	98.559	0.000	0.000	0.000	0.000	0.000	0.000
11.2500	90.278	98.690	0.000	0.000	0.000	0.000	0.000	0.000
11.5000	90.646	98.821	0.000	0.000	0.000	0.000	0.000	0.000
11.7500	91.012	98.937	0.000	0.000	0.000	0.000	0.000	0.000
12.0000	91.360	99.049	0.000	0.000	0.000	0.000	0.000	0.000
12.2500	91.747	99.127	0.000	0.000	0.000	0.000	0.000	0.000
12.5000	92.066	99.206	0.000	0.000	0.000	0.000	0.000	0.000

OVERALL VAPORIZATION EFFICIENCIES  
 FUEL= 92.00% OX= 98.21%

# MASS DISTRIBUTION PROFILE

X (IN)	CORE (lbm/s)		BARRIER (lbm/s)		LOCAL VAPOR MIXTURE RATIO	ETA-C*
	FUEL	OX	FUEL	OX		
0.0000	0.000	0.000	0.000	0.000	0.00	0.0000
0.2850	0.000	0.000	0.000	0.000	0.00	0.0000
0.5700	0.000	0.000	0.000	0.000	0.00	0.0000
0.8550	0.000	0.000	0.000	0.000	0.00	0.0000
1.1400	0.000	0.000	0.000	0.000	0.00	0.0000
1.4250	4.066	0.000	0.000	0.000	0.00	0.0039
1.7100	8.616	0.000	0.000	0.000	0.00	0.0118
1.9950	12.223	0.000	0.000	0.000	0.00	0.0167
2.2800	15.023	0.000	0.000	0.000	0.00	0.0205
2.5650	17.634	3.732	0.000	0.000	0.21	0.0441
2.8500	19.908	38.654	0.000	0.000	1.64	0.3092
3.1350	21.660	55.928	0.000	0.000	2.83	0.4391
3.4200	23.414	70.454	0.000	0.000	3.01	0.5175
3.7050	24.906	80.616	0.000	0.000	3.24	0.5756
3.9900	26.224	88.430	0.000	0.000	3.37	0.6213
4.2750	27.639	94.460	0.000	0.000	3.43	0.6594
4.5600	28.596	98.718	0.000	0.000	3.48	0.6915
4.8450	29.608	104.011	0.000	0.000	3.61	0.7192
5.1300	30.693	107.327	0.000	0.000	3.51	0.7425
5.4150	31.369	110.433	0.000	0.000	3.52	0.7629
5.7000	32.146	112.607	0.000	0.000	3.51	0.7798
5.9850	32.663	114.961	0.000	0.000	3.50	0.7964
6.2700	33.556	117.163	0.000	0.000	3.49	0.8120
6.5550	34.233	118.666	0.000	0.000	3.47	0.8249
6.8400	34.632	120.183	0.000	0.000	3.45	0.8369
7.1250	35.389	121.362	0.000	0.000	3.43	0.8472
7.4100	35.946	122.549	0.000	0.000	3.41	0.8573
7.6950	36.362	123.437	0.000	0.000	3.39	0.8652
7.9800	36.603	124.326	0.000	0.000	3.38	0.8729
8.2650	37.222	125.190	0.000	0.000	3.36	0.8805
8.5500	37.627	125.858	0.000	0.000	3.34	0.8871
8.8350	38.032	126.522	0.000	0.000	3.33	0.8937
9.1200	38.368	127.174	0.000	0.000	3.31	0.8998
9.4050	38.675	127.706	0.000	0.000	3.30	0.9047
9.6900	38.961	128.226	0.000	0.000	3.29	0.9096
9.9750	39.248	128.628	0.000	0.000	3.28	0.9139
10.2600	39.635	129.024	0.000	0.000	3.26	0.9189
10.5450	39.822	129.410	0.000	0.000	3.26	0.9226
10.8300	40.109	129.678	0.000	0.000	3.23	0.9263
11.1150	40.395	129.842	0.000	0.000	3.22	0.9301
11.4000	40.664	130.208	0.000	0.000	3.20	0.9337
11.6850	40.967	130.467	0.000	0.000	3.19	0.9387
11.9700	41.084	130.767	0.000	0.000	3.18	0.9398
12.2550	41.301	130.990	0.000	0.000	3.17	0.9428
12.5400	41.618	131.191	0.000	0.000	3.16	0.9468
12.8250	41.725	131.366	0.000	0.000	3.15	0.9482
13.1100	41.895	131.639	0.000	0.000	3.14	0.9505
13.3950	42.064	131.693	0.000	0.000	3.13	0.9527
13.6800	42.234	131.842	0.000	0.000	3.12	0.9548
13.9650	42.404	131.947	0.000	0.000	3.11	0.9568
14.2500	42.560	132.051	0.000	0.000	3.10	0.9586

# AXIAL PRESSURE PROFILE

X (in)	MACH #	Ptotal (psia)	Pstatic (psia)	Ttotal (°R)	Tstatic (°R)	Wdot (Lbm/s)	Local Radius (in)
1.20	0.000	2141.79	2142.78	1892.28	1886.78	0.59	3.354
1.30	0.001	2141.79	2142.78	1892.28	1886.78	2.00	3.354
1.41	0.006	2141.79	2142.71	1897.15	1890.65	7.08	3.354
1.51	0.008	2141.72	2142.62	1892.28	1886.78	11.10	3.354
2.21	0.010	2141.86	2142.50	1754.02	1747.31	14.73	3.354
2.51	0.012	2141.61	2142.39	1692.28	1686.77	17.83	3.354
2.81	0.064	2136.76	2132.69	5152.58	5141.79	51.60	3.354
3.11	0.112	2128.84	2112.78	6787.38	6767.68	77.20	3.354
3.41	0.137	2119.51	2097.96	6831.43	6831.87	93.48	3.354
3.71	0.158	2113.13	2085.07	6859.00	6828.16	105.81	3.354
4.01	0.171	2107.87	2074.00	6844.88	6811.00	115.34	3.354
4.32	0.183	2102.92	2064.34	6837.28	6802.45	122.98	3.354
4.62	0.193	2098.64	2055.61	6828.59	6792.91	129.44	3.354
4.92	0.202	2094.90	2047.98	6828.13	6789.67	134.78	3.354
5.22	0.208	2091.67	2041.38	6828.24	6789.08	139.18	3.354
5.52	0.216	2088.78	2035.47	6828.31	6787.54	142.99	3.354
5.82	0.220	2086.35	2030.45	6828.00	6786.67	146.09	3.354
6.12	0.228	2083.79	2028.20	6828.70	6789.80	149.28	3.354
6.42	0.230	2081.69	2020.97	6831.13	6791.73	151.94	3.354
6.72	0.234	2078.68	2016.74	6834.50	6794.65	154.18	3.354
7.02	0.238	2077.97	2013.21	6837.33	6797.08	156.17	3.354
7.33	0.241	2076.34	2009.65	6840.75	6800.11	158.01	3.354
7.63	0.244	2075.00	2007.08	6843.28	6802.31	159.82	3.354
7.93	0.248	2073.77	2004.55	6844.92	6803.66	160.88	3.354
8.23	0.249	2072.62	2001.97	6846.44	6804.89	162.26	3.354
8.53	0.253	2071.46	1999.78	6848.37	6806.55	163.41	3.354
8.83	0.255	2070.41	1997.61	6850.29	6808.22	164.54	3.354
9.13	0.256	2069.42	1995.68	6851.75	6809.44	165.80	3.354
9.43	0.258	2068.61	1993.88	6852.88	6810.37	166.48	3.354
9.73	0.258	2067.82	1992.23	6854.07	6811.38	167.30	3.354
10.04	0.259	2067.13	1990.61	6855.58	6812.69	168.01	3.354
10.34	0.260	2066.43	1988.34	6856.88	6813.94	168.75	3.354
10.64	0.261	2065.77	1987.98	6858.63	6815.32	169.42	3.354
10.94	0.263	2065.20	1986.80	6860.29	6816.94	169.98	3.354
11.24	0.264	2064.82	1985.58	6862.04	6818.53	170.56	3.354
11.54	0.265	2064.10	1984.51	6863.00	6819.38	171.11	3.354
11.84	0.270	2063.59	1980.77	6863.01	6818.13	171.83	3.354
12.14	0.290	2063.07	1967.88	6863.05	6813.03	172.12	3.228
12.44	0.327	2062.49	1941.92	6863.08	6802.57	172.57	3.061
12.74	0.376	2061.75	1904.03	6863.12	6788.84	172.99	2.868
13.05	0.438	2060.88	1848.58	6863.15	6783.60	173.38	2.712
13.35	0.524	2059.88	1785.38	6863.17	6728.31	173.70	2.538
13.65	0.655	2058.22	1618.15	6863.20	6687.03	174.04	2.387
13.95	0.818	2056.17	1417.33	6863.23	6583.39	174.34	2.257
14.25	0.998	2053.23	1185.64	6863.28	6421.39	174.81	2.223

# PERFORMANCE SUMMARY

## C\* EFFICIENCY CALCULATIONS (ODK)

INJECTED MR= 2.8800 CSTAR=5880.40 CORE Em=0.8930 BARRIER Em=1.0000  
 CORE: OVERALL MR= 2.8800 VAPOR MR= 3.1027 CSTAR-MIX=5789.54 MASS FRACTION= 1.0000  
 BARRIER: OVERALL MR= 0.0000 VAPOR MR=99.9000 CSTAR-MIX= 0.00 MASS FRACTION= 0.0000  
 ENGINE: OVERALL MR= 2.8800 VAPOR MR= 3.1027 CSTAR-DEL=5817.83  
 C\* EFFICIENCY = 9.588E-01

## ISP EFFICIENCY CALCULATIONS

ISP-ODK, INJ = 2.880E+02 SEC.  
 ISP-ODK, M.Z. INJ = 2.882E+02 SEC.  
 VAPORIZATION EFFICIENCY = 9.650E-01  
 ENERGY RELEASE EFFICIENCY = 9.588E-01

ISP-ODK, M.Z. VAPOR = 2.828E+02 SEC.  
 MIXING EFFICIENCY = 9.935E-01

NOTE: ISP-DEL = ISP-ODK, INJ. \* ERE \* ETADIV \* DELISP-BL

Cohem, In.=9.09E-03	FUEL Cohem, In.=2.293E+02	OX Cohem, In.=8.194E+01
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FUEL: CIn, In=1.874E-02 Lvap, In= 0.868 ATOMIZATION LENGTH USED, In.= 1.078E+00
Temp=3.127E-02 Tatom=2.541E-01 Tvp=1.335E-01 Total=4.187E-01
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OX:  Cln], ln.=1.607E-02  Lvap, ln.= 0.206  ATOMIZATION LENGTH USED, ln.= 2.369E+00
      Timp=5.668E-02      Tatom=6.908E-01      Tvp=6.767E-02      Total=7.780E-01

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FUEL: C10, In.=1.574E-02 Lvap, In.= 0.565  
Timp=3.127E-02 Tatom=2.645E-01 Tvap=1.393E-01  
Total=4.187E-01

DX: Clnj. In.=1.67E-02 Lvap. In.= 0.206  
Timp=5.69E-02 Tailom=6.60E-01 Tvp=5.79E-02 Total=7.76E-01

# CHAMBER-NOZZLE OPTIMIZATION RESULTS

CHAMBER LENGTH (FEET)	ETA-C*	ETA-NOZ	OVERALL EFFICIENCY
0.0000	0.0000	0.8607	0.0000
0.1667	0.0166	0.8726	0.0146
0.3333	0.6226	0.8646	0.8363
0.5000	0.7972	0.8637	0.8606
0.6667	0.8734	0.8419	0.7364
0.8333	0.9143	0.8302	0.7591
1.0000	0.9402	0.8186	0.7386
1.1667	0.9670	0.8057	0.7711
1.3333	0.9899	0.7915	0.7677
1.5000	0.9927	0.7773	0.7639
1.6667	0.9956	0.7631	0.7597
1.8333	1.0000	0.7485	0.7485
2.0000	1.0000	0.7299	0.7299
2.1667	1.0000	0.7114	0.7114
2.3333	1.0000	0.6928	0.6928
2.5000	1.0000	0.6742	0.6742
2.6667	1.0000	0.6461	0.6461
2.8333	1.0000	0.6142	0.6142
3.0000	1.0000	0.5823	0.5823
3.1667	1.0000	0.5504	0.5504
3.3333	1.0000	0.5071	0.5071
3.5000	1.0000	0.4439	0.4439
3.6667	1.0000	0.3808	0.3808
3.8333	1.0000	0.3176	0.3176
4.0000	1.0000	0.2544	0.2544

OPTIMUM CHAMBER LENGTH= 1.1667 FT  
 MAXIMUM OVERALL EFFICIENCY= 0.7711

BEGIN STEADY STATE COMBUSTION ANALYSIS  
PC=1551.40 PSIA

PROPELLANT PROPERTIES

FUEL=RP-1	Phase=Liquid	T <sub>man.</sub> , F= 71.00	
	Injected Density= 49.69 Lbm/Cu. Ft	Viscosity=1.360E-03 Lbm/Ft-s	Surface Tension=1.657E-03 Lbf/Ft
OX=LOX	Phase=Liquid	T <sub>man.</sub> , F=-279.00	
	Injected Density= 69.62 Lbm/Cu. Ft	Viscosity=1.143E-04 Lbm/Ft-s	Surface Tension=7.326E-04 Lbf/Ft

OPERATING CONDITIONS

PC FACE=1551.40 PSIA    PC THROAT=1497.56    MIXTURE RATIO= 2.660  
 FUEL INJECTION PRESSURE DROP= 353.97 Psia    FUEL INJECTION VELOCITY= 256.39 Ft/s  
 OX INJECTION PRESSURE DROP= 356.78 Psia    OX INJECTION VELOCITY= 217.92 Ft/s  
 FUEL FLOWRATE= 33.533    OX FLOWRATE= 96.676

ATOMIZATION OUTPUT

DROPSIZE MODEL=AEROJET

ELEMENT TYPE 1 IS LOL			
FUEL:	ATOMIZATION LENGTH, In.=1.04042	ATOMIZATION LENGTH FOR VAPORIZATION, In.=1.04042	DROPLET RADIUS, Microns= 77.80
OX:	ATOMIZATION LENGTH, In.=2.26933	ATOMIZATION LENGTH FOR VAPORIZATION, In.=2.26933	DROPLET RADIUS, Microns= 85.85

# VAPORIZATION CALCULATIONS

X (in.)	CORE=LOL		BAFFLE=		BARRIER=		FFC=	
	%FUEL VAP	%OX VAP	%FUEL VAP	%OX VAP	%FUEL VAP	%OX VAP	%FUEL VAP	%OX VAP
0.0000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
0.2850	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
0.5700	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
0.8550	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.1400	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.4250	10.171	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.7100	19.343	0.000	0.000	0.000	0.000	0.000	0.000	0.000
2.0000	26.639	0.000	0.000	0.000	0.000	0.000	0.000	0.000
2.2850	32.768	0.000	0.000	0.000	0.000	0.000	0.000	0.000
2.5650	38.255	10.299	0.000	0.000	0.000	0.000	0.000	0.000
2.8500	43.050	31.289	0.000	0.000	0.000	0.000	0.000	0.000
3.1350	46.758	44.877	0.000	0.000	0.000	0.000	0.000	0.000
3.4200	50.467	54.332	0.000	0.000	0.000	0.000	0.000	0.000
3.7050	53.681	61.504	0.000	0.000	0.000	0.000	0.000	0.000
3.9900	56.462	67.049	0.000	0.000	0.000	0.000	0.000	0.000
4.2750	59.244	71.422	0.000	0.000	0.000	0.000	0.000	0.000
4.5600	61.559	75.203	0.000	0.000	0.000	0.000	0.000	0.000
4.8450	63.697	78.300	0.000	0.000	0.000	0.000	0.000	0.000
5.1300	65.636	80.780	0.000	0.000	0.000	0.000	0.000	0.000
5.4150	67.515	82.995	0.000	0.000	0.000	0.000	0.000	0.000
5.7000	69.155	84.857	0.000	0.000	0.000	0.000	0.000	0.000
5.9850	70.770	86.316	0.000	0.000	0.000	0.000	0.000	0.000
6.2700	72.196	87.954	0.000	0.000	0.000	0.000	0.000	0.000
6.5550	73.622	89.076	0.000	0.000	0.000	0.000	0.000	0.000
6.8400	74.970	90.202	0.000	0.000	0.000	0.000	0.000	0.000
7.1250	76.147	91.081	0.000	0.000	0.000	0.000	0.000	0.000
7.4100	77.323	91.961	0.000	0.000	0.000	0.000	0.000	0.000
7.6950	78.355	92.623	0.000	0.000	0.000	0.000	0.000	0.000
7.9800	79.246	93.274	0.000	0.000	0.000	0.000	0.000	0.000
8.2650	80.137	93.926	0.000	0.000	0.000	0.000	0.000	0.000
8.5500	81.004	94.439	0.000	0.000	0.000	0.000	0.000	0.000
8.8350	81.859	94.922	0.000	0.000	0.000	0.000	0.000	0.000
9.1200	82.715	95.410	0.000	0.000	0.000	0.000	0.000	0.000
9.4050	83.346	95.819	0.000	0.000	0.000	0.000	0.000	0.000
9.6900	83.952	96.210	0.000	0.000	0.000	0.000	0.000	0.000
9.9750	84.556	96.529	0.000	0.000	0.000	0.000	0.000	0.000
10.2600	85.165	96.819	0.000	0.000	0.000	0.000	0.000	0.000
10.5450	85.771	97.112	0.000	0.000	0.000	0.000	0.000	0.000
10.8300	86.377	97.337	0.000	0.000	0.000	0.000	0.000	0.000
11.1150	86.983	97.533	0.000	0.000	0.000	0.000	0.000	0.000
11.4000	87.589	97.728	0.000	0.000	0.000	0.000	0.000	0.000
11.6850	88.118	97.932	0.000	0.000	0.000	0.000	0.000	0.000
11.9700	88.577	98.151	0.000	0.000	0.000	0.000	0.000	0.000
12.2550	89.036	98.327	0.000	0.000	0.000	0.000	0.000	0.000
12.5400	89.494	98.481	0.000	0.000	0.000	0.000	0.000	0.000
12.8250	89.953	98.621	0.000	0.000	0.000	0.000	0.000	0.000
13.1100	90.368	98.748	0.000	0.000	0.000	0.000	0.000	0.000
13.3950	90.725	98.872	0.000	0.000	0.000	0.000	0.000	0.000
13.6800	91.084	98.981	0.000	0.000	0.000	0.000	0.000	0.000
13.9650	91.442	99.079	0.000	0.000	0.000	0.000	0.000	0.000
14.2500	91.801	99.155	0.000	0.000	0.000	0.000	0.000	0.000

OVERALL VAPORIZATION EFFICIENCIES  
 FUEL= 91.80% OX= 99.16%

MASS DISTRIBUTION PROFILE

X (IN)	CORE (lbm/s)		BARRIER (lbm/s)		LOCAL VAPOR MIXTURE RATIO		ETA-C*
	FUEL	OX	FUEL	OX			
0.0000	0.000	0.000	0.000	0.000	0.00	0.0000	0.0000
0.2650	0.000	0.000	0.000	0.000	0.00	0.0000	0.0000
0.5700	0.000	0.000	0.000	0.000	0.00	0.0000	0.0000
0.8650	0.000	0.000	0.000	0.000	0.00	0.0000	0.0000
1.1400	0.000	0.000	0.000	0.000	0.00	0.0000	0.0000
1.4250	3.411	0.000	0.000	0.000	0.00	0.0064	0.0064
1.7100	6.487	0.000	0.000	0.000	0.00	0.0122	0.0122
1.9950	9.000	0.000	0.000	0.000	0.00	0.0169	0.0169
2.2800	10.986	0.000	0.000	0.000	0.00	0.0207	0.0207
2.5650	12.828	8.946	0.000	0.000	0.78	0.1202	0.1202
2.8500	14.438	30.216	0.000	0.000	2.09	0.3430	0.3430
3.1350	15.690	43.341	0.000	0.000	2.78	0.4526	0.4526
3.4200	16.923	52.471	0.000	0.000	3.10	0.5261	0.5261
3.7050	18.001	59.396	0.000	0.000	3.30	0.5801	0.5801
3.9900	18.934	64.753	0.000	0.000	3.42	0.6236	0.6236
4.2750	19.666	68.976	0.000	0.000	3.47	0.6604	0.6604
4.5600	20.642	72.628	0.000	0.000	3.52	0.6917	0.6917
4.8450	21.380	75.620	0.000	0.000	3.54	0.7185	0.7185
5.1300	22.077	77.966	0.000	0.000	3.53	0.7415	0.7415
5.4150	22.640	80.154	0.000	0.000	3.54	0.7615	0.7615
5.7000	23.190	81.759	0.000	0.000	3.53	0.7780	0.7780
5.9850	23.732	83.363	0.000	0.000	3.51	0.7945	0.7945
6.2700	24.210	84.943	0.000	0.000	3.51	0.8099	0.8099
6.5550	24.688	86.028	0.000	0.000	3.48	0.8224	0.8224
6.8400	25.140	87.113	0.000	0.000	3.47	0.8347	0.8347
7.1250	25.535	87.983	0.000	0.000	3.44	0.8447	0.8447
7.4100	25.928	88.612	0.000	0.000	3.43	0.8548	0.8548
7.6950	26.275	89.452	0.000	0.000	3.40	0.8630	0.8630
7.9800	26.574	90.081	0.000	0.000	3.39	0.8705	0.8705
8.2650	26.873	90.710	0.000	0.000	3.38	0.8780	0.8780
8.5500	27.169	91.200	0.000	0.000	3.36	0.8846	0.8846
8.8350	27.450	91.672	0.000	0.000	3.34	0.8911	0.8911
9.1200	27.737	92.144	0.000	0.000	3.32	0.8978	0.8978
9.4050	27.949	92.539	0.000	0.000	3.31	0.9028	0.9028
9.6900	28.152	92.916	0.000	0.000	3.30	0.9074	0.9074
9.9750	28.355	93.221	0.000	0.000	3.29	0.9116	0.9116
10.2600	28.559	93.504	0.000	0.000	3.27	0.9169	0.9169
10.5450	28.762	93.767	0.000	0.000	3.26	0.9203	0.9203
10.8300	28.965	94.004	0.000	0.000	3.25	0.9241	0.9241
11.1150	29.168	94.193	0.000	0.000	3.23	0.9278	0.9278
11.4000	29.372	94.382	0.000	0.000	3.21	0.9315	0.9315
11.6850	29.548	94.578	0.000	0.000	3.20	0.9349	0.9349
11.9700	29.709	94.780	0.000	0.000	3.19	0.9381	0.9381
12.2550	29.857	94.981	0.000	0.000	3.18	0.9410	0.9410
12.5400	30.010	95.119	0.000	0.000	3.17	0.9438	0.9438
12.8250	30.164	95.244	0.000	0.000	3.16	0.9465	0.9465
13.1100	30.303	95.367	0.000	0.000	3.15	0.9489	0.9489
13.3950	30.423	95.487	0.000	0.000	3.14	0.9511	0.9511
13.6800	30.543	95.593	0.000	0.000	3.13	0.9532	0.9532
13.9650	30.664	95.686	0.000	0.000	3.12	0.9552	0.9552
14.2500	30.784	95.760	0.000	0.000	3.11	0.9572	0.9572

# AXIAL PRESSURE PROFILE

X (in)	MACH $\phi$	Ptotal (psia)	Pstatic (psia)	Ttotal (R)	Tstatic (R)	Wdot (Lbm/s)	Local Radius (in)
1.20	0.001	1551.85	1552.56	1677.21	1670.62	0.53	3.354
1.30	0.002	1551.85	1552.56	1677.24	1670.65	1.78	3.354
1.61	0.005	1551.83	1552.51	1677.16	1670.79	5.45	3.354
1.81	0.008	1551.80	1552.44	1677.83	1671.53	6.42	3.354
2.21	0.010	1551.77	1552.40	1677.16	1670.79	9.99	3.354
2.51	0.021	1551.45	1551.75	1677.94	1670.82	19.32	3.354
2.61	0.077	1546.67	1542.17	1679.97	1676.25	41.78	3.354
3.11	0.117	1540.07	1528.90	1679.86	1674.14	58.05	3.354
3.41	0.140	1535.06	1518.79	1676.10	1673.01	69.13	3.354
3.71	0.158	1530.64	1509.85	1675.82	1672.58	77.80	3.354
4.01	0.172	1526.85	1502.16	1674.97	1671.75	84.16	3.354
4.32	0.183	1523.51	1495.36	1673.28	1670.17	89.51	3.354
4.62	0.193	1520.48	1488.19	1671.91	1668.96	94.06	3.354
4.92	0.202	1517.66	1483.63	1670.28	1668.61	97.80	3.354
5.22	0.209	1515.55	1479.12	1670.94	1668.59	100.94	3.354
5.52	0.215	1513.52	1474.95	1670.45	1668.51	103.63	3.354
5.82	0.220	1511.79	1471.40	1672.73	1668.27	105.64	3.354
6.12	0.226	1509.95	1467.83	1673.49	1668.48	108.13	3.354
6.42	0.230	1508.40	1464.43	1675.26	1668.77	110.02	3.354
6.72	0.234	1507.02	1461.60	1676.14	1669.22	111.65	3.354
7.02	0.237	1505.78	1459.07	1674.84	1670.24	113.07	3.354
7.33	0.241	1504.65	1456.70	1673.01	1670.35	114.39	3.354
7.63	0.243	1503.65	1454.65	1674.56	1670.55	115.51	3.354
7.93	0.246	1502.78	1452.85	1674.47	1670.20	116.46	3.354
8.23	0.248	1501.59	1451.00	1676.93	1670.36	117.47	3.354
8.53	0.250	1501.12	1448.41	1670.61	1671.01	118.31	3.354
8.83	0.252	1500.37	1447.87	1672.71	1671.66	119.11	3.354
9.13	0.254	1499.63	1446.33	1674.55	1673.26	119.91	3.354
9.43	0.256	1498.46	1445.09	1675.68	1674.20	120.55	3.354
9.73	0.257	1498.46	1443.91	1676.82	1675.18	121.15	3.354
10.04	0.259	1497.96	1442.66	1676.22	1676.22	121.68	3.354
10.34	0.260	1497.46	1441.83	1676.82	1677.82	122.20	3.354
10.64	0.261	1496.98	1440.88	1678.08	1678.92	122.69	3.354
10.94	0.262	1496.56	1439.86	1672.78	1678.45	123.12	3.354
11.24	0.263	1496.16	1438.08	1674.46	1672.00	123.53	3.354
11.54	0.264	1495.75	1436.27	1676.07	1673.48	123.94	3.354
11.84	0.270	1495.36	1435.52	1676.87	1673.06	124.33	3.328
12.14	0.290	1494.97	1426.04	1677.16	1671.36	124.69	3.225
12.44	0.327	1494.54	1407.41	1677.54	1670.56	125.03	3.061
12.74	0.376	1494.02	1380.00	1677.94	1669.65	125.33	2.866
13.05	0.436	1493.36	1340.56	1678.31	1667.14	125.61	2.712
13.35	0.524	1492.58	1279.61	1678.61	1665.50	125.87	2.539
13.65	0.655	1491.46	1173.28	1678.91	1668.56	126.11	2.367
13.95	0.816	1489.78	1027.76	1679.22	1666.21	126.34	2.257
14.25	0.999	1487.55	869.65	1679.55	1640.07	126.54	2.223

# PERFORMANCE SUMMARY

## C\* EFFICIENCY CALCULATIONS (ODK)

INJECTED MR= 2.8800 CSTAR=8880.40 CORE Em=0.8820 BARRIER Em=1.0000  
 CORE: OVERALL MR= 2.8800 VAPOR MR= 3.1107 CSTAR-MIX=6787.38 MASS FRACTION= 1.0000  
 BARRIER: OVERALL MR= 0.0000 VAPOR MR=99.9000 CSTAR-MIX= 0.00 MASS FRACTION= 0.0000  
 ENGINE: OVERALL MR= 2.8800 VAPOR MR= 3.1107 CSTAR-DEL=8889.32  
 C\* EFFICIENCY = 9.572E-01

## ISP EFFICIENCY CALCULATIONS

ISP-ODK, INJ = 2.689E+02 SEC.  
 ISP-ODK, M.Z. INJ = 2.882E+02 SEC. ISP-ODK, M.Z. VAPOR = 2.627E+02 SEC.  
 VAPORIZATION EFFICIENCY = 9.698E-01 MIXING EFFICIENCY = 9.936E-01  
 ENERGY RELEASE EFFICIENCY = 9.572E-01

NOTE: ISP-DEL = ISP-ODK, INJ. \* ERE \* STADIV \* DELISP-BL

TIME-LAG CALCULATIONS, Milliseconds

Cohem, In.=1.126E-02 FUEL Cohem, In.=2.293E+02 OX Cohem, In.=6.194E+01

ELEMENT 1 IS TYPE=LOL

FUEL: CInJ, In.=1.302E-02 Lvap, In.= 0.579 ATOMIZATION LENGTH USED, In.= 1.040E+00  
Timp=4.313E-02 Tatom=3.382E-01 Tvp=1.881E-01 Total=5.694E-01

OX: CInJ, In.=1.330E-02 Lvap, In.= 0.211 ATOMIZATION LENGTH USED, In.= 2.299E+00  
Timp=7.780E-02 Tatom=8.755E-01 Tvp=8.074E-02 Total=1.034E+00

EFFECTIVE TIMELAGS, Milliseconds

FUEL: CInJ, In.=1.302E-02 Lvap, In.= 0.579 Total=5.694E-01  
Timp=4.313E-02 Tatom=3.382E-01 Tvp=1.881E-01

OX: CInJ, In.=1.330E-02 Lvap, In.= 0.211 Total=1.034E+00  
Timp=7.780E-02 Tatom=8.755E-01 Tvp=8.074E-02

BEGIN STEADY STATE COMBUSTION ANALYSIS  
 PC= 767.34 PSIA

PROPELLANT PROPERTIES					
FUEL=RP-1	Phase=Liquid		Tman., F= 71.00		
	Injected Density= 49.69 Lbm/Cu. Ft		Viscosity=1.300E-03 Lbm/Ft-S	Surface Tension=1.857E-03 Lbf/Ft	
OX=LOX	Phase=Liquid		Tman., F=-279.00		
	Injected Density= 69.77 Lbm/Cu. Ft		Viscosity=1.082E-04 Lbm/Ft-S	Surface Tension=7.326E-04 Lbf/Ft	
OPERATING CONDITIONS					
PC FACE= 767.34 PSIA	PC THROAT= 735.66	MIXTURE RATIO= 2.880			
FUEL INJECTION PRESSURE DROP= 87.21 Psia	FUEL INJECTION VELOCITY= 127.27 Ft/S				
OX INJECTION PRESSURE DROP= 88.07 Psia	OX INJECTION VELOCITY= 109.55 Ft/S				
FUEL FLOWRATE= 16.644	OX FLOWRATE= 47.936				
ATOMIZATION OUTPUT					
DROPSIZE MODEL=AEROJET					
ELEMENT TYPE 1 IS LOL					
FUEL:	ATOMIZATION LENGTH, In.=0.98262	ATOMIZATION LENGTH FOR VAPORIZATION, In.=0.98262		DROPLET RADIUS, Microns=	84.09
OX:	ATOMIZATION LENGTH, In.=2.13096	ATOMIZATION LENGTH FOR VAPORIZATION, In.=2.13096		DROPLET RADIUS, Microns=	92.23

VAPORIZATION CALCULATIONS

X (in.)	CORE=LOL			BAFFLE=			BARRIER=			FFC=		
	%FUEL VAP	%OX VAP	%H <sub>2</sub> O VAP	%FUEL VAP	%OX VAP	%H <sub>2</sub> O VAP	%FUEL VAP	%OX VAP	%H <sub>2</sub> O VAP	%FUEL VAP	%OX VAP	%H <sub>2</sub> O VAP
0.0000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
0.2500	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
0.5000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
0.7500	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.0000	2.136	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.2500	12.618	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.5000	20.569	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.7500	27.488	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
2.0000	33.106	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
2.2500	38.298	22.485	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
2.5000	42.882	38.290	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
2.7500	46.412	49.009	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
3.0000	49.931	57.039	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
3.2500	53.126	63.286	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
3.5000	55.778	68.163	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
3.7500	58.418	72.207	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
4.0000	60.914	75.688	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
4.2500	62.845	78.540	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
4.5000	64.876	80.840	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
4.7500	66.685	82.968	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
5.0000	68.251	84.564	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
5.2500	69.908	86.140	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
5.5000	71.285	87.716	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
5.7500	72.619	88.642	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
6.0000	73.973	89.908	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
6.2500	75.200	90.606	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
6.5000	76.318	91.641	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
6.7500	77.433	92.352	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
7.0000	78.393	92.970	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
7.2500	79.239	93.688	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
7.5000	80.085	94.155	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
7.7500	80.910	94.618	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
8.0000	81.722	95.082	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
8.2500	82.534	95.536	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
8.5000	83.187	95.907	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
8.7500	83.762	96.278	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
9.0000	84.338	96.562	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
9.2500	84.913	96.840	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
9.5000	85.489	97.119	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
9.7500	86.064	97.331	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
10.0000	86.639	97.518	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
10.2500	87.240	97.710	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
10.5000	87.884	97.917	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
10.7500	88.325	98.125	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
11.0000	88.760	98.299	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
11.2500	89.185	98.455	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
11.5000	89.631	98.588	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
11.7500	90.066	98.707	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
12.0000	90.438	98.826	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
12.2500	90.777	98.935	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
12.5000	91.117	99.038	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000

OVERALL VAPORIZATION EFFICIENCIES  
FUEL= 91.12% OX= 98.04%

MASS DISTRIBUTION PROFILE

X (IN)	CORE (lbm/s)		BARRIER (lbm/s)		LOCAL VAPOR MIXTURE RATIO	ETA-C*
	FUEL	OX	FUEL	OX		
0.0000	0.000	0.000	0.000	0.000	0.00	0.0000
0.2600	0.000	0.000	0.000	0.000	0.00	0.0000
0.5700	0.000	0.000	0.000	0.000	0.00	0.0000
0.8550	0.000	0.000	0.000	0.000	0.00	0.0000
1.1400	0.356	0.000	0.000	0.000	0.00	0.0013
1.4250	2.100	0.000	0.000	0.000	0.00	0.0060
1.7100	3.427	0.000	0.000	0.000	0.00	0.0130
1.9950	4.575	0.000	0.000	0.000	0.00	0.0174
2.2800	5.510	0.000	0.000	0.000	0.00	0.0209
2.5650	6.374	10.783	0.000	0.000	1.89	0.2556
2.8500	7.138	18.355	0.000	0.000	2.57	0.3961
3.1350	7.725	23.493	0.000	0.000	3.04	0.4772
3.4200	8.311	27.342	0.000	0.000	3.29	0.5386
3.7050	8.844	30.336	0.000	0.000	3.43	0.5879
3.9900	9.264	32.664	0.000	0.000	3.52	0.6270
4.2750	9.723	34.613	0.000	0.000	3.56	0.6611
4.5600	10.122	36.280	0.000	0.000	3.58	0.6910
4.8450	10.460	37.649	0.000	0.000	3.60	0.7159
5.1300	10.786	38.751	0.000	0.000	3.59	0.7377
5.4150	11.101	39.780	0.000	0.000	3.58	0.7578
5.7000	11.360	40.636	0.000	0.000	3.57	0.7735
5.9850	11.519	41.291	0.000	0.000	3.55	0.7892
6.2700	11.662	42.047	0.000	0.000	3.54	0.8044
6.5550	12.067	42.567	0.000	0.000	3.52	0.8167
6.8400	12.312	43.096	0.000	0.000	3.50	0.8286
7.1250	12.518	43.528	0.000	0.000	3.48	0.8391
7.4100	12.702	43.928	0.000	0.000	3.46	0.8486
7.6950	12.868	44.269	0.000	0.000	3.43	0.8575
7.9800	13.046	44.565	0.000	0.000	3.42	0.8651
8.2650	13.169	44.862	0.000	0.000	3.40	0.8729
8.5500	13.329	45.138	0.000	0.000	3.39	0.8791
8.8350	13.467	45.365	0.000	0.000	3.37	0.8853
9.1200	13.602	45.578	0.000	0.000	3.35	0.8914
9.4050	13.737	45.766	0.000	0.000	3.33	0.8975
9.6900	13.846	45.978	0.000	0.000	3.32	0.9024
9.9750	13.942	46.151	0.000	0.000	3.31	0.9070
10.2600	14.037	46.297	0.000	0.000	3.30	0.9110
10.5450	14.133	46.420	0.000	0.000	3.28	0.9151
10.8300	14.229	46.554	0.000	0.000	3.27	0.9191
11.1150	14.325	46.656	0.000	0.000	3.26	0.9226
11.4000	14.420	46.743	0.000	0.000	3.24	0.9253
11.6850	14.520	46.837	0.000	0.000	3.23	0.9298
11.9700	14.626	46.937	0.000	0.000	3.21	0.9338
12.2550	14.701	47.036	0.000	0.000	3.20	0.9369
12.5400	14.778	47.120	0.000	0.000	3.19	0.9397
12.8250	14.848	47.194	0.000	0.000	3.18	0.9424
13.1100	14.916	47.267	0.000	0.000	3.17	0.9449
13.3950	14.991	47.315	0.000	0.000	3.16	0.9474
13.6800	15.052	47.373	0.000	0.000	3.15	0.9497
13.9650	15.109	47.425	0.000	0.000	3.14	0.9517
14.2500	15.166	47.474	0.000	0.000	3.13	0.9537

# AXIAL PRESSURE PROFILE

X (in)	MACH #	Ptotal (psia)	Pstatic (psia)	Ttotal (R)	Tstatic (R)	Wdot (Lbm/s)	Local Radius (in)
1.00	0.000	767.56	767.51	1636.06	1629.81	0.03	3.354
1.30	0.003	767.56	767.51	1642.51	1636.21	1.32	3.354
1.61	0.006	767.56	767.59	1636.08	1629.81	2.93	3.354
1.91	0.009	767.53	767.88	1723.80	1717.29	4.49	3.354
2.21	0.008	767.53	767.85	1636.06	1629.81	4.40	3.354
2.51	0.048	766.55	766.69	4446.27	4427.77	14.69	3.354
2.81	0.099	768.40	768.56	9468.84	9431.18	24.56	3.354
3.11	0.125	760.88	764.48	8580.30	8551.07	30.79	3.354
3.41	0.145	758.63	749.95	8563.62	8533.30	35.54	3.354
3.71	0.181	758.66	745.93	8547.50	8516.23	39.27	3.354
4.01	0.174	754.95	742.48	8533.95	8501.85	42.18	3.354
4.32	0.185	753.40	739.33	8528.37	8485.48	44.65	3.354
4.62	0.194	751.99	736.45	8524.67	8491.04	46.77	3.354
4.92	0.202	750.79	733.99	8523.30	8489.02	48.49	3.354
5.22	0.209	749.69	731.75	8524.97	8490.08	49.98	3.354
5.52	0.215	749.71	729.75	8526.07	8490.93	51.28	3.354
5.82	0.220	747.90	728.08	8526.71	8492.80	52.31	3.354
6.12	0.225	747.03	726.30	8520.34	8493.23	53.41	3.354
6.42	0.229	746.27	724.74	8532.63	8495.78	54.34	3.354
6.72	0.233	746.62	723.39	8536.28	8499.01	55.11	3.354
7.02	0.237	745.00	722.12	8539.84	8502.21	55.83	3.354
7.33	0.240	744.45	720.99	8542.94	8504.97	56.46	3.354
7.63	0.243	743.94	719.94	8546.43	8506.14	57.04	3.354
7.93	0.245	743.50	719.02	8549.63	8511.06	57.53	3.354
8.23	0.247	743.08	718.15	8551.91	8513.08	58.00	3.354
8.53	0.250	742.68	717.33	8553.66	8514.57	58.44	3.354
8.83	0.251	742.35	716.60	8555.64	8516.24	58.82	3.354
9.13	0.253	741.98	715.97	8557.95	8517.84	59.20	3.354
9.43	0.255	741.63	715.18	8559.15	8519.41	59.56	3.354
9.73	0.257	741.36	714.88	8560.49	8520.57	59.86	3.354
10.04	0.258	741.09	714.02	8561.88	8521.50	60.14	3.354
10.34	0.259	740.86	713.55	8562.98	8522.75	60.38	3.354
10.64	0.261	740.62	713.08	8564.31	8523.94	60.63	3.354
10.94	0.262	740.40	712.69	8565.72	8525.20	60.86	3.354
11.24	0.263	740.20	712.17	8567.34	8526.87	61.06	3.354
11.54	0.264	740.00	711.78	8568.97	8528.19	61.26	3.354
11.84	0.270	739.79	710.37	8570.78	8528.83	61.47	3.328
12.14	0.260	739.57	706.95	8572.12	8526.45	61.67	3.325
12.44	0.327	739.36	696.47	8572.73	8519.50	61.84	3.081
12.74	0.375	739.06	682.96	8573.26	8503.85	62.00	2.866
13.05	0.436	738.76	663.51	8573.88	8492.01	62.15	2.712
13.35	0.523	738.32	633.41	8574.48	8448.84	62.28	2.539
13.65	0.655	737.72	590.91	8574.95	8385.77	62.41	2.387
13.95	0.817	736.67	509.26	8575.40	8291.66	62.53	2.257
14.25	0.989	735.66	425.76	8575.85	8171.04	62.64	2.223

# PERFORMANCE SUMMARY

## C\* EFFICIENCY CALCULATIONS (ODK)

INJECTED MR= 2.8800 CSTAR=5880.40 CORE Em=0.8930 BARRIER Em=1.0000  
 CORE: OVERALL MR= 2.8800 VAPOR MR= 3.1304 CSTAR-MIX=5782.01 MASS FRACTION= 1.0000  
 BARRIER: OVERALL MR= 0.0000 VAPOR MR=99.9000 CSTAR-MIX= 0.00 MASS FRACTION= 0.0000  
 ENGINE: OVERALL MR= 2.8800 VAPOR MR= 3.1304 CSTAR-DEL=5588.98  
 C\* EFFICIENCY = 9.537E-01

## ISP EFFICIENCY CALCULATIONS

ISP-ODK, INJ = 2.888E+02 SEC.  
 ISP-ODK, M.Z. INJ = 2.852E+02 SEC. ISP-ODK, M.Z. VAPOR = 2.824E+02 SEC.  
 VAPORIZATION EFFICIENCY = 9.588E-01 MIXING EFFICIENCY = 9.935E-01  
 ENERGY RELEASE EFFICIENCY = 9.538E-01

NOTE: ISP-DEL = ISP-ODK, INJ. \* ERE \* ETADIV - DELISP-BL

TIME-LAG CALCULATIONS, Milliseconds

Cohem, In.=1.791E-02      FUEL Cohem, In.=2.293E+02      OX Cohem, In.=8.184E+01

ELEMENT 1 IS TYPE=LOL

FUEL:      CInj, In.=8.821E-03      Lvap, In.= 0.010      ATOMIZATION LENGTH USED, In.= 9.826E-01  
              Timp=8.888E-02      Tatom=8.303E-01      Tvp=3.893E-01      Total=1.118E+00

OX:      CInj, In.=8.808E-03      Lvap, In.= 0.223      ATOMIZATION LENGTH USED, In.= 2.131E+00  
              Timp=1.548E-01      Tatom=1.821E+00      Tvp=1.893E-01      Total=1.945E+00

EFFECTIVE TIMELAGS, Milliseconds

FUEL:      CInj, In.=8.821E-03      Lvap, In.= 0.010      Total=1.118E+00  
              Timp=8.888E-02      Tatom=8.303E-01      Tvp=3.893E-01

OX:      CInj, In.=8.808E-03      Lvap, In.= 0.223      Total=1.945E+00  
              Timp=1.548E-01      Tatom=1.821E+00      Tvp=1.893E-01

# LOW FREQUENCY COMBUSTION STABILITY CALCULATIONS

ROCKET COMBUSTOR INTERACTIVE DESIGN METHODOLOGY  
Version 23-FEB-91

ROCCID

DIRECT INPUT ECHO FROM SUBROUTINE SINPUT

ROCCID POINT DESIGN TEST CASE 1  
LOX/RP-1 LIKE DOUBLET PAIR WITH FIXED PC  
APPROXIMATES -0100 SUBSCALE DOUBLET

\$MODELS

MCHAM= 1

MBURN= 2

MINJ= 1

\$END

\$OPCOND

PISPA=

8.9600E+01, 1.0940E+02, 1.3940E+02, 1.5860E+02, 1.7780E+02,  
2.0870E+02, 2.3650E+02, 2.6470E+02, 2.8700E+02, 2.7040E+02,  
2.7180E+02, 2.7190E+02, 2.7190E+02, 2.8980E+02, 2.8920E+02,  
2.8860E+02, 2.8500E+02, 2.8210E+02, 2.5920E+02, 2.5630E+02,  
2.5080E+02, 2.3840E+02, 2.2750E+02, 2.1100E+02, 2.0140E+02,  
1.5660E+02, 1.0970E+02, 6.0000E+01,

PCSA=

1.7760E+03, 2.4560E+03, 3.4870E+03, 4.0740E+03, 4.6800E+03,  
5.0700E+03, 5.4540E+03, 5.7340E+03, 5.9080E+03, 5.9610E+03,  
5.9690E+03, 5.9540E+03, 5.9340E+03, 5.9090E+03, 5.8820E+03,  
5.8550E+03, 5.8260E+03, 5.7590E+03, 5.7000E+03, 5.6430E+03,  
5.5320E+03, 5.2760E+03, 5.0470E+03, 4.8760E+03, 4.4200E+03,  
3.4570E+03, 2.4940E+03, 1.4650E+03,

FUEL=

'RP-1'

,

'LOX'

PC=

2.1411E+03

XMR=

2.8800E+00

FTMAN=

7.1000E+01

XTMAN=

-2.7800E+02

EMMAN=

1.0000E+00

NPERFP=

28

PMRA=

1.0000E-01, 3.0000E-01, 6.0000E-01, 8.0000E-01, 1.0000E+00,  
1.2500E+00, 1.5000E+00, 1.7500E+00, 2.0000E+00, 2.2000E+00,  
2.4000E+00, 2.6000E+00, 2.8000E+00, 3.0000E+00, 3.2000E+00,  
3.4000E+00, 3.6000E+00, 3.8000E+00, 4.0000E+00, 4.2000E+00,  
4.4000E+00, 4.6000E+00, 4.8000E+00, 5.0000E+00, 5.2000E+00,  
5.4000E+00, 5.6000E+00, 5.8000E+00, 6.0000E+00, 6.2000E+00,  
6.4000E+00, 6.6000E+00, 6.8000E+00, 7.0000E+00, 7.2000E+00,  
7.4000E+00, 7.6000E+00, 7.8000E+00, 8.0000E+00, 8.2000E+00,  
8.4000E+00, 8.6000E+00, 8.8000E+00, 9.0000E+00, 9.2000E+00,  
9.4000E+00, 9.6000E+00, 9.8000E+00, 1.0000E+01,

\$SEND

\$GEOM

ROHAMB=

2.7949E-01

RTHRT=

1.8523E-01

RNE=

1.1100E-01

RTE=

1.1100E-01

ALPHA=

3.0000E+01

CHAMBL=

1.1875E+00

XC=

9.6475E-01

\$SEND

\$SCORE

TYPE=

'LOL'

NEL=

80

FCD=

6.6346E-02

FIH=

9.1000E-01

FIA=

1.1491E-01

3.0000E+01

FCANT=	1.6000E+01	
FFACET=	2.6728E-01	
XDJ=	1.0172E-01	
XCD=	9.4000E-01	
XIH=	1.7819E-01	
XIA=	3.0000E+01	
XCANT=	1.6000E+01	
XFACET=	4.4047E-01	
EMJN1=	6.5000E-01	
SEND		
\$BAFFLE		
\$END		
\$BARRIER		
\$END		
\$FFC		
\$END		
\$BURN		
GAMMA=	1.1462E+00	
AO=	4.0656E+03	
GMW=	2.3218E+01	
GPR=	5.9752E-01	
GK=	5.4160E-05	
GMU=	6.6324E-06	
RML=	7.5067E+01	
VJL=	3.5356E+02	
TJL=	6.3100E+02	
RHOL=	4.9695E+01	
CPL=	4.7321E-01	
PCRTIL=	3.1500E+02	
TCRTIL=	1.2180E+03	
TBOILL=	6.6200E+02	
XMWL=	1.7200E+02	
HVAPL=	1.2500E+02	
EN=	6.6793E-01	
TAUGEN=	1.3326E-04	
ISEN=	1	
\$END		
\$INJ		
FMAND=	6.7078E+00	
XMAND=	6.7078E+00	
FMANL=	3.3536E+00	
XMANL=	3.3536E+00	
PCA=	2.1411E+03	1.5514E+03, 7.6794E+02
FRA=	6.2681E-01	4.5633E-01, 2.2731E-01
FCAPA=	7.8914E-04	6.1781E-04, 8.8053E-04
NFE=	4	
FTLA=	4.1866E-04	5.8540E-04, 1.1164E-03
FINA=	4.4707E-06	4.4767E-06, 4.4924E-06
FFA=	1.0000E+00	1.0000E+00, 1.0000E+00
NXE=	4	
XRA=	6.2881E-01	4.5998E-01, 2.3216E-01
XCAPA=	6.9193E-04	6.1747E-04, 6.8616E-04
XTLA=	7.7504E-04	1.0340E-03, 1.9451E-03
XINA=	5.3932E-05	8.4084E-06, 8.4389E-06
XFA=	1.0000E+00	1.0000E+00, 1.0000E+00
XUOR=	6.2931E-03	1.0000E-01
AUOR=	3.4571E-03	3.1265E-03
RUOR=	3.3172E-02	5.0960E-02
XOR=	6.2931E-02	1.0000E-01
AOR=	3.4571E-03	6.1266E-03
ROR=	3.3172E-02	5.0960E-02
XDOR=	1.6586E-01	3.0861E-01
ADOR=	3.4571E-03	3.1265E-03
NDOR=	3.3172E-02	5.0960E-02

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WDOT= 4.8218E+01, 1.3311E+02, 3.3333E+01, 9.8578E+01, 1.8644E+01,
4.7939E+01
$END
$CHAMBER
XB= 2.6125E-01
ZS= 1.1875E-01
ZE= 4.9875E-01
MJB= 3
T= 3.3000E-02
CC1= 5*2.5000E+03
GAMC1= 5*1.2000E+00
CC2= 5*2.5000E+03
GAMC2= 5*1.2000E+00
$END
$FDORC
ZCONB= 9.5000E-01
NZON= 5
FTER= 2.1988E-02, 7.4060E-01, 8.9632E-01, 9.8388E-01, 1.0000E+00,
15*1.0000E+00
CCAV1= 20*2.5000E+03
CGAM1= 20*1.2000E+00
XMMC1= 20*1.7314E+01
RHOAP1= 20*1.8048E+00
CCAV2= 20*2.5000E+03
CGAM2= 20*1.2000E+00
XMMC2= 20*1.7314E+01
RHOAP2= 20*1.8048E+00
$END
$MIX
EM(1)= 8.930E-01, 1.000E+00
$END
$DEBUGC
DEBUG= F
$END
$HIFIC
SHORT= F
POC= 2.0000E-01, 2.0000E-01
$END
$DISTSDC
SHORT= F
POC= 2.0000E-01, 2.0000E-01
PAMP= 2.0000E-01
MC= 11
LC= 8
MB= 11
LB= 8
IDMAX= 10
$END
$CRPC
NDPC= 18
NDTFQ= 18
NDTLF= 1000
NPRINT= 80
NSLMS= 3500
PAMPC= 2.0000E-01
NRAD= 3
NCIRC= 5
$END
$LEINJC
IDMEM= 2
PAMPCH= 2.0000E-01
NTINJ= 18
USGF= 2.0000E-02

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OGF= 2.0000E-02
DSGF= 2.0000E-02
FGDO= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00
XCDO= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00
$END
$COMBUSTC
FALVM= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00
XALVM= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00
FALTM= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00
XALTM= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00
FRMM= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00
XRMM= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00
EMMULT= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00
COMBXS= 5.0000E-01
ACMULT= 1.0000E+00
CCMULT= 1.0000E+00
ENMULT= 1.0000E+00
TAMULT= 1.0000E+00
$END
$FDORCC
SHORT= F
EPSIL= 2.0000E+01
ERROR= 1.0000E-01
LTS= 5
NTS= 5
NTS= 10
ITMAX= 100
RELX= 5.5000E-01
NEET= 3
NXFST= 3
NYFST= 3
NAFST= 3
MORE= F
$END

```

END OF INPUT ECHO

# STABILITY MODEL INPUTS

## RUN DESCRIPTOR

ROCCID POINT DESIGN TEST CASE 1  
LOX/FP-1 LIKE DOUBLET PAIR WITH FIXED PC  
APPROXIMATES -0100 SUBSCALE DOUBLET

## SELECTED MODELS

BURNING MODEL=N-TAU INJECTION MODEL=INJ CHAMBER MODEL=HIFI

AXISYMETRIC=T DEBUG OUTPUT=F

## CHAMBER GEOMETRY AND OPERATING CONDITIONS

CHAMBER RADIUS, FT= 0.2795 THROAT RADIUS, FT= 0.1862  
CYLINDRICAL SECTION, FT= 0.9847 CONVERGENCE HALF-ANGLE, DEG=30.0000  
NOZZLE ENTRANCE RADIUS OF CURVATURE, FT= 0.1110 THROAT ENTRANCE RADIUS OF CURVATURE, FT= 0.1110  
CHAMBER PRESSURE, PSIA=2141.10 MIXTURE RATIO= 2.8800  
SOUND SPEED, FT/SEC=4086.80 GAMMA=1.1462

## N-TAU BURNING MODEL INPUTS

PRESSURE INTERACTION INDEX, EN= 0.8879 SENSITIVE TIMELAG, TAU, SEC= 1.338E-04 SENSITIVE CIRCUIT=FUEL

## LUMPED INJECTION MODEL INPUTS

PC, PSIA=2141.10

	% TOTAL FLOW	RESISTANCE	INERTANCE (SEC)	CAPACITANCE (SEC)	TIMELAG (SEC)
FUEL:	100.000	8.288E-01	4.471E-08	7.891E-04	4.187E-04
	0.000	8.288E-01	0.000E+00	7.891E-04	0.000E+00
	0.000	8.288E-01	0.000E+00	7.891E-04	0.000E+00
	0.000	8.288E-01	0.000E+00	7.891E-04	0.000E+00
OX:	100.000	8.288E-01	8.388E-08	8.918E-04	7.750E-04
	0.000	8.288E-01	0.000E+00	8.918E-04	0.000E+00
	0.000	8.288E-01	0.000E+00	8.918E-04	0.000E+00
	0.000	8.288E-01	0.000E+00	8.918E-04	0.000E+00

PC, PSIA=1851.40

% TOTAL FLOW	RESISTANCE	INERTANCE	CAPACITANCE	TIMELAG
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FUEL:	100.000	4.593E-01	4.477E-05	9.179E-04	5.604E-04
	0.000	4.593E-01	0.000E+00	9.179E-04	0.000E+00
	0.000	4.593E-01	0.000E+00	9.179E-04	0.000E+00
	0.000	4.593E-01	0.000E+00	9.179E-04	0.000E+00
OX:	100.000	4.600E-01	6.409E-05	9.175E-04	1.034E-03
	0.000	4.600E-01	0.000E+00	9.175E-04	0.000E+00
	0.000	4.600E-01	0.000E+00	9.175E-04	0.000E+00
	0.000	4.600E-01	0.000E+00	9.175E-04	0.000E+00

PC. PSIA= 767.34

% TOTAL FLOW	RESISTANCE	INERTANCE (SEC)	CAPACITANCE (SEC)	TIMELAG (SEC)
FUEL:	2.275E-01	4.492E-05	6.605E-04	1.116E-03
	0.000	0.000E+00	6.605E-04	0.000E+00
	0.000	0.000E+00	6.605E-04	0.000E+00
	0.000	0.000E+00	6.605E-04	0.000E+00
OX:	2.322E-01	6.439E-05	6.561E-04	1.945E-03
	0.000	0.000E+00	6.561E-04	0.000E+00
	0.000	0.000E+00	6.561E-04	0.000E+00
	0.000	0.000E+00	6.561E-04	0.000E+00

# HIFI CHAMBER MODEL INPUTS

COMBUSTION PLANE, FT= 2.612E-01      SHORT NOZZLE ASSUMED=F

## ACOUSTIC CAVITY INPUTS

CAVITY TYPE=NONE  
CAVITY TYPE 1:    NUMBER OF CAVITIES= 0      NUMBER OF PROPERTY SECTIONS= 1  
CAVITY TYPE 2:    NUMBER OF CAVITIES= 0      NUMBER OF PROPERTY SECTIONS= 1  
PARTITION THICKNESS, FT= 0.0000

PC (PSIA)	MAX. AMPLITUDE	FREQUENCY (HZ)	PHASE MARGIN (DEG)
2.141E+03	2.974E-01	5.450E+02	180.00
2.091E+03	3.076E-01	5.376E+02	180.00
2.041E+03	3.188E-01	5.316E+02	180.00
1.991E+03	3.267E-01	5.262E+02	180.00
1.941E+03	3.351E-01	5.180E+02	180.00
1.891E+03	3.449E-01	5.125E+02	180.00
1.841E+03	3.542E-01	5.060E+02	180.00
1.791E+03	3.640E-01	4.994E+02	180.00
1.741E+03	3.739E-01	4.926E+02	180.00
1.691E+03	3.837E-01	4.861E+02	180.00
1.641E+03	3.936E-01	4.793E+02	180.00
1.591E+03	4.034E-01	4.724E+02	180.00
1.541E+03	4.130E-01	4.653E+02	180.00
1.491E+03	4.224E-01	4.581E+02	180.00
1.441E+03	4.319E-01	4.508E+02	180.00
1.391E+03	4.409E-01	4.433E+02	180.00
1.341E+03	4.489E-01	4.355E+02	180.00
1.291E+03	4.563E-01	4.276E+02	180.00
1.241E+03	4.617E-01	4.194E+02	180.00
1.191E+03	4.671E-01	4.109E+02	180.00
1.141E+03	4.711E-01	4.021E+02	180.00
1.091E+03	4.739E-01	3.929E+02	180.00
1.041E+03	4.760E-01	3.833E+02	180.00
9.911E+02	4.744E-01	3.733E+02	180.00
9.411E+02	4.720E-01	3.627E+02	180.00
8.911E+02	4.676E-01	3.516E+02	180.00
8.411E+02	4.611E-01	3.399E+02	180.00
7.911E+02	4.524E-01	3.276E+02	180.00
7.411E+02	4.526E-01	3.156E+02	180.00
6.911E+02	5.050E-01	3.186E+02	180.00
6.411E+02	5.492E-01	3.909E+02	180.00
5.911E+02	5.741E-01	5.918E+02	180.00
5.411E+02	6.707E-01	6.306E+02	180.00
4.911E+02	7.656E-01	7.626E+02	-75.41
4.411E+02	1.087E+00	7.088E+02	24.25
4.578E+02	9.794E-01	7.252E+02	-5.97
4.522E+02	1.016E+00	7.190E+02	4.26
4.541E+02	1.003E+00	7.211E+02	0.89

MARGINAL CHUG POINT FOUND:  
PC= 484.06 PSIA FREQUENCY= 721.07 HZ

#### CHUG STABILITY ITERATION SUMMARY

THE CURRENT CONFIGURATION IS CHUG STABLE

DESIRED MARGINAL PC = 600.06 PSIA  
CURRENT MARGINAL PC = 484.06 PSIA  
CURRENT CHUG MARGIN = 345.94 PSI  
CHUG FREQUENCY = 721.07 HZ

# REDESIGNED CHAMBER RESULTS

NOMINAL CHAMBER PRESSURE = 2.141E+03 PSIA  
 THROTTLED CHAMBER PRESSURE = 1.826E+03 PSIA  
 FUEL INJECTION PRESSURE DROP = 6.090E+02 PSI  
 OX INJECTION PRESSURE DROP = 6.090E+02 PSI  
 CHAMBER RADIUS = 2.796E-01 FT  
 THROAT RADIUS = 1.852E-01 FT  
 NOZZLE ENTRANCE RADIUS OF CURVATURE = 1.110E-01 FT  
 THROAT ENTRANCE RADIUS OF CURVATURE = 1.110E-01 FT  
 NOZZLE CONVERGENCE HALF-ANGLE = 3.000E+01 DEG  
 INJECTOR-TO-THROAT CHAMBER LENGTH = 1.188E+00 FT  
 BARREL SECTION LENGTH = 9.848E-01 FT

## IMPINGING ELEMENT SIZING RESULTS

ELEMENT TYPE = -LOL  
 NO. OF ELEMENTS = 69  
 FUEL ORIFICE DIAMETER = 6.634E-02 IN  
 OX ORIFICE DIAMETER = 1.017E-01 IN

## CORE ELEMENT SPACING RESULTS

ELEMENT TYPE = LOL  
 NUMBER OF ELEMENTS = 69  
 FUEL ORIFICE/ANNULUS DIAMETER = 6.634E-02 IN  
 OX'DIZER ORIFICE DIAMETER = 1.017E-01 IN  
 FUEL INJECTION VELOCITY = 3.076E+02 FT/S  
 OXIDIZER INJECTION VELOCITY = 2.592E+02 FT/S

ROW	# ELEMENTS	MID-ROW RADIUS (IN)
1	6	9.115E-01
2	12	1.809E+00
3	21	2.907E+00
4	30	3.996E+00

# LOW FREQUENCY COMBUSTION STABILITY CALCULATIONS

ROCCID  
Rocket Combustor Interactive Design Methodology  
Version 23-FEB-91

DIRECT INPUT ECHO FROM SUBROUTINE SINPUT

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ROCCID POINT DESIGN TEST CASE 1
LOX/RP-1 LIKE DOUBLET PAIR WITH FIXED PC
APPROXIMATES .0100 SUBSCALE DOUBLET
$MODELS
  $CHAM= 1
  $BURN= 2
  $INJ= 1
  $END
$OPCOND
  PISPA= 8.9800E+01, 1.0940E+02, 1.3940E+02, 1.5880E+02, 1.7780E+02,
    2.0970E+02, 2.3650E+02, 2.6470E+02, 2.8700E+02, 2.7040E+02,
    2.7180E+02, 2.7190E+02, 2.7190E+02, 2.9880E+02, 2.6820E+02,
    2.8860E+02, 2.6500E+02, 2.8210E+02, 2.8920E+02, 2.5830E+02,
    2.5080E+02, 2.3640E+02, 2.2760E+02, 2.1100E+02, 2.0140E+02,
    1.5560E+02, 1.0970E+02, 8.0000E+01,
  PCSA= 1.7780E+03, 2.4680E+03, 3.4870E+03, 4.0740E+03, 4.6800E+03,
    5.0700E+03, 5.4540E+03, 5.7840E+03, 5.9060E+03, 5.9610E+03,
    5.9690E+03, 5.9540E+03, 5.9340E+03, 5.9090E+03, 5.8820E+03,
    5.8550E+03, 5.8280E+03, 5.7990E+03, 5.7590E+03, 5.7000E+03,
    5.6430E+03, 5.5320E+03, 5.2760E+03, 5.0470E+03, 4.8760E+03,
    4.4200E+03,
    3.4570E+03, 2.4940E+03, 1.4850E+03,
  FUEL= 'RP-1',
  OX= 'LOX',
  PC= 2.1411E+03
  XMR= 2.9800E+00
  FTMAN= 7.1000E+01
  XTMAN= -2.7900E+02
  EMMAN= 1.0000E+00
  NPERFP= 28
  PMRA= 1.0000E-01, 9.0000E-01, 9.0000E-01, 9.0000E-01, 1.0000E+00,
    1.2800E+00, 1.5000E+00, 1.7800E+00, 2.0000E+00, 2.2000E+00,
    2.4000E+00, 2.5000E+00, 2.6000E+00, 2.7000E+00, 2.8000E+00,
    2.9000E+00, 3.0000E+00, 3.2000E+00, 3.4000E+00, 3.6000E+00,
    4.0000E+00, 5.0000E+00, 6.0000E+00, 8.0000E+00, 1.0000E+01,
    1.5000E+01, 2.0000E+01, 5.0000E+01,
  $END
$GECM
  $CHAMB= 2.7949E-01
  $THRT= 1.8828E-01
  $RNE= 1.1100E-01
  $RTE= 1.1100E-01
  $ALPHA= 3.0000E+01
  $CHAMBL= 1.1878E+00
  $XC= 8.8477E-01
  $END
$SCORE
  $TYPE= 'LOL',
  $NEL= 80
  $FDJ= 8.8348E-02
  $FCD= 9.1000E-01
  $FIH= 1.1481E-01
  $FIA= 3.0000E+01

```

FCANT=	1.0000E+01	
FFACET=	2.8728E-01	
XDJ=	1.0178E-01	
XCD=	9.4000E-01	
XIH=	1.7618E-01	
XIA=	3.0000E+01	
XCANT=	1.0000E+01	
XFACET=	4.4047E-01	
EMUNI=	8.5000E-01	
\$END		
\$BAFFLE		
\$END		
\$BARRIER		
\$END		
\$FFC		
\$END		
\$BURN		
GAMMA=	1.1482E+00	
AO=	4.0666E+03	
GAM=	2.3218E+01	
GPR=	5.9752E-01	
GK=	5.4180E-05	
GMU=	6.6324E-05	
RML=	7.8242E+01	
VJL=	3.0748E+02	
TJL=	5.3100E+02	
RHOL=	4.9895E+01	
CPL=	4.7321E-01	
PCRITL=	3.1500E+02	
TCRITL=	1.2180E+03	
TBOILL=	8.6200E+02	
XMWL=	1.7200E+02	
HVAPL=	1.2500E+02	
EN=	8.6788E-01	
TAUSEN=	1.4114E-04	
ISEN=	1	
\$END		
\$INJ		
FMAND=	6.7076E+00	
XMAND=	6.7076E+00	
FMANL=	3.3636E+00	
XMANL=	3.3636E+00	
PCA=	2.1411E+03	1.5514E+03, 7.0734E+02
FRA=	4.7647E-01	3.4543E-01, 1.7208E-01
FCAPA=	7.6914E-04	6.1791E-04, 8.6053E-04
NFE=	4	
FTLA=	4.6146E-04	6.6481E-04, 1.2839E-03
FINA=	3.8876E-05	3.8928E-05, 3.9064E-05
FFA=	1.0000E+00	1.0000E+00, 1.0000E+00
NXE=	4	
XRA=	4.7647E-01	3.4543E-01, 1.7205E-01
XCAPA=	8.9183E-04	6.1747E-04, 8.6615E-04
XTLA=	8.9130E-04	1.1991E-03, 2.2389E-03
XINA=	7.3028E-05	7.3125E-05, 7.3382E-05
XFA=	1.0000E+00	1.0000E+00, 1.0000E+00
XUOR=	4.7680E-02	7.9412E-02
AUOR=	3.4570E-03	6.1269E-03
RUOR=	3.3172E-02	6.0861E-02
XOR=	4.7660E-02	7.9412E-02
AOR=	3.4570E-03	6.1269E-03
ROR=	3.3172E-02	5.0661E-02
XOOR=	9.5761E-02	1.4882E-01
ADOR=	3.4570E-03	8.1269E-03
RDOR=	3.3172E-02	5.0661E-02

WDOT= 4.8218E+01, 1.3311E+02, 3.3533E+01, 9.8578E+01, 1.8644E+01,  
 4.7935E+01  
 \$END  
 \$CHAMBER  
 XB= 2.6125E-01  
 ZS= 1.1875E-01  
 ZE= 4.9875E-01  
 MJB= 3  
 T= 3.3000E-02  
 CC1= 5\*2.5000E+03  
 GAMC1= 5\*1.2000E+00  
 CC2= 5\*2.5000E+03  
 GAMC2= 5\*1.2000E+00  
 \$END  
 \$FDORC  
 NZON= 6  
 FTER= 20\*1.0000E+00  
 \$END  
 \$MIX  
 EM(1)= 9.003E-01, 1.000E+00  
 \$END  
 \$DEBUGC  
 DEBUG= F  
 \$END  
 \$HIFIC  
 SHORT= F  
 POC= 2.0000E-01, 2.0000E-01  
 \$END  
 \$DIST3DC  
 SHORT= F  
 POC= 2.0000E-01, 2.0000E-01  
 PAMP= 2.0000E-01  
 MC= 11  
 LG= 8  
 MB= 11  
 LB= 6  
 IDMAX= 10  
 \$END  
 \$CRPC  
 NDCP= 16  
 NDTFQ= 16  
 NDTLF= 1000  
 NPRINT= 50  
 NSUMS= 3500  
 PAMP= 2.0000E-01  
 NRAD= 3  
 NCIRC= 5  
 \$END  
 \$LEINJC  
 IDMEM= 2  
 PAMPCH= 2.0000E-01  
 NTINJ= 16  
 USGF= 2.0000E-02  
 OGF= 2.0000E-02  
 DSGF= 2.0000E-02  
 FCDO= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00  
 XCDO= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00  
 \$END  
 \$COMBUSTC  
 FALVM= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00  
 XALVM= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00  
 FALTM= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00  
 XALTM= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00  
 FRMM= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00

```

XMM= 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00
ENMULT= 1.000E+00, 1.000E+00
COMB= 5.000E-01
ACMULT= 1.000E+00
CCMULT= 1.000E+00
ENMULT= 1.000E+00
TAMULT= 1.000E+00
$END
$FORCC
SHORT= F
EPIL= 2.000E+01
ERROR= 1.000E-01
LTS= 5
MTS= 5
NTS= 10
ITMAX= 100
RELX= 6.500E-01
NEET= 3
NXFT= 3
NYFT= 3
NAFT= 3
MORE= F
$END
END OF INPUT ECHO

```

# STABILITY MODEL INPUTS

## RUN DESCRIPTOR

ROCCID POINT DESIGN TEST CASE 1  
LOX/JP-1 LIKE DOUBLET PAIR WITH FIXED PC  
APPROXIMATES .0100 SUBSCALE DOUBLET

## SELECTED MODELS

BURNING MODEL=N-TAU INJECTION MODEL=INJ CHAMBER MODEL=HIFI

AXISYMETRIC=T DEBUG OUTPUT=F

## CHAMBER GEOMETRY AND OPERATING CONDITIONS

CHAMBER RADIUS, FT= 0.2795 THROAT RADIUS, FT= 0.1852  
CYLINDRICAL SECTION, FT= 0.9848 CONVERGENCE HALF-ANGLE, DEG=30.0000  
NOZZLE ENTRANCE RADIUS OF CURVATURE, FT= 0.1110 THROAT ENTRANCE RADIUS OF CURVATURE, FT= 0.1110  
CHAMBER PRESSURE, PSIA=2141.10 MIXTURE RATIO= 2.8800  
SOUND SPEED, FT/SEC=4085.80 GAMMA=1.1462

## N-TAU BURNING MODEL INPUTS

PRESSURE INTERACTION INDEX, EN= 0.8879 SENSITIVE TIMELAG, TAU, SEC= 1.411E-04 SENSITIVE CIRCUIT=FUEL

## LUMPED INJECTION MODEL INPUTS

PC, PSIA=2141.10

	% TOTAL FLOW	RESISTANCE	INERTANCE (SEC)	CAPACITANCE (SEC)	TIMELAG (SEC)
FUEL:	100.000	4.788E-01	9.888E-06	7.891E-04	4.818E-04
	0.000	4.788E-01	0.000E+00	7.891E-04	0.000E+00
	0.000	4.788E-01	0.000E+00	7.891E-04	0.000E+00
	0.000	4.788E-01	0.000E+00	7.891E-04	0.000E+00
OX:	100.000	4.788E-01	7.303E-06	5.918E-04	8.913E-04
	0.000	4.788E-01	0.000E+00	5.918E-04	0.000E+00
	0.000	4.788E-01	0.000E+00	5.918E-04	0.000E+00
	0.000	4.788E-01	0.000E+00	5.918E-04	0.000E+00

PC, PSIA=1551.40

% TOTAL FLOW	RESISTANCE	INERTANCE	CAPACITANCE	TIMELAG
--------------	------------	-----------	-------------	---------

		(SEC)	(SEC)	(SEC)
FUEL:	100.000	3.454E-01	3.000E-05	6.540E-04
	0.000	3.454E-01	0.000E+00	0.000E+00
	0.000	3.454E-01	0.000E+00	0.000E+00
	0.000	3.454E-01	0.000E+00	0.000E+00
OX:	100.000	3.454E-01	7.312E-05	1.100E-03
	0.000	3.454E-01	0.000E+00	0.000E+00
	0.000	3.454E-01	0.000E+00	0.000E+00
	0.000	3.454E-01	0.000E+00	0.000E+00

PC, PSIA= 787.34

	% TOTAL FLOW	RESISTANCE	INERTANCE (SEC)	CAPACITANCE (SEC)	TIME LAG (SEC)
FUEL:	100.000	1.721E-01	3.000E-05	6.605E-04	1.204E-03
	0.000	1.721E-01	0.000E+00	6.605E-04	0.000E+00
	0.000	1.721E-01	0.000E+00	6.605E-04	0.000E+00
	0.000	1.721E-01	0.000E+00	6.605E-04	0.000E+00
OX:	100.000	1.720E-01	7.398E-05	6.561E-04	2.237E-03
	0.000	1.720E-01	0.000E+00	6.561E-04	0.000E+00
	0.000	1.720E-01	0.000E+00	6.561E-04	0.000E+00
	0.000	1.720E-01	0.000E+00	6.561E-04	0.000E+00

# HIFI CHAMBER MODEL INPUTS

COMBUSTION PLANE, FT= 2.612E-01      SHORT NOZZLE ASSUMED=F

## ACOUSTIC CAVITY INPUTS

CAVITY TYPE=NONE  
CAVITY TYPE 1:    NUMBER OF CAVITIES= 0      NUMBER OF PROPERTY SECTIONS= 1  
CAVITY TYPE 2:    NUMBER OF CAVITIES= 0      NUMBER OF PROPERTY SECTIONS= 1  
PARTITION THICKNESS, FT= 0.0000

PC (PSIA)	MAX. AMPLITUDE	FREQUENCY (HZ)	PHASE MARGIN (DEG)
2.141E+03	3.710E-01	5.199E+02	180.00
2.091E+03	3.619E-01	5.131E+02	180.00
2.041E+03	3.598E-01	5.073E+02	180.00
1.991E+03	3.977E-01	5.014E+02	180.00
1.941E+03	4.054E-01	4.955E+02	180.00
1.891E+03	4.130E-01	4.894E+02	180.00
1.841E+03	4.204E-01	4.833E+02	180.00
1.791E+03	4.276E-01	4.770E+02	180.00
1.741E+03	4.345E-01	4.707E+02	180.00
1.691E+03	4.408E-01	4.641E+02	180.00
1.641E+03	4.469E-01	4.574E+02	180.00
1.591E+03	4.524E-01	4.506E+02	180.00
1.541E+03	4.573E-01	4.438E+02	180.00
1.491E+03	4.616E-01	4.363E+02	180.00
1.441E+03	4.650E-01	4.286E+02	180.00
1.391E+03	4.676E-01	4.211E+02	180.00
1.341E+03	4.693E-01	4.131E+02	180.00
1.291E+03	4.699E-01	4.048E+02	180.00
1.241E+03	4.695E-01	3.963E+02	180.00
1.191E+03	4.679E-01	3.873E+02	180.00
1.141E+03	4.651E-01	3.780E+02	180.00
1.091E+03	4.611E-01	3.683E+02	180.00
1.041E+03	4.558E-01	3.582E+02	180.00
9.911E+02	4.491E-01	3.476E+02	180.00
9.411E+02	4.412E-01	3.366E+02	180.00
8.911E+02	4.754E-01	6.716E+02	180.00
8.411E+02	5.129E-01	6.463E+02	180.00
7.911E+02	5.419E-01	6.243E+02	180.00
7.411E+02	5.690E-01	5.991E+02	180.00
6.911E+02	5.619E-01	5.725E+02	180.00
6.411E+02	6.223E-01	6.394E+02	180.00
5.911E+02	6.399E-01	7.913E+02	-49.18
5.411E+02	1.009E+00	7.444E+02	26.49
5.678E+02	1.009E+00	7.589E+02	2.58

MARGINAL CHUG POINT FOUND:  
PC= 557.77 PSIA      FREQUENCY= 759.90 HZ

# CHUG STABILITY ITERATION SUMMARY

THE CURRENT CONFIGURATION IS CHUG STABLE

DESIRED MARGINAL PC = 800.00 PSIA  
CURRENT MARGINAL PC = 557.77 PSIA  
CURRENT CHUG MARGIN = 242.28 PSI  
CHUG FREQUENCY = 759.90 HZ

# LOW FREQUENCY COMBUSTION STABILITY CALCULATIONS

ROCKET COMBUSTOR INTERACTIVE DESIGN METHODOLOGY  
Version 23-FEB-91

ROCCID

DIRECT INPUT ECHO FROM SUBROUTINE SINPUT

ROCCID POINT DESIGN TEST CASE 1

LOX/RP-1 LIKE DOUBLET PAIR WITH FIXED PC

APPROXIMATES -0100 SUBSCALE DOUBLET

\$MODELS

MCHAM= 1

MBURN= 2

MINJ= 1

\$END

\$OPCOND

P:SPA=

5.9600E+01, 1.0940E+02, 1.3940E+02, 1.5980E+02, 1.7780E+02,  
2.0970E+02, 2.3550E+02, 2.5470E+02, 2.6700E+02, 2.7040E+02,  
2.7180E+02, 2.7190E+02, 2.7130E+02, 2.8980E+02, 2.8820E+02,  
2.8800E+02, 2.8500E+02, 2.8210E+02, 2.5920E+02, 2.5830E+02,  
2.5080E+02, 2.3840E+02, 2.2760E+02, 2.1100E+02, 2.0140E+02,  
1.5580E+02, 1.0970E+02, 8.0000E+01, 4.0740E+03, 4.5800E+03,  
1.7780E+03, 2.4880E+03, 3.4870E+03, 5.0880E+03, 5.9610E+03,  
5.0750E+03, 5.4840E+03, 5.7340E+03, 5.9090E+03, 5.8820E+03,  
5.9880E+03, 5.9540E+03, 5.9340E+03, 5.7000E+03, 5.6430E+03,  
5.8580E+03, 5.8260E+03, 5.7880E+03, 5.7000E+03, 5.6430E+03,  
5.5320E+03, 5.2750E+03, 5.0470E+03, 4.6780E+03, 4.4200E+03,  
3.4570E+03, 2.4940E+03, 1.4650E+03,

'RP-1'

'LOX'

'LOX'

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'LOX'

FCANT=	1.6000E+01	
FFACET=	2.6726E-01	
XDJ=	1.0172E-01	
XCD=	9.4000E-01	
XIN=	1.7618E-01	
XIA=	3.0000E+01	
XCANT=	1.6000E+01	
XFACET=	4.4047E-01	
EMUNI=	6.5000E-01	
\$END		
\$BAFFLE		
\$END		
\$BARRIER		
\$END		
\$FFC		
\$END		
\$BURN		
GAMMA=	1.1462E+00	
AO=	4.0696E+03	
GAM=	2.3216E+01	
GPR=	5.9752E-01	
GK=	5.4180E-05	
GMU=	6.6324E-06	
RML=	7.6242E+01	
VJL=	3.0746E+02	
TJL=	5.3100E+02	
RHOL=	4.9895E+01	
CPL=	4.7821E-01	
PCRTIL=	3.1500E+02	
TCRITL=	1.2180E+03	
TBOILL=	8.8200E+02	
XMWL=	1.7200E+02	
HVAPL=	1.2500E+02	
EN=	6.6783E-01	
TAUSEN=	1.4114E-04	
ISEN=	1	
\$END		
8INJ		
FMAND=	6.7076E+00	
XMAND=	6.7076E+00	
FMANL=	3.5538E+00	
XMANL=	3.5538E+00	
PCA=	2.1411E+03,	1.5514E+03, 7.6784E+02
FRA=	4.7847E-01,	3.4543E-01, 1.7208E-01
FCAPA=	7.6914E-04,	6.1781E-04, 6.6053E-04
NFE=	4	
FTLA=	6.0142E-04,	6.8196E-04, 1.3371E-03,
FINA=	3.8978E-05,	3.8925E-05, 3.9084E-05,
FFA=	1.0000E+00,	1.0000E+00, 1.0000E+00,
NXE=	4	
XRA=	4.7847E-01,	3.4543E-01, 1.7208E-01
XCAPA=	6.9169E-04,	6.1747E-04, 6.6018E-04
XTLA=	9.3298E-04,	1.2436E-03, 2.3388E-03,
XINA=	7.3028E-05,	7.3125E-05, 7.3862E-05,
XFA=	1.0000E+00,	1.0000E+00, 1.0000E+00,
XUOR=	4.7860E-02,	7.3412E-02,
AUOR=	3.4570E-03,	6.1269E-03,
RUOR=	3.3172E-02,	6.0861E-02,
XOR=	4.7860E-02,	7.3412E-02,
AOR=	3.4570E-03,	6.1269E-03,
ROR=	3.3172E-02,	6.0861E-02,
XDOR=	9.5761E-02,	1.4682E-01,
ADOR=	3.4570E-03,	6.1269E-03,
ROR=	3.3172E-02,	6.0861E-02,

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WOOT= 4.6218E+01, 1.3311E+02, 3.3533E+01, 9.6576E+01, 1.6644E+01,
4.7939E+01

$END
$CHAMBER
XB= 2.6125E-01
ZB= 1.1675E-01
ZE= 4.9875E-01
XUB= 3
YUB= 3
T= 3.3000E-02
CO1= 5*2.5000E+03
GAMC1= 5*1.2000E+00
CC2= 5*2.5000E+03
GAMC2= 5*1.2000E+00
$END
$FDORC
NZON= 5
FTER= 20*1.0000E+00
$END
$MIX
EM(1)= 9.003E-01, 1.000E+00
$END
$DEBUG
F
$END
$HIFIC
SHORT= F
POC= 2.0000E-01, 2.0000E-01
$END
$DIST3DC
SHORT= F
POC= 2.0000E-01, 2.0000E-01
PAMP= 2.0000E-01
MC= 11
LC= 8
MB= 11
LB= 8
IDMAX= 10
$END
$CRPC
NDPC= 16
NDTFQ= 16
NDTLF= 1000
NPRINT= 50
NBLMS= 3500
PAMP= 2.0000E-01
NRAD= 3
NGIRC= 5
$END
$LEINJC
IDMEM= 2
PAMPCH= 2.0000E-01
NTINJ= 16
USGF= 2.0000E-02
OGF= 2.0000E-02
D8GF= 2.0000E-02
FCDO= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00
XCDO= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00
$END
$COMBUSTC
FALVM= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00
XALVM= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00
FALTM= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00
XALTM= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00
FMM= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00

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XNMA= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00
ENMULT= 1.0000E+00, 1.0000E+00, 1.0000E+00
COMBX= 5.0000E-01
ACMULT= 1.0000E+00
CCMULT= 1.0000E+00
ENMULT= 1.0000E+00
TAUMULT= 1.0000E+00
$END
$FDMCC
SHORT= F
EPSIL= 2.0000E+01
ERROR= 1.0000E-01
LTS= 5
MTS= 5
NTS= 10
ITMAX= 100
RELX= 5.5000E-01
NEET= 3
MXFST= 3
MYFST= 3
NAFST= 3
MORE= F
$END

```

END OF INPUT ECHO

# STABILITY MODEL INPUTS

## RUN DESCRIPTOR

ROCCID POINT DESIGN TEST CASE 1  
LOX/RP-1 LIKE DOUBLET PAIR WITH FIXED PC  
APPROXIMATES -0100 SUBSCALE DOUBLET

## SELECTED MODELS

BURNING MODEL=N-TAU INJECTION MODEL=INJ CHAMBER MODEL=HIFI

AXISYMETRIC=T DEBUG OUTPUT=F

## CHAMBER GEOMETRY AND OPERATING CONDITIONS

CHAMBER RADIUS, FT= 0.2798 THROAT RADIUS, FT= 0.1852  
CYLINDRICAL SECTION, FT= 0.9648 CONVERGENCE HALF-ANGLE, DEG=30.0000  
NOZZLE ENTRANCE RADIUS OF CURVATURE, FT= 0.1110 THROAT ENTRANCE RADIUS OF CURVATURE, FT= 0.1110  
CHAMBER PRESSURE, PSIA=2141.10 MIXTURE RATIO= 2.8800  
SOUND SPEED, FT/SEC=4086.80 GAMMA=1.1462

## N-TAU BURNING MODEL INPUTS

PRESSURE INTERACTION INDEX, EN= 0.8879 SENSITIVE TIMELAG, TAU, SEC= 1.411E-04 SENSITIVE CIRCUIT=FUEL

## LUMPED INJECTION MODEL INPUTS

PC, PSIA=2141.10

	% TOTAL FLOW	RESISTANCE	INERTANCE (SEC)	CAPACITANCE (SEC)	TIMELAG (SEC)
FUEL:	100.000	4.788E-01	3.888E-08	7.881E-04	8.014E-04
	0.000	4.788E-01	0.000E+00	7.881E-04	0.000E+00
	0.000	4.788E-01	0.000E+00	7.881E-04	0.000E+00
	0.000	4.788E-01	0.000E+00	7.881E-04	0.000E+00
OX:	100.000	4.788E-01	7.903E-08	8.916E-04	9.924E-04
	0.000	4.788E-01	0.000E+00	8.916E-04	0.000E+00
	0.000	4.788E-01	0.000E+00	8.916E-04	0.000E+00
	0.000	4.788E-01	0.000E+00	8.916E-04	0.000E+00

PC, PSIA=1551.40

% TOTAL FLOW	RESISTANCE	INERTANCE	CAPACITANCE	TIMELAG
--------------	------------	-----------	-------------	---------

		(SEC)	(SEC)	(SEC)
FUEL:	100.000	3.464E-01	3.893E-05	9.175E-04
	0.000	3.464E-01	0.000E+00	9.175E-04
	0.000	3.464E-01	0.000E+00	9.175E-04
	0.000	3.464E-01	0.000E+00	9.175E-04
OX:	100.000	3.464E-01	7.312E-05	9.175E-04
	0.000	3.464E-01	0.000E+00	9.175E-04
	0.000	3.464E-01	0.000E+00	9.175E-04
	0.000	3.464E-01	0.000E+00	9.175E-04

PC, PSIA= 767.34

	% TOTAL FLOW	RESISTANCE	INERTANCE (SEC)	CAPACITANCE (SEC)	TIMELAG (SEC)
FUEL:	100.000	1.721E-01	3.906E-05	9.605E-04	1.337E-03
	0.000	1.721E-01	0.000E+00	9.605E-04	0.000E+00
	0.000	1.721E-01	0.000E+00	9.605E-04	0.000E+00
	0.000	1.721E-01	0.000E+00	9.605E-04	0.000E+00
OX:	100.000	1.720E-01	7.336E-05	9.561E-04	2.340E-03
	0.000	1.720E-01	0.000E+00	9.561E-04	0.000E+00
	0.000	1.720E-01	0.000E+00	9.561E-04	0.000E+00
	0.000	1.720E-01	0.000E+00	9.561E-04	0.000E+00

# HIFI CHAMBER MODEL INPUTS

COMBUSTION PLANE, FT= 2.612E-01      SHORT NOZZLE ASSUMED=F

## ACOUSTIC CAVITY INPUTS

CAVITY TYPE=NONE  
 CAVITY TYPE 1:      NUMBER OF CAVITIES= 0      NUMBER OF PROPERTY SECTIONS= 1  
 CAVITY TYPE 2:      NUMBER OF CAVITIES= 0      NUMBER OF PROPERTY SECTIONS= 1  
 PARTITION THICKNESS, FT= 0.0000

PC (PSIA)	MAX. AMPLITUDE	FREQUENCY (HZ)	PHASE MARGIN (DEG)
2.141E+03	3.725E-01	5.066E+02	180.00
2.091E+03	3.835E-01	4.992E+02	180.00
2.041E+03	3.909E-01	4.935E+02	180.00
1.991E+03	3.981E-01	4.878E+02	180.00
1.941E+03	4.053E-01	4.819E+02	180.00
1.891E+03	4.125E-01	4.760E+02	180.00
1.841E+03	4.189E-01	4.700E+02	180.00
1.791E+03	4.253E-01	4.636E+02	180.00
1.741E+03	4.315E-01	4.575E+02	180.00
1.691E+03	4.370E-01	4.511E+02	180.00
1.641E+03	4.421E-01	4.445E+02	180.00
1.591E+03	4.468E-01	4.377E+02	180.00
1.541E+03	4.509E-01	4.307E+02	180.00
1.491E+03	4.541E-01	4.236E+02	180.00
1.441E+03	4.566E-01	4.162E+02	180.00
1.391E+03	4.584E-01	4.086E+02	180.00
1.341E+03	4.592E-01	4.008E+02	180.00
1.291E+03	4.590E-01	3.925E+02	180.00
1.241E+03	4.577E-01	3.840E+02	180.00
1.191E+03	4.553E-01	3.751E+02	180.00
1.141E+03	4.519E-01	3.659E+02	180.00
1.091E+03	4.472E-01	3.564E+02	180.00
1.041E+03	4.413E-01	3.464E+02	180.00
9.911E+02	4.343E-01	3.360E+02	180.00
9.411E+02	4.205E-01	3.272E+02	180.00
8.911E+02	4.677E-01	6.504E+02	180.00
8.411E+02	5.173E-01	6.275E+02	180.00
7.911E+02	5.377E-01	6.037E+02	180.00
7.411E+02	5.467E-01	5.786E+02	180.00
6.911E+02	5.434E-01	5.521E+02	180.00
6.411E+02	7.171E-01	8.109E+02	-119.37
5.911E+02	9.370E-01	7.654E+02	-19.25
5.411E+02	1.179E+00	7.210E+02	61.09
5.378E+02	1.098E+00	7.367E+02	28.36
5.744E+02	1.017E+00	7.506E+02	5.04
5.800E+02	9.901E-01	7.555E+02	-2.92

MARGINAL CHUG POINT FOUND:  
PC= 579.99 PSIA      FREQUENCY= 755.47 HZ

# CHUG STABILITY ITERATION SUMMARY

THE CURRENT CONFIGURATION IS CHUG STABLE

DESIRED MARGINAL PC = 900.00 PSIA  
CURRENT MARGINAL PC = 579.99 PSIA  
CURRENT CHUG MARGIN = 220.01 PSI  
CHUG FREQUENCY = 755.47 HZ

# HIGH FREQUENCY COMBUSTION STABILITY CALCULATIONS

ROCCID  
 Rocket Combuster Interactive Design Methodology  
 Version 23-FEB-91

DIRECT INPUT ECHO FROM SUBROUTINE SINPUT

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ROCCID POINT DESIGN TEST CASE 1
LOX/RP-1 LIKE DOUBLET PAIR WITH FIXED PC
APPROXIMATES -0100 SUBSCALE DOUBLET
$MODELS
MCHAM=      1
MBURN=      2
MINJ=       1
SEND
$OPCOND
PISPA=      6.9600E+01, 1.0940E+02, 1.3940E+02, 1.5860E+02, 1.7780E+02,
2.0970E+02, 2.3550E+02, 2.5470E+02, 2.6700E+02, 2.7040E+02,
2.7180E+02, 2.7180E+02, 2.7130E+02, 2.6980E+02, 2.6820E+02,
2.6680E+02, 2.6500E+02, 2.6210E+02, 2.5920E+02, 2.5830E+02,
2.5680E+02, 2.540E+02, 2.3760E+02, 2.1100E+02, 2.0140E+02,
1.5560E+02, 1.0970E+02, 8.0000E+01,
PCSA=      1.7780E+03, 2.4560E+03, 3.4870E+03, 4.0740E+03, 4.5800E+03,
5.0700E+03, 5.4540E+03, 5.7840E+03, 5.9080E+03, 5.9610E+03,
5.9680E+03, 5.9440E+03, 5.9340E+03, 5.9080E+03, 5.8320E+03,
5.6550E+03, 5.8280E+03, 5.7590E+03, 5.7000E+03, 5.6430E+03,
5.5320E+03, 5.2760E+03, 5.0470E+03, 4.6760E+03, 4.4200E+03,
3.4570E+03, 2.4940E+03, 1.4650E+03,
FUEL=      'RP-1'
OX=      'LOX'
PC=      2.1411E+03
XMR=      2.6600E+00
FTMAN=     7.1000E+01
XTMAN=     -2.7900E+02
EMMAN=     1.0000E+00
NPERFP=    28
PWRA=      1.0000E-01, 3.0000E-01, 6.0000E-01, 9.0000E-01, 1.0000E+00,
1.2500E+00, 1.5000E+00, 1.7500E+00, 2.0000E+00, 2.2000E+00,
2.4000E+00, 2.5000E+00, 2.6000E+00, 2.7000E+00, 2.8000E+00,
2.9000E+00, 3.0000E+00, 3.2000E+00, 3.4000E+00, 3.6000E+00,
4.0000E+00, 5.0000E+00, 6.0000E+00, 8.0000E+00, 1.0000E+01,
1.5000E+01, 2.0000E+01, 5.0000E+01,
$GEOM
RCHAMB=     2.7948E-01
RTHRT=      1.6523E-01
RNE=        1.100E-01
RTE=        1.1100E-01
ALPHA=      3.0000E+01
CHAMBL=     1.1675E+00
XC=         9.6477E-01
$END
$SCORE
TYPE=      'LOL'
NEL=      88
FDJ=      6.6948E-02
FCD=      9.1000E-01
FIN=      1.1491E-01
FIA=      3.0000E+01
    
```

FCANT=	1.6000E+01	
FFACET=	2.8728E-01	
XDJ=	1.0172E-01	
XOD=	8.4000E-01	
XIH=	1.7819E-01	
XIA=	3.0000E+01	
XCANT=	1.6000E+01	
XFACET=	4.4047E-01	
EMUNI=	6.6000E-01	
SEND		
8BAFFLE		
SEND		
8BARRIER		
SEND		
8FFC		
SEND		
8BURN		
GAMMA=	1.1462E+00	
AO=	4.0668E+03	
GAM=	2.3218E+01	
GPR=	5.9752E-01	
GK=	5.4180E-05	
GAU=	6.8324E-05	
RML=	7.8242E+01	
VJL=	3.0748E+02	
TJL=	6.3100E+02	
RHOL=	4.9885E+01	
CPL=	4.7321E-01	
PCRITL=	3.1500E+02	
TCRITL=	1.2180E+03	
TBOILL=	8.8200E+02	
XMWL=	1.7200E+02	
HVAPL=	1.2500E+02	
EN=	8.8783E-01	
TAUSEN=	1.4114E-04	
ISEN=	1	
SEND		
\$INJ		
FMAND=	6.7076E+00	
XMAND=	6.7076E+00	
FMANL=	3.3538E+00	
XMANL=	3.3538E+00	
PCA=	2.1411E+03,	1.5514E+03, 7.6734E+02
FRA=	4.7647E-01,	3.4543E-01, 1.7206E-01
FCAPA=	7.8914E-04,	8.1781E-04, 8.6053E-04
NFE=	4	
FTLA=	5.0142E-04,	8.8188E-04, 1.3371E-03,
FINA=	3.8876E-06,	3.8928E-06, 3.9084E-06,
FFA=	1.0000E+00,	1.0000E+00, 1.0000E+00,
NXE=	4	
XRA=	4.7647E-01,	3.4543E-01, 1.7206E-01
XCAPA=	5.8188E-04,	8.1747E-04, 8.6615E-04
XTLA=	9.3236E-04,	1.2439E-03, 2.3395E-03,
XINA=	7.3028E-06,	7.3126E-06, 7.3382E-06,
XFA=	1.0000E+00,	1.0000E+00, 1.0000E+00,
XUOR=	4.7880E-02,	7.3412E-02,
AUOR=	3.4570E-03,	8.1289E-03,
RUOR=	3.3172E-02,	5.0881E-02,
XOR=	4.7880E-02,	7.3412E-02,
AOR=	3.4570E-03,	8.1289E-03,
ROR=	3.3172E-02,	5.0881E-02,
XDOR=	9.5781E-02,	1.4882E-01,
ADOR=	3.4570E-03,	8.1289E-03,
RDOR=	3.3172E-02,	8.0861E-02,

WOOT= 4.9210E+01, 1.3311E+02, 3.3333E+01, 9.9970E+01, 1.0044E+01,  
 4.7936E+01  
 \$END  
 \$CHAMBER  
 XB= 2.6126E-01  
 ZB= 1.1878E-01  
 ZE= 4.9878E-01  
 MJB= 3  
 T= 3.3000E-02  
 CC1= 5\*2.5000E+03  
 GAMC1= 5\*1.2000E+00  
 CC2= 5\*2.5000E+03  
 GAMC2= 5\*1.2000E+00  
 \$END  
 \$FDORG  
 NZON= 5  
 FTER= 20\*1.0000E+00  
 \$END  
 \$MIX  
 EM(1)= 9.003E-01, 1.000E+00  
 \$END  
 \$DEBUGC  
 DEBUG= F  
 \$END  
 \$HIFIC  
 SHORT= F  
 POC= 2.0000E-01, 2.0000E-01  
 \$END  
 \$DIST3DC  
 SHORT= F  
 POC= 2.0000E-01, 2.0000E-01  
 PAMP= 2.0000E-01  
 MC= 11  
 LC= 8  
 MB= 11  
 LB= 8  
 IDMAX= 10  
 \$END  
 \$CRPC  
 NDPC= 16  
 NDTFQ= 16  
 NDTLF= 1000  
 NPRINT= 50  
 NSUMS= 3500  
 PAMPC= 2.0000E-01  
 NRAD= 3  
 NCIRC= 6  
 \$END  
 \$LEINJC  
 IDOMEM= 2  
 PAMPCH= 2.0000E-01  
 NTINJ= 16  
 USGF= 2.0000E-02  
 OGF= 2.0000E-02  
 DSGF= 2.0000E-02  
 FGOO= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00  
 XCDO= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00  
 \$END  
 \$COMBUSTC  
 FALVM= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00  
 XALVM= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00  
 FALTM= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00  
 XALTM= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00  
 FRMM= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00

```

XPMW= 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00
ENMULT= 1.0000E+00, 1.0000E+00
COMBXS= 8.0000E-01
ACMULT= 1.0000E+00
CCMULT= 1.0000E+00
ENMULT= 1.0000E+00
TAMULT= 1.0000E+00
$END
$FDORC
SHORT= F
EP81L= 2.0000E+01
ERROR= 1.0000E-01
LTS= 5
MTS= 5
NTS= 10
ITMAX= 100
RELX= 8.5000E-01
NEET= 3
NXFST= 3
NYFST= 3
NAFST= 3
MORE= F
$END
END OF INPUT ECHO

```

# STABILITY MODEL INPUTS

## RUN DESCRIPTOR

ROCCID POINT DESIGN TEST CASE 1  
LOX/RP-1 LIKE DOUBLET PAIR WITH FIXED PC  
APPROXIMATES -0100 SUBSCALE DOUBLET

## SELECTED MODELS

BURNING MODEL=N-TAU      INJECTION MODEL=INJ      CHAMBER MODEL=HIFI  
AXISYMETRIC=T      DEBUG OUTPUT=F

## CHAMBER GEOMETRY AND OPERATING CONDITIONS

CHAMBER RADIUS, FT= 0.2795      THROAT RADIUS, FT= 0.1962  
CYLINDRICAL SECTION, FT= 0.9646      CONVERGENCE HALF-ANGLE, DEG=20.0000  
NOZZLE ENTRANCE RADIUS OF CURVATURE, FT= 0.1110      THROAT ENTRANCE RADIUS OF CURVATURE, FT= 0.1110  
CHAMBER PRESSURE, PSIA=2141.10      MIXTURE RATIO= 2.8800  
SOUND SPEED, FT/SEC=4065.60      GAMMA=1.1462

## N-TAU BURNING MODEL INPUTS

PRESSURE INTERACTION INDEX, EN= 0.8679      SENSITIVE TIMELAG, TAU, SEC= 1.411E-04      SENSITIVE CIRCUIT=FUEL

## LUMPED INJECTION MODEL INPUTS

PC, PSIA=2141.10

	% TOTAL FLOW	RESISTANCE	INERTANCE (SEC)	CAPACITANCE (SEC)	TIMELAG (SEC)
FUEL:	100.000	4.766E-01	3.886E-06	7.891E-04	6.014E-04
	0.000	4.766E-01	0.000E+00	7.891E-04	0.000E+00
	0.000	4.766E-01	0.000E+00	7.891E-04	0.000E+00
	0.000	4.766E-01	0.000E+00	7.891E-04	0.000E+00
OX:	100.000	4.766E-01	7.303E-06	6.916E-04	9.324E-04
	0.000	4.766E-01	0.000E+00	6.916E-04	0.000E+00
	0.000	4.766E-01	0.000E+00	6.916E-04	0.000E+00
	0.000	4.766E-01	0.000E+00	6.916E-04	0.000E+00

PC, PSIA=1551.40

% TOTAL FLOW	RESISTANCE	INERTANCE	CAPACITANCE	TIMELAG
--------------	------------	-----------	-------------	---------

FUEL:	100.000	3.454E-01	3.693E-08	(SEC)	(SEC)	(SEC)
	0.000	3.454E-01	0.000E+00		6.178E-04	6.820E-04
	0.000	3.454E-01	0.000E+00		6.178E-04	0.000E+00
	0.000	3.454E-01	0.000E+00		6.178E-04	0.000E+00
OX:	100.000	3.454E-01	7.312E-06		6.178E-04	1.244E-03
	0.000	3.454E-01	0.000E+00		6.178E-04	0.000E+00
	0.000	3.454E-01	0.000E+00		6.178E-04	0.000E+00
	0.000	3.454E-01	0.000E+00		6.178E-04	0.000E+00

PC, PSIA= 767.34

	% TOTAL FLOW	RESISTANCE	INERTANCE (SEC)	CAPACITANCE (SEC)	TIMELAG (SEC)
FUEL:	100.000	1.721E-01	3.906E-06	6.605E-04	1.337E-03
	0.000	1.721E-01	0.000E+00	6.605E-04	0.000E+00
	0.000	1.721E-01	0.000E+00	6.605E-04	0.000E+00
	0.000	1.721E-01	0.000E+00	6.605E-04	0.000E+00
OX:	100.000	1.720E-01	7.336E-06	6.561E-04	2.340E-03
	0.000	1.720E-01	0.000E+00	6.561E-04	0.000E+00
	0.000	1.720E-01	0.000E+00	6.561E-04	0.000E+00
	0.000	1.720E-01	0.000E+00	6.561E-04	0.000E+00

# HIFI CHAMBER MODEL INPUTS

COMBUSTION PLANE, FT= 2.612E-01    SHORT NOZZLE ASSUMED=F

## ACOUSTIC CAVITY INPUTS

CAVITY TYPE=NONE  
 CAVITY TYPE 1:    NUMBER OF CAVITIES= 0    NUMBER OF PROPERTY SECTIONS= 1  
 CAVITY TYPE 2:    NUMBER OF CAVITIES= 0    NUMBER OF PROPERTY SECTIONS= 1  
 PARTITION THICKNESS, FT= 0.0000

## HIGH FREQUENCY INSTABILITY IS BURNING-COUPLED

TANGENTIAL MODE	RADIAL MODE	FREQUENCY (HZ)	GAIN
1	0	4196.57	1.0233

# HIGH FREQUENCY COMBUSTION STABILITY CALCULATIONS

ROCKET COMBUSTOR INTERACTIVE DESIGN METHODOLOGY  
Version 23-FEB-91

ROCCID

DIRECT INPUT ECHO FROM SUBROUTINE SINPUT

```

ROCCID POINT DESIGN TEST CASE 1
LOX/RP-1 LIKE DOUBLET PAIR WITH FIXED PC
APPROXIMATES .0100 SUBSCALE DOUBLET
$MODELS
MCHAM=      1
MBURN=      2
MINJ=       1
$END
$OPCOND
PISPA= 8.9800E+01, 1.0940E+02, 1.3940E+02, 1.5880E+02, 1.7780E+02,
2.0870E+02, 2.3650E+02, 2.5470E+02, 2.6700E+02, 2.7040E+02,
2.7180E+02, 2.7190E+02, 2.7190E+02, 2.6980E+02, 2.6920E+02,
2.6860E+02, 2.6500E+02, 2.6210E+02, 2.5920E+02, 2.5630E+02,
2.5080E+02, 2.3840E+02, 2.2780E+02, 2.1100E+02, 2.0140E+02,
1.5660E+02, 1.0970E+02, 8.0000E+01,
1.7760E+03, 2.4680E+03, 3.4870E+03, 4.0740E+03, 4.5800E+03,
5.0700E+03, 5.4540E+03, 5.7340E+03, 5.9080E+03, 5.9610E+03,
5.9890E+03, 5.9540E+03, 5.9340E+03, 5.9080E+03, 5.8820E+03,
5.8550E+03, 5.8260E+03, 5.7590E+03, 5.7000E+03, 5.6430E+03,
5.5820E+03, 5.2750E+03, 5.0470E+03, 4.8760E+03, 4.4200E+03,
3.4570E+03, 2.4940E+03, 1.4650E+03,
FUEL= 'RP-1',
OX= 'LOX',
PC= 2.1411E+03
XMR= 2.8800E+00
FTMAN= 7.1000E+01
XTMAN= -2.7800E+02
EMMAN= 1.0000E+00
NPERFP= 28
PMRA= 1.0000E-01, 3.0000E-01, 6.0000E-01, 8.0000E-01, 1.0000E+00,
1.2600E+00, 1.5000E+00, 1.7500E+00, 2.0000E+00, 2.2000E+00,
2.4000E+00, 2.5000E+00, 2.6000E+00, 2.7000E+00, 2.8000E+00,
2.9000E+00, 3.0000E+00, 3.2000E+00, 3.4000E+00, 3.6000E+00,
4.0000E+00, 5.0000E+00, 6.0000E+00, 8.0000E+00, 1.0000E+01,
1.5000E+01, 2.0000E+01, 5.0000E+01,
$END
$GECM
RCHAMB= 2.7949E-01
RTHRT= 1.8523E-01
RNE= 1.1100E-01
RTE= 1.1100E-01
ALPHA= 3.0000E+01
CHAMBL= 1.1875E+00
XC= 9.6477E-01
$END
$SCORE
TYPE= 'LOL',
NEL= 69
FDJ= 6.8345E-02
FCO= 9.1000E-01
FIH= 1.1491E-01
FIA= 3.0000E+01

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FCANT=	1.6000E+01	
FFACET=	2.8728E-01	
XDJ=	1.0172E-01	
XCD=	9.4000E-01	
XIH=	1.7619E-01	
XIA=	3.0000E+01	
XCANT=	1.6000E+01	
XFACET=	4.4047E-01	
EMUNI=	6.5000E-01	
\$END		
\$BAFFLE		
\$END		
\$BARRIER		
\$FFC		
\$END		
\$BURN		
GAMMA=	1.1452E+00	
AO=	4.0656E+03	
GMW=	2.3218E+01	
GPR=	5.9752E-01	
GK=	5.4180E-06	
GMU=	6.6324E-05	
RML=	7.6242E+01	
VJL=	3.0748E+02	
TJL=	6.3100E+02	
RHOL=	4.9896E+01	
CPL=	4.7321E-01	
PCRITL=	3.1500E+02	
TCRITL=	1.2180E+03	
TBOILL=	8.6200E+02	
XMWL=	1.7200E+02	
HVAPL=	1.2500E+02	
EN=	6.6793E-01	
TAUSEN=	1.4114E-04	
ISEN=	1	
\$END		
\$INJ		
FMAND=	6.7076E+00	
XMAND=	6.7076E+00	
FMANL=	3.3636E+00	
XMANL=	3.3636E+00	
PCA=	2.1411E+03,	1.6614E+03,
FRA=	4.7647E-01,	3.4643E-01,
FCAPA=	7.6914E-04,	9.1791E-04,
NFE=	4	
FTLA=	5.0142E-04,	9.8196E-04,
FINA=	3.6879E-05,	3.8928E-06,
FFA=	1.0000E+00,	1.0000E+00,
NXE=	4	
XRA=	4.7647E-01,	3.4643E-01,
XCAPA=	6.9163E-04,	9.1747E-04,
XTLA=	9.3239E-04,	1.2439E-03,
XINA=	7.3023E-05,	7.3126E-06,
XFA=	1.0000E+00,	1.0000E+00,
XUOR=	4.7680E-02,	7.3412E-02,
AUOR=	3.4570E-03,	6.1299E-03,
RUOR=	3.3172E-02,	6.0651E-02,
XOR=	4.7680E-02,	7.3412E-02,
AOR=	3.4570E-03,	6.1299E-03,
NOR=	3.3172E-02,	6.0651E-02,
XDOR=	9.5761E-02,	1.4982E-01,
ADOR=	3.4670E-03,	6.1299E-03,
NDOR=	3.3172E-02,	6.0651E-02,

```

WDOT= 4.9218E+01, 1.3311E+02, 3.3633E+01, 9.8576E+01, 1.8644E+01,
4.7936E+01
$END
$CHAMBER
NCAV= 1,
XB= 2.6125E-01
ZB= 1.1875E-01
ZE= 4.9875E-01
MJB= 3
T= 3.3000E-02
WC= 2.0861E-02,
AC= 3.8809E-02,
ARATIO= 1.0000E+00,
IDCAV= 1,
DC1= 1.4000E-01,
CC1= 5*2.5000E+03
GAMC1= 5*1.2000E+00
CC2= 5*2.5000E+03
GAMC2= 5*1.2000E+00
ICAV= 1
$END
$FDORC
NZON= 5
FTEF= 20*1.0000E+00
NCAV1= 1
ICTVP1= 4,
ZE1= 2.0861E-02,
AE1= 3.6000E+02,
APL1= 1.4000E-01,
WC1= 1.9186E-01,
ZLOW1= -1.2315E-02,
ZUP1= 2.0417E-01,
$END
$MIX
EM(1)= 9.003E-01, 1.000E+00
$END
$DEBUGC
DEBUG= F
$END
$HIFIG
SHORT= F
POC= 2.0000E-01, 2.0000E-01
$END
$DISTDC
SHORT= F
POC= 2.0000E-01, 2.0000E-01
PAMP= 2.0000E-01
MC= 11
LC= 8
MS= 11
LB= 8
IDMAX= 10
$END
$CRPC
NDPC= 16
NDTFQ= 16
NDTLF= 1000
NPRINT= 60
NSUMS= 3500
PAMPC= 2.0000E-01
NRAD= 3
NCIRC= 6
$END
$LEINJC

```

```

1
IDOMEU=      2
PAMPCH= 2.000E-01
NTINJ=      16
USGF= 2.000E-02
OGF= 2.000E-02
DSGF= 2.000E-02
FCDO= 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00
XCDO= 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00
$END
$COMBUSTC
FALVM= 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00
XALVM= 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00
FALTM= 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00
XALTM= 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00
FRMM= 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00
XRMM= 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00
ENMUL= 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00
COMBXB= 5.000E-01
ACMUL= 1.000E+00
CCMUL= 1.000E+00
ENMUL= 1.000E+00
TAMUL= 1.000E+00
$END
$FDORCC
F
SHORT=
EPSIL= 2.000E+01
ERROR= 1.000E-01
LTS= 5
MTS= 5
NTS= 10
ITMAX= 100
RELX= 6.500E-01
NEET= 3
NXFST= 3
NYFST= 3
NAFST= 3
MORE= F
$END

```

END OF INPUT ECHO

# STABILITY MODEL INPUTS

## RUN DESCRIPTOR

ROCCID POINT DESIGN TEST CASE 1  
LOX/JP-1 LIKE DOUBLET PAIR WITH FIXED PC  
APPROXIMATES 0100 SUBSCALE DOUBLET

## SELECTED MODELS

CHAMBER MODEL=HIFI

BURNING MODEL=N-TAU INJECTION MODEL=INJ

AXISYMETRIC=T DEBUG OUTPUT=F

## CHAMBER GEOMETRY AND OPERATING CONDITIONS

CHAMBER RADIUS, FT= 0.2795 THROAT RADIUS, FT= 0.1652  
CYLINDRICAL SECTION, FT= 0.9848 CONVERGENCE HALF-ANGLE, DEG=30.0000  
NOZZLE ENTRANCE RADIUS OF CURVATURE, FT= 0.1110 THROAT ENTRANCE RADIUS OF CURVATURE, FT= 0.1110  
CHAMBER PRESSURE, PSIA=2141.10 MIXTURE RATIO= 2.6800  
SOUND SPEED, FT/SEC=4065.80 GAMMA=1.1482

## N-TAU BURNING MODEL INPUTS

PRESSURE INTERACTION INDEX, EN= 0.8879 SENSITIVE TIMELAG, TAU, SEC= 1.411E-04 SENSITIVE CIRCUIT=FUEL

## LUMPED INJECTION MODEL INPUTS

PC, PSIA=2141.10

	% TOTAL FLOW	RESISTANCE	INERTANCE (SEC)	CAPACITANCE (SEC)	TIMELAG (SEC)
FUEL:	100.000	4.788E-01	3.888E-05	7.881E-04	8.014E-04
	0.000	4.788E-01	0.000E+00	7.881E-04	0.000E+00
	0.000	4.788E-01	0.000E+00	7.881E-04	0.000E+00
	0.000	4.788E-01	0.000E+00	7.881E-04	0.000E+00
OX:	100.000	4.788E-01	7.303E-05	5.916E-04	9.324E-04
	0.000	4.788E-01	0.000E+00	5.916E-04	0.000E+00
	0.000	4.788E-01	0.000E+00	5.916E-04	0.000E+00
	0.000	4.788E-01	0.000E+00	5.916E-04	0.000E+00

PC, PSIA=1851.40

% TOTAL FLOW	RESISTANCE	INERTANCE	CAPACITANCE	TIMELAG
--------------	------------	-----------	-------------	---------

		(SEC)	(SEC)	(SEC)
FUEL:	100.000	3.454E-01	3.698E-06	6.179E-04
	0.000	3.454E-01	0.000E+00	0.000E+00
	0.000	3.454E-01	0.000E+00	0.000E+00
	0.000	3.454E-01	0.000E+00	0.000E+00
OX:	100.000	3.454E-01	7.312E-06	6.179E-04
	0.000	3.454E-01	0.000E+00	0.000E+00
	0.000	3.454E-01	0.000E+00	0.000E+00
	0.000	3.454E-01	0.000E+00	0.000E+00

PC. PSIA= 767.34

	% TOTAL FLOW	RESISTANCE	INERTANCE (SEC)	CAPACITANCE (SEC)	TIMELAG (SEC)
FUEL:	100.000	1.721E-01	3.906E-06	6.605E-04	1.337E-03
	0.000	1.721E-01	0.000E+00	6.605E-04	0.000E+00
	0.000	1.721E-01	0.000E+00	6.605E-04	0.000E+00
	0.000	1.721E-01	0.000E+00	6.605E-04	0.000E+00
OX:	100.000	1.720E-01	7.336E-06	6.561E-04	2.340E-03
	0.000	1.720E-01	0.000E+00	6.561E-04	0.000E+00
	0.000	1.720E-01	0.000E+00	6.561E-04	0.000E+00
	0.000	1.720E-01	0.000E+00	6.561E-04	0.000E+00

# HIFI CHAMBER MODEL INPUTS

COMBUSTION PLANE, FT= 2.612E-01      SHORT NOZZLE ASSUMED=F

## ACOUSTIC CAVITY INPUTS

CAVITY TYPE=1/4 WAVE  
 CAVITY TYPE 1: NUMBER OF CAVITIES= 1      NUMBER OF PROPERTY SECTIONS= 1  
 CAVITY TYPE 2: NUMBER OF CAVITIES= 0      NUMBER OF PROPERTY SECTIONS= 1  
 PARTITION THICKNESS, FT= 0.0000

	TOTAL DEPTH, FT	WIDTH, FT	AREA, SQ. FT	P'/PC	INLET TYPE
TYPE 1:	0.14000	0.02096	3.681E-02	0.200	0
TYPE 2:	0.00000	0.00000	0.000E+00	0.200	0

CAVITY TYPE ORIENTATION= 1

## CAVITY GAS PROPERTIES

	SECT	LENGTH	VSONIC	GAMMA
TYPE 1:	1	0.1400	2500.00	1.2000

## NO HIGH FREQUENCY INSTABILITY MODES OBSERVED

TANGENTIAL MODE	RADIAL MODE	GAIN MAGNITUDE	FREQUENCY (HZ)	YB / YJ
0	0	0.4805	1651.69	1.03
1	0	0.7916	4733.72	2.75

# HIGH FREQUENCY COMBUSTION STABILITY CALCULATIONS

ROCKET COMBUSTOR INTERACTIVE DESIGN METHODOLOGY  
Version 23-FEB-91

ROCCID

DIRECT INPUT ECHO FROM SUBROUTINE SINPUT

```

ROCCID POINT DESIGN TEST CASE 1
LOX/FP-1 LIKE DOUBLET PAIR WITH FIXED PC
APPROXIMATES .0100 SUBSCALE DOUBLET
$MODELS
MCHAM= 1
MBURN= 2
MINJ= 1
$END
$OPCOND
PISPA= 8.9800E+01, 1.0940E+02, 1.3940E+02, 1.5880E+02, 1.7780E+02,
2.0970E+02, 2.3650E+02, 2.5470E+02, 2.6700E+02, 2.7040E+02,
2.7180E+02, 2.7190E+02, 2.7130E+02, 2.8980E+02, 2.8820E+02,
2.8660E+02, 2.8500E+02, 2.8210E+02, 2.8920E+02, 2.5680E+02,
2.8080E+02, 2.3840E+02, 2.2760E+02, 2.1100E+02, 2.0140E+02,
1.5680E+02, 1.0970E+02, 0.0000E+01,
PCSA= 1.7780E+03, 2.4600E+03, 3.4870E+03, 4.0740E+03, 4.5800E+03,
5.0700E+03, 5.4840E+03, 5.7340E+03, 5.9080E+03, 5.9610E+03,
5.9690E+03, 5.9540E+03, 6.9340E+03, 6.9090E+03, 5.6020E+03,
5.8550E+03, 5.8280E+03, 5.7590E+03, 5.7000E+03, 5.6430E+03,
5.5320E+03, 5.2780E+03, 5.0470E+03, 4.6760E+03, 4.4200E+03,
3.4570E+03, 2.4840E+03, 1.4680E+03,
FUEL= 'RP-1'
OX= 'LOX'
PC= 2.1411E+03
XMR= 2.8600E+00
FTMAN= 7.1000E+01
XTMAN= -2.7900E+02
EMMAN= 1.0000E+00
NPERFP= 28
PMRA= 1.0000E-01, 3.0000E-01, 8.0000E-01, 8.0000E-01, 1.0000E+00,
1.2600E+00, 1.5000E+00, 1.7500E+00, 2.0000E+00, 2.2000E+00,
2.4000E+00, 2.5000E+00, 2.6000E+00, 2.8000E+00, 2.7000E+00, 2.8000E+00,
2.9000E+00, 3.0000E+00, 3.2000E+00, 3.4000E+00, 3.6000E+00,
4.0000E+00, 5.0000E+00, 8.0000E+00, 8.0000E+00, 8.0000E+00,
1.5000E+01, 2.0000E+01, 5.0000E+01,
$END
$GEOM
RCHAMB= 2.7949E-01
RTHRT= 1.8528E-01
RNE= 1.1100E-01
RTE= 1.1100E-01
ALPHA= 3.0000E+01
CHAMBL= 1.1876E+00
XC= 9.6477E-01
$END
$SCORE
TYPE= 'LOL'
NEL= 88
FDJ= 8.8345E-02
FCD= 8.1000E-01
FIH= 1.1481E-01
FIA= 3.0000E+01

```

FCANT=	1.6000E+01	
FFACET=	2.8728E-01	
XDJ=	1.0172E-01	
XCD=	9.4000E-01	
XIH=	1.7619E-01	
XIA=	3.0000E+01	
XCANT=	1.6000E+01	
XFACET=	4.4047E-01	
EMUNI=	6.6000E-01	
\$END		
\$BAFFLE		
\$END		
\$BARRIER		
\$END		
\$FFC		
\$END		
\$BURN		
GAMMA=	1.1462E+00	
AO=	4.0656E+03	
GMW=	2.3218E+01	
GPR=	5.9762E-01	
GK=	5.4180E-06	
GMU=	6.6324E-06	
RML=	7.6242E+01	
VJL=	9.0746E+02	
TJL=	5.3100E+02	
RHOL=	4.9895E+01	
CPL=	4.7921E-01	
PCRITL=	3.1600E+02	
TCRITL=	1.2180E+03	
TBOILL=	8.8200E+02	
XMWL=	1.7200E+02	
HVAPL=	1.2600E+02	
EN=	8.6789E-01	
TAUSEN=	1.4114E-04	
ISEN=	1	
\$END		
\$INJ		
FMAND=	6.7078E+00	
XMAND=	6.7078E+00	
FMANL=	3.3638E+00	
XMANL=	3.3638E+00	
PCA=	2.1411E+03,	1.6614E+03,
FRA=	4.7647E-01,	3.4643E-01,
FCAPA=	7.6914E-04,	8.1791E-04,
NFE=	4	8.6053E-04
FTLA=	6.0142E-04,	6.8198E-04,
FINA=	3.8878E-06,	3.8928E-06,
FFA=	1.0000E+00,	1.0000E+00,
NXE=	4	1.0000E+00,
XRA=	4.7647E-01,	3.4643E-01,
XCAPA=	6.8198E-04,	6.1747E-04,
XTLA=	9.3288E-04,	1.2438E-03,
XINA=	7.3028E-06,	7.3126E-06,
XFA=	1.0000E+00,	1.0000E+00,
XUOR=	4.7680E-02,	7.3412E-02,
AUOR=	3.4670E-03,	8.1288E-03,
RUOR=	3.3172E-02,	5.0881E-02,
XOR=	4.7680E-02,	7.3412E-02,
AOR=	3.4670E-03,	8.1288E-03,
ROR=	3.3172E-02,	5.0881E-02,
XDOR=	9.5761E-02,	1.4682E-01,
ADOR=	3.4670E-03,	8.1288E-03,
RDOR=	3.3172E-02,	5.0881E-02,

```

WDOT= 4.6216E+01, 1.3511E+02, 3.3533E+01, 9.6576E+01, 1.6644E+01,
4.7935E+01

$END
$CHAMBER
NCAV= 1,
XB= 2.6125E-01
ZB= 1.1875E-01
ZE= 4.9876E-01
MUB= 3
T= 3.3000E-02
WC= 2.0961E-02,
AC= 3.6906E-02,
ARATIO= 1.0000E+00,
IDCAV= 1,
DC1= 1.4000E-01,
CC1= 6*2.5000E+03
GAMC1= 5*1.2000E+00
CC2= 6*2.5000E+03
GAMC2= 5*1.2000E+00
ICAV= 1
$END
$FDORC
NZON= 5
FTER= 20*1.0000E+00
NCAV1= 1
ICTYP1= 4,
ZET= 2.0961E-02,
AE1= 3.6000E+02,
APL1= 1.4000E-01,
WC1= 1.9166E-01,
ZLOW1= -1.2316E-02,
ZUP1= 2.0417E-01,
$END
$MIX
EW(1)= 9.003E-01, 1.000E+00
$END
$DEBUGC
DEBUG= F
$END
$HIFIC
SHORT= F
POC= 2.0000E-01, 2.0000E-01
$END
$DISTDC
SHORT= F
POC= 2.0000E-01, 2.0000E-01
PAMP= 2.0000E-01
MC= 11
LC= 8
MS= 11
LB= 6
IDMAX= 10
$END
$CRPC
NDPC= 16
NDTFQ= 16
NDTLF= 1000
NPRINT= 60
NSUMB= 3600
PAMPC= 2.0000E-01
NRAD= 3
NCIRC= 5
$END
$LEINJC

```

```

IDMEM=      2
PAMPCN=    2.000E-01
NTINJ=      18
USGF=      2.000E-02
OGF=        2.000E-02
DSGF=      2.000E-02
FCDO=      1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00
XCDO=      1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00
$END
$COMBUSTC
FALVM=      1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00
XALVM=      1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00
FALTM=      1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00
XALTM=      1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00
FRNM=      1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00
XRNM=      1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00
ENMULT=     1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00
COMBX8=     5.000E-01
ACMULT=     1.000E+00
CCMULT=     1.000E+00
ENMULT=     1.000E+00
TAMULT=     1.000E+00
$END
$FDORCC
SHORT=      F
EPSIL=      2.000E+01
ERROR=      1.000E-01
LTS=        5
MTS=        5
NTS=        10
ITMAX=      100
RELX=      5.500E-01
NEET=       3
NXF8T=      3
NYF8T=      3
NAF8T=      3
MORE=       F
$END

```

END OF INPUT ECHO

# STABILITY MODEL INPUTS

## RUN DESCRIPTOR

ROCCID POINT DESIGN TEST CASE 1  
LOX/RP-1 LIKE DOUBLET PAIR WITH FIXED PC  
APPROXIMATES .0100 SUBSCALE DOUBLET

## SELECTED MODELS

BURNING MODEL=N-TAU INJECTION MODEL=INJ CHAMBER MODEL=HIFI

AXISYMETRIC=T DEBUG OUTPUT=F

## CHAMBER GEOMETRY AND OPERATING CONDITIONS

CHAMBER RADIUS, FT= 0.2785 THROAT RADIUS, FT= 0.1852  
CYLINDRICAL SECTION, FT= 0.9048 CONVERGENCE HALF ANGLE, DEG=30.0000  
NOZZLE ENTRANCE RADIUS OF CURVATURE, FT= 0.1110 THROAT ENTRANCE RADIUS OF CURVATURE, FT= 0.1110  
CHAMBER PRESSURE, PSIA=2141.10 MIXTURE RATIO= 2.8800  
SOUND SPEED, FT/SEC=4885.80 GAMMA=1.1462

## N-TAU BURNING MODEL INPUTS

PRESSURE INTERACTION INDEX, EN= 0.8878 SENSITIVE TIMELAG, TAU, SEC= 1.411E-04 SENSITIVE CIRCUIT=FUEL

## LUMPED INJECTION MODEL INPUTS

PC, PSIA=2141.10

	% TOTAL FLOW	RESISTANCE	INERTANCE (SEC)	CAPACITANCE (SEC)	TIMELAG (SEC)
FUEL:	100.000	4.788E-01	3.888E-08	7.881E-04	5.014E-04
	0.000	4.788E-01	0.000E+00	7.881E-04	0.000E+00
	0.000	4.788E-01	0.000E+00	7.881E-04	0.000E+00
	0.000	4.788E-01	0.000E+00	7.881E-04	0.000E+00
OX:	100.000	4.788E-01	7.303E-05	5.918E-04	9.324E-04
	0.000	4.788E-01	0.000E+00	5.918E-04	0.000E+00
	0.000	4.788E-01	0.000E+00	5.918E-04	0.000E+00
	0.000	4.788E-01	0.000E+00	5.918E-04	0.000E+00

PC, PSIA=1651.40

% TOTAL FLOW	RESISTANCE	INERTANCE	CAPACITANCE	TIMELAG
--------------	------------	-----------	-------------	---------

		(SEC)	(SEC)	(SEC)
FUEL:	100.000	3.454E-01	3.895E-05	6.179E-04
	0.000	3.454E-01	0.000E+00	0.000E+00
	0.000	3.454E-01	0.000E+00	0.000E+00
	0.000	3.454E-01	0.000E+00	0.000E+00
OX:	100.000	3.454E-01	7.312E-05	6.175E-04
	0.000	3.454E-01	0.000E+00	0.000E+00
	0.000	3.454E-01	0.000E+00	0.000E+00
	0.000	3.454E-01	0.000E+00	0.000E+00

PC, PSIA= 767.34

	% TOTAL FLOW	RESISTANCE	INERTANCE (SEC)	CAPACITANCE (SEC)	TIMELAG (SEC)
FUEL:	100.000	1.721E-01	3.905E-05	8.605E-04	1.337E-03
	0.000	1.721E-01	0.000E+00	8.605E-04	0.000E+00
	0.000	1.721E-01	0.000E+00	8.605E-04	0.000E+00
	0.000	1.721E-01	0.000E+00	8.605E-04	0.000E+00
OX:	100.000	1.720E-01	7.338E-05	8.561E-04	2.340E-03
	0.000	1.720E-01	0.000E+00	8.561E-04	0.000E+00
	0.000	1.720E-01	0.000E+00	8.561E-04	0.000E+00
	0.000	1.720E-01	0.000E+00	8.561E-04	0.000E+00

# HIFI CHAMBER MODEL INPUTS

COMBUSTION PLANE, FT= 2.612E-01      SHORT NOZZLE ASSUMED=F

## ACOUSTIC CAVITY INPUTS

CAVITY TYPE=1/4 WAVE  
 CAVITY TYPE 1:    NUMBER OF CAVITIES= 1      NUMBER OF PROPERTY SECTIONS= 1  
 CAVITY TYPE 2:    NUMBER OF CAVITIES= 0      NUMBER OF PROPERTY SECTIONS= 1  
 PARTITION THICKNESS, FT= 0.0000

	TOTAL DEPTH, FT	WIDTH, FT	AREA, SQ.FT	P'/PC	INLET TYPE
TYPE 1:	0.14000	0.02088	3.861E-02	0.200	0
TYPE 2:	0.00000	1.00000	0.000E+00	0.200	0

CAVITY TYPE ORIENTATION= 1

## CAVITY GAS PROPERTIES

	SECT	LENGTH	VSONIC	GAMMA
TYPE 1:	1	0.1400	2500.00	1.2000

BEGIN CALCULATIONS FOR 0 TANGENTIAL + 0 RADIAL MODE

AL (1/SEC)	MAX. AMPLITUDE	FREQUENCY (HZ)	Yb / Y	PHASE MARGIN (DEG)
0.000E+00	4.905E-01	1.552E+03	1.029E+00	180.00
-1.000E+02	5.466E-01	1.553E+03	1.029E+00	180.00
-2.000E+02	6.177E-01	1.555E+03	1.029E+00	180.00
-3.000E+02	7.096E-01	1.556E+03	1.029E+00	180.00
-4.000E+02	8.316E-01	1.557E+03	1.029E+00	-32.00
-6.000E+02	9.866E-01	1.559E+03	1.029E+00	-0.82

BEGIN CALCULATIONS FOR 1 TANGENTIAL + 0 RADIAL MODE					
AL (1/SEC)	MAX. AMPLITUDE	FREQUENCY (HZ)	YB / YJ	PHASE MARGIN (DEG)	
0.000E+00	7.91E-01	4.734E+03	2.745E+00	180.00	
-1.000E+02	6.23E-01	4.740E+03	2.745E+00	180.00	
-2.000E+02	6.801E-01	4.745E+03	2.745E+00	180.00	
-3.000E+02	9.010E-01	4.751E+03	2.745E+00	180.00	
-4.000E+02	9.471E-01	4.756E+03	2.745E+00	9.86	
-5.000E+02	9.991E-01	4.764E+03	2.926E+00	0.14	

# HIGH FREQUENCY STABILITY RESULTS

TANGENTIAL MODE	RADIAL MODE	GROWTH COEF. (1/S)	FREQUENCY (HZ)	YB / YJ
0	0	-500.0	1559.01	1.03
1	0	-500.0	4763.92	2.83

ROCCID  
Rocket Combustor Interactive Design Methodology  
Version 23-FEB-91

I-109

XIH=	1.7618E-01	
XIA=	3.0000E+01	
XCANT=	1.6000E+01	
XFACET=	4.4047E-01	
EMUNI=	6.6000E-01	
SEND		
\$BAFFLE		
SEND		
\$BARRIER		
SEND		
\$FFC		
SEND		
\$END		
\$BURN		
GAMMA=	1.1462E+00	
AO=	4.0656E+03	
GMM=	2.3218E+01	
GPR=	6.9752E-01	
GK=	6.4160E-06	
GMU=	6.6324E-06	
RML=	7.6242E+01	
VJL=	3.0746E+02	
TJL=	5.3100E+02	
RHOL=	4.8695E+01	
CPL=	4.7921E-01	
PCRITL=	3.1500E+02	
TCRITL=	1.2180E+03	
TBOILL=	6.6200E+02	
XMWL=	1.7200E+02	
HVAPL=	1.2500E+02	
EN=	6.6793E-01	
TAUSEN=	1.4114E-04	
ISEN=	1	
SEND		
\$INJ		
FMAND=	6.7076E+00	
XMAND=	6.7076E+00	
FWANL=	3.3636E+00	
XWANL=	3.3536E+00	
PCA=	2.1411E+03,	1.5514E+03, 7.6734E+02
FRA=	4.7547E-01,	3.4543E-01, 1.7206E-01
FCAPA=	7.6614E-04,	6.1761E-04, 6.6053E-04
NFE=	4	
FTLA=	6.0142E-04,	6.6196E-04, 1.3371E-03,
FINA=	3.6676E-06,	3.6626E-06, 3.9064E-06,
FFA=	1.0000E+00,	1.0000E+00, 1.0000E+00,
NXE=	4	
XRA=	4.7547E-01,	3.4543E-01, 1.7206E-01
XCAPA=	6.9169E-04,	6.1747E-04, 6.6615E-04
XTLA=	6.3236E-04,	1.2436E-03, 2.3386E-03,
XINA=	7.3026E-06,	7.3126E-06, 7.3382E-06,
XFA=	1.0000E+00,	1.0000E+00, 1.0000E+00,
XUOR=	4.7660E-02,	7.3412E-02,
AUOR=	3.4570E-03,	6.1269E-03,
RUOR=	3.3172E-02,	5.0661E-02,
XOR=	4.7660E-02,	7.3412E-02,
ACR=	3.4570E-03,	6.1269E-03,
ROR=	3.3172E-02,	5.0661E-02,
XDOR=	6.6761E-02,	1.4662E-01,
ADOR=	3.4570E-03,	6.1269E-03,
RDOR=	3.3172E-02,	5.0661E-03,
WDOT=	4.6216E+01,	1.3311E+02, 3.3533E+01, 6.6676E+01, 1.6644E+01,
SEND		
\$CHAMBER		

NCAV=	1,	
XB=	2.6126E-01	
ZB=	1.1675E-01	
ZE=	4.9875E-01	
MJB=	3	
T=	3.3000E-02	
WC=	2.0861E-02	
AC=	3.6600E-02	
ARATIO=	1.0000E+00	
IDCAV=	1,	
DC1=	1.4000E-01	
CC1=	5*2.5000E+03	
GAMC1=	5*1.2000E+00	
CC2=	5*2.5000E+03	
GAMC2=	5*1.2000E+00	
ICAV=	1	
\$END		
\$FDORC		
NZON=	5	
FTER=	20*1.0000E+00	
NCAV1=	1	
ICTYP1=	4,	
ZE1=	2.0861E-02	
AE1=	3.6000E+02	
APL1=	1.4000E-01	
WC1=	1.9166E-01	
ZLOW1=	-1.2315E-02	
ZUP1=	2.0417E-01	
\$END		
\$MIX		
EM(1)=	9.003E-01,	1.000E+00
\$END		
\$DEBUGC		
DEBUG=	F	
\$END		
\$HIFIC		
SHORT=	F	
POC=	2.0000E-01,	2.0000E-01
\$END		
\$DIST3DG		
SHORT=	F	
POC=	2.0000E-01,	2.0000E-01
PAMP=	2.0000E-01	
MC=	11	
LC=	8	
MB=	11	
LB=	8	
IDMAX=	10	
\$END		
\$CRPC		
NOPC=	16	
NOTFO=	16	
NOTLF=	1000	
NPRINT=	50	
NSUMB=	3500	
PAMPC=	2.0000E-01	
NRAD=	3	
NCIRC=	5	
\$END		
\$LEINJC		
IDMEM=	2	
PAMPCH=	2.0000E-01	
NTINJ=	16	
USGF=	2.0000E-02	

```

CGF=      2.0000E-02
DBGF=     2.0000E-02
FDDO=     1.0000E+00,
XCDO=     1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00
$END
$COMBUSTC
FALVM=     1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00
XALVM=     1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00
FALTM=     1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00
XALTM=     1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00
FRMA=     1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00
XRMA=     1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00, 1.0000E+00
ENMULT=    1.0000E+00, 1.0000E+00, 1.0000E+00
COMBXS=    5.0000E-01
ACMULT=    1.0000E+00
CCMULT=    1.0000E+00
ENMULT=    1.0000E+00
TAUMULT=    1.0000E+00
$END
$FDORCC
SHORT=      F
EPSIL=     2.0000E+01
ERROR=     1.0000E-01
LTS=        5
MTS=        5
NTS=        10
ITMAX=      100
RELX=       6.5000E-01
NEET=       3
NXFST=      3
NYFST=      3
NAFST=      3
MORE=       F
$END

```

END OF INPUT ECHO

## RUN DESCRIPTORS

ROCCID POINT DESIGN TEST CASE 1  
LOX/RP-1 LIKE DOUBLET PAIR WITH FIXED PC  
APPROXIMATES -0100 SUBSCALE DOUBLET

## PROPELLENT DESCRIPTION

FUEL=AP-1	Tmax., F=	71.00
OX=LOX	Tmax., F=	279.00

## CHAMBER GEOMETRY

CHAMBER RADIUS = 3.3539 IN.  
CYLINDRICAL SECTION = 11.8770 IN.  
NOZZLE ENTRANCE RADIUS OF CURVATURE = 1.3320 IN.  
CONVERGENCE HALF-ANGLE = 30.0000 DEG.  
THROAT RADIUS = 2.2326 IN.  
CONVERGENT SECTION LENGTH = 2.6730 IN.  
THROAT ENTRANCE RADIUS OF CURVATURE = 1.3320 IN.  
CONTRACTION RATIO = 2.28

## INJECTOR DATA

INJECTOR CORE CONTAINS 89 LOL ELEMENTS

ITEM	DESCRIPTION	UNIT	VALUE	UNIT	VALUE
FUEL SIDE:	Orifice Diam.	=	6.634E-02	in.	
	Impingement Half-angle	=	30.00	Deg.	
	Orifice Diam.	=	1.017E-01	in.	
	Impingement Half-angle	=	30.00	Deg.	
OX SIDE:	Orifice Diam.	=	6.634E-02	in.	
	Impingement Half-angle	=	30.00	Deg.	
	Orifice Diam.	=	1.017E-01	in.	
	Impingement Half-angle	=	30.00	Deg.	
FUEL SIDE:	Orifice Diam.	=	6.634E-02	in.	
	Impingement Half-angle	=	30.00	Deg.	
	Orifice Diam.	=	1.017E-01	in.	
	Impingement Half-angle	=	30.00	Deg.	
OX SIDE:	Orifice Diam.	=	6.634E-02	in.	
	Impingement Half-angle	=	30.00	Deg.	
	Orifice Diam.	=	1.017E-01	in.	
	Impingement Half-angle	=	30.00	Deg.	
FUEL SIDE:	Orifice Diam.	=	6.634E-02	in.	
	Impingement Half-angle	=	30.00	Deg.	
	Orifice Diam.	=	1.017E-01	in.	
	Impingement Half-angle	=	30.00	Deg.	
OX SIDE:	Orifice Diam.	=	6.634E-02	in.	
	Impingement Half-angle	=	30.00	Deg.	
	Orifice Diam.	=	1.017E-01	in.	
	Impingement Half-angle	=	30.00	Deg.	

## MIXING EFFICIENCIES

**CORE MIXING EFFICIENCY=0.9003**

## COMBUST CONTROL PARAMETERS

MULTIPLIERS:		CORE	BAFFLE	BARRIER	FFC
FUEL ATOMIZATION LENGTH FOR VAPORIZATION:		1.000	1.000	1.000	1.000
OX ATOMIZATION LENGTH FOR VAPORIZATION:		1.000	1.000	1.000	1.000
FUEL ATOMIZATION LENGTH FOR TIMELAGS:		1.000	1.000	1.000	1.000
OX ATOMIZATION LENGTH FOR TIMELAGS:		1.000	1.000	1.000	1.000
FUEL DROPSIZE:		1.000	1.000	1.000	1.000
OX DROPSIZE:		1.000	1.000	1.000	1.000
MIXING (Em):		1.000	1.000	1.000	
AO-Multiplier=1.000 Eta-C* for XB=0.500		CC-Multiplier=1.000	N-Multiplier=1.000		Tau-Multiplier=1.000

BEGIN STEADY STATE COMBUSTION ANALYSIS  
PC=2141.10 PSIA

PROPELLANT PROPERTIES

FUEL=RP-1	Phase=Liquid	T <sub>man</sub> .. F= 71.00	
	Injected Density= 49.89 Lbm/Cu. Ft	Viscosity=1.380E-03 Lbm/Ft-S	Surface Tension=1.067E-09 Lbf/Ft
OX=LOX	Phase=Liquid	T <sub>man</sub> .. F=279.00	
	Injected Density= 70.20 Lbm/Cu. Ft	Viscosity=1.167E-04 Lbm/Ft-S	Surface Tension=7.326E-04 Lbf/Ft

OPERATING CONDITIONS

PC FACE=2141.10 PSIA PC THROAT=2064.41 MIXTURE RATIO= 2.880  
FUEL INJECTION PRESSURE DROP= 502.74 Psia FUEL INJECTION VELOCITY= 309.56 Ft/S  
OX INJECTION PRESSURE DROP= 501.92 Psia OX INJECTION VELOCITY= 287.36 Ft/S  
FUEL FLOWRATE= 45.953 OX FLOWRATE= 132.345

ATOMIZATION OUTPUT

DROPSIZE MODEL=AEROJET

ELEMENT TYPE 1 IS LOL				
FUEL:	ATOMIZATION LENGTH, In.=1.06090	ATOMIZATION LENGTH FOR VAPORIZATION, In.=1.06090	DROPLET RADIUS, Microns= 76.29	
OX:	ATOMIZATION LENGTH, In.=2.32433	ATOMIZATION LENGTH FOR VAPORIZATION, In.=2.32433	DROPLET RADIUS, Microns= 84.56	

# VAPORIZATION CALCULATIONS

X (In.)	CORE=LOL		BAFFLE=		BARRIER=		FFC=	
	%FUEL VAP	%OX VAP	%FUEL VAP	%OX VAP	%FUEL VAP	%OX VAP	%FUEL VAP	%OX VAP
0.0000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
0.2850	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
0.5700	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
0.8550	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.1400	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.4250	10.524	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.7100	20.465	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.9950	28.379	0.000	0.000	0.000	0.000	0.000	0.000	0.000
2.2800	34.918	0.000	0.000	0.000	0.000	0.000	0.000	0.000
2.5650	40.682	7.081	0.000	0.000	0.000	0.000	0.000	0.000
2.8500	45.300	31.805	0.000	0.000	0.000	0.000	0.000	0.000
3.1350	49.437	48.257	0.000	0.000	0.000	0.000	0.000	0.000
3.4200	53.230	58.252	0.000	0.000	0.000	0.000	0.000	0.000
3.7050	55.333	63.892	0.000	0.000	0.000	0.000	0.000	0.000
3.9900	56.435	68.225	0.000	0.000	0.000	0.000	0.000	0.000
4.2750	61.962	73.768	0.000	0.000	0.000	0.000	0.000	0.000
4.5600	64.339	77.514	0.000	0.000	0.000	0.000	0.000	0.000
4.8450	65.558	80.335	0.000	0.000	0.000	0.000	0.000	0.000
5.1300	66.388	82.910	0.000	0.000	0.000	0.000	0.000	0.000
5.4150	70.215	84.764	0.000	0.000	0.000	0.000	0.000	0.000
5.7000	71.859	86.619	0.000	0.000	0.000	0.000	0.000	0.000
5.9850	73.448	88.288	0.000	0.000	0.000	0.000	0.000	0.000
6.2700	74.982	89.542	0.000	0.000	0.000	0.000	0.000	0.000
6.5550	76.274	90.867	0.000	0.000	0.000	0.000	0.000	0.000
6.8400	77.567	91.649	0.000	0.000	0.000	0.000	0.000	0.000
7.1250	78.857	92.457	0.000	0.000	0.000	0.000	0.000	0.000
7.4100	79.852	93.194	0.000	0.000	0.000	0.000	0.000	0.000
7.6950	80.836	93.922	0.000	0.000	0.000	0.000	0.000	0.000
7.9800	81.591	94.487	0.000	0.000	0.000	0.000	0.000	0.000
8.2650	82.545	95.032	0.000	0.000	0.000	0.000	0.000	0.000
8.5500	83.296	95.562	0.000	0.000	0.000	0.000	0.000	0.000
8.8350	83.872	95.998	0.000	0.000	0.000	0.000	0.000	0.000
9.1200	84.646	96.401	0.000	0.000	0.000	0.000	0.000	0.000
9.4050	85.324	96.728	0.000	0.000	0.000	0.000	0.000	0.000
9.6900	86.001	97.058	0.000	0.000	0.000	0.000	0.000	0.000
9.9750	86.677	97.322	0.000	0.000	0.000	0.000	0.000	0.000
10.2600	87.383	97.540	0.000	0.000	0.000	0.000	0.000	0.000
10.5450	87.987	97.768	0.000	0.000	0.000	0.000	0.000	0.000
10.8300	88.446	97.976	0.000	0.000	0.000	0.000	0.000	0.000
11.1150	88.802	98.185	0.000	0.000	0.000	0.000	0.000	0.000
11.4000	89.360	98.380	0.000	0.000	0.000	0.000	0.000	0.000
11.6850	89.836	98.524	0.000	0.000	0.000	0.000	0.000	0.000
11.9700	90.317	98.686	0.000	0.000	0.000	0.000	0.000	0.000
12.2550	91.118	98.938	0.000	0.000	0.000	0.000	0.000	0.000
12.5400	91.618	99.058	0.000	0.000	0.000	0.000	0.000	0.000
12.8250	91.919	99.141	0.000	0.000	0.000	0.000	0.000	0.000
13.1100	92.236	99.227	0.000	0.000	0.000	0.000	0.000	0.000
13.3950	92.533	99.312	0.000	0.000	0.000	0.000	0.000	0.000
13.6800	92.830	99.388	0.000	0.000	0.000	0.000	0.000	0.000
13.9650	93.126	99.447	0.000	0.000	0.000	0.000	0.000	0.000

OVERALL VAPORIZATION EFFICIENCIES  
 FUEL= 93.13% OX= 98.48%

MASS DISTRIBUTION PROFILE					
X (IN)	CORE (lbm/s)		BARRIER (lbm/s)		LOCAL VAPOR MIXTURE RATIO
	FUEL	OX	FUEL	OX	
0.0000	0.000	0.000	0.000	0.000	0.0000
0.2500	0.000	0.000	0.000	0.000	0.0000
0.5000	0.000	0.000	0.000	0.000	0.0000
0.7500	0.000	0.000	0.000	0.000	0.0000
1.0000	0.000	0.000	0.000	0.000	0.0000
1.2500	1.4250	0.000	0.000	0.000	0.0000
1.5000	9.404	0.000	0.000	0.000	0.0129
1.7500	13.041	0.000	0.000	0.000	0.0179
2.0000	16.046	0.000	0.000	0.000	0.0221
2.2500	18.787	9.365	0.000	0.000	0.0856
2.5000	20.917	41.628	0.000	0.000	0.3498
2.7500	22.718	61.219	0.000	0.000	0.4712
3.0000	24.461	74.446	0.000	0.000	0.6479
3.2500	26.887	84.253	0.000	0.000	0.8040
3.5000	27.312	91.617	0.000	0.000	0.8490
3.7500	28.469	97.629	0.000	0.000	0.8656
4.0000	29.586	102.586	0.000	0.000	0.7171
4.2500	30.585	108.319	0.000	0.000	0.7427
4.5000	31.426	109.727	0.000	0.000	0.7652
4.7500	32.266	112.162	0.000	0.000	0.7836
5.0000	33.020	114.636	0.000	0.000	0.8012
5.2500	33.751	116.645	0.000	0.000	0.8175
5.5000	34.447	118.505	0.000	0.000	0.8312
5.7500	35.061	119.954	0.000	0.000	0.8432
6.0000	35.654	121.293	0.000	0.000	0.8545
6.2500	36.146	122.376	0.000	0.000	0.8638
6.5000	36.602	123.338	0.000	0.000	0.8722
6.7500	37.055	124.301	0.000	0.000	0.8806
7.0000	37.484	125.049	0.000	0.000	0.8879
7.2500	37.932	125.770	0.000	0.000	0.8951
7.5000	38.277	126.472	0.000	0.000	0.9013
7.7500	38.568	127.049	0.000	0.000	0.9067
8.0000	38.898	127.582	0.000	0.000	0.9119
8.2500	39.208	128.016	0.000	0.000	0.9168
8.5000	39.520	128.448	0.000	0.000	0.9213
8.7500	39.831	128.801	0.000	0.000	0.9257
9.0000	40.141	129.090	0.000	0.000	0.9298
9.2500	40.433	129.378	0.000	0.000	0.9337
9.5000	40.643	129.667	0.000	0.000	0.9369
9.7500	40.863	129.858	0.000	0.000	0.9401
10.0000	41.084	130.174	0.000	0.000	0.9429
10.2500	41.283	130.391	0.000	0.000	0.9459
10.5000	41.503	130.580	0.000	0.000	0.9487
10.7500	41.687	130.768	0.000	0.000	0.9512
11.0000	41.871	130.938	0.000	0.000	0.9536
11.2500	42.056	131.098	0.000	0.000	0.9559
11.5000	42.239	131.209	0.000	0.000	0.9580
11.7500	42.388	131.322	0.000	0.000	0.9596
12.0000	42.522	131.436	0.000	0.000	0.9615
12.2500	42.656	131.546	0.000	0.000	0.9632
12.5000	42.784	131.614	0.000	0.000	0.9647

# AXIAL PRESSURE PROFILE

X (in)	MACH #	Ptotal (psia)	Pstatic (psia)	Ttotal (R)	Tstatic (R)	Wdot (Lbm/s)	Local Radius (in)
1.20	0.001	2141.80	2142.78	1692.28	1696.27	0.74	3.354
1.30	0.002	2141.80	2142.75	1692.26	1696.27	2.48	3.354
1.61	0.006	2141.76	2142.69	1694.00	1696.01	7.88	3.354
1.91	0.008	2141.71	2142.69	1692.26	1696.27	12.04	3.354
2.21	0.011	2141.66	2142.49	1714.30	1708.23	15.13	3.354
2.51	0.018	2141.40	2141.97	1987.60	1980.46	23.75	3.354
2.81	0.076	2134.79	2126.73	5659.05	5637.47	57.98	3.354
3.11	0.120	2124.47	2107.97	6839.73	6811.23	82.67	3.354
3.41	0.145	2116.91	2092.70	6861.86	6831.71	98.53	3.354
3.71	0.163	2110.42	2078.57	6856.87	6825.41	110.41	3.354
4.01	0.178	2104.90	2068.35	6847.00	6814.45	119.59	3.354
4.32	0.190	2100.08	2056.55	6837.35	6803.85	127.02	3.354
4.62	0.199	2095.85	2049.90	6832.15	6797.78	133.17	3.354
4.92	0.207	2092.31	2042.67	6830.93	6795.61	138.04	3.354
5.22	0.214	2089.14	2036.16	6829.57	6793.77	142.25	3.354
5.52	0.220	2086.49	2030.73	6831.98	6795.58	145.62	3.354
5.82	0.226	2083.81	2025.22	6832.58	6795.58	148.96	3.354
6.12	0.231	2081.48	2020.40	6835.22	6797.88	151.78	3.354
6.42	0.235	2078.50	2016.35	6838.29	6800.30	154.09	3.354
6.72	0.238	2077.66	2012.57	6841.42	6803.00	156.20	3.354
7.02	0.242	2076.08	2009.31	6843.92	6805.12	158.00	3.354
7.33	0.244	2074.73	2006.62	6846.64	6806.42	159.82	3.354
7.63	0.247	2073.37	2003.70	6847.15	6807.71	161.03	3.354
7.93	0.249	2072.18	2001.25	6849.08	6809.35	162.33	3.354
8.23	0.251	2071.04	1998.90	6851.21	6811.21	163.56	3.354
8.53	0.253	2070.00	1996.74	6852.50	6812.25	164.68	3.354
8.83	0.255	2069.12	1994.90	6853.72	6813.25	165.82	3.354
9.13	0.257	2068.27	1993.16	6855.05	6814.38	166.51	3.354
9.43	0.258	2067.52	1991.60	6856.64	6815.78	167.30	3.354
9.73	0.260	2066.77	1990.04	6858.20	6817.18	168.08	3.354
10.04	0.261	2066.10	1988.66	6859.97	6818.76	168.76	3.354
10.34	0.262	2065.48	1987.37	6861.91	6820.53	169.39	3.354
10.64	0.263	2064.90	1986.17	6863.12	6821.60	169.98	3.354
10.94	0.264	2064.40	1985.13	6862.98	6821.35	170.50	3.354
11.24	0.265	2063.92	1984.09	6863.06	6821.30	171.00	3.354
11.54	0.266	2063.48	1983.22	6863.08	6821.23	171.45	3.354
11.84	0.267	2063.04	1979.56	6863.12	6820.03	171.90	3.328
12.14	0.291	2062.60	1966.57	6863.15	6814.89	172.31	3.225
12.44	0.326	2062.11	1940.65	6863.18	6804.38	172.69	3.061
12.74	0.377	2061.49	1902.92	6863.21	6788.55	173.06	2.886
13.05	0.439	2060.76	1848.36	6863.24	6863.20	173.39	2.712
13.35	0.528	2059.89	1764.18	6863.26	6727.74	173.67	2.538
13.65	0.687	2058.68	1617.13	6863.29	6856.19	173.93	2.387
13.95	0.920	2056.71	1415.52	6863.31	6853.31	174.19	2.257
14.25	0.999	2054.41	1186.32	6863.33	6421.46	174.41	2.223

# PERFORMANCE SUMMARY

## C\* EFFICIENCY CALCULATIONS (ODK)

INJECTED MR= 2.8800 CSTAR=5880.40 CORE Em=0.9003 BARRIER Em=1.0000  
 CORE: OVERALL MR= 2.8800 VAPOR MR= 3.0755 CSTAR-MIX=5779.55 MASS FRACTION= 1.0000  
 BARRIER: OVERALL MR= 0.0000 VAPOR MR=99.9000 CSTAR-MIX= 0.00 MASS FRACTION= 0.0000  
 ENGINE: OVERALL MR= 2.8800 VAPOR MR= 3.0755 CSTAR-DEL=5653.55  
 C\* EFFICIENCY = 9.847E-01

## ISP EFFICIENCY CALCULATIONS

ISP-ODK, INJ = 2.889E+02 SEC.  
 ISP-ODK, M.Z. INJ = 2.855E+02 SEC.  
 VAPORIZATION EFFICIENCY = 9.701E-01  
 ENERGY RELEASE EFFICIENCY = 9.850E-01

ISP-ODK, M.Z. VAPOR = 2.833E+02 SEC.  
 MIXING EFFICIENCY = 9.847E-01

NOTE: ISP-DEL = ISP-ODK, INJ. \* ERE \* ETADIV - DELISP-BL

TIME-LAG CALCULATIONS, Milliseconds

Cohem, In.=9.099E-03      FUEL Cohem, In.=2.293E+02      OX Cohem, In.=8.194E+01

ELEMENT 1 IS TYPE=LOL

FUEL:      CinJ, In.=1.444E-02      Lvap, In.= 0.519      ATOMIZATION LENGTH USED, In.= 1.081E+00  
             Timp=9.619E-02      Tatom=2.893E-01      Tvap=1.415E-01      Total=4.670E-01

OX:      CinJ, In.=1.474E-02      Lvap, In.= 0.189      ATOMIZATION LENGTH USED, In.= 2.324E+00  
             Timp=6.598E-02      Tatom=7.526E-01      Tvap=6.126E-02      Total=8.798E-01

EFFECTIVE TIMELAGS, Milliseconds

FUEL:      CinJ, In.=1.444E-02      Lvap, In.= 0.519      Total=4.670E-01  
             Timp=9.619E-02      Tatom=2.893E-01      Tvap=1.415E-01

OX:      CinJ, In.=1.474E-02      Lvap, In.= 0.189      Total=8.798E-01  
             Timp=6.598E-02      Tatom=7.526E-01      Tvap=6.126E-02

# CHAMBER-NOZZLE OPTIMIZATION RESULTS

CHAMBER LENGTH (FEET)	ETA-C*	ETA-NOZ	OVERALL EFFICIENCY
0.0000	0.0000	0.8807	0.0000
0.1667	0.0160	0.8726	0.0187
0.3333	0.0503	0.8646	0.0629
0.5000	0.1182	0.8537	0.0985
0.6667	0.1884	0.8419	0.1740
0.8333	0.2681	0.8302	0.2689
1.0000	0.3489	0.8186	0.3767
1.1667	0.4234	0.8087	0.4762
1.3333	0.4739	0.7916	0.5504
1.5000	0.5044	0.7773	0.6071
1.6667	0.5246	0.7631	0.6439
1.8333	1.0000	0.7486	0.6689
2.0000	1.0000	0.7299	0.6762
2.1667	1.0000	0.7114	0.6770
2.3333	1.0000	0.6928	0.6761
2.5000	1.0000	0.6742	0.6751
2.6667	1.0000	0.6461	0.6746
2.8333	1.0000	0.6142	0.6729
3.0000	1.0000	0.5823	0.7114
3.1667	1.0000	0.5504	0.8028
3.3333	1.0000	0.5071	0.8742
3.5000	1.0000	0.4439	0.8461
3.6667	1.0000	0.3809	0.8142
3.8333	1.0000	0.3176	0.5823
4.0000	1.0000	0.2544	0.5504

OPTIMUM CHAMBER LENGTH= 1.0000 FT  
 MAXIMUM OVERALL EFFICIENCY= 0.7767

# REDESIGNED CHAMBER RESULTS

NOMINAL CHAMBER PRESSURE = 2.141E+03 PSIA  
 THROTTLED CHAMBER PRESSURE = 1.636E+03 PSIA  
 FUEL INJECTION PRESSURE DROP = 6.027E+02 PSI  
 OX INJECTION PRESSURE DROP = 6.018E+02 PSI  
 CHAMBER RADIUS = 2.795E-01 FT  
 THROAT RADIUS = 1.862E-01 FT  
 NOZZLE ENTRANCE RADIUS OF CURVATURE = 1.110E-01 FT  
 THROAT ENTRANCE RADIUS OF CURVATURE = 1.110E-01 FT  
 NOZZLE CONVERGENCE HALF-ANGLE = 3.000E+01 DEG  
 INJECTOR-TO-THROAT CHAMBER LENGTH = 1.188E+00 FT  
 BARREL SECTION LENGTH = 9.648E-01 FT

## IMPINGING ELEMENT SIZING RESULTS

ELEMENT TYPE = LOL  
 NO. OF ELEMENTS = 80  
 FUEL ORIFICE DIAMETER = 7.119E-02 IN  
 OX ORIFICE DIAMETER = 1.091E-01 IN

## CORE ELEMENT SPACING RESULTS

ELEMENT TYPE = LOL  
 NUMBER OF ELEMENTS = 80  
 FUEL ORIFICE/ANNULUS DIAMETER = 7.119E-02 IN  
 OXIDIZER ORIFICE DIAMETER = 1.091E-01 IN  
 FUEL INJECTION VELOCITY = 3.056E+02 FT/S  
 OXIDIZER INJECTION VELOCITY = 2.574E+02 FT/S

ROW	# ELEMENTS	MID-ROW RADIUS (IN)
1	8	7.347E-01
2	12	1.493E+00
3	18	2.291E+00
4	24	2.900E+00

DIRECT INPUT ECHO FROM SUBROUTINE PINPUT

ROCCID  
Rocket Combustor Interactive Design Methodology  
Version 23-FEB-91

ROCCID POINT DESIGN TEST CASE 1  
LOX/RP-1 LIKE DOUBLET PAIR WITH FIXED PC  
APPROXIMATES -0100 SUBSCALE DOUBLET

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$MODELS
MCHAM= 1
MBURN= 2
MINJ= 1
$END

$OPCOND
PISPA= 3.9800E+01, 1.0840E+02, 1.3840E+02, 1.5880E+02, 1.7780E+02,
2.0970E+02, 2.3850E+02, 2.6470E+02, 2.6700E+02, 2.7040E+02,
2.7180E+02, 2.7180E+02, 2.7130E+02, 2.6800E+02, 2.6820E+02,
2.6800E+02, 2.6500E+02, 2.6210E+02, 2.5920E+02, 2.5830E+02,
2.5080E+02, 2.3840E+02, 2.2750E+02, 2.1100E+02, 2.0140E+02,
1.5580E+02, 1.0870E+02, 8.0000E+01,
PCBA= 1.7780E+03, 2.4860E+03, 3.4870E+03, 4.0740E+03, 4.8800E+03,
5.0700E+03, 5.4840E+03, 5.7340E+03, 5.8080E+03, 5.8810E+03,
5.8890E+03, 5.9540E+03, 5.9840E+03, 5.9080E+03, 5.820E+03,
5.6550E+03, 5.8280E+03, 5.7590E+03, 5.7000E+03, 5.6430E+03,
5.5320E+03, 5.2750E+03, 5.0470E+03, 4.6760E+03, 4.4200E+03,
2.4670E+03, 2.4940E+03, 1.4850E+03,
FUEL= 'RP-1',
OX= 'LOX',
PC= 2.1411E+03
XMR= 2.8800E+00
FTMAN= 7.1000E+01
XTMAN= -2.7900E+02
EMMAN= 1.0000E+00
NPERFP= 28
PMRA= 1.0000E-01, 3.0000E-01, 6.0000E-01, 8.0000E-01, 1.0000E+00,
1.2500E+00, 1.5000E+00, 1.7800E+00, 2.0000E+00, 2.2000E+00,
2.4000E+00, 2.5000E+00, 2.6000E+00, 2.7000E+00, 2.8000E+00,
2.8000E+00, 3.0000E+00, 3.2000E+00, 3.4000E+00, 3.6000E+00,
4.0000E+00, 5.0000E+00, 6.0000E+00, 8.0000E+00, 1.0000E+01,
1.5000E+01, 2.0000E+01, 5.0000E+01, 1.0000E+02, 2.0000E+02,
$END

$GEOM
RCHAMB= 2.7849E-01
RTHRT= 1.8523E-01
RNE= 1.1100E-01
RTE= 1.1100E-01
ALPHA= 3.0000E+01
CHAMBL= 1.1878E+00
XC= 8.6477E-01
$END

$SCORE
TYPE= 'LOL'
NEL= 80
FDJ= 7.1147E-02
FOD= 9.1000E-01
FIH= 1.2329E-01
FIA= 3.0000E+01
FCANT= 1.6000E+01
FFACET= 3.0808E-01
XDJ= 1.0909E-01
XCD= 9.4000E-01

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XIH=	1.8894E-01		
XIA=	3.0000E+01		
XCANT=	1.6000E+01		
XFACET=	4.7285E-01		
EMUNI=	6.5000E-01		
SEND			
\$BAFFLE			
SEND			
\$BARRIER			
SEND			
\$FFC			
SEND			
\$BURN			
GAMMA=	1.1482E+00		
AO=	4.0656E+03		
GMM=	2.3218E+01		
GPR=	5.9732E-01		
GK=	5.4160E-05		
GMU=	6.6924E-05		
RML=	7.6242E+01		
VJL=	3.0746E+02		
TJL=	5.3100E+02		
RHOL=	4.9895E+01		
CPL=	4.7321E-01		
PCRTL=	3.1500E+02		
TCRTL=	1.2180E+03		
TBOILL=	6.8200E+02		
XMWL=	1.7200E+02		
HVAPL=	1.2500E+02		
EN=	8.6783E-01		
TAUSEN=	1.4114E-04		
ISEN=	1		
SEND			
\$INJ			
FMAND=	6.7076E+00		
FMAND=	6.7076E+00		
FMANL=	3.3538E+00		
XMANL=	3.3538E+00		
PCA=	2.1411E+03,	1.8814E+03,	7.6734E+02
FRA=	4.7847E-01,	3.4843E-01,	1.7206E-01
FCAPA=	7.8914E-04,	8.1781E-04,	8.6053E-04
NFE=	4		
FTLA=	5.0142E-04,	6.8186E-04,	1.3371E-03,
FINA=	3.8976E-06,	3.8928E-06,	3.9084E-06,
FFA=	1.0000E+00,	1.0000E+00,	1.0000E+00,
NXE=	4		
XRA=	4.7847E-01,	3.4843E-01,	1.7206E-01
XCAPA=	5.9168E-04,	6.1747E-04,	6.5615E-04
XTLA=	9.3296E-04,	1.2439E-03,	2.3399E-03,
XINA=	7.3026E-06,	7.3126E-06,	7.3382E-06,
XFA=	1.0000E+00,	1.0000E+00,	1.0000E+00,
XUOR=	4.7880E-02,	7.3412E-02,	
AUOR=	3.4570E-03,	8.1269E-03,	
RUOR=	3.3172E-02,	6.0861E-02,	
XOR=	4.7880E-02,	7.3412E-02,	
AOR=	3.4570E-03,	8.1269E-03,	
ROR=	3.3172E-02,	6.0861E-02,	
XDOR=	8.6761E-02,	1.4682E-01,	
ADOR=	3.4570E-03,	6.1269E-03,	
RDOR=	3.3172E-02,	6.0861E-02,	
WDOT=	4.8218E+01,	1.3911E+02,	3.3633E+01, 9.8576E+01, 1.8844E+01,
	4.7936E+01		
SEND			
\$CHAMBER			

NCAV=	1,
XB=	2.8126E-01
ZB=	1.1675E-01
ZE=	4.8875E-01
MJB=	3
T=	3.3000E-02
WC=	2.0881E-02,
AC=	3.8808E-02,
ARATIO=	1.0000E+00,
IDCAV=	1,
DC1=	1.4000E-01,
CG1=	5*2.5000E+03
GAMC1=	5*1.2000E+00
CC2=	5*2.5000E+03
GAMC2=	5*1.2000E+00
ICAV=	1
\$END	
\$FDORG	
NZON=	5
FTER=	20*1.0000E+00
NCAV1=	1
ICTYP1=	4,
ZE1=	2.0881E-02,
AE1=	3.8000E+02,
APL1=	1.4000E-01,
WC1=	1.8168E-01,
ZLOW1=	-1.2315E-02,
ZUP1=	2.0417E-01,
\$END	
\$MIX	
EM(1)=	8.830E-01, 1.000E+00
\$END	
\$DEBUG	
DEBUG=	F
\$END	
\$SHIFIG	
SHORT=	F
POC=	2.0000E-01, 2.0000E-01
\$END	
\$DIST3DC	
SHORT=	F
POC=	2.0000E-01, 2.0000E-01
PAMP=	2.0000E-01
MC=	11
LG=	8
MB=	11
LB=	8
IDMAX=	10
\$END	
\$CRPG	
NOPC=	16
NOTFG=	16
NDTLF=	1000
NPRINT=	50
NSUMS=	3600
PAMP=	2.0000E-01
NRAD=	3
NCIRC=	5
\$END	
\$LEINJC	
IDMEM=	2
PAMPCH=	2.0000E-01
NTINJ=	16
USGF=	2.0000E-02

QGF= 2.000E-02  
 DGGF= 2.000E-02  
 FCDO= 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00  
 XCDO= 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00  
 \$END  
 \$COMBUSTC  
 FALVM= 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00  
 XALVM= 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00  
 FALTM= 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00  
 XALTM= 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00  
 FRMM= 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00  
 XFRMM= 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00  
 ENMULT= 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00  
 COMBXB= 5.000E-01  
 ACMULT= 1.000E+00  
 CCMULT= 1.000E+00  
 ENMULT= 1.000E+00  
 TALMULT= 1.000E+00  
 \$END  
 \$FDORCC  
 SHORT= F  
 EPSIL= 2.000E+01  
 ERROR= 1.000E-01  
 LTS= 5  
 MTS= 5  
 NTS= 10  
 ITMAX= 100  
 RELX= 8.500E-01  
 NEET= 3  
 NXFST= 3  
 NYFST= 3  
 NAFST= 3  
 MORE= F  
 \$END

END OF INPUT ECHO

# STEADY STATE COMBUSTION ANALYSIS PROGRAM

## RUN DESCRIPTORS

ROCCID POINT DESIGN TEST CASE 1  
LOX/RP-1 LIKE DOUBLET PAIR WITH FIXED PC  
APPROXIMATES 0100 SUBSCALE DOUBLET

## PROPELLENT DESCRIPTION

FUEL=RP-1 Tman., F= 71.00  
OX=LOX Tman., F=-279.00

## CHAMBER GEOMETRY

CHAMBER RADIUS = 3.3639 IN.  
CYLINDRICAL SECTION =11.5770 IN.  
NOZZLE ENTRANCE RADIUS OF CURVATURE = 1.3320 IN.  
CONVERGENCE HALF-ANGLE =30.0000 DEG.  
THROAT RADIUS = 2.2228 IN.  
CONVERGENT SECTION LENGTH = 2.9730 IN.  
THROAT ENTRANCE RADIUS OF CURVATURE = 1.3320 IN.  
CONTRACTION RATIO = 2.28

## INJECTOR DATA

## INJECTOR CORE CONTAINS 6.0 LOL ELEMENTS

FUEL SIDE: Orifice Diam. =7.115E-02 In.  
Impingement Half-angle =30.00 Deg.  
Cd =0.9100  
OX SIDE: Orifice Diam. =1.091E-01 In.  
Impingement Half-angle =30.00 Deg.  
Cd =0.9400  
Unlike Cant Angle =16.00 Deg.  
Unlike Cant Angle =16.00 Deg.  
Impingement Height =0.123 in.  
Faceplate Thickness = 0.3081 in.  
Impingement Height =0.169 in.  
Faceplate Thickness = 0.4724 in.

## MIXING EFFICIENCIES

CORE MIXING EFFICIENCY=0.9930 BARRIER MIXING EFFICIENCY=1.0000

## COMBUST CONTROL PARAMETERS

MULTIPLIERS:  
FUEL ATOMIZATION LENGTH FOR VAPORIZATION: 1.000  
OX ATOMIZATION LENGTH FOR VAPORIZATION: 1.000  
FUEL ATOMIZATION LENGTH FOR TIMELAGS: 1.000  
OX ATOMIZATION LENGTH FOR TIMELAGS: 1.000  
FUEL DROPSIZE: 1.000  
OX DROPSIZE: 1.000  
MIXING (Em): 1.000  
AO-Multiplier=1.000 CO-Multiplier=1.000 N-Multiplier=1.000  
Eta-C\* for XB=0.500 Tau-Multiplier=1.000

BEGIN STEADY STATE COMBUSTION ANALYSIS  
PC=2141.10 PSIA

PROPELLANT PROPERTIES			
FUEL=RP-1	Phase=Liquid	Tman., F= 71.00	
	Injected Density= 49.69 Lbm/Cu. Ft	Viscosity=1.360E-03 Lbm/Ft-S	Surface Tension=1.657E-03 Lbf/Ft
OX=LOX	Phase=Liquid	Tman., F=-278.00	
	Injected Density= 70.20 Lbm/Cu. Ft	Viscosity=1.167E-04 Lbm/Ft-S	Surface Tension=7.326E-04 Lbf/Ft
OPERATING CONDITIONS			
PC FACE=2141.10 PSIA	PC THROAT=2092.78	MIXTURE RATIO= 2.890	
FUEL INJECTION PRESSURE DROP= 510.07 Psia	FUEL INJECTION VELOCITY= 307.78 Ft/S		
OX INJECTION PRESSURE DROP= 509.14 Psia	OX INJECTION VELOCITY= 259.23 Ft/S		
FUEL FLOWRATE= 49.277	OX FLOWRATE= 139.277		
ATOMIZATION OUTPUT			
DROPSIZE MODEL=AEROJET			
ELEMENT TYPE 1 IS LOL			
FUEL:	ATOMIZATION LENGTH, In.=1.14748	ATOMIZATION LENGTH FOR VAPORIZATION, In.=1.14748	DROPLET RADIUS, Microns= 81.12
OX:	ATOMIZATION LENGTH, In.=2.51421	ATOMIZATION LENGTH FOR VAPORIZATION, In.=2.51421	DROPLET RADIUS, Microns= 89.91

# VAPORIZATION CALCULATIONS

X (in.)	CORE-LOL		BAFFLE-		BARRIER-		FFC-	
	%FUEL VAP	%OX VAP	%FUEL VAP	%OX VAP	%FUEL VAP	%OX VAP	%FUEL VAP	%OX VAP
0.0000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
0.2850	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
0.5700	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
0.8550	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.1400	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.4250	5.403	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.7100	16.248	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.9950	24.267	0.000	0.000	0.000	0.000	0.000	0.000	0.000
2.2800	30.676	0.000	0.000	0.000	0.000	0.000	0.000	0.000
2.5650	36.456	0.000	0.000	0.000	0.000	0.000	0.000	0.000
2.8500	41.859	15.086	0.000	0.000	0.000	0.000	0.000	0.000
3.1350	45.624	34.360	0.000	0.000	0.000	0.000	0.000	0.000
3.4200	49.368	46.985	0.000	0.000	0.000	0.000	0.000	0.000
3.7050	52.914	56.016	0.000	0.000	0.000	0.000	0.000	0.000
3.9900	56.797	62.691	0.000	0.000	0.000	0.000	0.000	0.000
4.2750	58.561	68.162	0.000	0.000	0.000	0.000	0.000	0.000
4.5600	61.064	72.468	0.000	0.000	0.000	0.000	0.000	0.000
4.8450	63.238	76.117	0.000	0.000	0.000	0.000	0.000	0.000
5.1300	65.408	79.030	0.000	0.000	0.000	0.000	0.000	0.000
5.4150	67.211	81.467	0.000	0.000	0.000	0.000	0.000	0.000
5.7000	68.676	83.543	0.000	0.000	0.000	0.000	0.000	0.000
5.9850	70.641	85.231	0.000	0.000	0.000	0.000	0.000	0.000
6.2700	71.996	86.916	0.000	0.000	0.000	0.000	0.000	0.000
6.5550	73.444	88.377	0.000	0.000	0.000	0.000	0.000	0.000
6.8400	74.941	89.519	0.000	0.000	0.000	0.000	0.000	0.000
7.1250	76.026	90.560	0.000	0.000	0.000	0.000	0.000	0.000
7.4100	77.230	91.454	0.000	0.000	0.000	0.000	0.000	0.000
7.6950	78.287	92.257	0.000	0.000	0.000	0.000	0.000	0.000
7.9800	79.202	92.919	0.000	0.000	0.000	0.000	0.000	0.000
8.2650	80.107	93.560	0.000	0.000	0.000	0.000	0.000	0.000
8.5500	80.987	94.182	0.000	0.000	0.000	0.000	0.000	0.000
8.8350	81.866	94.678	0.000	0.000	0.000	0.000	0.000	0.000
9.1200	82.726	95.174	0.000	0.000	0.000	0.000	0.000	0.000
9.4050	83.562	95.636	0.000	0.000	0.000	0.000	0.000	0.000
9.6900	83.377	96.033	0.000	0.000	0.000	0.000	0.000	0.000
9.9750	84.593	96.396	0.000	0.000	0.000	0.000	0.000	0.000
10.2600	86.208	96.896	0.000	0.000	0.000	0.000	0.000	0.000
10.5450	86.823	96.996	0.000	0.000	0.000	0.000	0.000	0.000
10.8300	86.436	97.261	0.000	0.000	0.000	0.000	0.000	0.000
11.1150	87.064	97.466	0.000	0.000	0.000	0.000	0.000	0.000
11.4000	87.669	97.666	0.000	0.000	0.000	0.000	0.000	0.000
11.6850	88.178	97.868	0.000	0.000	0.000	0.000	0.000	0.000
11.9700	88.644	98.067	0.000	0.000	0.000	0.000	0.000	0.000
12.2550	89.110	98.262	0.000	0.000	0.000	0.000	0.000	0.000
12.5400	89.576	98.449	0.000	0.000	0.000	0.000	0.000	0.000
12.8250	90.041	98.690	0.000	0.000	0.000	0.000	0.000	0.000
13.1100	90.440	98.719	0.000	0.000	0.000	0.000	0.000	0.000
13.3950	90.804	98.949	0.000	0.000	0.000	0.000	0.000	0.000
13.6800	91.169	99.060	0.000	0.000	0.000	0.000	0.000	0.000
13.9650	91.533	99.065	0.000	0.000	0.000	0.000	0.000	0.000
14.2500	91.898	99.142	0.000	0.000	0.000	0.000	0.000	0.000

OVERALL VAPORIZATION EFFICIENCIES  
 FUEL= 91.00% OX= 99.14%

# MASS DISTRIBUTION PROFILE

X (IN)	CORE (lbm/s)		BARRIER (lbm/s)		LOCAL VAPOR MIXTURE RATIO	ETA-C*
	FUEL	OX	FUEL	OX		
0.0000	0.000	0.000	0.000	0.000	0.00	0.0000
0.2850	0.000	0.000	0.000	0.000	0.00	0.0000
0.5700	0.000	0.000	0.000	0.000	0.00	0.0000
0.8550	0.000	0.000	0.000	0.000	0.00	0.0000
1.1400	0.000	0.000	0.000	0.000	0.00	0.0000
1.4250	2.500	0.000	0.000	0.000	0.00	0.0034
1.7100	7.619	0.000	0.000	0.000	0.00	0.0103
1.9950	11.225	0.000	0.000	0.000	0.00	0.0153
2.2800	14.197	0.000	0.000	0.000	0.00	0.0194
2.5650	16.671	0.000	0.000	0.000	0.00	0.0230
2.8500	19.371	20.106	0.000	0.000	1.04	0.1760
3.1350	21.113	45.784	0.000	0.000	2.17	0.3732
3.4200	22.665	82.533	0.000	0.000	2.74	0.4752
3.7050	24.467	74.656	0.000	0.000	3.05	0.5449
3.9900	25.793	83.616	0.000	0.000	3.25	0.5988
4.2750	27.100	90.671	0.000	0.000	3.35	0.6390
4.5600	28.259	96.583	0.000	0.000	3.42	0.6742
4.8450	29.284	101.446	0.000	0.000	3.47	0.7042
5.1300	30.289	105.329	0.000	0.000	3.48	0.7301
5.4150	31.103	108.677	0.000	0.000	3.49	0.7516
5.7000	31.674	111.344	0.000	0.000	3.49	0.7706
5.9850	32.644	113.553	0.000	0.000	3.48	0.7874
6.2700	33.316	115.642	0.000	0.000	3.48	0.8032
6.5550	33.986	117.787	0.000	0.000	3.47	0.8177
6.8400	34.634	119.306	0.000	0.000	3.44	0.8303
7.1250	35.167	120.996	0.000	0.000	3.43	0.8413
7.4100	35.736	121.686	0.000	0.000	3.41	0.8515
7.6950	36.233	122.957	0.000	0.000	3.39	0.8607
7.9800	36.652	123.839	0.000	0.000	3.38	0.8683
8.2650	37.071	124.721	0.000	0.000	3.36	0.8759
8.5500	37.476	125.622	0.000	0.000	3.35	0.8831
8.8350	37.860	126.184	0.000	0.000	3.33	0.8897
9.1200	38.232	126.846	0.000	0.000	3.31	0.8962
9.4050	38.577	127.461	0.000	0.000	3.30	0.9016
9.6900	38.862	127.990	0.000	0.000	3.29	0.9066
9.9750	39.147	128.476	0.000	0.000	3.28	0.9112
10.2600	39.431	128.873	0.000	0.000	3.27	0.9155
10.5450	39.716	129.270	0.000	0.000	3.25	0.9198
10.8300	40.001	129.626	0.000	0.000	3.24	0.9240
11.1150	40.286	129.881	0.000	0.000	3.22	0.9277
11.4000	40.570	130.155	0.000	0.000	3.21	0.9314
11.6850	40.806	130.432	0.000	0.000	3.20	0.9348
11.9700	41.022	130.728	0.000	0.000	3.19	0.9380
12.2550	41.237	130.987	0.000	0.000	3.18	0.9410
12.5400	41.453	131.209	0.000	0.000	3.17	0.9439
12.8250	41.668	131.397	0.000	0.000	3.15	0.9468
13.1100	41.883	131.670	0.000	0.000	3.14	0.9490
13.3950	42.021	131.742	0.000	0.000	3.14	0.9513
13.6800	42.180	131.691	0.000	0.000	3.13	0.9534
13.9650	42.359	132.030	0.000	0.000	3.12	0.9555
14.2500	42.527	132.134	0.000	0.000	3.11	0.9575

# AXIAL PRESSURE PROFILE

X (in)	MACH $\phi$	Ptotal (psia)	Pstatic (psia)	Ttotal (R)	Tstatic (R)	Wdot (Lbm/s)	Local Radius (in)
1.20	0.000	2141.73	2142.70	1002.26	1005.71	0.27	3.354
1.30	0.001	2141.73	2142.70	1002.26	1005.71	1.02	3.354
1.40	0.004	2141.71	2142.66	1002.07	1005.51	5.69	3.354
1.50	0.007	2141.66	2142.58	1002.26	1005.70	10.12	3.354
1.61	0.010	2141.61	2142.47	1002.26	1005.92	13.85	3.354
1.71	0.011	2141.58	2142.43	1002.26	1005.70	15.13	3.354
1.81	0.031	2140.55	2140.34	1002.19	1002.00	35.47	3.354
1.91	0.089	2132.16	2132.53	1002.49	1002.17	64.88	3.354
2.01	0.124	2125.41	2125.81	1002.35	1002.26	85.01	3.354
2.11	0.146	2118.46	2118.84	1002.41	1002.26	99.48	3.354
2.21	0.168	2110.52	2110.52	1002.41	1002.26	110.39	3.354
2.31	0.178	2105.36	2105.36	1002.41	1002.26	119.02	3.354
2.41	0.188	2100.84	2100.84	1002.41	1002.26	126.07	3.354
2.51	0.197	2096.79	2096.79	1002.41	1002.26	132.04	3.354
2.61	0.206	2093.29	2093.29	1002.41	1002.26	136.93	3.354
2.71	0.212	2090.24	2090.24	1002.41	1002.26	141.04	3.354
2.81	0.218	2087.55	2087.55	1002.41	1002.26	144.53	3.354
2.91	0.223	2085.07	2085.07	1002.41	1002.26	147.65	3.354
3.01	0.228	2082.66	2082.66	1002.41	1002.26	150.62	3.354
3.11	0.232	2080.56	2080.56	1002.41	1002.26	153.10	3.354
3.21	0.236	2078.77	2078.77	1002.41	1002.26	155.22	3.354
3.31	0.239	2077.11	2077.11	1002.41	1002.26	157.13	3.354
3.41	0.242	2075.59	2075.59	1002.41	1002.26	158.84	3.354
3.51	0.245	2074.33	2074.33	1002.41	1002.26	160.26	3.354
3.61	0.247	2073.10	2073.10	1002.41	1002.26	161.63	3.354
3.71	0.250	2071.92	2071.92	1002.41	1002.26	162.92	3.354
3.81	0.252	2070.67	2070.67	1002.41	1002.26	164.05	3.354
3.91	0.254	2069.63	2069.63	1002.41	1002.26	165.17	3.354
4.01	0.255	2068.64	2068.64	1002.41	1002.26	166.12	3.354
4.11	0.257	2067.78	2067.78	1002.41	1002.26	166.98	3.354
4.21	0.258	2067.06	2067.06	1002.41	1002.26	167.77	3.354
4.31	0.260	2066.58	2066.58	1002.41	1002.26	168.48	3.354
4.41	0.261	2066.00	2066.00	1002.41	1002.26	169.20	3.354
4.51	0.262	2065.37	2065.37	1002.41	1002.26	169.84	3.354
4.61	0.263	2064.80	2064.80	1002.41	1002.26	170.42	3.354
4.71	0.264	2064.24	2064.24	1002.41	1002.26	170.98	3.354
4.81	0.265	2063.71	2063.71	1002.41	1002.26	171.52	3.354
4.91	0.266	2063.18	2063.18	1002.41	1002.26	172.04	3.354
5.01	0.267	2062.65	2062.65	1002.41	1002.26	172.52	3.354
5.11	0.268	2062.11	2062.11	1002.41	1002.26	173.00	3.354
5.21	0.269	2061.58	2061.58	1002.41	1002.26	173.44	3.354
5.31	0.270	2061.05	2061.05	1002.41	1002.26	173.71	3.354
5.41	0.271	2060.52	2060.52	1002.41	1002.26	174.05	3.354
5.51	0.272	2060.00	2060.00	1002.41	1002.26	174.37	3.354
5.61	0.273	2059.47	2059.47	1002.41	1002.26	174.66	3.354
5.71	0.274	2058.94	2058.94	1002.41	1002.26	174.91	3.354
5.81	0.275	2058.41	2058.41	1002.41	1002.26	175.13	3.354
5.91	0.276	2057.88	2057.88	1002.41	1002.26	175.31	3.354
6.01	0.277	2057.35	2057.35	1002.41	1002.26	175.48	3.354
6.11	0.278	2056.82	2056.82	1002.41	1002.26	175.63	3.354
6.21	0.279	2056.29	2056.29	1002.41	1002.26	175.77	3.354
6.31	0.280	2055.76	2055.76	1002.41	1002.26	175.89	3.354
6.41	0.281	2055.23	2055.23	1002.41	1002.26	176.00	3.354
6.51	0.282	2054.70	2054.70	1002.41	1002.26	176.10	3.354
6.61	0.283	2054.17	2054.17	1002.41	1002.26	176.19	3.354
6.71	0.284	2053.64	2053.64	1002.41	1002.26	176.27	3.354
6.81	0.285	2053.11	2053.11	1002.41	1002.26	176.34	3.354
6.91	0.286	2052.58	2052.58	1002.41	1002.26	176.40	3.354
7.01	0.287	2052.05	2052.05	1002.41	1002.26	176.45	3.354
7.11	0.288	2051.52	2051.52	1002.41	1002.26	176.49	3.354
7.21	0.289	2050.99	2050.99	1002.41	1002.26	176.52	3.354
7.31	0.290	2050.46	2050.46	1002.41	1002.26	176.55	3.354
7.41	0.291	2049.93	2049.93	1002.41	1002.26	176.57	3.354
7.51	0.292	2049.40	2049.40	1002.41	1002.26	176.59	3.354
7.61	0.293	2048.87	2048.87	1002.41	1002.26	176.60	3.354
7.71	0.294	2048.34	2048.34	1002.41	1002.26	176.61	3.354
7.81	0.295	2047.81	2047.81	1002.41	1002.26	176.62	3.354
7.91	0.296	2047.28	2047.28	1002.41	1002.26	176.63	3.354
8.01	0.297	2046.75	2046.75	1002.41	1002.26	176.64	3.354
8.11	0.298	2046.22	2046.22	1002.41	1002.26	176.65	3.354
8.21	0.299	2045.69	2045.69	1002.41	1002.26	176.66	3.354
8.31	0.300	2045.16	2045.16	1002.41	1002.26	176.67	3.354
8.41	0.301	2044.63	2044.63	1002.41	1002.26	176.68	3.354
8.51	0.302	2044.10	2044.10	1002.41	1002.26	176.69	3.354
8.61	0.303	2043.57	2043.57	1002.41	1002.26	176.70	3.354
8.71	0.304	2043.04	2043.04	1002.41	1002.26	176.71	3.354
8.81	0.305	2042.51	2042.51	1002.41	1002.26	176.72	3.354
8.91	0.306	2041.98	2041.98	1002.41	1002.26	176.73	3.354
9.01	0.307	2041.45	2041.45	1002.41	1002.26	176.74	3.354
9.11	0.308	2040.92	2040.92	1002.41	1002.26	176.75	3.354
9.21	0.309	2040.39	2040.39	1002.41	1002.26	176.76	3.354
9.31	0.310	2039.86	2039.86	1002.41	1002.26	176.77	3.354
9.41	0.311	2039.33	2039.33	1002.41	1002.26	176.78	3.354
9.51	0.312	2038.80	2038.80	1002.41	1002.26	176.79	3.354
9.61	0.313	2038.27	2038.27	1002.41	1002.26	176.80	3.354
9.71	0.314	2037.74	2037.74	1002.41	1002.26	176.81	3.354
9.81	0.315	2037.21	2037.21	1002.41	1002.26	176.82	3.354
9.91	0.316	2036.68	2036.68	1002.41	1002.26	176.83	3.354
10.01	0.317	2036.15	2036.15	1002.41	1002.26	176.84	3.354
10.11	0.318	2035.62	2035.62	1002.41	1002.26	176.85	3.354
10.21	0.319	2035.09	2035.09	1002.41	1002.26	176.86	3.354
10.31	0.320	2034.56	2034.56	1002.41	1002.26	176.87	3.354
10.41	0.321	2034.03	2034.03	1002.41	1002.26	176.88	3.354
10.51	0.322	2033.50	2033.50	1002.41	1002.26	176.89	3.354
10.61	0.323	2032.97	2032.97	1002.41	1002.26	176.90	3.354
10.71	0.324	2032.44	2032.44	1002.41	1002.26	176.91	3.354
10.81	0.325	2031.91	2031.91	1002.41	1002.26	176.92	3.354
10.91	0.326	2031.38	2031.38	1002.41	1002.26	176.93	3.354
11.01	0.327	2030.85	2030.85	1002.41	1002.26	176.94	3.354
11.11	0.328	2030.32	2030.32	1002.41	1002.26	176.95	3.354
11.21	0.329	2029.79	2029.79	1002.41	1002.26	176.96	3.354
11.31	0.330	2029.26	2029.26	1002.41	1002.26	176.97	3.354
11.41	0.331	2028.73	2028.73	1002.41	1002.26	176.98	3.354
11.51	0.332	2028.20	2028.20	1002.41	1002.26	176.99	3.354
11.61	0.333	2027.67	2027.67	1002.41	1002.26	177.00	3.354
11.71	0.334	2027.14	2027.14	1002.41	1002.26	177.01	3.354
11.81	0.335	2026.61	2026.61	1002.41	1002.26	177.02	3.354
11.91	0.336	2026.08	2026.08	1002.41	1002.26	177.03	3.354
12.01	0.337	2025.55	2025.55	1002.41	1002.26	177.04	3.354
12.11	0.338	2025.02	2025.02	1002.41	1002.26	177.05	3.354
12.21	0.339	2024.49	2024.49	1002.41	1002.26	177.06	3.354
12.31	0.340	2023.96	2023.96	1002.41	1002.26	177.07	3.354
12.41	0.341	2023.43	2023.43	1002.41	1002.26	177.08	3.354
12.51	0.342	2022.90	2022.90	1002.41	1002.26	177.09	3.354
12.61	0.343	2022.37	2022.37	1002.41	1002.26	177.10	3.354
12.71	0.344	2021.84	2021.84	1002.41	1002.26	177.11	3.354
12.81	0.345	2021.31	2021.31	1002.41	1002.26	177.12	3.354
12.91	0.346	2020.78	2020.78	1002.41	1002.26	177.13	3.354
13.01	0.347	2020.25	2020.25	1002.41	1002.26	177.14	3.354
13.11	0.348	2019.72	2019.72	1002.41	1002.26	177.15	3.354
13.21	0.349	2019.19	2019.19	1002.41	1002.26	177.16	3.354
13.31	0.350	2018.66	2018.66	1002.41	1002.26	177.17	3.354
13.41	0.351	2018.13	2018.13	1002.41	1002.26	177.18	3.354
13.51	0.352	2017.60	2017.60	1002.41	1002.26	177.19	3.354
13.61	0.353	2017.07	2017.07	1002.41	1002.26	177.20	3.354
13.71	0.354	2016.54	2016.54	1002.41	1002.26	177.21	3.354
13.81	0.355	2016.01	2016.01	1002.41	1002.26	177.22	3.354
13.91	0.356	2015.48	2015.48	1002.41	1002.26	177.23	3.354
14.01	0.357	2014.95	2014.95	1002.41	1002.26	177.24	3.354
14.11	0.358	2014.42	2014.42	1002.41	1002.26	177.25	3.354
14.21	0.359	2013.89	2013.89	1002.41	1002.26	177.26	3.354
14.31							

# PERFORMANCE SUMMARY

## C\* EFFICIENCY CALCULATIONS (ODK)

INJECTED MR= 2.8800 CSTAR=6880.40 CORE Em=0.8930 BARRIER Em=1.0000  
 CORE: OVERALL MR= 2.8800 VAPOR MR= 3.1070 CSTAR-MIX=5788.35 MASS FRACTION= 1.0000  
 BARRIER: OVERALL MR= 0.0000 VAPOR MR=99.9000 CSTAR-MIX= 0.00 MASS FRACTION= 0.0000  
 ENGINE: OVERALL MR= 2.8800 VAPOR MR= 3.1070 CSTAR-DEL=5811.17  
 C\* EFFICIENCY = 9.675E-01

## ISP EFFICIENCY CALCULATIONS

ISP-ODK, INJ = 2.669E+02 SEC.  
 ISP-ODK, M.Z. INJ = 2.662E+02 SEC. ISP-ODK, M.Z. VAPOR = 2.828E+02 SEC.  
 VAPORIZATION EFFICIENCY = 9.638E-01 MIXING EFFICIENCY = 9.935E-01  
 ENERGY RELEASE EFFICIENCY = 9.676E-01

NOTE: ISP-DEL = ISP-ODK, INJ. \* ERE \* ETADIV \* DELISP-BL

# TIME-LAG CALCULATIONS, Milliseconds

Cohem, In.=9.090E-03      FUEL Cohem, In.=2.293E+02      OX Cohem, In.=8.194E+01

## ELEMENT 1 IS TYPE=LOL

FUEL:      Cinj, In.=1.587E-02      Lvap, In.= 0.570      ATOMIZATION LENGTH USED, In.= 1.147E+00  
              Timp=3.863E-02      Tatom=3.107E-01      Tvp=1.544E-01      Total=5.036E-01

OX:      Cinj, In.=1.620E-02      Lvap, In.= 0.206      ATOMIZATION LENGTH USED, In.= 2.514E+00  
              Timp=7.013E-02      Tatom=6.082E-01      Tvp=6.884E-02      Total=9.452E-01

## EFFECTIVE TIMELAGS, Milliseconds

FUEL:      Cinj, In.=1.587E-02      Lvap, In.= 0.570      Total=5.036E-01  
              Timp=3.863E-02      Tatom=3.107E-01      Tvp=1.544E-01

OX:      Cinj, In.=1.620E-02      Lvap, In.= 0.206      Total=9.452E-01  
              Timp=7.013E-02      Tatom=6.082E-01      Tvp=6.884E-02

# CHAMBER-NOZZLE OPTIMIZATION RESULTS

CHAMBER LENGTH (FEET)	ETA-C*	ETA-NOZ	OVERALL EFFICIENCY
0.0000	0.0000	0.8807	0.0000
0.1667	0.0154	0.8726	0.0134
0.3333	0.5983	0.8646	0.5173
0.5000	0.7682	0.8537	0.6728
0.6667	0.8686	0.8419	0.7315
0.8333	0.9116	0.8302	0.7568
1.0000	0.9383	0.8185	0.7680
1.1667	0.9556	0.8057	0.7701
1.3333	0.9694	0.7915	0.7673
1.5000	0.9830	0.7773	0.7641
1.6667	0.9966	0.7631	0.7606
1.8333	1.0000	0.7465	0.7465
2.0000	1.0000	0.7299	0.7299
2.1667	1.0000	0.7114	0.7114
2.3333	1.0000	0.6926	0.6926
2.5000	1.0000	0.6742	0.6742
2.6667	1.0000	0.6461	0.6461
2.8333	1.0000	0.6142	0.6142
3.0000	1.0000	0.5823	0.5823
3.1667	1.0000	0.5504	0.5504
3.3333	1.0000	0.5071	0.5071
3.5000	1.0000	0.4439	0.4439
3.6667	1.0000	0.3608	0.3608
3.8333	1.0000	0.3176	0.3176
4.0000	1.0000	0.2544	0.2544

OPTIMUM CHAMBER LENGTH= 1.1667 FT  
 MAXIMUM OVERALL EFFICIENCY= 0.7701

# REDESIGNED CHAMBER RESULTS

NOMINAL CHAMBER PRESSURE	=	2.141E+03	PSIA
THROTTLED CHAMBER PRESSURE	=	1.538E+03	PSIA
FUEL INJECTION PRESSURE DROP	=	5.101E+02	PSI
OX INJECTION PRESSURE DROP	=	5.091E+02	PSI
CHAMBER RADIUS	=	2.786E-01	FT
THROAT RADIUS	=	1.852E-01	FT
NOZZLE ENTRANCE RADIUS OF CURVATURE	=	1.110E-01	FT
THROAT ENTRANCE RADIUS OF CURVATURE	=	1.110E-01	FT
NOZZLE CONVERGENCE HALF-ANGLE	=	3.000E+01	DEG
INJECTOR-TO-THROAT CHAMBER LENGTH	=	1.190E+00	FT
BARREL SECTION LENGTH	=	9.573E-01	FT

## IMPINGING ELEMENT SIZING RESULTS

ELEMENT TYPE	=	LOL
NO. OF ELEMENTS	=	80
FUEL ORIFICE DIAMETER	=	7.116E-02 IN
OX ORIFICE DIAMETER	=	1.091E-01 IN

ROCKET COMBUSTOR INTERACTIVE DESIGN METHODOLOGY  
Version 23-FEB-91

I-135

XIH=	1.8894E-01		
XIA=	3.0000E+01		
XCANT=	1.6000E+01		
XFACET=	4.7298E-01		
EMUNI=	6.5000E-01		
SEND			
\$BAFFLE			
\$END			
\$BARRIER			
\$END			
\$FFC			
\$END			
\$BURN			
\$GAMMA=			
AO=	1.1482E+00		
GMW=	4.0658E+03		
GPR=	2.3218E+01		
GK=	5.9752E-01		
GMU=	6.4180E-05		
GK=	6.6324E-05		
RML=	7.6242E+01		
VJL=	3.0748E+02		
TJL=	5.3100E+02		
RHOL=	4.9895E+01		
CPL=	4.7321E-01		
PCRITL=	3.1500E+02		
TCRITL=	1.2180E+03		
TBOILL=	6.6200E+02		
XMWL=	1.7200E+02		
HVAPL=	1.2500E+02		
EN=	6.6793E-01		
TAUSEN=	1.4114E-04		
ISEN=	1		
\$END			
\$INJ			
\$MAND=	6.7076E+00		
\$XMAND=	6.7076E+00		
\$FMANL=	3.3535E+00		
\$XMANL=	3.3535E+00		
PCA=	2.1411E+03,	1.5514E+03,	7.6734E+02
FRA=	4.7547E-01,	3.4543E-01,	1.7206E-01
FCAPA=	7.6914E-04,	8.1791E-04,	8.6053E-04
NFE=	4		
FTLA=	5.0142E-04,	6.6166E-04,	1.3371E-03,
FINA=	3.8678E-05,	3.8928E-05,	3.9064E-05,
FFA=	1.0000E+00,	1.0000E+00,	1.0000E+00,
NXE=	4		
XRA=	4.7547E-01,	3.4543E-01,	1.7206E-01
XCAPA=	5.9193E-04,	6.1747E-04,	6.5815E-04
XTLA=	9.3238E-04,	1.2439E-03,	2.3399E-03,
XINA=	7.3025E-05,	7.3125E-05,	7.3382E-05,
XFA=	1.0000E+00,	1.0000E+00,	1.0000E+00,
XUOR=	4.7805E-02,	7.3412E-02,	
AUOR=	3.4570E-03,	8.1288E-03,	
RUOR=	3.3172E-02,	6.0881E-02,	
XOR=	4.7805E-02,	7.3412E-02,	
ACR=	3.4570E-03,	8.1288E-03,	
ROR=	3.3172E-02,	6.0881E-02,	
XDOR=	9.5781E-02,	1.4682E-01,	
ADOR=	3.4570E-03,	6.1288E-03,	
RDOR=	3.3172E-02,	5.0861E-02,	
WOOT=	4.6218E+01,	1.3311E+02,	3.3533E+01,
			0.6578E+01,
			1.6644E+01,
\$END			
\$CHAMBER			

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NCAV=      1,
XB=      2.6126E-01
ZE=      1.1675E-01
ZE=      4.9676E-01
MJB=      3
T=      3.3000E-02
WC=      2.0961E-02,
AC=      3.6609E-02,
ARATIO=    1.0000E+00,
IDCAV=      1,
DC1=      1.4000E-01,
CC1=      5*2.5000E+03
GAMC1=      5*1.2000E+00
CC2=      5*2.5000E+03
GAMC2=      5*1.2000E+00
ICAV=      1
$END
$FDORG
NZON=      5
FTER=      20*1.0000E+00
NCAV1=      1
ICTYP1=      4,
ZE1=      2.0961E-02,
AE1=      3.6000E+02,
APL1=      1.4000E-01,
WC1=      1.9166E-01,
ZLOW1=     -1.2316E-02,
ZUP1=      2.0417E-01,
$END
$MIX
EM(1)= 8.930E-01, 1.000E+00
$END
$DEBUGC
DEBUG=      F
$END
$SHIFIG
SHORT=      F
POC=      2.0000E-01, 2.0000E-01
$END
$D18T8DC
SHORT=      F
POC=      2.0000E-01, 2.0000E-01
PAMP=      2.0000E-01
MC=      11
LC=      8
MB=      11
LB=      8
IDMAX=      10
$END
$CRPG
NDPC=      16
NOTFO=      16
NOTLF=      1000
NPRINT=      50
NSUMS=      3500
PAMP=      2.0000E-01
NRAD=      3
NCING=      5
$END
$LEINJC
IDMEM=      2
PAMPCH=      2.0000E-01
NTINJ=      16
U99F=      2.0000E-02

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OSF= 2.000E-02
DSOF= 2.000E-02
PCDO= 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00
XCDO= 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00
$END
$COMBUSTC
FALVM= 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00
XALVM= 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00
FALTM= 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00
XALTM= 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00
FRMA= 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00
XRMA= 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00
EMULT= 1.000E+00, 1.000E+00, 1.000E+00, 1.000E+00
COMBXS= 5.000E-01
ACMULT= 1.000E+00
COMULT= 1.000E+00
ENMULT= 1.000E+00
TAUMULT= 1.000E+00
$END
$FDOROC
SHORT= F
EPSIL= 2.000E+01
ERROR= 1.000E-01
LTS= 5
WTS= 5
NTS= 10
ITMAX= 100
RELX= 6.600E-01
NEET= 3
NXFST= 3
NYFST= 3
NAFST= 3
MORE= F
$END

```

END OF INPUT ECHO

# STEADY STATE COMBUSTION ANALYSIS PROGRAM

## RUN DESCRIPTORS

ROCCID POINT DESIGN TEST CASE 1  
LOX/RP-1 LIKE DOUBLET PAIR WITH FIXED PC  
APPROXIMATES .0100 SUBSCALE DOUBLET

## PROPELLANT DESCRIPTION

FUEL=RP-1 Tman., F= 71.00  
OX=LOX Tman., F=-278.00

## CHAMBER GEOMETRY

CHAMBER RADIUS = 3.3530 IN.  
CYLINDRICAL SECTION =11.6070 IN.  
NOZZLE ENTRANCE RADIUS OF CURVATURE = 1.3320 IN.  
CONVERGENCE HALF-ANGLE =30.0000 DEG.  
THROAT RADIUS = 2.2228 IN.  
CONVERGENT SECTION LENGTH = 2.8730 IN.  
THROAT ENTRANCE RADIUS OF CURVATURE = 1.3320 IN.  
CONTRACTION RATIO = 2.28

## INJECTOR DATA

INJECTOR CORE CONTAINS 60 LOL ELEMENTS

FUEL SIDE: Orifice Diam. =7.115E-02 In.  
Impingement Half-angle =30.00 Deg.  
OX SIDE: Orifice Diam. =1.091E-01 In.  
Impingement Half-angle =30.00 Deg.  
Cd =0.9100  
Unlike Cant Angle =16.00 Deg.  
Cd =0.9400  
Unlike Cant Angle =16.00 Deg.  
Impingement Height =0.123 In.  
Faceplate Thickness = 0.3081 In.  
Impingement Height =0.189 In.  
Faceplate Thickness = 0.4724 In.

## MIXING EFFICIENCIES

CORE MIXING EFFICIENCY=0.8930 BARRIER MIXING EFFICIENCY=1.0000

## COMBUST CONTROL PARAMETERS

MULTIPLIERS:  
FUEL ATOMIZATION LENGTH FOR VAPORIZATION: 1.000 CORE 1.000 BARRIER 1.000 FFC 1.000  
OX ATOMIZATION LENGTH FOR VAPORIZATION: 1.000 1.000 1.000 1.000  
FUEL ATOMIZATION LENGTH FOR TIMELAGS: 1.000 1.000 1.000 1.000  
OX ATOMIZATION LENGTH FOR TIMELAGS: 1.000 1.000 1.000 1.000  
FUEL DROPSIZE: 1.000 1.000 1.000 1.000  
OX DROPSIZE: 1.000 1.000 1.000 1.000  
MIXING (Em): 1.000 1.000 1.000 1.000  
AO-Multiplier=1.000 CC-Multiplier=1.000 N-Multiplier=1.000 Tau-Multiplier=1.000  
Eta-C\* for XB=0.500

BEGIN STEADY STATE COMBUSTION ANALYSIS  
PC=2141.10 PSIA

PROPELLANT PROPERTIES

FUEL=RP-1	Phase=Liquid	Tman., F= 71.00	
	Injected Density= 49.89 Lbm/Cu. Ft	Viscosity=1.380E-03 Lbm/Ft-S	Surface Tension=1.867E-03 Lb/Ft
OX=LOX	Phase=Liquid	Tman., F=-279.00	
	Injected Density= 70.20 Lbm/Cu. Ft	Viscosity=1.187E-04 Lbm/Ft-S	Surface Tension=7.326E-04 Lb/Ft

OPERATING CONDITIONS

PC FACE=2141.10 PSIA PC THROAT=2092.82 MIXTURE RATIO= 2.880  
FUEL INJECTION PRESSURE DROP= 609.96 Psia FUEL INJECTION VELOCITY= 307.74 Ft/S  
OX INJECTION PRESSURE DROP= 607.62 Psia OX INJECTION VELOCITY= 268.90 Ft/S

FUEL FLOWRATE= 46.262 OX FLOWRATE= 133.206

ATOMIZATION OUTPUT

DROPSIZE MODEL=AEROJET

ELEMENT TYPE 1 IS LOL			
FUEL:	ATOMIZATION LENGTH, In.=1.14748	ATOMIZATION LENGTH FOR VAPORIZATION, In.=1.14748	DROPLET RADIUS, Microns= 81.12
OX:	ATOMIZATION LENGTH, In.=2.51385	ATOMIZATION LENGTH FOR VAPORIZATION, In.=2.51385	DROPLET RADIUS, Microns= 88.92

VAPORIZATION CALCULATIONS

X (In.)	CORE=LOL		BAFFLE=		BARRIER=		FFC=	
	%FUEL VAP	%OX VAP	%FUEL VAP	%OX VAP	%FUEL VAP	%OX VAP	%FUEL VAP	%OX VAP
0.0000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
0.2856	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
0.5712	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
0.8568	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.1424	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.4280	6.496	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.7136	16.369	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.9992	24.874	0.000	0.000	0.000	0.000	0.000	0.000	0.000
2.2848	30.781	0.000	0.000	0.000	0.000	0.000	0.000	0.000
2.5704	36.561	0.000	0.000	0.000	0.000	0.000	0.000	0.000
2.8560	41.940	16.850	0.000	0.000	0.000	0.000	0.000	0.000
3.1416	45.713	34.742	0.000	0.000	0.000	0.000	0.000	0.000
3.4272	49.485	47.258	0.000	0.000	0.000	0.000	0.000	0.000
3.7128	52.998	56.256	0.000	0.000	0.000	0.000	0.000	0.000
3.9984	55.822	63.085	0.000	0.000	0.000	0.000	0.000	0.000
4.2840	58.662	68.351	0.000	0.000	0.000	0.000	0.000	0.000
4.5696	61.136	72.627	0.000	0.000	0.000	0.000	0.000	0.000
4.8552	63.316	76.257	0.000	0.000	0.000	0.000	0.000	0.000
5.1408	65.492	79.144	0.000	0.000	0.000	0.000	0.000	0.000
5.4264	67.279	81.583	0.000	0.000	0.000	0.000	0.000	0.000
5.7120	68.948	83.630	0.000	0.000	0.000	0.000	0.000	0.000
5.9976	70.614	85.322	0.000	0.000	0.000	0.000	0.000	0.000
6.2832	72.065	87.015	0.000	0.000	0.000	0.000	0.000	0.000
6.5688	73.516	88.446	0.000	0.000	0.000	0.000	0.000	0.000
6.8544	74.903	89.591	0.000	0.000	0.000	0.000	0.000	0.000
7.1400	76.100	90.919	0.000	0.000	0.000	0.000	0.000	0.000
7.4256	77.297	91.515	0.000	0.000	0.000	0.000	0.000	0.000
7.7112	78.350	92.304	0.000	0.000	0.000	0.000	0.000	0.000
7.9968	79.257	92.968	0.000	0.000	0.000	0.000	0.000	0.000
8.2824	80.164	93.632	0.000	0.000	0.000	0.000	0.000	0.000
8.5680	81.044	94.221	0.000	0.000	0.000	0.000	0.000	0.000
8.8536	81.914	94.719	0.000	0.000	0.000	0.000	0.000	0.000
9.1392	82.785	95.217	0.000	0.000	0.000	0.000	0.000	0.000
9.4248	83.406	95.672	0.000	0.000	0.000	0.000	0.000	0.000
9.7104	84.023	96.070	0.000	0.000	0.000	0.000	0.000	0.000
9.9960	84.639	96.426	0.000	0.000	0.000	0.000	0.000	0.000
10.2816	85.256	96.726	0.000	0.000	0.000	0.000	0.000	0.000
10.5672	85.872	97.023	0.000	0.000	0.000	0.000	0.000	0.000
10.8528	86.489	97.261	0.000	0.000	0.000	0.000	0.000	0.000
11.1384	87.106	97.480	0.000	0.000	0.000	0.000	0.000	0.000
11.4240	87.722	97.679	0.000	0.000	0.000	0.000	0.000	0.000
11.7096	88.215	97.867	0.000	0.000	0.000	0.000	0.000	0.000
11.9952	88.681	98.110	0.000	0.000	0.000	0.000	0.000	0.000
12.2808	89.148	98.299	0.000	0.000	0.000	0.000	0.000	0.000
12.5664	89.615	98.466	0.000	0.000	0.000	0.000	0.000	0.000
12.8520	90.081	98.604	0.000	0.000	0.000	0.000	0.000	0.000
13.1376	90.472	98.734	0.000	0.000	0.000	0.000	0.000	0.000
13.4232	90.897	98.862	0.000	0.000	0.000	0.000	0.000	0.000
13.7088	91.203	98.973	0.000	0.000	0.000	0.000	0.000	0.000
13.9944	91.568	99.074	0.000	0.000	0.000	0.000	0.000	0.000
14.2800	91.933	99.162	0.000	0.000	0.000	0.000	0.000	0.000

OVERALL VAPORIZATION EFFICIENCIES  
FUEL= 91.98% OX= 99.15%

MASS DISTRIBUTION PROFILE					
X (IN)	CORE (lbm/s)		BARRIER (lbm/s)		LOCAL VAPOR MIXTURE RATIO
	FUEL	OX	FUEL	OX	
0.0000	0.000	0.000	0.000	0.000	0.0000
0.2056	0.000	0.000	0.000	0.000	0.0000
0.5712	0.000	0.000	0.000	0.000	0.0000
0.6568	0.000	0.000	0.000	0.000	0.0000
1.1424	0.000	0.000	0.000	0.000	0.0000
1.4280	2.542	0.000	0.000	0.000	0.0035
1.7136	7.568	0.000	0.000	0.000	0.0103
1.9992	11.274	0.000	0.000	0.000	0.0164
2.2848	14.237	0.000	0.000	0.000	0.0194
2.5704	16.910	0.000	0.000	0.000	0.0231
2.8560	19.398	0.000	0.000	0.000	0.0281
3.1416	21.143	20.847	0.000	0.000	1.07
3.4272	22.868	48.278	0.000	0.000	0.3764
3.7128	24.610	82.852	0.000	0.000	2.75
3.9984	26.619	74.940	0.000	0.000	3.08
4.2840	28.278	84.046	0.000	0.000	0.5467
4.5696	27.126	91.048	0.000	0.000	0.5983
4.8552	29.285	96.743	0.000	0.000	0.6404
5.1408	30.292	101.579	0.000	0.000	0.6754
5.4264	31.116	106.424	0.000	0.000	3.47
5.7120	31.890	111.401	0.000	0.000	0.7059
5.9976	32.661	113.655	0.000	0.000	0.7311
6.2832	33.332	115.809	0.000	0.000	0.7526
6.5688	34.003	117.818	0.000	0.000	0.7714
6.8544	34.644	119.340	0.000	0.000	0.7862
7.1400	35.198	120.710	0.000	0.000	0.8040
7.4256	36.751	121.904	0.000	0.000	0.8164
7.7112	38.239	122.955	0.000	0.000	0.8309
7.9968	36.656	123.839	0.000	0.000	0.8419
8.2824	37.077	124.729	0.000	0.000	0.8622
8.5680	37.484	125.609	0.000	0.000	0.8611
8.8536	37.887	126.172	0.000	0.000	0.8866
9.1392	38.290	126.635	0.000	0.000	0.8765
9.4248	38.577	127.441	0.000	0.000	0.8938
9.7104	38.862	127.971	0.000	0.000	0.8802
9.9960	39.147	128.445	0.000	0.000	0.9020
10.2816	39.433	128.643	0.000	0.000	0.9089
10.5672	39.716	129.241	0.000	0.000	0.9116
10.8528	40.003	129.650	0.000	0.000	0.9159
11.1384	40.288	129.855	0.000	0.000	0.9202
11.4240	40.574	130.115	0.000	0.000	0.9243
11.7096	40.801	130.392	0.000	0.000	0.9280
11.9952	41.017	130.688	0.000	0.000	3.21
12.2808	41.233	130.941	0.000	0.000	0.9350
12.5664	41.449	131.163	0.000	0.000	0.9383
12.8520	41.665	131.346	0.000	0.000	0.9413
13.1376	41.845	131.519	0.000	0.000	0.9442
13.4232	42.014	131.690	0.000	0.000	0.9469
13.7088	42.183	131.838	0.000	0.000	0.9483
13.9944	42.352	131.973	0.000	0.000	0.9515
14.2800	42.521	132.077	0.000	0.000	0.9526
					0.9557
					0.9577

# AXIAL PRESSURE PROFILE

X (in)	MACH $\phi$	Ptotal (psia)	Pstatic (psia)	Ttotal (R)	Tstatic (R)	Wdot (Lbm/s)	Local Radius (in)
1.21	0.000	2141.71	2142.69	1692.26	1686.71	0.26	3.354
1.31	0.001	2141.71	2142.69	1692.26	1686.71	1.06	3.354
1.81	0.004	2141.69	2142.65	1693.22	1686.68	5.74	3.354
1.91	0.007	2141.65	2142.56	1692.26	1686.71	10.16	3.354
2.21	0.010	2141.60	2142.45	1704.47	1697.87	13.91	3.354
2.51	0.011	2141.58	2142.42	1692.26	1686.71	15.10	3.354
2.82	0.032	2140.45	2140.17	2665.03	2654.61	36.18	3.354
3.12	0.090	2131.91	2132.04	6222.30	6195.78	65.43	3.354
3.42	0.126	2123.20	2105.51	6649.21	6616.06	65.41	3.354
3.72	0.147	2116.27	2091.50	6663.38	6630.74	99.78	3.354
4.02	0.163	2110.34	2079.49	6655.10	6621.29	110.64	3.354
4.32	0.177	2105.19	2069.04	6646.78	6611.93	119.22	3.354
4.63	0.188	2100.68	2059.85	6636.05	6602.32	126.25	3.354
4.93	0.197	2096.64	2051.61	6631.69	6595.55	132.18	3.354
5.23	0.206	2093.16	2044.46	6630.43	6593.15	137.05	3.354
5.53	0.212	2090.11	2036.26	6626.75	6590.82	141.14	3.354
5.83	0.218	2087.44	2032.77	6629.64	6591.32	144.60	3.354
6.13	0.223	2084.94	2027.66	6631.26	6592.18	147.74	3.354
6.44	0.228	2082.55	2022.74	6632.10	6592.49	150.66	3.354
6.74	0.232	2080.49	2018.49	6635.08	6594.99	153.14	3.354
7.04	0.236	2078.68	2014.76	6637.74	6597.22	155.25	3.354
7.34	0.239	2077.02	2011.34	6640.65	6599.74	157.16	3.354
7.64	0.242	2075.52	2008.25	6643.26	6601.99	158.85	3.354
7.94	0.246	2074.26	2005.54	6644.91	6603.34	160.26	3.354
8.25	0.247	2073.02	2003.08	6646.41	6604.55	161.64	3.354
8.55	0.250	2071.85	2000.67	6648.04	6605.90	162.91	3.354
8.85	0.252	2070.81	1998.51	6649.93	6607.54	164.04	3.354
9.15	0.254	2069.76	1996.34	6651.78	6609.14	165.17	3.354
9.45	0.255	2068.69	1994.53	6652.72	6609.86	166.10	3.354
9.75	0.257	2068.08	1992.86	6653.87	6610.81	166.96	3.354
10.06	0.258	2067.39	1991.30	6655.15	6611.92	167.74	3.354
10.36	0.260	2066.64	1989.67	6656.60	6613.19	168.49	3.354
10.66	0.261	2065.95	1988.44	6658.04	6614.48	169.17	3.354
10.96	0.262	2065.39	1987.16	6659.64	6615.91	169.80	3.354
11.26	0.263	2064.76	1985.93	6661.44	6617.54	170.38	3.354
11.56	0.264	2064.21	1984.82	6662.76	6618.74	170.94	3.354
11.87	0.270	2063.68	1981.14	6663.04	6617.61	171.48	3.329
12.17	0.289	2063.13	1968.16	6663.09	6612.73	171.99	3.228
12.47	0.326	2062.62	1942.54	6663.07	6603.34	172.47	3.089
12.77	0.376	2061.90	1904.69	6663.11	6586.66	172.90	2.867
13.07	0.437	2060.94	1850.34	6663.14	6569.50	173.29	2.713
13.37	0.523	2059.82	1766.26	6663.16	6528.32	173.66	2.540
13.68	0.654	2058.26	1618.42	6663.19	6467.24	173.98	2.367
13.98	0.817	2055.91	1416.06	6663.22	6403.62	174.31	2.257
14.28	0.999	2052.62	1166.41	6663.25	6421.38	174.60	2.228

# PERFORMANCE SUMMARY

## C\* EFFICIENCY CALCULATIONS (CDK)

INJECTED MR= 2.8800 CSTAR=5860.40 CORE Em=0.8930 BARRIER Em=1.0000  
 CORE: OVERALL MR= 2.8800 VAPOR MR= 3.1082 CSTAR-MIX=5786.60 MASS FRACTION= 1.0000  
 BARRIER: OVERALL MR= 0.0000 VAPOR MR=99.9000 CSTAR-MIX= 0.00 MASS FRACTION= 0.0000  
 ENGINE: OVERALL MR= 2.8800 VAPOR MR= 3.1082 CSTAR-DEL=5612.36  
 C\* EFFICIENCY = 9.577E-01

## ISP EFFICIENCY CALCULATIONS

ISP-CDK, INJ = 2.889E+02 SEC.  
 ISP-CDK, M.Z. INJ = 2.862E+02 SEC.  
 VAPORIZATION EFFICIENCY = 9.640E-01  
 ENERGY RELEASE EFFICIENCY = 9.678E-01

NOTE: ISP-DEL = ISP-CDK, INJ. \* ERE \* ETADIV - DELISP-BL

TIME-LAG CALCULATIONS, Milliseconds

Cohem, In.=0.000E-03	FUEL Cohem, In.=2.293E+02	OX Cohem, In.=0.194E+01
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ELEMENT 1 IS TYPE-LOL

FUEL:	CinJ, In.=1.687E-02	Lvap, In.= 0.870	ATOMIZATION LENGTH USED, In.= 1.147E+00
	Temp=3.653E-02	Tatom=3.107E-01	Tvap=1.544E-01
			Total=5.036E-01
OX:	CinJ, In.=1.618E-02	Lvap, In.= 0.208	ATOMIZATION LENGTH USED, In.= 2.514E+00
	Temp=7.022E-02	Tatom=6.092E-01	Tvap=6.667E-02
			Total=9.463E-01

EFFECTIVE TIMELAGS, Milliseconds

FUEL:	CinJ, In.=1.587E-02	Lvap, In.= 0.870	
	Temp=3.653E-02	Tatom=3.107E-01	Tvap=1.544E-01
			Total=6.036E-01
OX:	CinJ, In.=1.618E-02	Lvap, In.= 0.208	
	Temp=7.022E-02	Tatom=6.092E-01	Tvap=6.667E-02
			Total=9.463E-01

# CHAMBER-NOZZLE OPTIMIZATION RESULTS

CHAMBER LENGTH (FEET)	ETA-C*	ETA-NOZ	OVERALL EFFICIENCY
0.0000	0.0000	0.8807	0.0000
0.1667	0.0154	0.8726	0.0184
0.3333	0.5986	0.8646	0.8176
0.5000	0.7683	0.8537	0.6730
0.6667	0.6889	0.8419	0.7316
0.8333	0.9116	0.8302	0.7569
1.0000	0.9383	0.8185	0.7680
1.1667	0.9556	0.8067	0.7701
1.3333	0.9694	0.7915	0.7673
1.5000	0.9830	0.7773	0.7641
1.6667	0.9866	0.7631	0.7605
1.8333	1.0000	0.7485	0.7485
2.0000	1.0000	0.7299	0.7299
2.1667	1.0000	0.7114	0.7114
2.3333	1.0000	0.6928	0.6928
2.5000	1.0000	0.6742	0.6742
2.6667	1.0000	0.6481	0.6481
2.8333	1.0000	0.6142	0.6142
3.0000	1.0000	0.5823	0.5823
3.1667	1.0000	0.5504	0.5504
3.3333	1.0000	0.5071	0.5071
3.5000	1.0000	0.4439	0.4439
3.6667	1.0000	0.3808	0.3808
3.8333	1.0000	0.3176	0.3176
4.0000	1.0000	0.2544	0.2544

OPTIMUM CHAMBER LENGTH= 1.1667 FT  
 MAXIMUM OVERALL EFFICIENCY= 0.7701

BEGIN STEADY STATE COMBUSTION ANALYSIS  
PC=1661.40 PSIA

PROPELLANT PROPERTIES			
FUEL=RP-1	Phase=Liquid	Tman., F= 71.00	
	Injected Density= 49.60 Lbm/Cu. Ft	Viscosity=1.360E-03 Lbm/Ft-S	Surface Tension=1.657E-03 Lbf/Ft
OX=LOX	Phase=Liquid	Tman., F=-279.00	
	Injected Density= 69.62 Lbm/Cu. Ft	Viscosity=1.143E-04 Lbm/Ft-S	Surface Tension=7.326E-04 Lbf/Ft

OPERATING CONDITIONS

PC FACE=1661.40 PSIA    PC THROAT=1497.36    MIXTURE RATIO= 2.660  
 FUEL INJECTION PRESSURE DROP= 268.15 Psia    FUEL INJECTION VELOCITY= 223.16 Ft/S  
 OX INJECTION PRESSURE DROP= 270.43 Psia    OX INJECTION VELOCITY= 189.72 Ft/S  
 FUEL FLOWRATE= 33.666    OX FLOWRATE= 96.669

ATOMIZATION OUTPUT

DROPSIZE MODEL=AEROJET

ELEMENT TYPE 1 IS LOL			
FUEL:	ATOMIZATION LENGTH, In.=1.10721	ATOMIZATION LENGTH FOR VAPORIZATION, In.=1.10721	DROPLET RADIUS, Microns= 84.07
OX:	ATOMIZATION LENGTH, In.=2.43656	ATOMIZATION LENGTH FOR VAPORIZATION, In.=2.43656	DROPLET RADIUS, Microns= 92.76

# VAPORIZATION CALCULATIONS

X (in.)	CORE=LOL		BAFFLE=		BARRIER=		FFC=	
	%FUEL VAP	%OX VAP	%FUEL VAP	%OX VAP	%FUEL VAP	%OX VAP	%FUEL VAP	%OX VAP
0.0000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
0.2868	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
0.5712	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
0.8566	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.1424	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.4280	7.034	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.7136	17.267	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.9992	24.996	0.000	0.000	0.000	0.000	0.000	0.000	0.000
2.2848	31.116	0.000	0.000	0.000	0.000	0.000	0.000	0.000
2.5704	36.738	0.000	0.000	0.000	0.000	0.000	0.000	0.000
2.8560	41.978	21.297	0.000	0.000	0.000	0.000	0.000	0.000
3.1416	46.861	36.123	0.000	0.000	0.000	0.000	0.000	0.000
3.4272	49.347	49.950	0.000	0.000	0.000	0.000	0.000	0.000
3.7128	52.624	57.637	0.000	0.000	0.000	0.000	0.000	0.000
3.9984	55.566	64.009	0.000	0.000	0.000	0.000	0.000	0.000
4.2840	58.352	68.940	0.000	0.000	0.000	0.000	0.000	0.000
4.5696	60.866	73.041	0.000	0.000	0.000	0.000	0.000	0.000
4.8552	62.985	76.519	0.000	0.000	0.000	0.000	0.000	0.000
5.1408	65.111	79.281	0.000	0.000	0.000	0.000	0.000	0.000
5.4264	66.946	81.957	0.000	0.000	0.000	0.000	0.000	0.000
5.7120	68.576	83.641	0.000	0.000	0.000	0.000	0.000	0.000
5.9976	70.208	85.292	0.000	0.000	0.000	0.000	0.000	0.000
6.2832	71.677	86.943	0.000	0.000	0.000	0.000	0.000	0.000
6.5688	73.094	88.389	0.000	0.000	0.000	0.000	0.000	0.000
6.8544	74.512	89.486	0.000	0.000	0.000	0.000	0.000	0.000
7.1400	75.697	90.515	0.000	0.000	0.000	0.000	0.000	0.000
7.4256	76.666	91.369	0.000	0.000	0.000	0.000	0.000	0.000
7.7112	78.003	92.195	0.000	0.000	0.000	0.000	0.000	0.000
7.9968	78.666	92.842	0.000	0.000	0.000	0.000	0.000	0.000
8.2824	79.774	93.490	0.000	0.000	0.000	0.000	0.000	0.000
8.5680	80.650	94.103	0.000	0.000	0.000	0.000	0.000	0.000
8.8536	81.600	94.666	0.000	0.000	0.000	0.000	0.000	0.000
9.1392	82.351	95.074	0.000	0.000	0.000	0.000	0.000	0.000
9.4248	83.084	95.548	0.000	0.000	0.000	0.000	0.000	0.000
9.7104	83.667	95.926	0.000	0.000	0.000	0.000	0.000	0.000
9.9960	84.288	96.318	0.000	0.000	0.000	0.000	0.000	0.000
10.2816	84.891	96.810	0.000	0.000	0.000	0.000	0.000	0.000
10.5672	85.494	96.901	0.000	0.000	0.000	0.000	0.000	0.000
10.8528	86.086	97.192	0.000	0.000	0.000	0.000	0.000	0.000
11.1384	86.666	97.369	0.000	0.000	0.000	0.000	0.000	0.000
11.4240	87.301	97.563	0.000	0.000	0.000	0.000	0.000	0.000
11.7096	87.920	97.788	0.000	0.000	0.000	0.000	0.000	0.000
11.9952	88.376	98.003	0.000	0.000	0.000	0.000	0.000	0.000
12.2808	88.831	98.216	0.000	0.000	0.000	0.000	0.000	0.000
12.5664	89.287	98.378	0.000	0.000	0.000	0.000	0.000	0.000
12.8520	89.749	98.632	0.000	0.000	0.000	0.000	0.000	0.000
13.1376	90.199	98.889	0.000	0.000	0.000	0.000	0.000	0.000
13.4232	90.566	98.786	0.000	0.000	0.000	0.000	0.000	0.000
13.7088	90.913	98.903	0.000	0.000	0.000	0.000	0.000	0.000
13.9944	91.289	99.012	0.000	0.000	0.000	0.000	0.000	0.000
14.2800	91.626	99.099	0.000	0.000	0.000	0.000	0.000	0.000

OVERALL VAPORIZATION EFFICIENCIES  
 FUEL= 91.63% OX= 99.10%

# MASS DISTRIBUTION PROFILE

X (IN)	CORE (lbm/s)		BARRIER (lbm/s)		LOCAL VAPOR MIXTURE RATIO	ETA-C*
	FUEL	OX	FUEL	OX		
0.0000	0.000	0.000	0.000	0.000	0.00	0.0000
0.2056	0.000	0.000	0.000	0.000	0.00	0.0000
0.5712	0.000	0.000	0.000	0.000	0.00	0.0000
0.8568	0.000	0.000	0.000	0.000	0.00	0.0000
1.1424	0.000	0.000	0.000	0.000	0.00	0.0000
1.4280	2.361	0.000	0.000	0.000	0.00	0.0044
1.7136	5.803	0.000	0.000	0.000	0.00	0.0109
1.9992	8.390	0.000	0.000	0.000	0.00	0.0158
2.2848	10.444	0.000	0.000	0.000	0.00	0.0197
2.5704	12.331	0.000	0.000	0.000	0.00	0.0232
2.8560	14.090	20.587	0.000	0.000	1.48	0.2446
3.1416	15.327	36.853	0.000	0.000	2.40	0.4028
3.4272	16.664	47.706	0.000	0.000	2.88	0.4906
3.7128	17.731	55.717	0.000	0.000	3.14	0.5542
3.9984	18.658	61.877	0.000	0.000	3.32	0.6026
4.2840	19.586	66.644	0.000	0.000	3.40	0.6425
4.5696	20.428	70.608	0.000	0.000	3.46	0.6765
4.8552	21.141	73.970	0.000	0.000	3.50	0.7064
5.1408	21.855	76.641	0.000	0.000	3.51	0.7302
5.4264	22.472	78.937	0.000	0.000	3.51	0.7516
5.7120	23.019	80.955	0.000	0.000	3.51	0.7698
5.9976	23.586	82.451	0.000	0.000	3.50	0.7862
6.2832	24.069	84.047	0.000	0.000	3.49	0.8019
6.5688	24.535	85.426	0.000	0.000	3.46	0.8182
6.8544	25.010	86.505	0.000	0.000	3.46	0.8288
7.1400	25.408	87.500	0.000	0.000	3.44	0.8398
7.4256	25.601	88.315	0.000	0.000	3.42	0.8486
7.7112	26.182	89.124	0.000	0.000	3.40	0.8590
7.9968	26.479	89.750	0.000	0.000	3.38	0.8686
8.2824	26.777	90.376	0.000	0.000	3.36	0.8740
8.5680	27.071	90.968	0.000	0.000	3.36	0.8812
8.8536	27.358	91.436	0.000	0.000	3.34	0.8877
9.1392	27.642	91.807	0.000	0.000	3.32	0.8941
9.4248	27.886	92.365	0.000	0.000	3.31	0.8999
9.7104	28.080	92.741	0.000	0.000	3.30	0.9047
9.9960	28.292	93.110	0.000	0.000	3.29	0.9084
10.2816	28.484	93.392	0.000	0.000	3.28	0.9137
10.5672	28.667	93.673	0.000	0.000	3.26	0.9179
10.8528	28.899	93.955	0.000	0.000	3.25	0.9221
11.1384	29.101	94.146	0.000	0.000	3.24	0.9259
11.4240	29.303	94.333	0.000	0.000	3.22	0.9284
11.7096	29.511	94.529	0.000	0.000	3.20	0.9332
11.9952	29.684	94.739	0.000	0.000	3.19	0.9384
12.2808	29.817	94.944	0.000	0.000	3.18	0.9395
12.5664	29.970	95.101	0.000	0.000	3.17	0.9423
12.8520	30.123	95.250	0.000	0.000	3.16	0.9461
13.1376	30.276	95.373	0.000	0.000	3.15	0.9477
13.4232	30.396	95.485	0.000	0.000	3.14	0.9498
13.7088	30.516	95.609	0.000	0.000	3.13	0.9521
13.9944	30.635	95.714	0.000	0.000	3.12	0.9541
14.2800	30.755	95.799	0.000	0.000	3.11	0.9561

# AXIAL PRESSURE PROFILE

X (in)	MACH #	Ptotal (psia)	Pstatic (psia)	Ttotal (R)	Tstatic (R)	Wdot (Lbm/s)	Local Radius (in)
1.21	0.000	1551.57	1552.55	1577.15	1570.71	0.31	3.354
1.31	0.001	1551.57	1552.55	1577.15	1570.71	1.08	3.354
1.41	0.004	1551.55	1552.54	1581.33	1574.84	4.51	3.354
1.51	0.007	1551.52	1552.48	1577.15	1570.70	7.55	3.354
1.61	0.010	1551.75	1552.40	1729.52	1722.53	10.39	3.354
2.01	0.010	1551.75	1552.40	1577.15	1570.70	10.51	3.354
2.52	0.047	1549.95	1545.75	3545.95	3531.73	31.25	3.354
3.12	0.100	1543.17	1535.14	5492.51	5464.45	51.01	3.354
3.42	0.125	1537.52	1523.75	6771.12	6740.25	53.97	3.354
3.72	0.145	1532.75	1514.12	7559.45	7537.30	72.57	3.354
4.02	0.155	1525.59	1505.55	7559.15	7519.97	81.05	3.354
4.32	0.175	1525.14	1495.55	7455.51	7411.55	85.55	3.354
4.53	0.185	1521.55	1492.17	7355.70	7303.70	91.90	3.354
4.93	0.195	1515.13	1485.42	6734.53	6695.75	95.03	3.354
5.23	0.205	1515.57	1461.40	6733.92	6697.43	95.44	3.354
5.53	0.212	1514.50	1475.55	6733.20	6695.10	102.35	3.354
5.83	0.215	1512.52	1473.10	6734.05	6695.40	104.50	3.354
6.13	0.223	1510.55	1455.45	6735.43	6697.24	107.03	3.354
6.44	0.225	1505.15	1455.55	6735.15	6697.45	109.14	3.354
6.74	0.232	1507.57	1452.52	6735.55	6695.50	110.51	3.354
7.04	0.235	1505.37	1450.24	6740.53	6701.37	112.44	3.354
7.34	0.235	1505.20	1457.53	6743.05	6703.15	113.75	3.354
7.64	0.242	1504.05	1455.54	6745.72	6705.43	115.05	3.354
7.94	0.245	1503.15	1453.55	6747.50	6705.92	115.07	3.354
8.25	0.247	1502.31	1451.57	6745.55	6705.14	117.03	3.354
8.55	0.245	1501.45	1450.05	6750.54	6705.41	117.55	3.354
8.85	0.251	1500.71	1445.55	6752.41	6711.04	118.75	3.354
9.15	0.253	1495.57	1447.02	6754.25	6712.54	119.55	3.354
9.45	0.255	1495.25	1445.50	6755.50	6713.75	120.31	3.354
9.75	0.257	1495.71	1444.42	6755.55	6714.55	120.52	3.354
10.05	0.255	1495.15	1443.25	6757.55	6715.55	121.51	3.354
10.35	0.255	1497.57	1442.25	6755.30	6715.93	122.01	3.354
10.65	0.261	1497.17	1441.22	6750.57	6715.13	122.53	3.354
10.95	0.253	1495.70	1440.25	6752.15	6715.47	123.01	3.354
11.25	0.253	1495.31	1435.40	6753.55	6721.01	123.41	3.354
11.55	0.254	1495.55	1435.55	6755.54	6722.55	123.54	3.354
11.85	0.255	1495.45	1435.55	6755.75	6722.50	124.24	3.354
12.15	0.255	1495.05	1425.45	6757.05	6717.55	124.52	3.354
12.45	0.255	1494.53	1407.52	6757.43	6705.23	124.97	3.053
12.75	0.255	1494.05	1350.54	6757.50	6693.34	125.25	2.557
13.05	0.253	1493.42	1341.12	6755.21	6671.23	125.55	2.713
13.35	0.253	1492.50	1280.31	6755.51	6655.41	125.55	2.540
13.65	0.254	1491.45	1174.05	6755.50	6555.51	125.10	2.557
13.95	0.257	1485.55	1025.35	6755.10	6455.25	125.34	2.257
14.25	0.255	1487.55	555.53	6755.42	5355.55	125.55	2.223

# PERFORMANCE SUMMARY

## C\* EFFICIENCY CALCULATIONS (ODK)

INJECTED MR= 2.8800 CSTAR=5880.40 CORE Em=0.8930 BARRIER Em=1.0000  
 CORE: OVERALL MR= 2.8800 VAPOR MR= 3.1149 CSTAR-MIX=5786.22 MASS FRACTION= 1.0000  
 BARRIER: OVERALL MR= 0.0000 VAPOR MR=99.9000 CSTAR-MIX= 0.00 MASS FRACTION= 0.0000  
 ENGINE: OVERALL MR= 2.8800 VAPOR MR= 3.1149 CSTAR-DEL=5803.22  
 C\* EFFICIENCY = 9.561E-01

## ISP EFFICIENCY CALCULATIONS

ISP-ODK, INJ = 2.880E+02 SEC.  
 ISP-ODK, M.Z. INJ = 2.882E+02 SEC.  
 VAPORIZATION EFFICIENCY = 9.624E-01  
 ENERGY RELEASE EFFICIENCY = 9.602E-01  
 ISP-ODK, M.Z. VAPOR = 2.826E+02 SEC.  
 MIXING EFFICIENCY = 9.988E-01

NOTE: ISP-DEL = ISP-ODK, INJ. \* ERE \* ETADIV \* DELISP-BL

TIME-LAG CALCULATIONS, Milliseconds

Cohem, in.=1.125E-02      FUEL Cohem, in.=2.293E+02      OX Cohem, in.=8.194E+01

ELEMENT 1 IS TYPE=LOL  
 FUEL:      Cinj, in.=1.315E-02      Lvap, in.= 0.594      ATOMIZATION LENGTH USED, in.= 1.107E+00  
             Timp=5.314E-02      Tatom=4.195E-01      Tvap=2.179E-01      Total=6.945E-01  
 OX:      Cinj, in.=1.341E-02      Lvap, in.= 0.213      ATOMIZATION LENGTH USED, in.= 2.437E+00  
             Timp=9.595E-02      Tatom=1.070E+00      Tvap=9.354E-02      Total=1.280E+00

EFFECTIVE TIMELAGS, Milliseconds

FUEL:      Cinj, in.=1.315E-02      Lvap, in.= 0.594      Total=6.945E-01  
             Timp=5.314E-02      Tatom=4.195E-01      Tvap=2.179E-01  
 OX:      Cinj, in.=1.341E-02      Lvap, in.= 0.213      Total=1.280E+00  
             Timp=9.595E-02      Tatom=1.070E+00      Tvap=9.354E-02

BEGIN STEADY STATE COMBUSTION ANALYSIS  
PC= 767.34 PSIA

PROPELLANT PROPERTIES

FUEL=RP-1	Phase=Liquid	T <sub>man.</sub> , F= 71.00	
	Injected Density= 49.69 Lbm/Cu. Ft	Viscosity=1.360E-03 Lbm/Ft-S	Surface Tension=1.857E-03 Lbf/Ft
OX=LOX	Phase=Liquid	T <sub>man.</sub> , F=-279.00	
	Injected Density= 68.77 Lbm/Cu. Ft	Viscosity=1.082E-04 Lbm/Ft-S	Surface Tension=7.326E-04 Lbf/Ft

OPERATING CONDITIONS

PC FACE= 767.34 PSIA PC THROAT= 735.69 MIXTURE RATIO= 2.690  
FUEL INJECTION PRESSURE DROP= 66.11 Psia FUEL INJECTION VELOCITY= 110.80 Ft/S  
OX INJECTION PRESSURE DROP= 67.34 Psia OX INJECTION VELOCITY= 95.25 Ft/S  
FUEL FLOWRATE= 16.662 OX FLOWRATE= 47.988

ATOMIZATION OUTPUT

DROPSIZE MODEL=AEROJET

ELEMENT TYPE 1 IS LOL

FUEL:	ATOMIZATION LENGTH, in.=1.02434	ATOMIZATION LENGTH FOR VAPORIZATION, in.=1.02434	DROPLET RADIUS, Microns= 90.87
OX:	ATOMIZATION LENGTH, in.=2.26767	ATOMIZATION LENGTH FOR VAPORIZATION, in.=2.26767	DROPLET RADIUS, Microns= 99.68

# VAPORIZATION CALCULATIONS

X (in.)	CORE=LOL		BAFFLE=		BARRIER=		FFO=	
	%FUEL VAP	%OX VAP	%FUEL VAP	%OX VAP	%FUEL VAP	%OX VAP	%FUEL VAP	%OX VAP
0.0000	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
0.2558	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
0.5712	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
0.8566	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.1424	0.000	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.4280	8.927	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.7136	16.746	0.000	0.000	0.000	0.000	0.000	0.000	0.000
1.9992	26.046	0.000	0.000	0.000	0.000	0.000	0.000	0.000
2.2848	31.628	0.000	0.000	0.000	0.000	0.000	0.000	0.000
2.5704	36.931	11.188	0.000	0.000	0.000	0.000	0.000	0.000
2.8560	41.922	30.800	0.000	0.000	0.000	0.000	0.000	0.000
3.1416	45.419	43.985	0.000	0.000	0.000	0.000	0.000	0.000
3.4272	48.917	53.237	0.000	0.000	0.000	0.000	0.000	0.000
3.7128	52.361	60.329	0.000	0.000	0.000	0.000	0.000	0.000
3.9984	54.984	65.660	0.000	0.000	0.000	0.000	0.000	0.000
4.2840	57.607	70.133	0.000	0.000	0.000	0.000	0.000	0.000
4.5696	60.177	73.682	0.000	0.000	0.000	0.000	0.000	0.000
4.8552	62.195	77.050	0.000	0.000	0.000	0.000	0.000	0.000
5.1408	64.212	79.661	0.000	0.000	0.000	0.000	0.000	0.000
5.4264	66.176	81.807	0.000	0.000	0.000	0.000	0.000	0.000
5.7120	67.723	83.664	0.000	0.000	0.000	0.000	0.000	0.000
5.9976	69.270	85.231	0.000	0.000	0.000	0.000	0.000	0.000
6.2832	70.768	86.788	0.000	0.000	0.000	0.000	0.000	0.000
6.5688	72.134	88.215	0.000	0.000	0.000	0.000	0.000	0.000
6.8544	73.479	89.275	0.000	0.000	0.000	0.000	0.000	0.000
7.1400	74.785	90.308	0.000	0.000	0.000	0.000	0.000	0.000
7.4256	75.855	91.135	0.000	0.000	0.000	0.000	0.000	0.000
7.7112	77.005	91.865	0.000	0.000	0.000	0.000	0.000	0.000
7.9968	78.063	92.589	0.000	0.000	0.000	0.000	0.000	0.000
8.2824	79.003	93.203	0.000	0.000	0.000	0.000	0.000	0.000
8.5680	79.744	93.616	0.000	0.000	0.000	0.000	0.000	0.000
8.8536	80.577	94.324	0.000	0.000	0.000	0.000	0.000	0.000
9.1392	81.365	94.785	0.000	0.000	0.000	0.000	0.000	0.000
9.4248	82.192	95.248	0.000	0.000	0.000	0.000	0.000	0.000
9.7104	82.941	95.688	0.000	0.000	0.000	0.000	0.000	0.000
9.9960	83.512	96.034	0.000	0.000	0.000	0.000	0.000	0.000
10.2816	84.084	96.377	0.000	0.000	0.000	0.000	0.000	0.000
10.5672	84.656	96.654	0.000	0.000	0.000	0.000	0.000	0.000
10.8528	85.226	96.931	0.000	0.000	0.000	0.000	0.000	0.000
11.1384	85.788	97.205	0.000	0.000	0.000	0.000	0.000	0.000
11.4240	86.371	97.389	0.000	0.000	0.000	0.000	0.000	0.000
11.7096	86.967	97.581	0.000	0.000	0.000	0.000	0.000	0.000
11.9952	87.607	97.788	0.000	0.000	0.000	0.000	0.000	0.000
12.2808	88.134	97.984	0.000	0.000	0.000	0.000	0.000	0.000
12.5664	88.587	98.200	0.000	0.000	0.000	0.000	0.000	0.000
12.8520	88.999	98.365	0.000	0.000	0.000	0.000	0.000	0.000
13.1376	89.432	98.507	0.000	0.000	0.000	0.000	0.000	0.000
13.4232	89.865	98.626	0.000	0.000	0.000	0.000	0.000	0.000
13.7088	90.278	98.748	0.000	0.000	0.000	0.000	0.000	0.000
13.9944	90.615	98.866	0.000	0.000	0.000	0.000	0.000	0.000
14.2800	90.953	98.989	0.000	0.000	0.000	0.000	0.000	0.000

OVERALL VAPORIZATION EFFICIENCIES  
FUEL= 80.95% OX= 98.97%

# MASS DISTRIBUTION PROFILE

X (IN)	CORE (lmm/s)		BARRIER (lmm/s)		LOCAL VAPOR MIXTURE RATIO	ETA-C*
	FUEL	OX	FUEL	OX		
0.0000	0.000	0.000	0.000	0.000	0.00	0.0000
0.2856	0.000	0.000	0.000	0.000	0.00	0.0000
0.5712	0.000	0.000	0.000	0.000	0.00	0.0000
0.8568	0.000	0.000	0.000	0.000	0.00	0.0000
1.1424	0.000	0.000	0.000	0.000	0.00	0.0000
1.4280	1.654	0.000	0.000	0.000	0.00	0.0083
1.7136	3.124	0.000	0.000	0.000	0.00	0.0118
1.9992	4.340	0.000	0.000	0.000	0.00	0.0164
2.2848	5.270	0.000	0.000	0.000	0.00	0.0200
2.5704	6.183	5.368	0.000	0.000	0.87	0.1302
2.8560	6.986	14.780	0.000	0.000	2.12	0.3387
3.1416	7.588	21.112	0.000	0.000	2.78	0.4423
3.4272	8.150	28.548	0.000	0.000	3.13	0.5124
3.7128	8.724	36.950	0.000	0.000	3.32	0.5878
3.9984	9.161	31.604	0.000	0.000	3.46	0.6106
4.2840	9.598	33.854	0.000	0.000	3.51	0.6460
4.5696	10.027	36.483	0.000	0.000	3.54	0.6782
4.8552	10.383	36.873	0.000	0.000	3.57	0.7048
5.1408	10.699	36.178	0.000	0.000	3.57	0.7277
5.4264	11.026	36.266	0.000	0.000	3.58	0.7489
5.7120	11.284	40.147	0.000	0.000	3.59	0.7682
5.9976	11.542	40.699	0.000	0.000	3.64	0.7818
6.2832	11.796	41.661	0.000	0.000	3.63	0.7972
6.5688	12.019	42.331	0.000	0.000	3.62	0.8111
6.8544	12.243	42.840	0.000	0.000	3.60	0.8229
7.1400	12.461	43.394	0.000	0.000	3.48	0.8344
7.4256	12.646	43.732	0.000	0.000	3.48	0.8438
7.7112	12.830	44.131	0.000	0.000	3.44	0.8535
7.9968	13.007	44.430	0.000	0.000	3.42	0.8616
8.2824	13.147	44.725	0.000	0.000	3.40	0.8686
8.5680	13.287	45.020	0.000	0.000	3.39	0.8767
8.8536	13.426	45.263	0.000	0.000	3.37	0.8822
9.1392	13.560	45.454	0.000	0.000	3.35	0.8883
9.4248	13.695	45.705	0.000	0.000	3.34	0.8944
9.7104	13.820	45.906	0.000	0.000	3.32	0.9000
9.9960	13.915	46.053	0.000	0.000	3.31	0.9045
10.2816	14.010	46.246	0.000	0.000	3.30	0.9089
10.5672	14.105	46.381	0.000	0.000	3.29	0.9128
10.8528	14.201	46.513	0.000	0.000	3.28	0.9169
11.1384	14.296	46.645	0.000	0.000	3.26	0.9208
11.4240	14.391	46.733	0.000	0.000	3.25	0.9244
11.7096	14.490	46.828	0.000	0.000	3.23	0.9280
11.9952	14.587	46.925	0.000	0.000	3.21	0.9319
12.2808	14.685	47.024	0.000	0.000	3.20	0.9353
12.5664	14.767	47.123	0.000	0.000	3.19	0.9389
12.8520	14.829	47.197	0.000	0.000	3.18	0.9410
13.1376	14.901	47.270	0.000	0.000	3.17	0.9437
13.4232	14.973	47.326	0.000	0.000	3.16	0.9461
13.7088	15.042	47.385	0.000	0.000	3.15	0.9485
13.9944	15.098	47.442	0.000	0.000	3.14	0.9506
14.2800	15.155	47.491	0.000	0.000	3.13	0.9526

# AXIAL PRESSURE PROFILE

X (in)	MACH $\phi$	Ptotal (psia)	Pstatic (psia)	Ttotal (R)	Tstatic (R)	Wdot (Lbm/s)	Local Radius (in)
1.21	0.001	767.56	767.90	1629.16	1629.90	0.26	3.354
1.31	0.002	767.56	767.90	1629.23	1629.98	0.86	3.354
1.81	0.005	767.54	767.88	1629.08	1629.83	2.01	3.354
1.91	0.006	767.53	767.85	1629.11	1631.85	4.06	3.354
2.21	0.008	767.52	767.83	1629.08	1629.83	4.73	3.354
2.51	0.022	767.34	767.46	2187.25	2186.93	9.79	3.354
2.82	0.077	765.01	762.91	5744.05	5720.55	20.42	3.354
3.12	0.114	761.96	760.67	6675.67	6648.99	26.21	3.354
3.42	0.137	759.68	761.90	6673.80	6643.70	33.67	3.354
3.72	0.165	757.46	747.68	6659.48	6528.63	37.77	3.354
4.02	0.188	755.64	743.90	6644.49	6512.87	40.99	3.354
4.32	0.180	754.07	740.70	6635.91	6503.30	43.66	3.354
4.63	0.190	752.96	737.98	6631.48	6498.09	45.66	3.354
4.93	0.199	751.96	736.05	6627.25	6493.19	47.75	3.354
5.23	0.209	750.16	732.72	6626.10	6489.42	49.33	3.354
5.53	0.212	749.18	730.89	6626.84	6483.67	50.73	3.354
5.83	0.219	748.29	728.81	6629.82	6494.16	51.96	3.354
6.13	0.223	747.41	727.09	6632.30	6498.06	52.93	3.354
6.44	0.228	746.69	726.36	6633.67	6498.95	53.95	3.354
6.74	0.232	745.88	723.95	6636.32	6499.19	54.79	3.354
7.04	0.235	745.28	722.81	6639.92	6502.39	55.55	3.354
7.34	0.238	744.97	721.44	6643.00	6505.12	56.21	3.354
7.64	0.242	744.13	720.39	6645.88	6507.62	56.83	3.354
7.94	0.244	743.66	719.36	6649.45	6510.94	57.36	3.354
8.25	0.246	743.24	718.49	6651.68	6513.11	57.82	3.354
8.55	0.249	742.82	717.63	6653.43	6514.40	58.28	3.354
8.85	0.251	742.45	716.86	6655.20	6515.95	58.68	3.354
9.15	0.253	742.10	716.14	6657.01	6517.54	59.08	3.354
9.45	0.255	741.75	715.42	6658.78	6519.09	59.43	3.354
9.75	0.258	741.44	714.76	6660.36	6520.47	59.77	3.354
10.06	0.260	741.17	714.21	6661.39	6521.32	60.06	3.354
10.36	0.269	740.92	713.68	6662.69	6522.37	60.32	3.354
10.66	0.269	740.69	713.21	6663.97	6523.60	60.56	3.354
10.96	0.262	740.45	712.72	6665.28	6524.76	60.80	3.354
11.26	0.263	740.24	712.26	6666.73	6526.06	61.02	3.354
11.56	0.264	740.05	711.88	6668.40	6527.51	61.22	3.354
11.87	0.269	739.83	710.52	6670.15	6528.24	61.43	3.329
12.17	0.269	739.61	709.85	6671.74	6528.19	61.64	3.228
12.47	0.326	739.37	696.67	6672.55	6516.47	61.82	3.093
12.77	0.375	739.10	683.18	6673.07	6502.66	61.99	2.897
13.07	0.437	738.75	663.75	6673.63	6482.04	62.14	2.713
13.37	0.523	738.31	633.71	6674.22	6448.64	62.28	2.540
13.68	0.654	737.88	581.21	6674.80	6386.03	62.41	2.367
13.98	0.817	736.60	509.34	6675.23	6291.75	62.53	2.257
14.28	0.989	735.59	425.74	6675.67	6170.88	62.65	2.223

# PERFORMANCE SUMMARY

## C\* EFFICIENCY CALCULATIONS (OOK)

INJECTED MR= 2.8800 CSTAR=5880.40 CORE Em=0.8880 BARRIER Em=1.0000  
 CORE: OVERALL MR= 2.8800 VAPOR MR= 9.1396 CSTAR-MIX=8761.07 MASS FRACTION= 1.0000  
 BARRIER: OVERALL MR= 0.0000 VAPOR MR=98.0000 CSTAR-MIX= 0.00 MASS FRACTION= 0.0000  
 ENGINE: OVERALL MR= 2.8800 VAPOR MR= 9.1396 CSTAR-DEL=6562.64  
 C\* EFFICIENCY = 9.528E-01

## ISP EFFICIENCY CALCULATIONS

ISP-ODK, INJ = 2.068E+02 SEC.  
 ISP-ODK, M.Z. INJ = 2.852E+02 SEC.  
 VAPORIZATION EFFICIENCY = 9.597E-01  
 ENERGY RELEASE EFFICIENCY = 9.528E-01  
 ISP-ODK, M.Z. VAPOR = 2.824E+02 SEC.  
 MIXING EFFICIENCY = 9.936E-01

NOTE: ISP-DEL = ISP-ODK, INJ. \* EMS \* STADIV - DELISP-BL

TIME-LAG CALCULATIONS, Milliseconds

Cohem, In.=1.751E-02	FUEL Cohem, In.=2.299E+02	OX Cohem, In.=6.194E+01
ELEMENT 1 IS TYPE=LOL		
FUEL:	CinJ, In.=8.699E-03 Timp=1.070E-01	Lvap, In.= 0.616 Tatom=7.704E-01
		ATOMIZATION LENGTH USED, In.= 1.024E+00 Total=1.340E+00
OX:	CinJ, In.=8.876E-03 Timp=1.909E-01	Lvap, In.= 0.224 Tatom=1.994E+00
		ATOMIZATION LENGTH USED, In.= 2.266E+00 Total=2.371E+00
EFFECTIVE TIMELAGS, Milliseconds		
FUEL:	CinJ, In.=8.699E-03 Timp=1.070E-01	Lvap, In.= 0.616 Tatom=7.704E-01
		Total=1.340E+00
OX:	CinJ, In.=8.876E-03 Timp=1.909E-01	Lvap, In.= 0.224 Tatom=1.994E+00
		Total=2.371E+00

**APPENDIX J**  
**COMPONENT MODEL DOCUMENTATION**

**PART A**

**HIGH FREQUENCY ACOUSTIC CHAMBER RESPONSE MODEL  
(HIFI)**



THERMODYNAMIC ANALYSIS REPORT		NUMBER: 9980:1807
		DATE: 6 Feb. 1987
SUBJECT:  COMPUTER CODE FOR USE IN HIGH FREQUENCY COMBUSTION STABILITY ANALYSES	PAGE 1 OF	
	NO. OF ENCLOSURES	
	NO. OF APPENDICES	
ADDITIONAL INFORMATION AND WORK NOTES INCLUDED IN MICROFILM FILE CDN		

PREPARED FOR: J. L. Pieper

A computer code, HIFI has been developed for use in high frequency combustion stability analyses of rectangular or cylindrical cross-sectional chambers.

The code is capable of calculating the burning admittance and the  $n$ - $\gamma$  neutral stability curve. It is operational on the VAX computer system at ATC.

The attachment describes the theory, the computer code and the calculated results.

KEYWORDS: Misc (21), Chamber (52), Nozzle (53), Combustion Stability (105), LOX/HC (153), Model Development (209), Computer Program - New Develop. (210), 1986 (271), T. V. Nguyen (357)

DISTRIBUTION:  R. Hewitt, J. Hulka, J. Hyde, J. Ito, S. Mercer, J. Muss, K. Niiya, R. Schindler, R. Walker, 9980 File	PREPARED BY: <i>Thong Van Nguyen</i> THONG VAN NGUYEN
	REVIEWED BY: JAMES J. FANG <i>J. J. Fang</i>
	APPROVED BY: <i>M. F. Young</i> M. F. YOUNG, MANAGER
W.O. NO: KAE626	J-3

**COMPUTER CODE FOR USE  
IN  
HIGH FREQUENCY COMBUSTION STABILITY ANALYSES**

**by  
Thong Van Nguyen**

**Aerojet TechSystems Company  
Sacramento, CA 95813**

**February 5, 1987**

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- Figure 1.2 : Schematic diagram showing the relation between the pressure interaction index  $n$ , the insensitive time lag  $\tau_i$ , the sensitive time lag  $\tau_s$ , and the total time lag  $\tau_t$ .
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- Figure 4.4 : Comparisons between HIFI and IFAR predictions of burning admittance amplitude in a 1T mode.
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Table 3.1 : Description of variables, in namelist CNTRL.

Table 3.2 : Description of variables in namelist INPUT.

## I. INTRODUCTION

Aerojet TechSystems Company is currently conducting a program (contract F04611-85-C-0100) to formulate a procedure (Ref. 1) which can accurately characterize injector designs for large thrust (0.5 to 2.0 million pounds) high pressure (500 to 3000 psia) LOX/hydrocarbon engines. In this procedure, rectangular cross-sectional (hereafter will be referred to simply as rectangular) combustion chambers are to be used to simulate the lower transverse frequency modes of the large scale chamber. This requires the development of stability models for rectangular chambers.

As part of the development of models for use in combustion stability analyses of rectangular chambers, a computer code, High Frequency Intrinsic Stability Analysis (HIFI) has been developed to calculate the burning admittances and the  $n-\tau$  neutral stability curves. The code can be applied not only to rectangular chambers but also to cylindrical chambers.

### 1.1 High Frequency Intrinsic Combustion Stability

Combustion instability, characterized by organized pressure oscillations in rocket combustion chamber, can cause severe vibrations on various engine system components and payloads. In addition, combustion instabilities may cause excessive mechanical stresses and heat loads on the injector and combustion chamber walls.

Combustion instabilities have been generally classified

according to their frequency range: low, intermediate and high frequency. Significant efforts have been devoted to the understanding of high frequency instability because it is the most common in new engine developments and is the most destructive. High frequency instability results from the coupling between the combustion process and the acoustic waves in the chamber.

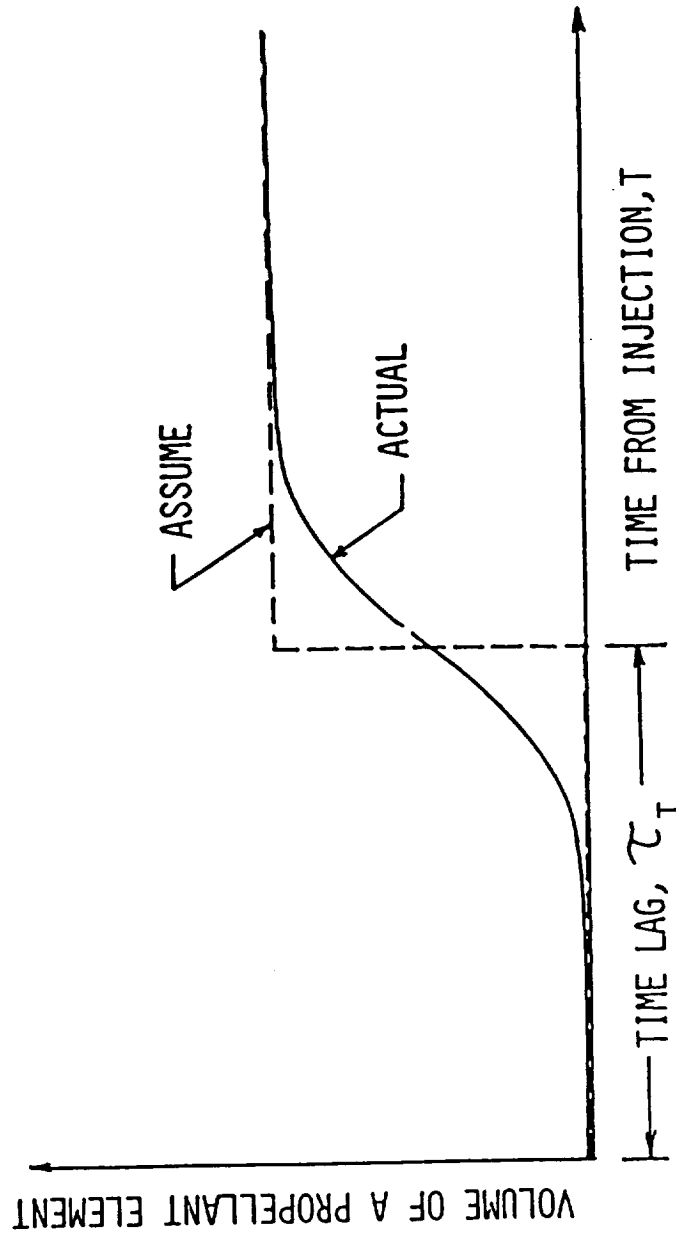
## 1.2 Concentrated Combustion and Sensitive Time-Lag Approach

Analytical models capable of characterizing combustion instability are obviously useful and valuable to engine designers during the development stage. Basic approaches in the modelling of high frequency combustion stability are described in reference 3. The concentrated combustion and sensitive time-lag approach developed by Crocco (Ref. 4) is discussed here since it is adopted in the present study. In this approach, the burning of propellant elements is assumed to occur instantaneously as shown in Figure 1.1. The time period between the instant of the injection and the burning of the propellant element is called the total time lag,  $\tau_r$ . All physical factors, e.g. pressure, temperature, that affect the burning process are assumed to correlate with the value of the local pressure. Consequently, the effects of these physical factors can be implicitly taken into accounts by relating the burning rate,  $\dot{m}_b$  to the instantaneous local pressure,  $p$ . The relation between  $\dot{m}_b$  and  $p$  is in the form:

$$\dot{m}_b \sim p^n$$

where  $n$  is called the (pressure) interaction index. The value

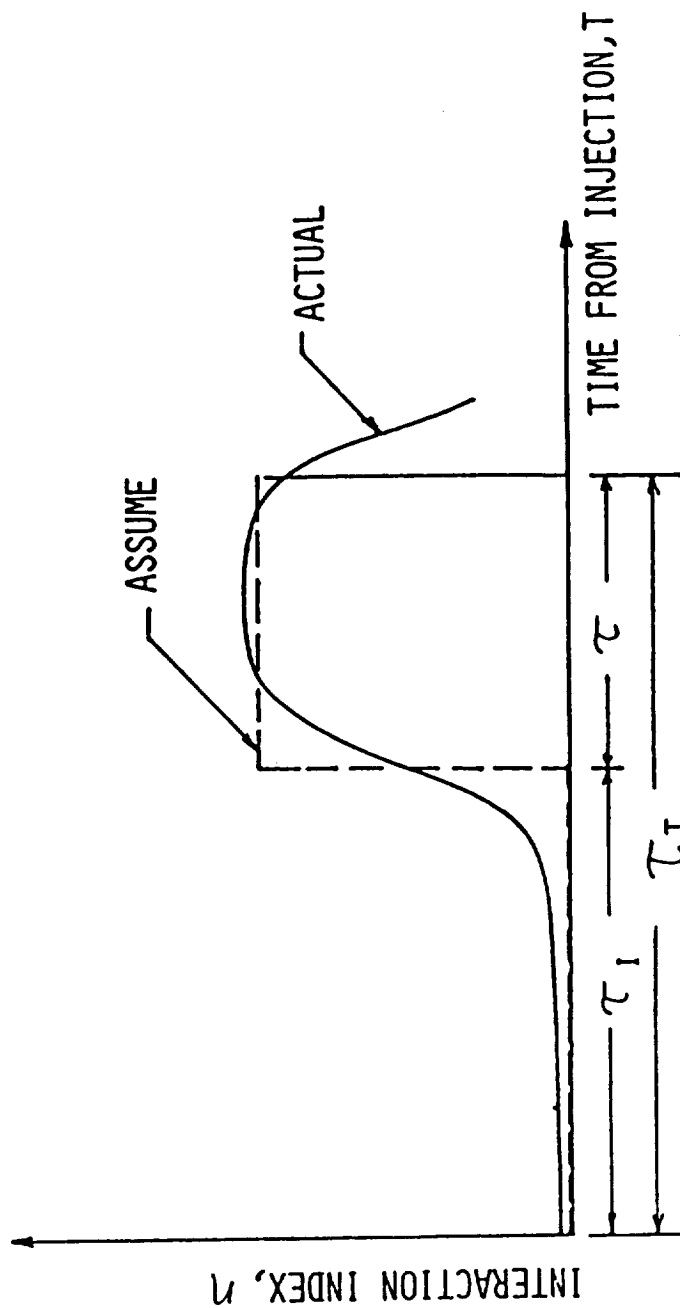
FIGURE 1.1 : DIAGRAM SHOWING COMBUSTION TIME LAG CONCEPT



of the interaction index is assumed to be zero during a time period called "insensitive" time lag,  $\tau_i$  and discontinuously becomes  $n$  during a time period called "sensitive" time lag,  $\tau_s$ . The sum of the insensitive time lag and the sensitive time lag equals to the total combustion time lag. Figure 1.2 is a schematic showing the relation between the interaction index, the insensitive time lag, the sensitive time lag and the total time lag. This sensitive time lag approach was first applied by Crocco (Ref. 4) in one-dimensional combustion stability analyses with both concentrated combustion and distributed combustion. As the names imply, the concentrated combustion approach assumes that the combustion concentrates in a plane at some distance from the injector face whereas in the distributed combustion approach the combustion distributes arbitrarily along the combustion chamber axis. In the concentrated-combustion approach, the combustion plane divides the chamber into two regions: the first region upstream of the combustion plane where the mean velocity is assumed to be zero and the second region downstream of the combustion plane where the velocity is non-zero and is assumed to be constant. This greatly simplifies the analysis since the equations which describe the flow dynamics in the two regions have no source terms. It is obvious that the concentrated combustion approach is not as realistic as the distributed combustion approach but it greatly simplifies the mathematical treatment of the analysis.

Crocco's original study was subsequently continued and

FIGURE 1.2 : SCHEMATIC SHOWING THE RELATION BETWEEN THE PRESSURE INTERACTION INDEX  $\eta$ , THE INSENSITIVE TIME LAG  $\tau_i$ , THE SENSITIVE TIME LAG  $\tau_s$  AND THE TOTAL TIME LAG  $\tau_t$



improved by several authors. Reference 3 describes subsequent studies following Crocco's original study. The reference also gives a brief history of the development of the sensitive time lag theory. The study of Crocco was first extended to tranverse modes by Scala (Ref. 5). Reardon then introduced the velocity interaction index to include the sensitivity of the burning rate to the tranverse components of the oscillating gas velocity (Ref. 6). This study also accounts for the effects of the non-uniform distribution of the propellant injection. The assumption of low Mach number in the chamber, which are used in all aforementioned models, was eliminated in the studies of longitudinal modes by Mitchell (Ref. 7) and by Harrje (Ref. 8.) and of tranverse modes by Smith (Ref. 2).

The concentrated-combustion and sensitive time-lag theory has been used extensively at Aerojet. Its evolution at Aerojet resulted into a computer code known as IFAR (Ref. 9). This computer code has been used in combustion stability analyses of virtually all liquid-propellant rocket engines developed in recent years at Aerojet. In general, the prediction capability of the code is satisfactory.

### 1.3 Objectives of the Present Study

The objective of the present work is to provide a computer code to predict the burning admittances and the  $n$  and  $\tau$  neutral stability curves for rectangular chambers. Although the objective is to provide a computer code for use in combustion stability analyses of rectangular chambers, the code

developed in the present study can also be used for cylindrical chambers. The code has many new features which are not available in IFAR. These features are:

- \* Values of nozzle admittances can be input in tabular forms as in IFAR or calculated internally by the program. If the nozzle admittances are calculated, the program will automatically generate<sup>is</sup> a table that ~~are~~ is to be used for any subsequent runs in which the values of the nozzle admittances do not change. The nozzle admittance values are calculated at the frequencies at which the burning admittance, its corresponding  $n$  and  $\tau$  are calculated, therefore no interpolation errors are introduced into the solutions.

- \* The user is not required to determine the Mach number a priori since it is calculated internally by the program given the specific heat ratio and the contraction ratio.

- \* Variables used in the theory and in the computer code are retained in the forms of complex variables. For users who wish to understand the theory or to make modifications to the code, this feature makes the theory described in the next section and the logic used in the code easy for them to follow.

- \* The code generates output files in the format which can be input to a computer graphic program, for example TELLEGRAF, to plot the calculated results.

\* The Mach number in the chamber is no longer assumed to be small.

#### 1.4 Approach

The present study follows the approach taken in the development of IFAR (Ref. 9). The difference between the present study and IFAR is in the calculation of chamber admittances. IFAR calculates the admittances by solving the pressure wave equation which has been derived on the assumption of low Mach number in the chamber. This assumption is also used to implement the boundary condition to the solutions of the equation and in the calculation of the burning admittance. The present study calculates chamber admittances by solving the wave equation for a velocity potential function (see Refs. 10 and 11). This has the advantage that the Mach number in the chamber is no longer assumed to be small, thus the code can be used for chambers having small contraction ratios.

## II. THEORY

The theory in the present study follows closely reference 9 to calculate the cavity admittance, the burning admittance and the  $n-\tau$  neutral stability curve. The nozzle-admittance model of reference 11 is extended to calculate the chamber admittances.

### 2.1 Theory Description

First, the continuity and momentum equations are written for an ideal gas. The thermodynamic variables, i.e. pressure density, etc., are decomposed into their mean and fluctuating components. These components are then normalized by the corresponding steady-state values. The mean components do not vary with time and are assumed to be uniform in the regions upstream and downstream of the combustion plane although they may be discontinuous at the combustion plane. The fluctuating components, however, vary in all directions and are functions of time. The velocity is also decomposed into a mean and a fluctuating component. The mean velocity is assumed to be only in the axial direction while the fluctuating component can vary in all directions. The fluctuating components of the velocity and the thermodynamic variables are assumed to be so small that the products of any two components can be neglected. As a result, equations for the fluctuating components are linear in time, thus their oscillations can be assumed to be sinusoidal. The flow is assumed to be irrotational and the fluctuating velocity components are defined to be the gradients of a

velocity potential function. The continuity and momentum equations and an isentropic relation are combined to yield a governing equation for the velocity potential function. The resulting equation is then written in <sup>the</sup> cartesian coordinate system for rectangular chambers and in <sup>the</sup> cylindrical coordinate system for cylindrical chambers.

Using <sup>the</sup> separation of variables technique, the partial differential equation governing the evolution of the potential function is separated into three second-degree ordinary differential equations. Using boundary conditions at the chamber walls two of the equations in the tranverse and lateral directions are solved explicitly to give the eigenvalues that correspond to tranverse and lateral resonance modes for rectangular chambers. Similarly for cylindrical chambers, boundary conditions at the chamber wall and at the axis of symmetry are used to calculate the eigenvalues that correspond to radial resonance modes. The eigenvalues that correspond to tangential resonance modes in cylindrical chambers are determined by requiring the solutions to the differential equation being single value functions. A general solution is obtained by solving the differential equation in the axial direction. The boundary condition at the injector face is then applied to calculate the chamber admittance upstream of the combustion plane. Using the nozzle admittance as the boundary equation, the chamber admittance downstream of the combustion plane is calculated. <sup>the</sup> Continuity condition is then applied at the combustion plane to relate the burning admittance to the

upstream and downstream chamber admittances. Finally, the pressure interaction index , n and the sensitive time lag,  $\tau$  for neutral stability condition is calculated using the expression derived by Crocco in reference 3 which relates n and  $\tau$  to the burning admittance.

## 2.2 Equation Derivations and Solutions

The continuity and momentum equations for an inviscid compressible gas are:

$$\frac{\partial \rho^*}{\partial t^*} + \nabla^* (\rho^* \vec{u}^*) = 0 , \quad (1)$$

and 
$$\rho^* \left( \frac{\partial \vec{u}^*}{\partial t^*} + \vec{u}^* \nabla^* \vec{u}^* \right) + \nabla^* p^* = 0 , \quad (2)$$

where t is the time,  $\rho$  is the gas density, u is the gas velocity and p is the gas pressure, the notation  $\vec{\phantom{x}}$  denotes vector quantities, and the supercript \* denotes demensional quantities. <sup>An</sup> Additional equation needed to close the above conservation equations is the following isentropic relation:

$$\frac{dp^*}{p^*} = \gamma \frac{d\rho^*}{\rho^*} \quad (3)$$

where  $\gamma$  is the gas specific heat ratio. Equations (1), (2) and (3) can be written in non-dimensional form as:

$$\frac{\partial \rho}{\partial t} + \nabla (\rho \vec{u}) = 0 , \quad (4)$$

$$\rho \left( \frac{\partial \vec{u}}{\partial t} + \vec{u} \nabla \vec{u} \right) + \frac{1}{\gamma} \nabla p = 0 , \quad (5)$$

and

$$dp = \gamma d\rho \quad (6)$$

In equations (4), (5), and (6) the density and the pressure are non-dimensionalized by their corresponding mean values; the velocity by the sound speed; length scales by some characteristic length, e.g. chamber radius or chamber half-width; and time by the sound speed and the characteristic length.

All non-dimensionalized dependent variables are then decomposed into the mean components which are time independent, and the perturbation components which are time dependent, i.e:

$$\vec{u} = \bar{\vec{u}} + \vec{u}', \quad p = 1 + p', \quad \rho = 1 + \rho', \quad (7)$$

where the bar and the superscript ' denote the mean and the perturbation components, respectively. It should be noted that the mean velocity  $\bar{\vec{u}}$  shown in the above equation has been non-dimensionalized by the sound speed and thus it is the same as the Mach number.

Assume there exists a velocity potential function,  $\phi$ , such that:

$$u' = \nabla \phi \quad (8)$$

and that the flow is irrotational, equations (4), (5) and (6) can be combined to yield the following relation:

$$\rho' = -\tau \left( \frac{\partial \phi}{\partial z} + (\bar{u} \cdot \nabla) \phi \right), \quad (9)$$

and the following equation governing the evolution of the velocity potential function:

$$\frac{\partial^2 \phi}{\partial t^2} - \nabla^2 \phi + \bar{u} \left[ \nabla (\bar{u} \cdot \nabla \phi) \right] + 2\bar{u} \frac{\partial \nabla \phi}{\partial t} = 0. \quad (10)$$

Assume that:

$$\phi = \phi e^{st} \quad (11)$$

where  $s$  is a complex quantity with its imaginary part representing the angular frequency of the oscillation and its real part representing the amplification coefficient of the oscillation. Equation (10) can then be written as:

$$s^2 \phi - \nabla^2 \phi + \bar{u} \left[ \nabla (\bar{u} \cdot \nabla \phi) \right] + 2\bar{u}s \nabla \phi = 0. \quad (12)$$

Assume the mean flow velocity exists only in the axial direction and its magnitude,  $\bar{u}$  is constant. Equation (12) can be written in <sup>the</sup> cylindrical coordinate system for cylindrical chambers and in <sup>the</sup> cartesian coordinate system for rectangular chambers. In each case, the separation of variables technique is used to separate the equation into three ordinary differential equations, of which the equation in the axial direction has the following form:

$$(1 - \bar{u}^2) \phi_x^{(n)} - 2\bar{u}s \phi_x^{(n)} - (s^2 + s_{mn}^2) \phi_x = 0, \quad (13)$$

where the superscripts (') and (") denote the first derivative and the second derivative of  $\phi_x$  with respect to  $x$ , the axial coordinate.  $\phi_x$  is the component of  $\phi$  that is dependent only on  $x$ . For rectangular chamber cases, the value of  $S_{mn}$  is:

$$S_{mn} = \frac{\pi}{2} \sqrt{m^2 + \frac{n^2}{b}} \quad (14)$$

where  $b$  is the ratio of the chamber thickness to the chamber width. The values of  $S_{mn}$  for cylindrical chamber cases are given in table 2.1 for selected values of  $m$  and  $n$ . The subscripts  $m$  and  $n$  correspond to the  $m^{\text{th}}$  tangential (transverse or width) and the  $n^{\text{th}}$  radial (lateral or thickness) resonance modes. In rectangular chamber cases, the values of  $S_{mn}$  are determined by applying appropriate boundary conditions at the chamber walls to the solutions of the differential equations in the transverse and the lateral directions. In cylindrical chamber cases, the values of  $S_{mn}$  are determined by requiring the solution to the differential equation in the circumferential direction being a single value function and by applying appropriate boundary conditions at the chamber wall and at the axis of symmetry to the solution of the differential equation in the radial direction.

The axial equation (Equation 12) is applicable to both regions upstream and downstream of the combustion plane in a cylindrical or a rectangular chamber. Applying the boundary condition at the injector face to the solution of the equation yields the following expression for the upstream chamber admittance:

m \ n					
	0	1	2	3	4
0	0.0000	3.8318	7.0155	10.1734	13.3238
1	1.8413	5.3313	8.5263	11.7059	14.8635
2	3.0543	6.7060	9.9695	13.1705	16.3476
3	4.2013	8.0151	11.3459	14.5858	17.7890
4	5.3175	9.2825	12.6820	15.9640	19.1961
5	6.4154	10.5199	13.9873	17.3127	20.5755
6	7.5012	11.7348	15.2681	18.6375	21.9318
7	8.5778	12.9324	16.5295	19.9419	23.2682
8	9.6475	14.1155	17.7739	21.2290	24.5874

Table 2.1: Selected values of  $S_{mn}$ .

*From [1] and [2].*

$$Y_I = \frac{u'_x}{p'} \Big|_I = \frac{\left(\frac{\alpha}{\gamma s}\right)(e^{\alpha x_p} - e^{-\alpha x_p}) - \psi_p \left(\frac{\alpha}{\gamma s}\right)(e^{\alpha x_p} + e^{-\alpha x_p})}{\psi_p (e^{\alpha x_p} - e^{-\alpha x_p}) - \left(\frac{\alpha}{\gamma s}\right)(e^{\alpha x_p} + e^{-\alpha x_p})}, \quad (15)$$

where the subscript I denotes the quantities evaluated at the location immediately upstream of the combustion plane,  $u'_x$  is the axial component of the local (non-dimensionalized) perturbation velocity, and  $x_p$  is the distance between the injector face and the combustion plane. Other quantities in the equation are defined as follows:

$$\alpha = \sqrt{S^2 + S_{mn}^2}, \quad (16)$$

and  $\psi_p$  is a quantity that is determined from the boundary conditions at the injector face. This quantity is described below.

For cylindrical chamber cases, suppose that the admittance, which is defined as the ratio of the local axial (non-dimensionalized) perturbation velocity to the local (non-dimensionalized) pressure perturbation, at the injector face can be expressed as:

$$Y \Big|_{x=0} = Y_r(r) Y_\theta(\theta), \quad (17)$$

where  $r$  and  $\theta$  are the radial and tangential coordinates, respectively. Then for a spinning  $m^{\text{th}}$  tangential mode, the expression for  $\psi_p$  can be written as:

$$\psi_p = \frac{\left(\int_0^1 Y_r(r) J_m^2(s_{mn}r) r dr\right) \left(\int_0^{2\pi} Y_\theta(\theta) d\theta\right)}{2\pi \int_0^1 J_m^2(s_{mn}r) r dr} \quad \text{J-23} \quad (18)$$

and for a standing  $m^{\text{th}}$  tangential mode, it can be written as:

$$\Psi_p = \frac{\left( \int_0^1 Y_r(r) J_m^2(s_{mn}r) r dr \right) \left( \int_0^{2\pi} Y_\theta(\theta) \cos^2(m\theta) d\theta \right)}{\pi \int_0^1 J_m^2(s_{mn}r) r dr} \quad (19)$$

In the above expressions for  $\Psi_p$ ,  $J_m$  is the  $m^{\text{th}}$  order Bessel function of the first kind.

For rectangular chamber cases, the expression for  $\Psi_p$  is written as:

$$\Psi_p = \int_{-1}^1 Y_y(y) \cos^2\left(\frac{m\pi}{2}[y+1]\right) dy \quad (20)$$

In equation (20), it has been assumed that the admittance at the injector face varies only in the  $y$  (width) direction and that oscillations in the  $z$  (thickness) direction do not exist. In the cases where oscillations in the  $z$  direction do exist, the value of  $\Psi_p$  will be different from that given by the above expression. In general cases, the difference and its effects on the overall solutions are believed to be small, thus the model is considered not to be limited to two-dimensional oscillations.

It should be noted that the admittances of acoustic cavities are included in  $Y_r(r)$ ,  $Y_\theta(\theta)$  and  $Y_y(y)$  in equations (18), (19) and (20). In the present study, the width of the cavities is assumed to be small compared to the radius or the width of the chamber. Furthermore, the cavities are assumed to locate at the circumference of the injector of the cylindrical chamber

or they are assumed to locate at the edge of the width of the rectangular chamber. These assumptions simplify the analysis since they allow the above expressions for  $\Psi_p$  being approximated analytically.

Applying the boundary condition at the nozzle entrance (nozzle admittance) to the solution of equation (13) yields the following expression for the downstream chamber admittance:

$$Y_{II} = \frac{u'_x}{p'} \bigg|_x = \frac{\alpha_1 e^{-\alpha_1 x_q} + \alpha_2 A e^{-\alpha_2 x_q}}{-\gamma (s e^{-\alpha_1 x_q} + s A e^{-\alpha_2 x_q} + \bar{u} \alpha_1 e^{-\alpha_1 x_q} + \bar{u} \alpha_2 A e^{-\alpha_2 x_q})}, \quad (21)$$

where the subscript II denotes the quantities evaluated at the location immediately downstream of the combustion plane, and

$$\alpha_1 = \frac{1}{1 - \bar{u}^2} \left( \bar{u} s - \sqrt{s^2 + s_{mn}^2 (1 - \bar{u}^2)} \right), \quad (22)$$

$$\alpha_2 = \frac{1}{1 - \bar{u}^2} \left( \bar{u} s + \sqrt{s^2 + s_{mn}^2 (1 - \bar{u}^2)} \right), \quad (23)$$

$$A = - \frac{\gamma s \Psi_q + \gamma \bar{u} \alpha_1 \Psi_q + \alpha_1}{\gamma s \Psi_q + \gamma \bar{u} \alpha_2 \Psi_q + \alpha_2}, \quad (24)$$

$\Psi_q$  is the nozzle admittance and  $x_q$  is the distance from the combustion plane to the nozzle entrance.

Continuity is then applied at the combustion plane to give the following expression for the burning admittance:

$$y = \frac{m'_b}{p'} = \frac{1}{\bar{u}} \left( Y_{II} - \frac{\bar{p}_I a_I}{\bar{p}_{II} a_{II}} Y_I \right) + \frac{1}{\gamma}, \quad (25)$$

where  $m'_b$  is the burning rate perturbation normalized by its mean value. In this expression, the non-dimensionalized pressure perturbations upstream and downstream of the combustion plane are assumed to be equal.

The concentrated-combustion analyses tend to give results that indicate the combustion is less stable than the more realistic distributed-combustion approach (private communication with J. Fang). In an attempt to compensate for this problem, a constant 1.0 is added to the right-hand side of equation (25) in reference 9; predictions using this practice appear to correlate better with test data (private communication with J. Fang). For these reasons, the present study follows the practice.

Finally,  $n$  and  $\tau$  can be related to the real part,  $y_R$  and the imaginary part,  $y_I$  of the burning admittance:

$$n = \frac{y_R^2 + y_I^2}{2y_R} \quad (26)$$

$$\tau = \frac{1}{\omega} \arctan \left( \frac{2y_I}{y_I^2/y_R - y_R} \right) \quad (27)$$

where  $\omega$  is the imaginary part of  $s$  which is the angular frequency of the oscillation.

### III. PROGRAM DESCRIPTION

The High Frequency Instability Analysis computer program HIFI consists of a main program and eight subroutines which are described in the next section. Program input and output are described in Sections 2 and 3. A listing of the computer code is provided in appendix A. Input and output for a sample case are provided in appendix B.

#### 3.1 Program Description

All input to the code are made in the main program. The input include chamber and nozzle geometry, location of the combustion plane, chamber gas properties, cavity geometry, cavity gas properties, chamber acoustic resonance mode, and frequency range of interests. For each frequency in the specified range, the main program calculates chamber admittance upstream of the combustion plane after calling subroutines TED and CAP2 to calculate cavity admittances. Next, it calls subroutine NOZADM to calculate nozzle admittance or it obtains the nozzle admittance value from a table generated by a previous run. If the nozzle admittance is calculated, its value is written to a file NOADTA.DAT for future runs in which the nozzle geometry and resonance modes are the same as the run that generates the nozzle admittance file. After the nozzle admittance value is determined, the main program calculates chamber admittance downstream of the combustion plane. Finally, it calculates and output the burning admittance and the corresponding values of  $n$  and  $\omega$ . In addition, it outputs

the value of the chamber admittances upstream and downstream of the combustion plane.

- Subroutine CALADM: called by subroutine NOZADM to calculate nozzle admittance.
- Subroutine CAP2: calculates cavity admittances. This subroutine is taken from the computer code IFAR (Ref. 9).
- Subroutine INTGRT: called by subroutine NOZADM to perform numerical integration. See reference 11 for more description of this subroutine.
- Subroutine MACH: calculates Mach number as function of area ratio and specific heat ratio using successive iteration techniques.
- Subroutine NOZADM: "main" program of the computer code for calculating nozzle admittance. See reference 11 for more description of this subroutine.
- Subroutine NOZINI: calculates values of variables that are independent of the frequency and are frequently used by subroutine NOZADM. This reduces computer time by avoiding repetitive calculations of these variables every time NOZADM is called.
- Subroutine NOZTAB: obtains the value of nozzle admittance from a previously generated table.
- Subroutine TED: calculates the effect of cavity distribution with respect to mode orientation. This

subroutine is taken from the computer code IFAR (Ref. 9)

### 3.2 Input Description

Input to the computer code is divided into four groups: the first group is the problem description; the second group is the namelist CNTRL which specifies chamber type (rectangular or cylindrical) and file generation options; the third group is the namelist INPUT which specifies the chamber and nozzle geometry, the chamber gas properties, the chamber acoustic resonance mode and the frequency range of interests; and the fourth group is the data specifying cavity geometries and cavity gas properties. A sample input file is provided in Appendix B.1.

The problem description can be specified using any number of lines but at least one line must be used although it can be a blank line. Following the problem description is the namelist CNTRL and subsequently the namelist INPUT. Variables in the namelists CNTRL and INPUT are described in tables 3.1 and 3.2, respectively. The last group of input data pertaining to cavity geometries and cavity gas properties immediately follows the namelist INPUT. This last group of data is described line-by-line here:

CARD 1: Variables NCAV1, NCAV2, NSEC1, NSEC2, NTESTM  
Format (5I10)

Variable name	Unit	Description
NCAV1	----	Number of group 1 cavities
NCAV2	----	Number of group 2 cavities

Name	Type	Unit	Description and Remarks
AXISYM	L		=TRUE for cylindrical chambers =FALSE for rectangular chambers
PLOT	L		=TRUE if plot files are to be generated =FALSE if plot files are NOT to be generated
TABLE	L		=TRUE if nozzle admittance is obtained from a file that has been generated by a previous run. =FALSE if nozzle admittance is to be calculated internally.

Table 3.1: Descriptions of namelist CNTRL variables

Name	Type	Unit	Description and Remarks (*)
GAMMA	R	None	Specific heat ratio
HST	R	ft	Throat radius (throat half-height)
RC	R	ft	Radius of curvature at the throat
XSSL	R	ft	Chamber straight-section length
XB	R	ft	Distance from injector face to combustion plane
RCHAMB	R	ft	Chamber radius (chamber half-height)
RWI	R	ft	Not used for axisymmetric case. (Chamber half-thickness)
RE	R	ft	Radius of curvature at nozzle entrance
ALPHA	R	deg.	Nozzle convergence half-angle
AO	R	ft/s	Speed of sound at stagnation condition
PCHAMB	R	psf	Chamber pressure
WS	R	Hertz	Initial frequency
DW	R	Hertz	Frequency increment
NW	I	None	Number of frequency values
M	I	None	Tangential (tranverse or width) mode number
N	I	None	Radial (lateral or thickness) mode number

Table 3.2: Description of namelist INPUT variables

(\*) Descriptions enclosed in parentheses are for rectangular chamber cases.

NSEC1	----	Number of sections of group 1 cavities
NSEC2	----	Number of sections of group 2 cavities
NTESTM	----	Maximum number of iterations when calculating cavity admittance.

```

*** The following input are not required ***
*** if NCAV1=0 and NCAV2=0 in card 1 ***

```

CARD 2: Variables WD, AC, LC, LOGIC1, LOGIC2, LOGIC3, LOGIC4  
Format (3E10.3, 4I10)

Variable name	Unit	Description
WD	ft.	Width of cavity section
AC	ft**2	Cross-sectional area of cavity section
LC	ft.	Length of cavity section
LOGIC1	----	Cavity inlet characteristic =0, square edged inlet =1, rounded inlet =2, well-rounded inlet
LOGIC2	----	Switch to specify whether or not to consider sound absorption coefficient. =0, not considered =1, considered
LOGIC3	----	Switch to specify data to be used =0, use cavity inlet data =1, use data inside cavity
LOGIC4	----	Switch to specify cavity type =0, circular cross-sectional cavity =1, rectangular cross-sectional cavity

CARD 3: Variables RHOC, CC, GAMMAC, PO, PRTL, VIS  
Format (6E10.3)

Variable name	Unit	Description
RHOC	lb/ft3	Cavity gas density

CC	ft/s	Sound speed in cavity
GAMMAC	----	Specific heat ratio of cavity gas
PO	----	Cavity inlet pressure amplitude normalized by mean chamber pressure
PRTL	----	Prandtl number of cavity gas
VIS	lbf-s/ft <sup>2</sup>	Viscosity of cavity gas

Cards 2 and 3 are repeated for NSEC1 sections of group 1 cavities, then they are repeated for NSEC2 sections of group 2 cavities. If NCAV1=0 in Card 1, then Cards 2 and 3 begin with the first section of group 2 cavities. Likewise, if NCAV2=0 Cards 2 and 3 end with last section of group 1 cavities.

For each group of cavity, the input for the sections must be in sequential order beginning with the section at the end of the cavity and ending with the section at the cavity entrance.

It should be noted here that for rectangular chamber cases, the cavities are assumed to be located at the edge of the chamber width. Although the effects of the cavities themselves are considered, the effects of the cavity distribution with respect to the lateral (thickness) modes are not considered. Minor modifications can be made to the code to account for the cavity distribution effects.

```

*** The following input are not required ***
*** if AXISYM=F in namelist CNTRL ***

```

CARD 4: Variables MO  
Format (I10)

Variable name	Unit	Descriptions
MO	----	Mode orientation number =1, mode orientation in which group 2 cavities are more effective than they are in other mode orientations =2, mode orientation in which group 1 cavities are more effective than they are in other mode orientations

CARD 5: Variables (IDCAV(I), I=1, NCAV1+NCAV2)  
Format (20I4)

Variable name	Unit	Descriptions
IDCAV	----	Cavity group number distribution. For example, there are 6 group-1 cavities and 3 group-2 cavities and the distribution such that there is one group-2 cavity followed by two group-1 cavities. Card 5 should be input as follow: 2 1 1 2 1 1 2 1 1

### 3.3 Output Description

Output from the code begin with the echo of input data which includes a problem description, variables in namelists CNTRL and INPUT, and cavity input data. Although the problem description can be input using any number of lines, only the first line is output. Following the echo of the input data is the chamber Mach number. The last section of the output is the calculated stability results which include the amplitude and phase of the burning admittance,  $n$  and  $\tau$ , the real parts and the imaginary parts of the chamber admittances upstream and downstream of the combustion plane, the real part and the

imaginary part of the burning admittance. These results are calculated and output for each frequency in the range specified in namelist INPUT. A sample output file is provided in Appendix B.2.

In addition to the output described above, a file NOADTA.DAT is also output by the code if the variable TABLE in the namelist CNTRL is equal to FALSE. This file contains a table of the calculated nozzle admittance vs. frequency. For future runs with the same nozzle geometry and resonance modes, the code obtains the values of nozzle admittances from the table instead of re-calculating them. This option saves computer time and can be selected by setting the variable TABLE in the namelist CNTRL to TRUE.

If the variable PLOT in the namelist CNTRL is equal to TRUE, three additional files are generated by the code. These files can be input to any x-y plotting package, for example TELLEGRAF, to plot the calculated results. The first file, AMPLD.PLT contains the amplitude of the burning admittance vs. frequency. The second file, PHASE.PLT contains the phase angle of the burning admittance vs. frequency. The third file, NTAU.PLT contains the pressure interaction index,  $n$  vs. the sensitive time lag,  $\tau$ .

#### IV. RESULTS AND DISCUSSION.

Calculations of the burning admittances and  $n-\tau$  neutral stability curves for a cylindrical chamber and a rectangular chamber were made using the computer code HIFI. Input parameters for the two cases are shown in figures 4.1 and 4.2, respectively. Results are compared with the IFAR predictions.

A typical run for 100 frequency values, in which the nozzle admittance are calculated internally, requires approximately 40 CPU seconds on the micro-VAX at Aerojet TechSystems. A similar run, in which nozzle admittance is provided as a table, requires approximately 4 CPU seconds.

Figures 4.3 shows the  $n-\tau$  neutral stability curves. Figures 4.4 and 4.5 show the burning-admittance amplitudes and the burning admittance phase angles versus frequency. Figure 4.6 is a replot of the burning admittance amplitude shown in Figure 4.4 but is shown on a larger scale to show the results near resonance frequencies. These results are calculated using the computer codes HIFI and IFAR for mixed 1T and longitudinal modes of a cylindrical chamber without acoustic cavities. The figures show that the differences between HIFI and IFAR predictions are small near resonance. The differences become larger at off-resonance frequencies but this is not important since we are more interested in the region near resonance.

Calculations are then made for the 1T mode of a cylindrical chamber with and without acoustic cavities using computer codes HIFI and IFAR. The calculations were made to

(a)

```

HIFI, AX, 1T
$CNTRL
  AXISYM=T, TABLE=F, PLOT= T,
$END
$INPUT
  RC=0.216, RE=0.216, ALPHA=45.0, HST=0.208, RCHAMB=0.623, RWI=0.2,
  GAMMA=1.14, AO=3850.0, WS=1600.0, DW=5.0, NW=100, M=1, N=0,
  XSSL=1.0, XB=0.21, PCHAMB=2.088E+05,
$END
      0          0          1          1          100

```

(b)

```

HIFI, AX, 1T
$CNTRL
  AXISYM=T, TABLE=T, PLOT= T,
$END
$INPUT
  RC=0.216, RE=0.216, ALPHA=45.0, HST=0.208, RCHAMB=0.623, RWI=0.2,
  GAMMA=1.14, AO=3850.0, WS=1600.0, DW=5.0, NW=100, M=1, N=0,
  XSSL=1.0, XB=0.21, PCHAMB=2.088E+05,
$END
      10          0          1          1          100
1.500E-01 0.750E-02 0.320E+00          2          1          0
0.514E-00 2.750E+03 1.250E+00 0.100E+00 0.472E+00 4.580E-05          1
      2
      1  1  1  1  1  1  1  1  1  1

```

Figure 4.1: Input data for a 1T mode in a cylindrical chamber.  
 (a) without cavities, (b) with cavities.

(a)

```
HIFI, 2D, 1W
$CNTRL
  AXISYM=F, TABLE=F, PLOT= T,
$END
$INPUT
  RC=0.216, RE=0.216, ALPHA=45.0, HST=0.208, RCHAMB=0.623, RWI=0.2,
  GAMMA=1.14, AO=3850.0, WS=1300.0, DW=5.0, NW=100, M=1, N=0,
  XSSL=1.0, XB=0.21, PCHAMB=2.088E+05,
$END
      0          0          1          1          100
```

(b)

```
HIFI, 2D, 1W
$CNTRL
  AXISYM=F, TABLE=T, PLOT= T,
$END
$INPUT
  RC=0.216, RE=0.216, ALPHA=45.0, HST=0.208, RCHAMB=0.623, RWI=0.2,
  GAMMA=1.14, AO=3850.0, WS=1300.0, DW=5.0, NW=100, M=1, N=0,
  XSSL=1.0, XB=0.21, PCHAMB=2.088E+05,
$END
      2          0          1          1          100
1.500E-01 0.750E-02 0.420E+00      2          1          0          1
0.514E-00 2.750E+03 1.250E+00 0.100E+00 0.472E+00 4.580E-05
```

Figure 4.2: Input data for a 1W mode in a rectangular chamber.  
(a) without cavities, (b) with cavities.

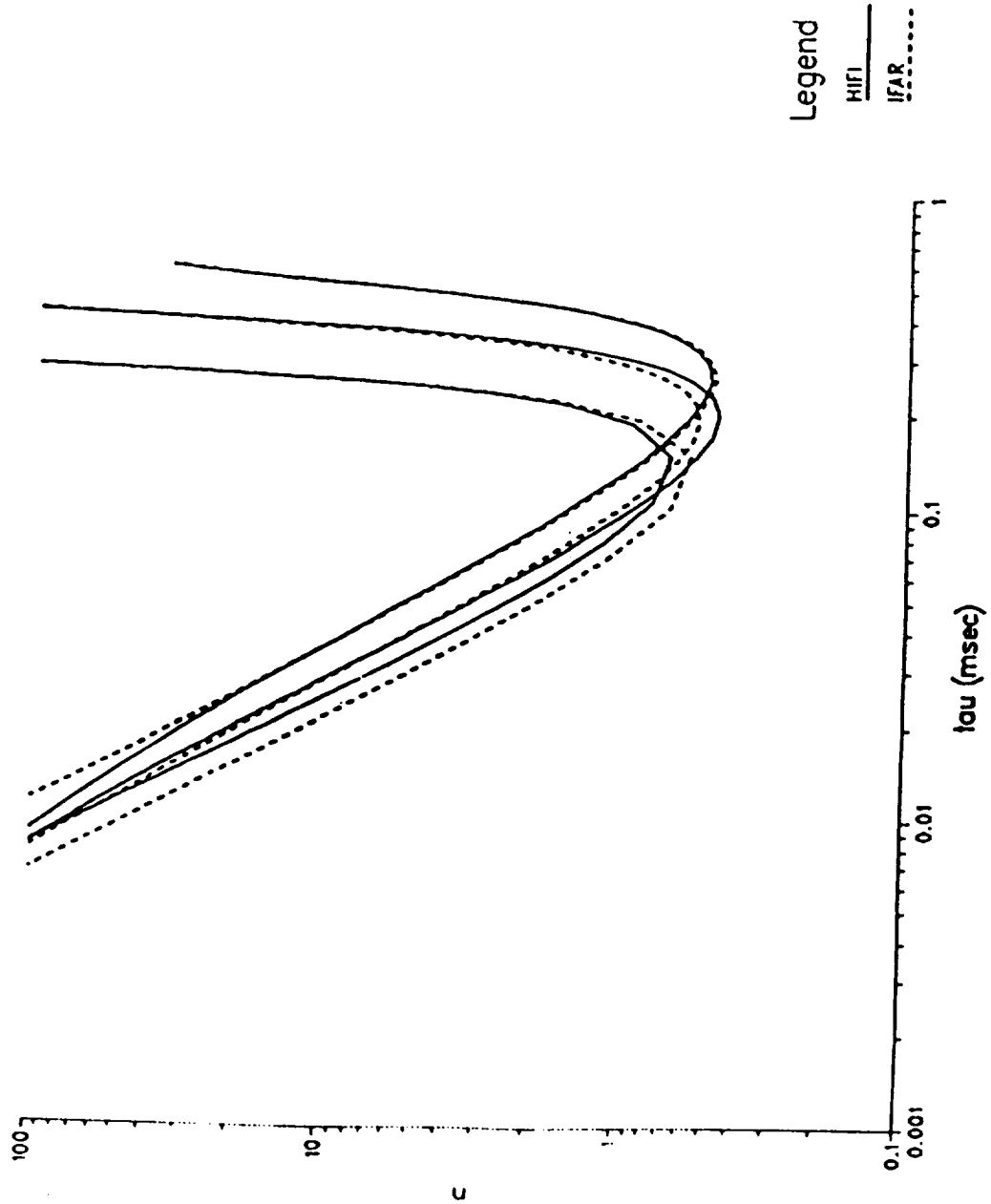


Figure 4.3 : Comparisons between HIFI and IFAR predictions of neutral stability curve in a IT mode.

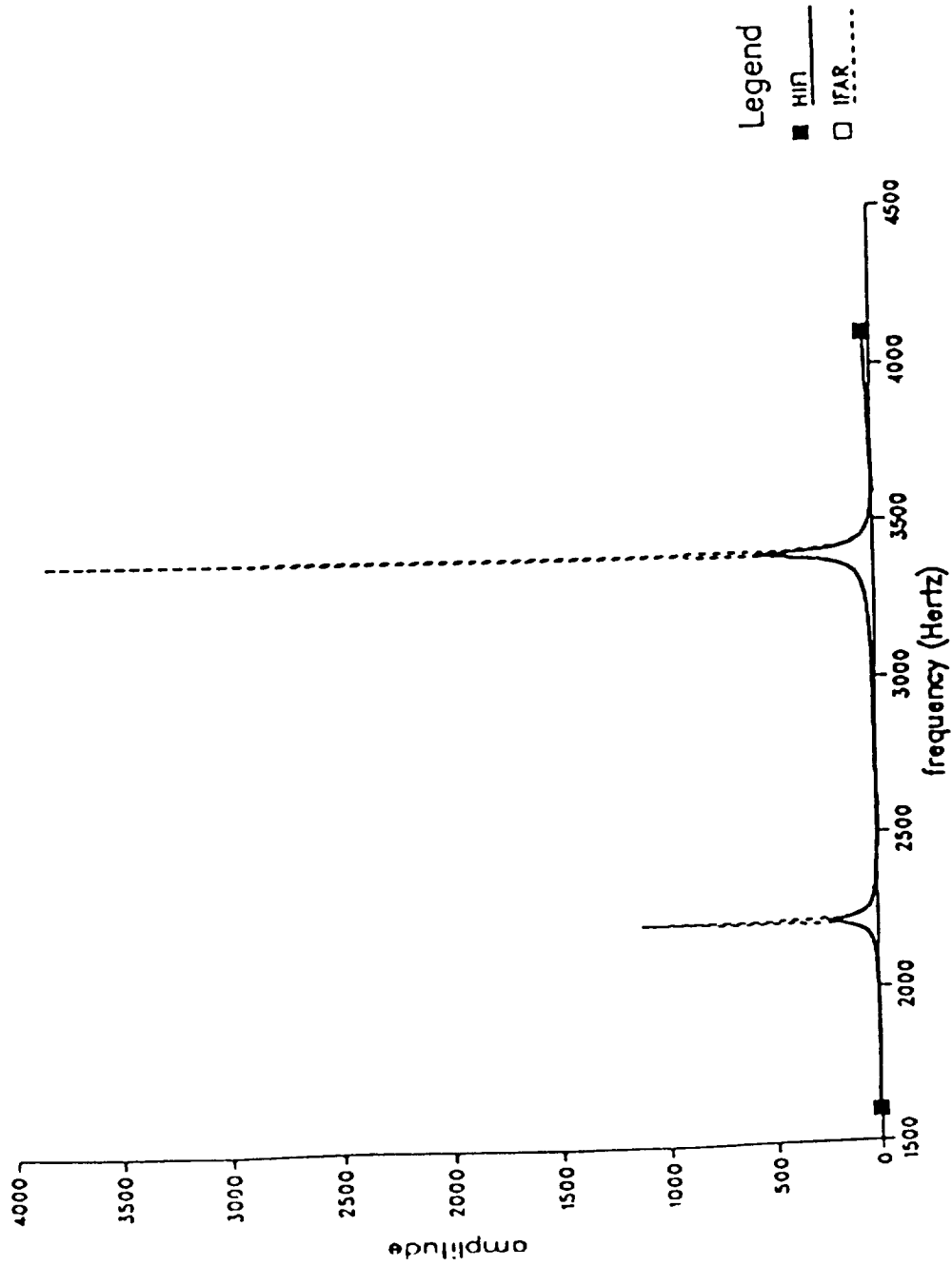


Figure 4.4 : Comparisons between HIFI and IFAR predictions of burning admittance amplitude in a 1T mode.

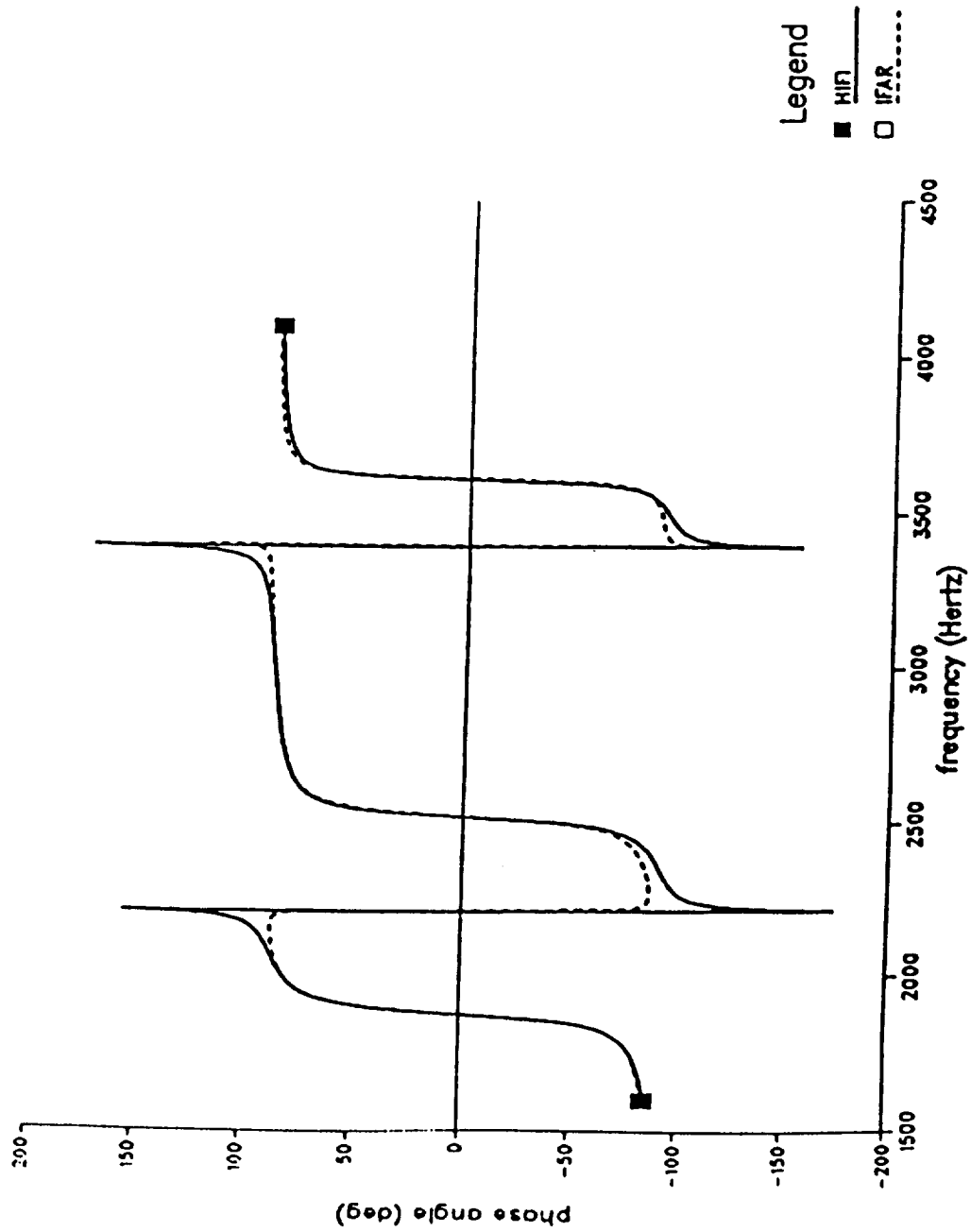


Figure 4.5 : Comparisons between HIFI and IFAR predictions of burning admittance phase angle in a 1T mode.

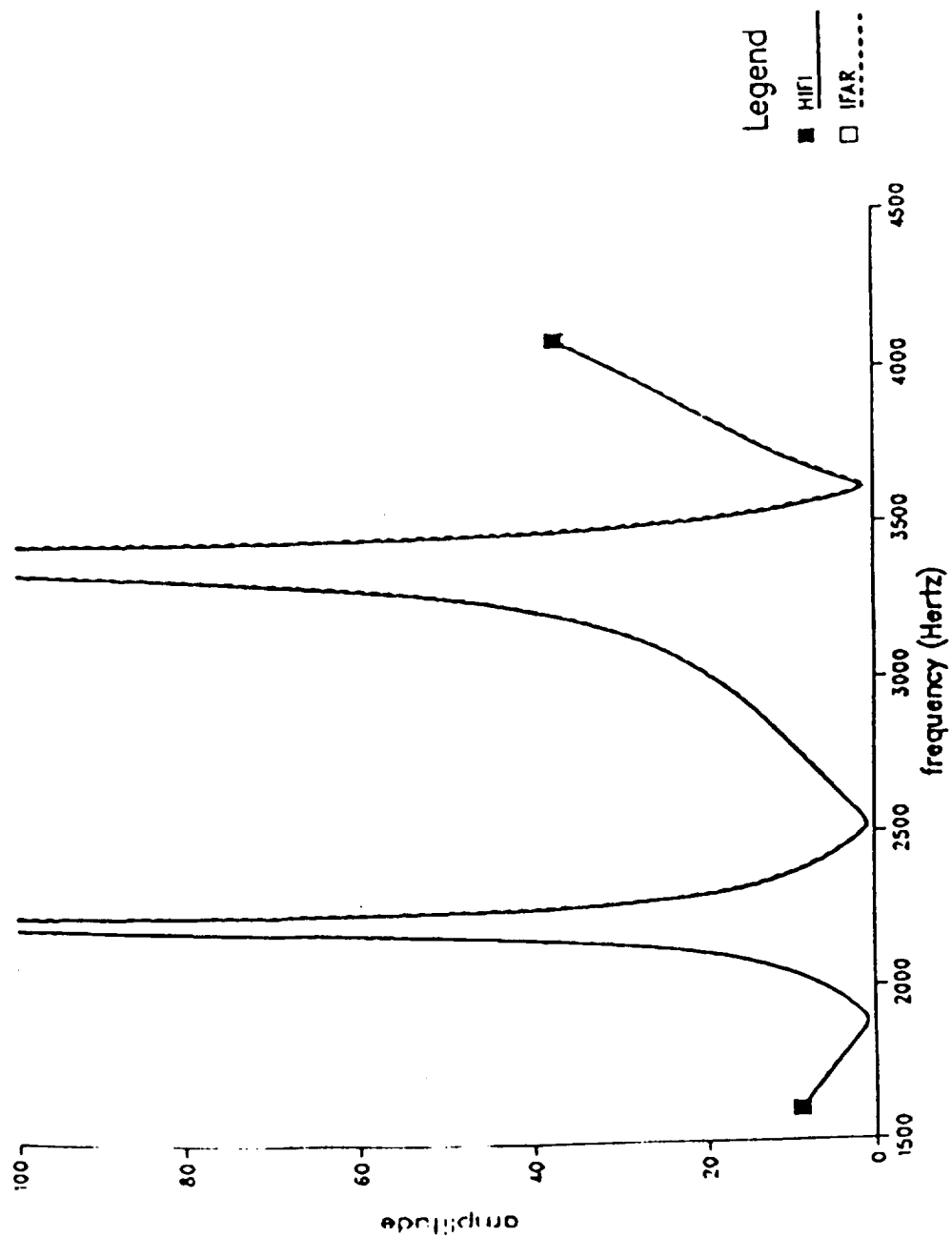


Figure 4.6 : Comparisons between HIFI and IFAR predictions of burning admittance amplitude in a 1T mode.

study the effects of acoustic cavities on combustion stability, and to provide further comparisons between HIFI and IFAR predictions. Figures 4.7, 4.8 and 4.9 show the calculated results. Again, the differences between HIFI and IFAR predictions are small. Both computer codes predict the stabilizing effects of the acoustic cavities as shown by the higher value of the  $n$  minimum, the minimum interaction index that can support linearly instability. Another effect of acoustic cavities predicted by the computer codes is to shift the value of  $\tau$  where the  $n$  minimum occurs to a higher value (Fig. 4.7) and to shift the resonance frequency to a lower value (Fig. 4.8).

Figures 4.10, 4.11 and 4.12 show the comparisons between HIFI and IFAR calculated results for a 1W mode of a rectangular chamber. Similar to the cylindrical case, the differences between the predictions are small.

Figures 4.13, 4.14 and 4.15 show the effects of acoustic cavities on the stability results for the 1W mode of a rectangular chamber. The effects are similar to those discussed above for the 1T mode of the cylindrical chamber.

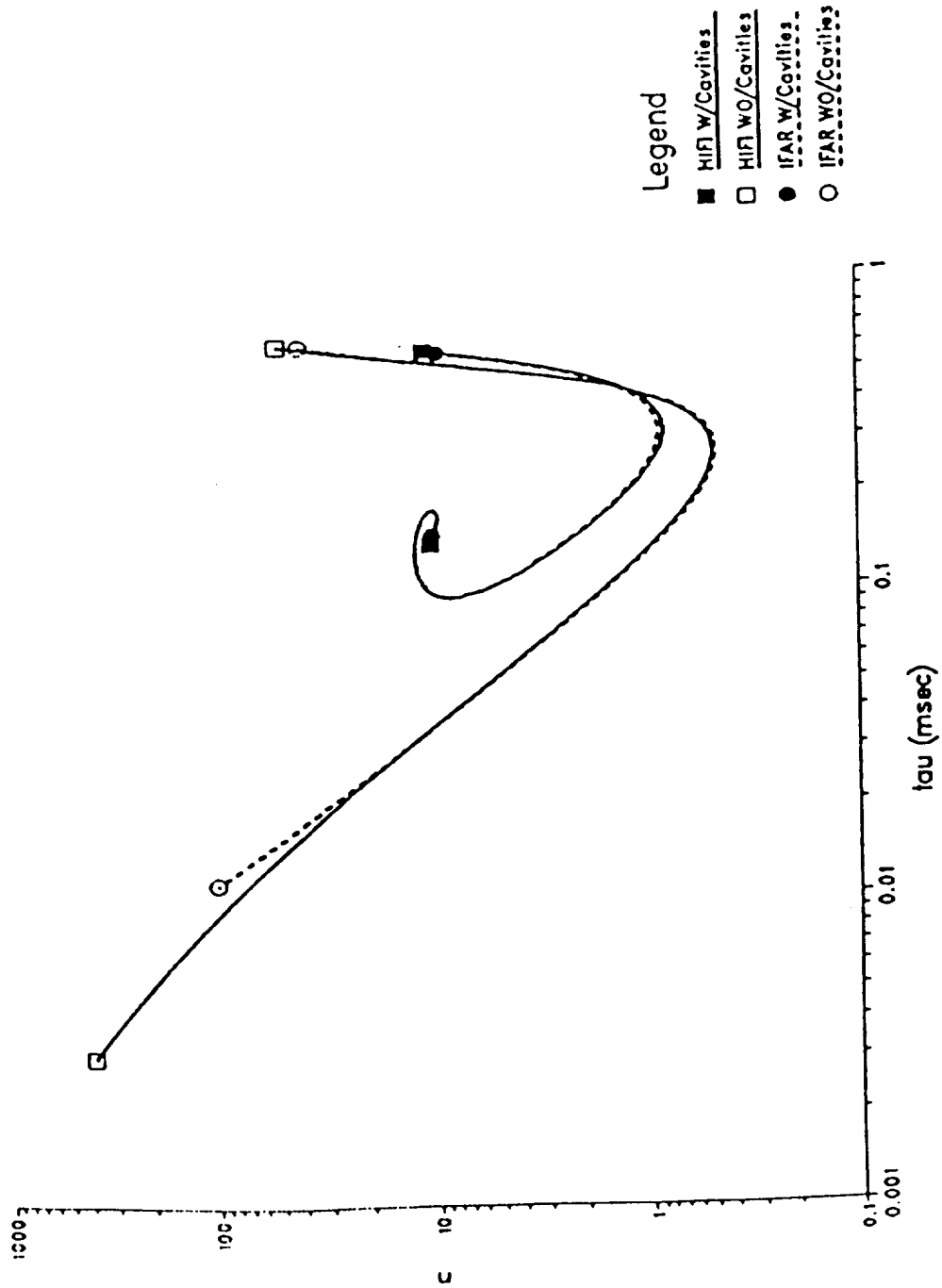


Figure 4.7 : Effects of acoustic cavities on neutral stability curve in a 1T mode. Comparisons between HIFI and IFAR predictions.

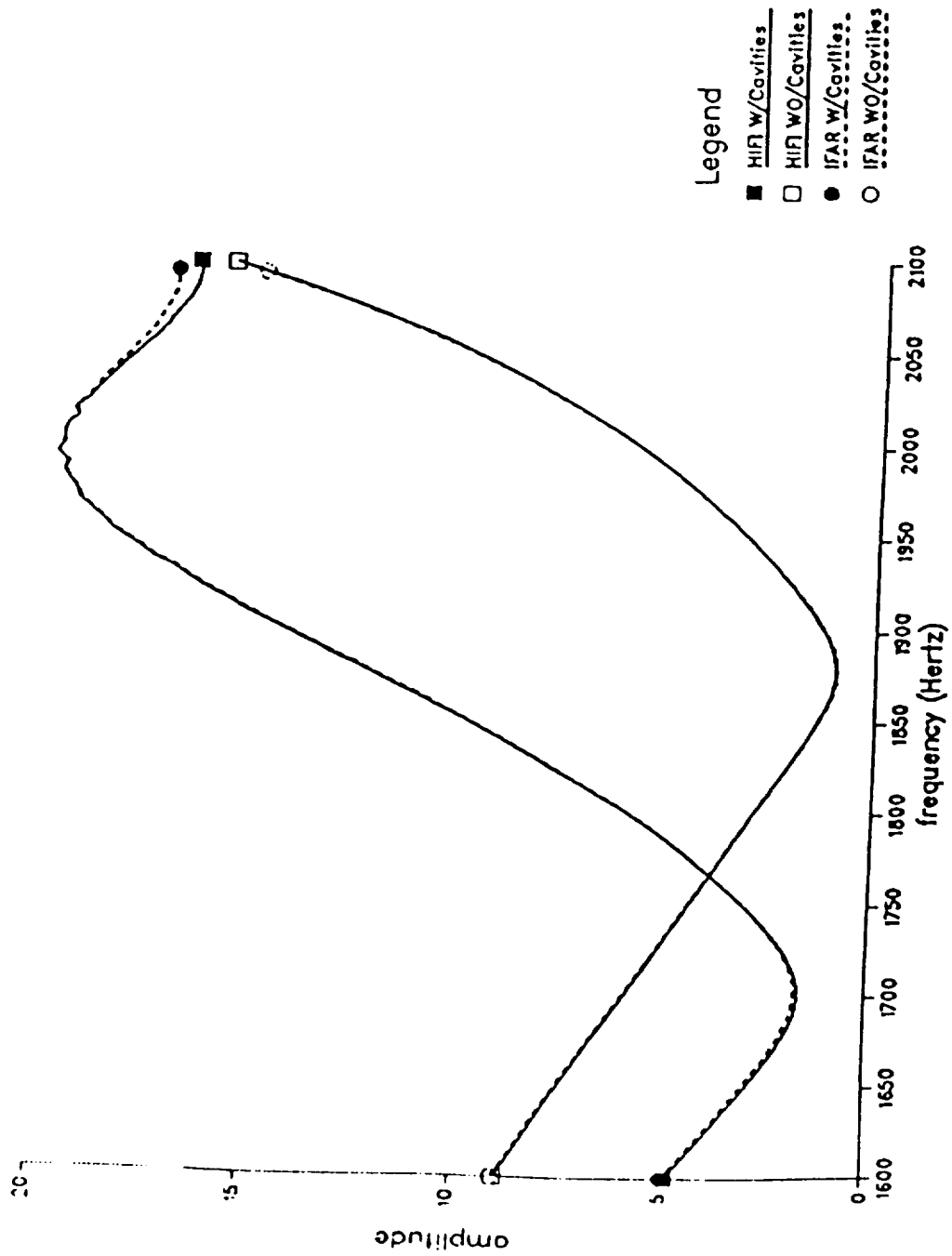


Figure 4.8 : Effects of acoustic cavities on burning admittance amplitude in a 1T mode. Comparisons between HIFI and IFAR predictions.

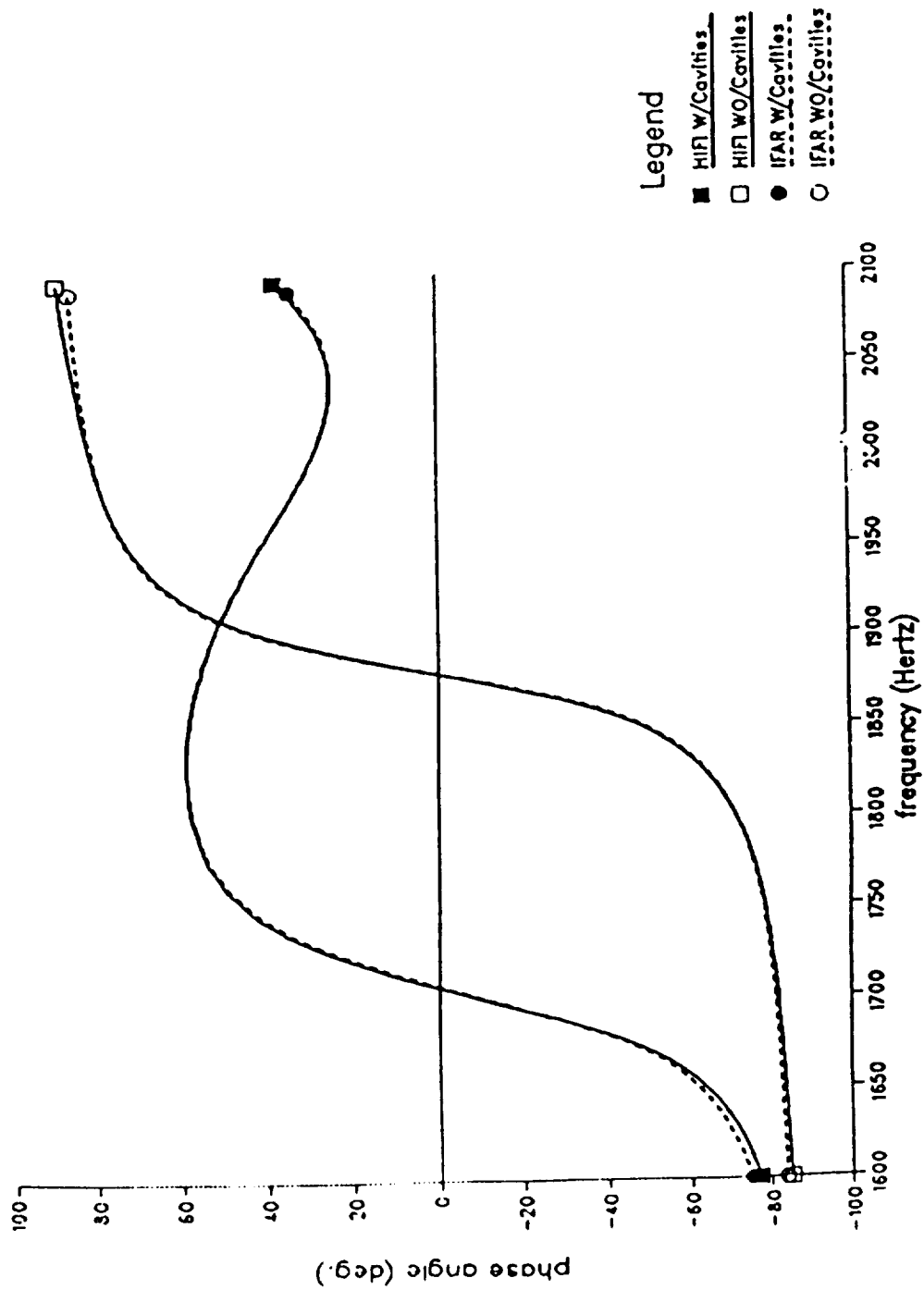


Figure 4.9 : Effects of acoustic cavities on burning admittance phase angle in a 1T mode. Comparisons between HIFI and IFAR predictions.

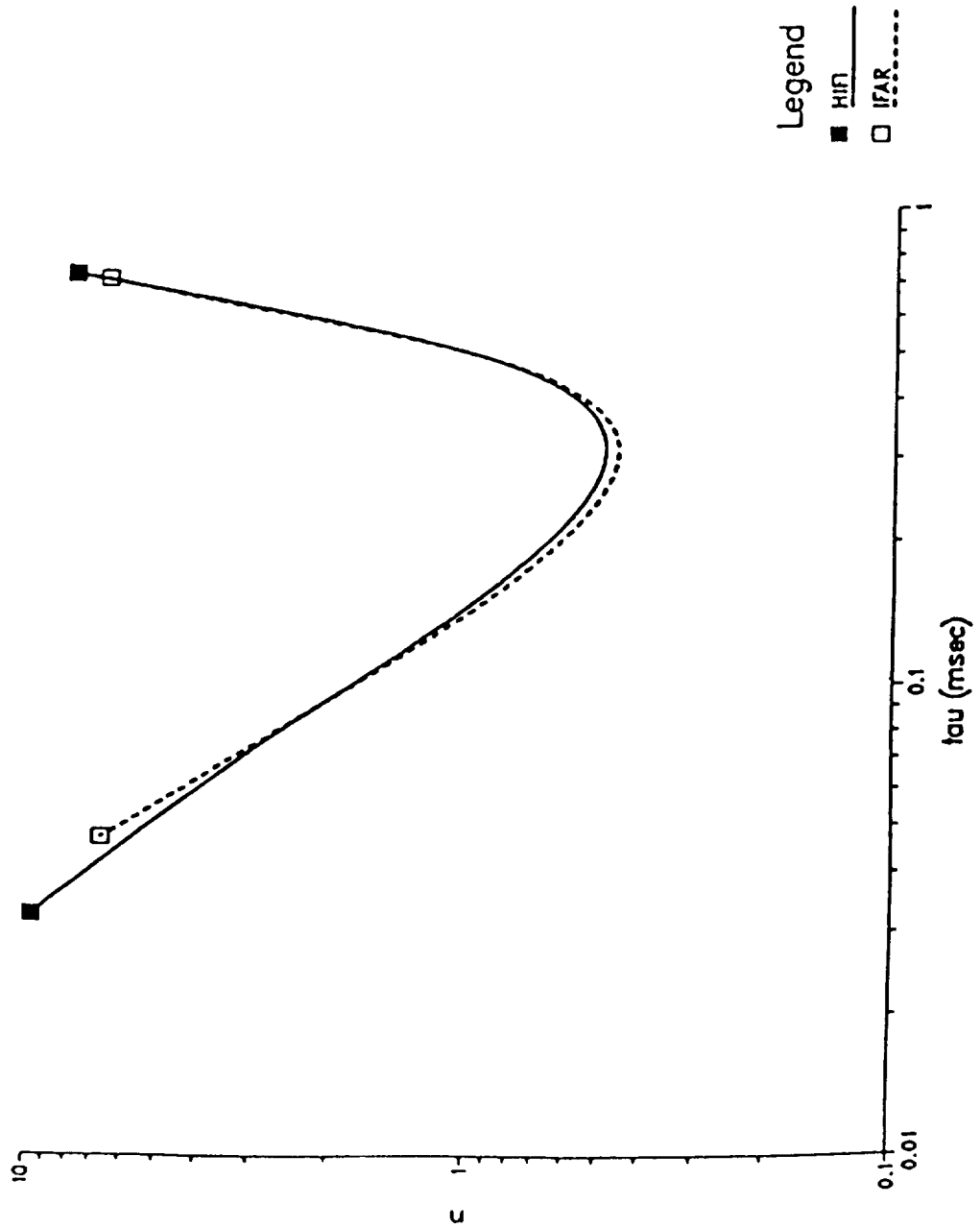


Figure 4.10: Comparisons between HIFI and IFAR predictions of neutral stability curve in a LW mode.

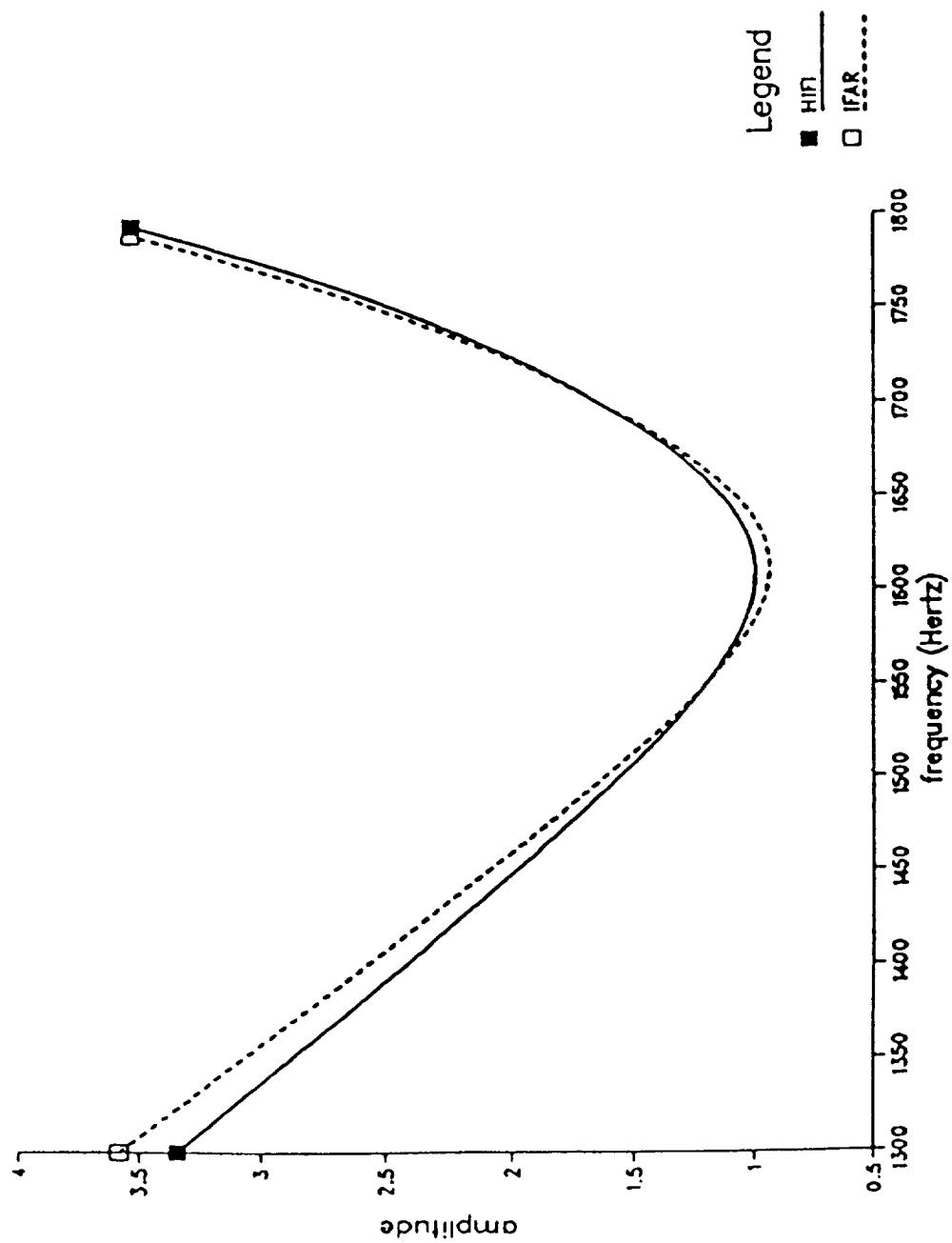


Figure 4.11: Comparisons between HIFI and IFAR predictions of burning admittance amplitude in a 1W mode.

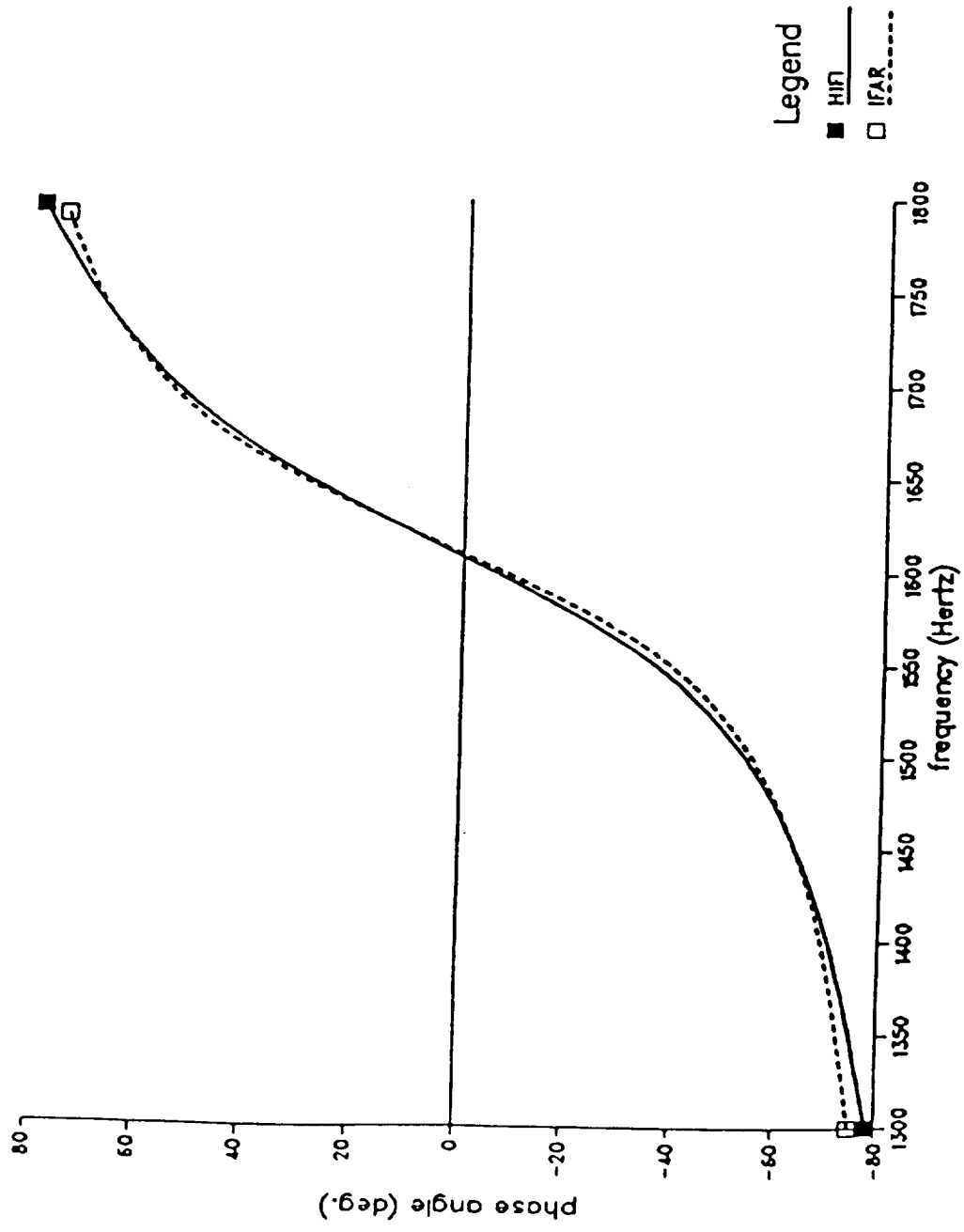


Figure 4.12: Comparisons between HIFI and IFAR predictions of burning admittance phase angle in a 1W mode.

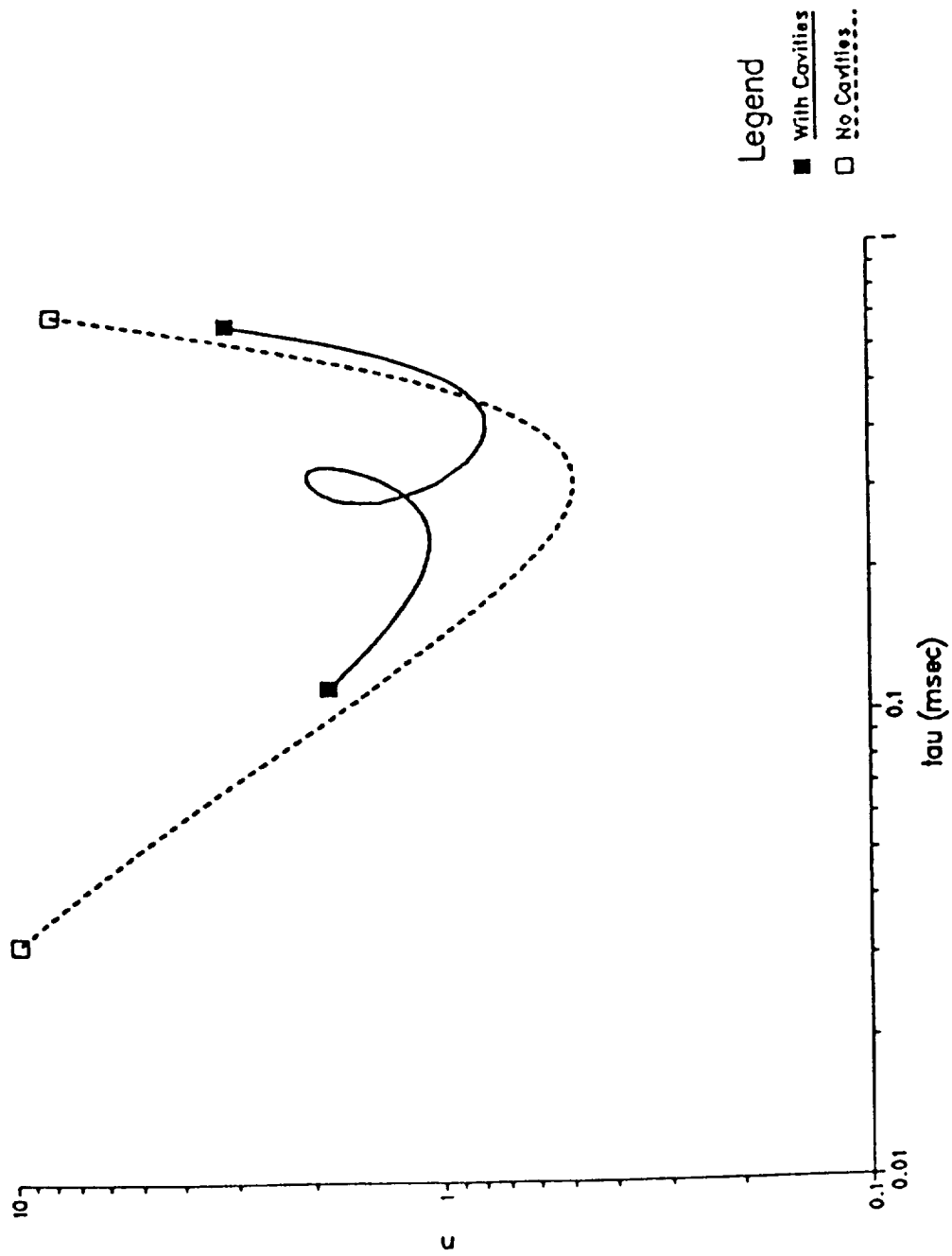


Figure 4.13: Effects of acoustic cavities on neutral stability curve in a 1W mode.

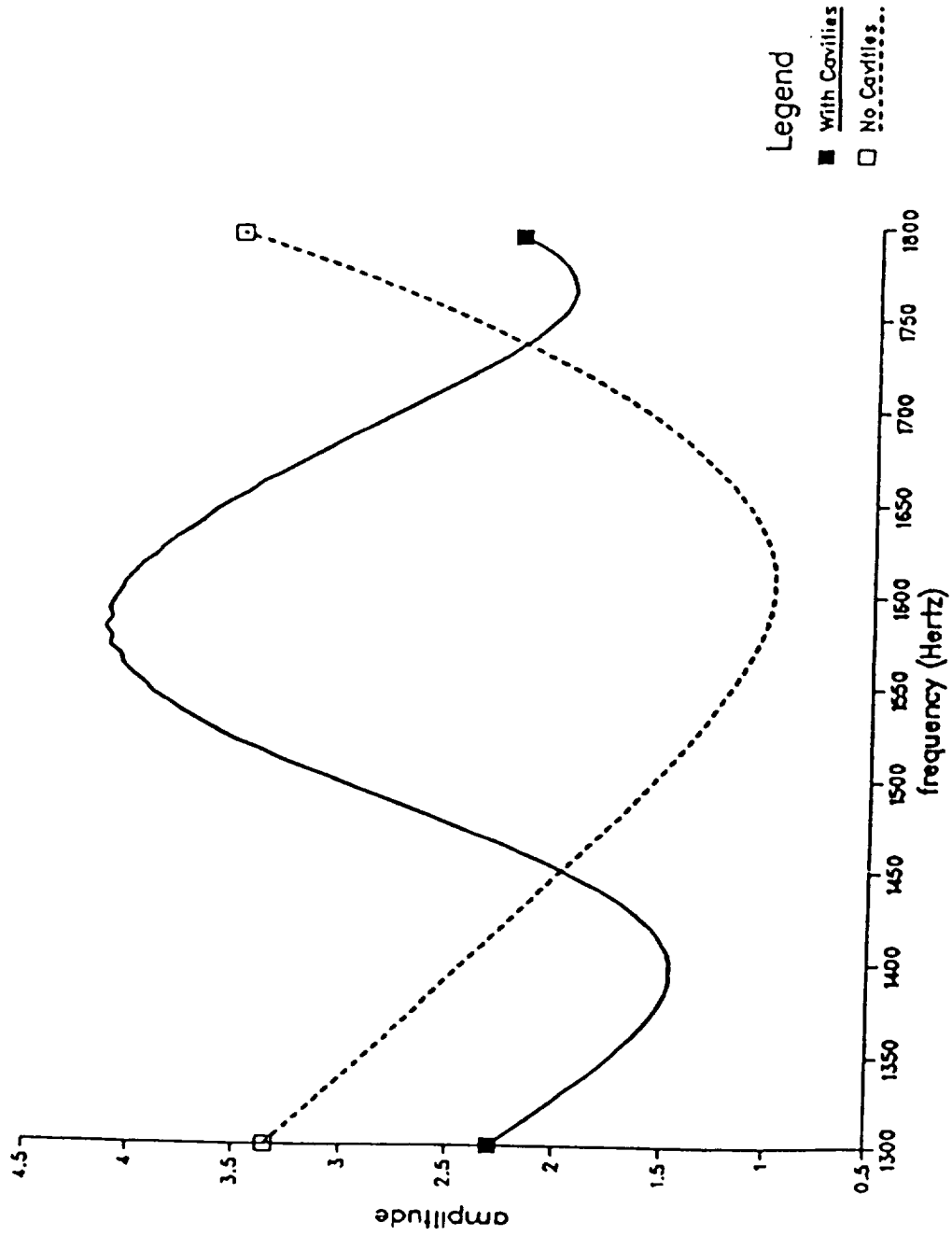


Figure 4.14: Effects of acoustic cavities on burning admittance amplitude in a 1W mode.

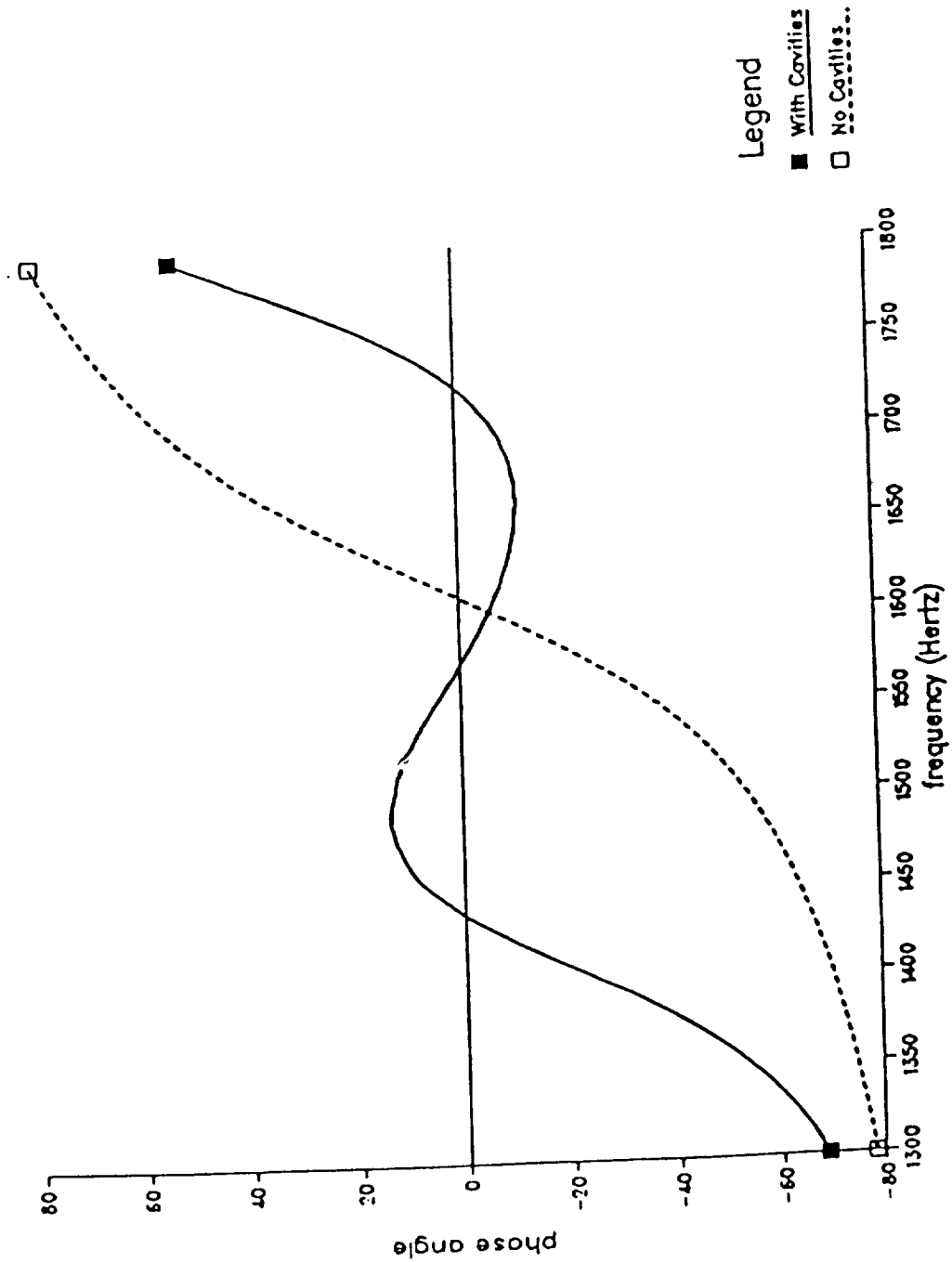


Figure 4.15: Effects of acoustic cavities on burning admittance phase angle in a 1W mode.

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**PART B**

**3-D DISTRIBUTED COMBUSTION BAFFLE MODEL  
(DIST3D)**

USER'S MANUAL FOR THE MULTIDIMENSIONAL  
BAFFLE MODEL COMPUTER PROGRAMS

by

C.E. Mitchell, D.J. Howell, F.E. Dodd  
T.L. Acker

Department of Mechanical Engineering  
Colorado State University  
Fort Collins, Colorado 80523

for

Aerojet TechSystems Company

31 July 1987

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## 1. Introduction

The overall goal toward which the analytical models and computer programs presented here are directed is the development of predictive tools for determining the stability behavior of two and three dimensional liquid propellant rocket thrust chambers. Two basic combustion distribution models are discussed. The first of these assumes that a concentrated zone of combustion exists at the injector face. This assumption implies that the remainder of the chamber is source free and consequently leads to a relatively simple analysis. This approach was employed by Baer and Mitchell in their original models and computer codes (Refs. 1 and 2). One objective of the current effort was the resurrection of these early programs in a form that would be immediately useful and convenient. This involved retrieving the codes from punched cards, rewriting portions of the code in standard Fortran 5, correcting minor errors originally present, adding a nozzle admittance code (Aerojet TechSystems NOZADM), generalization of the combustion response input so that arbitrary values of the interaction index ( $n$ ) and time lag ( $\tau$ ) could be used and, finally, conversion of most of the essential input and output data to dimensional form, consistent with Aerojet's stability models. Minor improvements to the code including modernization of the input and output modes and addition of comment statements were also made. The final result of this work was two computer programs: CON2D, for two dimensional thrust chambers and CON3D, for three dimensional thrust chambers.

The second combustion distribution model is considerably more realistic in that the zone of combustion is taken to be distributed over a significant fraction of the chamber axial length. The distribution is limited to a linear form (constant mean volumetric rate of mass and energy release over the length of the zone), but the beginning and end of the zone are arbitrary. This allows consideration of both relatively intense

(concentrated) zones of combustion as well as less intense (distributed) zones, and permits the determination of the impact of the concentration of combustion sources on stability. Moreover, the zone of combustion may be located completely in the baffle cavities, completely in the unbaffled main chamber, or partly in both regions. Thus, the influence of combustion zone location on stability can be assessed as well.

The analysis of the distributed combustion problem is both more sophisticated and more complex than that of the concentrated combustion model. The analytical approach employed follows in many respects recent work on combustion distributions in unbaffled combustion chambers done at Colorado State University (Ref. 3). The resulting computer codes are more flexible than those for the concentrated combustion mode. The main additional features are:

1. Output in terms of  $n$ , tau neutral stability data as well as in terms of frequency and decay rate for given combustion response input.
2. Inclusion of a radially oriented acoustic cavity in the stability model.
3. Output of local pressure amplitudes and phase angles at any spatial location in the chamber or baffle cavities.

For distributed combustion the two dimensional computer program is called DIST2D while the three dimensional program is DIST3D.

The remainder of this manual is devoted to a presentation of the analytical approaches involved and a description of the use of the resulting computer programs.

## 2. Theory

### 2.1 Analytical Approach

In many respects the general analytical approaches for treating the concentrated combustion model and the distributed combustion

model are quite similar. Consequently, a single description of the analytical development will be given. Specific differences between the analyses for the two combustion zone types (as well as for the differences between two and three dimensional chambers) will be noted as the presentation proceeds.

### 2.1.1 Model Assumptions

The following modeling assumptions are made in representing the major thrust chamber components in the analyses.

#### 1. Chamber Geometry

Three dimensional. A right circular cylinder is terminated by an axis-symmetric nozzle (see Fig. 1). Cylindrical coordinates  $(r, \Theta, z)$  are used in the analysis. The nozzle entrance is at  $Z^*=L^*$ , the injector is at  $Z^*=0$ , the cylindrical wall at  $r^*=R^*$  (asterisks indicate dimensional quantities). The radial absorber is located at  $r^*=R^*$ ,  $0 \leq Z^* \leq Z_A^*$ . Baffle blades are located at  $\Theta_j = \frac{2\pi j}{N}$ , where  $j$  is an integer,  $N$  is the total number of baffle cavities, and  $0 \leq j \leq N-1$ . Baffle blade length is  $Z_B^*$  thickness is  $T^*$ .

Two dimensional. A two dimensional (vertical pancake) chamber is terminated by a convergent two dimensional nozzle (see Fig. 1). Cartesian coordinates  $(Z^*, y^*)$  are used in the analysis. Chamber length is  $L^*$ , chamber height is  $R^*$ . The absorber is located at  $y^* = 0$  and  $y^* = R^*$ , with  $0 \leq Z^* \leq Z_A^*$ . Baffle blades are located at  $y_j^* = \frac{j R^*}{N}$  where again,  $N$  is the total number of baffle cavities and  $0 \leq j \leq N-1$ . Baffle blade length is  $Z_B^*$ .

#### 2. Gasdynamic Flowfield

- . uniform composition calorically perfect combustion product gas
- . irrotational flow outside of baffle boundary layers (exact for concentrated combustion, correct through

order mean Mach number squared in distributed combustion).

- . one dimensional axial flow in the steady state.
- . linear, nearly harmonic oscillations with small growth or decay rates (less than about 30% per cycle).
- . droplet volume and drag ignored (distributed combustion).
- . standing or traveling waves in the main chamber  
standing waves in the baffle compartments (standing waves only for two dimensional chamber).

### 3. Combustion Distribution

#### Concentrated combustion

- . all combustion concentrated at  $Z^*=0$
- . mass flow rate,  $\bar{m}^*$ , Mach number  $M$ , pressure  $\bar{p}^*$  constant for  $Z^*>0$  in the steady state (steady state indicated by superposed bars)

#### Distributed combustion

- . linear combustion distribution between  $Z^*=Z_s^*$  and  $Z^*=Z_e^*$

$$\bar{m}^* = \bar{m}_T^* \left( \frac{Z^* - Z_s^*}{Z_e^* - Z_s^*} \right) \quad Z_s^* \leq Z^* \leq Z_e^*$$

$$\bar{m}^* = 0 \quad Z^* < Z_s^*$$

$$\bar{m}^* = \bar{m}_T^* \quad Z^* > Z_e^*$$

$\bar{\dot{m}}_T^*$  is the total mean mass flow in the steady state

- $Z_s^*$  and  $Z_e^*$  are arbitrary, but  $Z_s^* - Z_e^*$  limited to at least 20% of the chamber diameter (or height for two dimensional chamber)

#### 4. Combustion Response

##### Concentrated combustion

$$\frac{\dot{m}'^*}{\bar{\dot{m}}^*} = \left[ n(1 - e^{-i\omega^* \tau^*}) - 1 \right] \frac{p'^*}{\bar{p}^*}$$

This form is corrected from earlier forms. The term -1 on the right-hand side is added, and can be shown formally to be the correct limit of distributed combustion.  $\dot{m}'^*$  is the unsteady perturbation in mass flow,  $p'^*$  the unsteady pressure perturbation,  $\omega^*$  is the dimensional angular frequency,  $\tau^*$  is the time lag.

##### Distributed combustion

$$\frac{\dot{Q}'^*}{\bar{Q}^*} = n(1 - e^{-i\omega^* \tau^*}) \frac{p'^*}{\bar{p}^*}$$

$\dot{Q}$  is the volumetric rate of local gas production in the combustion zone.

#### 5. Baffle Dissipation

- turbulent boundary layer in region near baffle blade tips

- . driven by inviscid outer flow in cavity and main chamber
- . uses Spalding's effective turbulent viscosity model
- . numerical integration over boundary layer near blade tips

#### 6. Acoustic Cavity

- . radial slot
- . located at injector
- .  $u_r^{*'} = \beta_c^* p^{*'}$  where  $u_r^{*'}$  is the radial unsteady velocity,  $p^{*'}$  is the local pressure perturbation,  $\beta_c^*$  is the cavity admittance (complex, must be supplied by user)

#### 7. Nozzle

- .  $u_z^{*' } = \beta_N^* p^{*'}$ ,  $u_z^{*'}$  axial velocity,  $p^{*'}$  pressure at  $z^*=L^*$
- .  $\beta_N^*$  supplied by NOZADM program

### 2.2 Basic Equations

Using the assumptions given above the conservation equations and equations of state can be reduced to a nondimensional set in terms of a velocity potential. The nondimensional scheme is defined in the table below the particular nondimensional variable is formed through division by the listed characteristic quality.

<u>Variable</u>	<u>Characteristic Quantity</u>		<u>Form</u>
	<u>3-D</u>	<u>2-D</u>	
p*(pressure)	$\bar{p}^* (z^* = L^*)$	same	p
$\rho^*$ (density)	$\bar{\rho}^* (z^* = L^*)$	same	$\rho$
a*(sonic speed)	$\bar{a}^* (z^* = L^*)$	same	a
Z*(axial coordinate)	$R^* (\text{chamber radius})$	$R^* (\text{chamber height})$	z
y*(vertical coordinate, 2-D)	—	$R^* (\text{chamber height})$	y
r*(radial coordinate, 3-D)	$R^* (\text{chamber radius})$	—	r
t*(time)	$R^* / \bar{a}_t^*$	same	t
$\omega^*$ (angular frequency)	$\bar{a}_t^* / R^*$	same	$\omega$
$\lambda^*$ (decay rate)	$\bar{a}_t^* / R^*$	same	$\lambda$
$\vec{q}^*$ (velocity vector)	$\bar{a}_t^*$	same	$\vec{q}$
$\phi^*$ (velocity potential)	$\bar{a}_t^* / R^*$	same	$\phi$

In this table it should be pointed out that the characteristic length for the two dimensional problem is the total chamber height  $R^*$ , rather than the half height which is used in some Aerojet programs. The appropriate conversion is made in our computer codes when NOZADM is called.

The dependent variables are all represented as the sum of a steady state (or time averaged) part (superposed bar) and an oscillatory part (prime). Thus,  $p = \bar{p} + p'$ , etc. The oscillatory parts are in turn represented as the product of a space dependent part and the factor  $e^{i\omega t}$ . Thus,  $p' = \bar{p} e^{i\omega t}$ , etc., where  $p = P(\bar{\eta}, z)$  or  $p = P(z, y)$ .  $\omega$  is the complex frequency ( $\omega = \omega_R + i\lambda$ ).

Since the flow is irrotational (at least through order  $u^2$ ) then

$$\vec{q} = \nabla \phi$$

and, from partial integration of the momentum equation

$$p = -\gamma \left( \bar{p} i \omega \phi + \frac{\partial}{\partial z} (\bar{u} \phi) \right)$$

where  $\gamma$  is the ratio of specific heats. Some manipulation finally results in the following basic equation for  $\phi$

$$\nabla^2 \phi + \omega^2 \phi = F_1(\phi, \bar{u}, \omega, n, z) \quad (1)$$

where for concentrated combustion,

$$F_1 = 2M i \omega \frac{\partial \phi}{\partial z} + M^2 \frac{\partial^2 \phi}{\partial z^2}$$

while for distributed combustion

$$F_1 = (1+\gamma) \left[ i \omega \frac{d\bar{u}}{dz} + \left( \frac{d\bar{u}}{dz} \right)^2 \right] \phi + \left( 2\bar{u} i \omega + (\gamma+2) \bar{u} \frac{d\bar{u}}{dz} \right) \frac{\partial \phi}{\partial z} \\ + \bar{u}^2 \frac{\partial^2 \phi}{\partial z^2} - n(1 - e^{-i\omega \tau}) \left[ \gamma i \omega \frac{d\bar{u}}{dz} + 2\gamma \bar{u} \frac{d\bar{u}}{dz} \right] \frac{\partial \phi}{\partial z}$$

The boundary conditions on Equation (1) are given by the general form  $\nabla \phi \cdot \vec{n} = \beta p$ , where  $\vec{n}$  is the unit outward normal on a given surface, and  $\beta$  the surface admittance is defined for the chamber bounding surfaces below.

Nozzle entrance plane ( $Z = L$ )       $\beta = \beta_N$  (Nozzle admittance)

Radial cavity entrance ( $r = 1$ )       $\beta = \beta_C$  (Cavity Admittance)

(or  $y = 1, y = 0$ )

Baffle blade surfaces       $\beta = \beta_B$  (Baffle admittance due to  
dissipation of energy in the  
boundary layer)

Injector surface ( $Z = 0$ )       $\beta = 0$  (Distributed combustion)

$$\beta = M \left( \frac{\gamma+1}{\gamma} - \eta (1 - e^{-i\omega\tau}) \right)$$

(Concentrated Combustion)

All other surfaces       $\beta = 0.$

### 2.3 Solution Technique

An integral technique is followed in order to predict either complex frequency ( $\omega_R$  and  $\lambda$ ) for a given  $n$  and  $\tau$ , or to predict  $n$  and  $\tau$  values required for neutral stability when the frequency ( $\omega_R$ ) is given. Consider Equation (1) written as a homogeneous equation with homogeneous boundary conditions

$$\nabla^2 \tilde{\phi} + \tilde{\omega}^2 \phi = 0$$

$$\nabla \tilde{\phi} \cdot \vec{n} = 0$$

all surfaces, including  
baffle blade surfaces

is the solution to this homogeneous problem, and  $\omega$  is the associated frequency. then, using Greens Theorem it can be shown that

$$\begin{aligned} (\omega^2 - \tilde{\omega}^2) \int_V \tilde{\phi} \phi dV &= - \int_S \tilde{\phi} \nabla \phi \cdot \vec{n} dS \\ &+ \int_V \tilde{\phi} F_1(\tilde{\phi}, \bar{u}, \omega, n, \tau) dV \end{aligned} \quad (2)$$

Equation (2) is exact and would determine either  $\omega$  or  $n$  and  $\tau$  exactly, if  $\tilde{\phi}$ ,  $\tilde{\omega}$ , and the functional form of  $\phi$  were known. It is possible

to determine  $\tilde{\phi}$  and  $\tilde{\omega}$  to arbitrary accuracy using an eigenfunction matching technique which will be described shortly. In determining an appropriate form for  $\phi$ , reliance is placed on two characteristics of the problem as posed. First, the function  $F_1$  on the right-hand side of Equation (1), as well as the values of  $\beta$  on the active surfaces are small, usually less than the mean flow Mach number in size. Second, the form of  $\phi$  is generated through integration of the  $\beta$  dependent terms over the chamber bounding surfaces. Taken together these encourage using  $\tilde{\phi}$  as the approximate form for  $\phi$  in the integrals appearing in Equation (2). Certainly an error no larger than the mean Mach number squared will occur upon this substitution. In practice, experience indicates that the error is usually considerably less. This is so, for example, for the concentrated combustion problem without an acoustic absorber, or boundary layer dissipation, for which an exact solution to Equation (1) exists. Whether this exact solution or  $\tilde{\phi}$  is used in the integrals, affects predictions of  $n$  and  $\tau$  only to an amount which is always considerably smaller than the Mach number squared in our calculations. If one accepts the substitution of  $\tilde{\phi}$  for  $\phi$  in the volume integrals and of  $\beta p(\tilde{\phi})$  for  $\nabla \phi \cdot \vec{n}$  in the surface integrals, then performing the indicated integrations leads to an algebraic relationship between  $n$ ,  $\tau$ , and  $\omega$ . This relationship is complex and can be solved for  $n$  and  $\tau$  given  $\omega$  or for  $\omega$  given  $n$  and  $\tau$ .

#### Solution for $\tilde{\phi}$

The function  $\tilde{\phi}$  is represented formally by eigenfunction expansions in each baffle cavity and in the main chamber. The form of these expansions is given below.

#### Two dimensional thrust chamber

$$\text{Main chamber: } \tilde{\phi}_c = \sum_{m=0}^{m_c} B_m \cos m\pi y \frac{\cosh(m^2\pi^2 - \tilde{\omega}^2)^{1/2}(z-L)}{\cosh(m^2\pi^2 - \tilde{\omega}^2)^{1/2}(z_B-L)}$$

baffle compartment  $\mu$ :

$$\Phi^\mu = \sum_{n=0}^{n_B} A_n^\mu \cos n\pi Ny \frac{\cosh(n^2\pi^2 N^2 - \tilde{\omega}^2)^{1/2} z}{\cosh(n^2\pi^2 N^2 - \tilde{\omega}^2)^{1/2} z_B}$$

where  $n$  and  $m$  are integers,  $N$  is the number of baffle compartments and  $n_B$  and  $m_C$  determine the number of terms in the eigenfunction representation. The coefficient vectors  $B_m$  and  $A_n^\mu$  (one  $A_n^\mu$  vector for each baffle compartment, not generally the same) as well as the frequency eigenvalue,  $\tilde{\omega}$  are determined by requiring that, at the interface between the main chamber and each baffle compartment,  $\tilde{\Phi}^\mu = \tilde{\Phi}^C$ , and  $\frac{\partial \tilde{\Phi}^\mu}{\partial z} = \frac{\partial \tilde{\Phi}^C}{\partial z}$ . In addition the dominant mode of oscillation in the main chamber is designated by setting  $B_{\hat{m}}^\Lambda = 1$ , where  $\hat{m}$  is arbitrary. Thus, the choice  $\hat{m} = 1$ , a first transverse type oscillation, would cause  $B_1 = 1$ . The other  $B_m$  would then represent corrections to the pure first transverse mode due to the presence of the baffle cavities. The coefficients are calculated using a successive approximation technique. It is first assumed that  $B_{\hat{m}}^\Lambda = 1$ , and  $\omega = \hat{m}\pi$ , while all other  $B_m = 0$ . The matching condition  $\tilde{\Phi}^\mu = \tilde{\Phi}^C$  then determines the  $A_n^\mu$  values for each baffle cavity. Using these values of  $A_n^\mu$ , values for all  $B_m$  (except  $B_{\hat{m}}^\Lambda$ , of course) are determined using the condition  $\frac{\partial \tilde{\Phi}^\mu}{\partial z} = \frac{\partial \tilde{\Phi}^C}{\partial z}$  at the matching plane. Finally the condition  $B_{\hat{m}}^\Lambda = 1$  is used to determine an improved approximation for  $\tilde{\omega}$ . The details of this type of matching solution are quite messy; an in-depth presentation is given in Reference 1. In general the convergence is rapid (less than 5 iterations). The number of iterations desired is specified by the program user, as are the values for  $m_C$  and  $n_B$ . Experience indicates that a choice of  $m_C = 30$  and  $n_B = 30$  give excellent matching and wave shape information.

### Three Dimensional Thrust Chamber

Main chamber:

$$\Phi_c = \sum_{m=0}^{m_c} \sum_{l=1}^{l_c} B_{lm} \cos m\theta J_m(\lambda_{lm}^c r) \frac{\cosh(\lambda_{lm}^c - \tilde{\omega}^2)^{1/2} (z-L)}{\cosh(\lambda_{lm}^c - \tilde{\omega}^2)^{1/2} (z_B-L)}$$

$J_m$  is Bessel function of order  $m$

$\lambda_{lm}^c$  is the  $l^{\text{th}}$  solution to  $\left(\frac{dJ_m}{dr}\right)_{r=1} = 0$

(Note: for a traveling wave in the main chamber  $\cos m\theta$  is replaced by  $e^{im\theta}$ )

Baffle compartment  $\mu$ :

$$\tilde{\Phi}^\mu = \sum_{m'=0}^{m_0} \sum_{l'=0}^{l_0} A_{l'm'}^\mu \cos \frac{m'N}{2} \theta J_{\frac{m'N}{2}}(\lambda_{l'm'}^B r) \frac{\cosh(\lambda_{l'm'}^B - \tilde{\omega}^2)^{1/2} z}{\cosh(\lambda_{l'm'}^B - \tilde{\omega}^2)^{1/2} z_B}$$

$J_{\frac{m'N}{2}}$  is Bessel function of order  $\frac{m'N}{2}$

$\lambda_{l'm'}^B$  is the  $l'^{\text{th}}$  solution to  $\left(\frac{dJ_{\frac{m'N}{2}}}{dr}\right)_{r=1} = 0$

The same matching requirements and successive approximation technique is used for the three dimensional thrust chamber as was used for the two dimensional chamber. In the three dimensional case  $B_{lm}$  and  $A_{l'm'}^\mu$  are matrices rather than vectors. In order to specify the main oscillation mode in the chamber both  $\hat{m}$  and  $\hat{l}$  must be chosen. For example  $\hat{m} = 1$ ,  $\hat{l} = 1$  leads to a 1st tangential type oscillation with  $\lambda_{11}^c = 1.8412$ ;  $\hat{m} = 2$ ,  $\hat{l}'$

= 1 a second tangential mode, etc. Choices for  $m_c$ ,  $l_c$ ,  $l'_B$ ,  $m'_B$  must be made as well. In the 3-D case computation time is a factor, and a compromise between accuracy and run time must be made. Choices of  $m_c = m'_B = 11$ ,  $l_c = l'_B = 4$  appear to give good results. The details and integrals involved in the 3-D matching problem are formidable; one is again referred to Reference 1.

A limit on the number of iterations must be specified by the user. Convergence is a little slower in the three dimensional model: approximately 10 iterations gives reliable convergence in most cases.

#### Asymptotic solution for $\tilde{\phi}$

Near the baffle blade tips it is desirable to have great accuracy in defining the local behavior of  $\tilde{\phi}$ . This is so because the outer inviscid solution will drive the boundary layer dissipation integral. An asymptotic solution which is valid near the blade tips was derived and explained in Reference 1. This is used without modification in the present work. Matching between the eigenfunction representations and the asymptotic solution is quite good over the region near the blade tips, if a matching at two points to determine constants in the asymptotic solution is made at a distance approximately equal to the baffle blade thickness  $T$ , on either side of the baffle blade.

#### Boundary layer dissipation integral

The portion of the boundary integral  $-\int_S \tilde{\phi} \nabla \phi \cdot \vec{n} ds$  of Equation (2) which lies on the baffle blade surfaces can be shown to be equivalent to  $E_{diss}^T$  the turbulent boundary layer dissipation integral.  $E_{diss}^T$  in turn is equal to

$$-\gamma \int_{S_{B.L.}} \left( \frac{\mu_{eff} \omega}{2} \right)^{1/2} |u'|^2 ds$$

where  $\mu_{eff} = C_{turb} \left[ \bar{u}^2 + \frac{|u'|^2}{2} \right]^{1/2}$

$$C_{turb} = .05 \text{ 3-Dimensional}$$

$$C_{turb} = .034 \text{ 2-Dimensional}$$

$|u'|$  is the modulus of the oscillatory velocity at the outer edge of the boundary layer as determined by the  $\tilde{\Phi}$  solution.

#### Application of Equation (2) and Final Solution

The  $\tilde{\Phi}$  representation for both main chamber and baffle cavities is now substituted for  $\phi$  in the volume integrals of Equation (2). It is to be recognized that these integrals extend over both the main chamber and baffle cavities. As can be imagined these integrals are somewhat involved. Solution is pursued analytically where possible, numerically where necessary. The surface integrals except for the baffle blade surfaces are evaluated by replacing  $\nabla\phi \cdot \hat{n}$  with  $\beta \tilde{P}$  and integrating over the appropriate surface (nozzle entrance plane, injector, acoustic cavity interface). As discussed above the surface integral for the baffle blades is replaced by the turbulent boundary layer integral  $E_{diss}^T$ . This integral is performed numerically (one dimension for two dimensional thrust chamber, two dimensions for three dimensional thrust chamber), over a streamline defining the edge of the boundary layer. Once the integrations are complete the resulting algebraic expression is solved for either  $\omega_R$  and  $\lambda$  (option 1) or  $n$  and  $\tau$  (option 2).

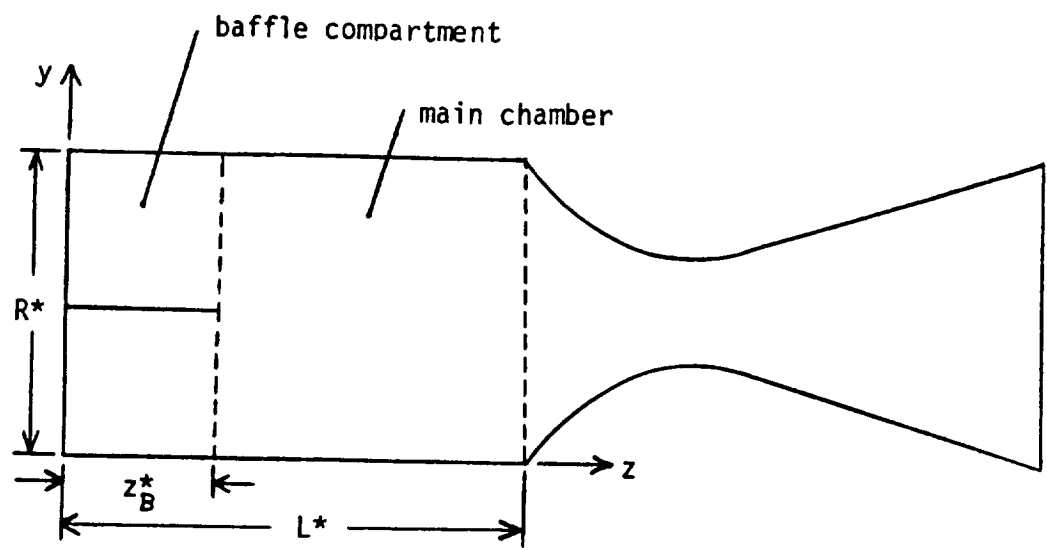


Figure 1a: Two-dimensional baffled chamber

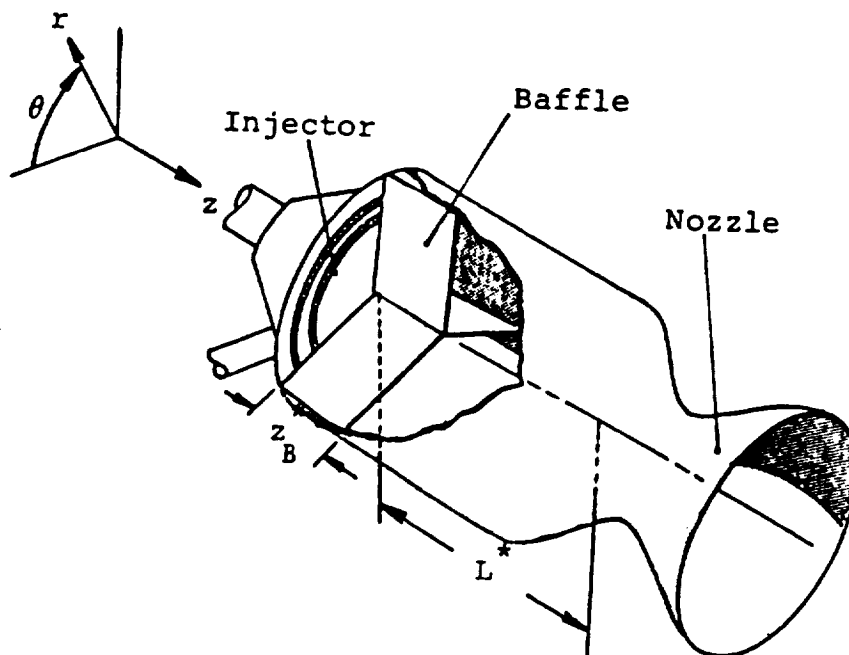


Figure 1b: Three-dimensional baffled chamber

### 3. - COMPUTER PROGRAMS

#### 3.1 - TWO-DIMENSIONAL PROGRAMS

This section will describe the two-dimensional distributed and concentrated combustion programs. A general overview of the distributed combustion program (DIST2D) will cover the program structure, input, output, and results. Distributed combustion sample runs and listing are located in the appendix. A brief discussion of the two-dimensional concentrated combustion program (CON2D) will be presented last. This discussion will cover the program input and output. The appendix includes a sample run from (CON2D) and listing.

##### 3.1.1 - DISTRIBUTED COMBUSTION PROGRAM DESCRIPTION

The computer program DIST2D consists of a main program and fourteen sub-programs which are listed in Table 1. The program has two main running options; Option one requires the user to input an interaction index ( $n$ ) and combustion time lag ( $\tau$ ), and the resulting complex frequency is calculated. Option two requires the input of a frequency range and increment and the program generates an  $n, \tau$  stability map. Additional program options are the capability of placing a radial acoustic absorber in the chamber and of making pressure amplitude calculations at any location.

A description of the program structure follows. (see Figure 2 for program flow chart.) After all input variables and program options are read in the program proceeds with the zero order (closed baffled chamber,  $\tilde{\phi}$ ) solution. The iteration counter is initialized along with the chamber coefficient vector and an initial approximation to the frequency,  $\tilde{\omega}$  is made. The program then proceeds into a loop that iterates on frequency,  $\tilde{\omega}$  until a correct solution to the matching condition equation (Eq. 11 Ref. 1) converges. The converged frequency and iteration counter are then printed out. At this point one full

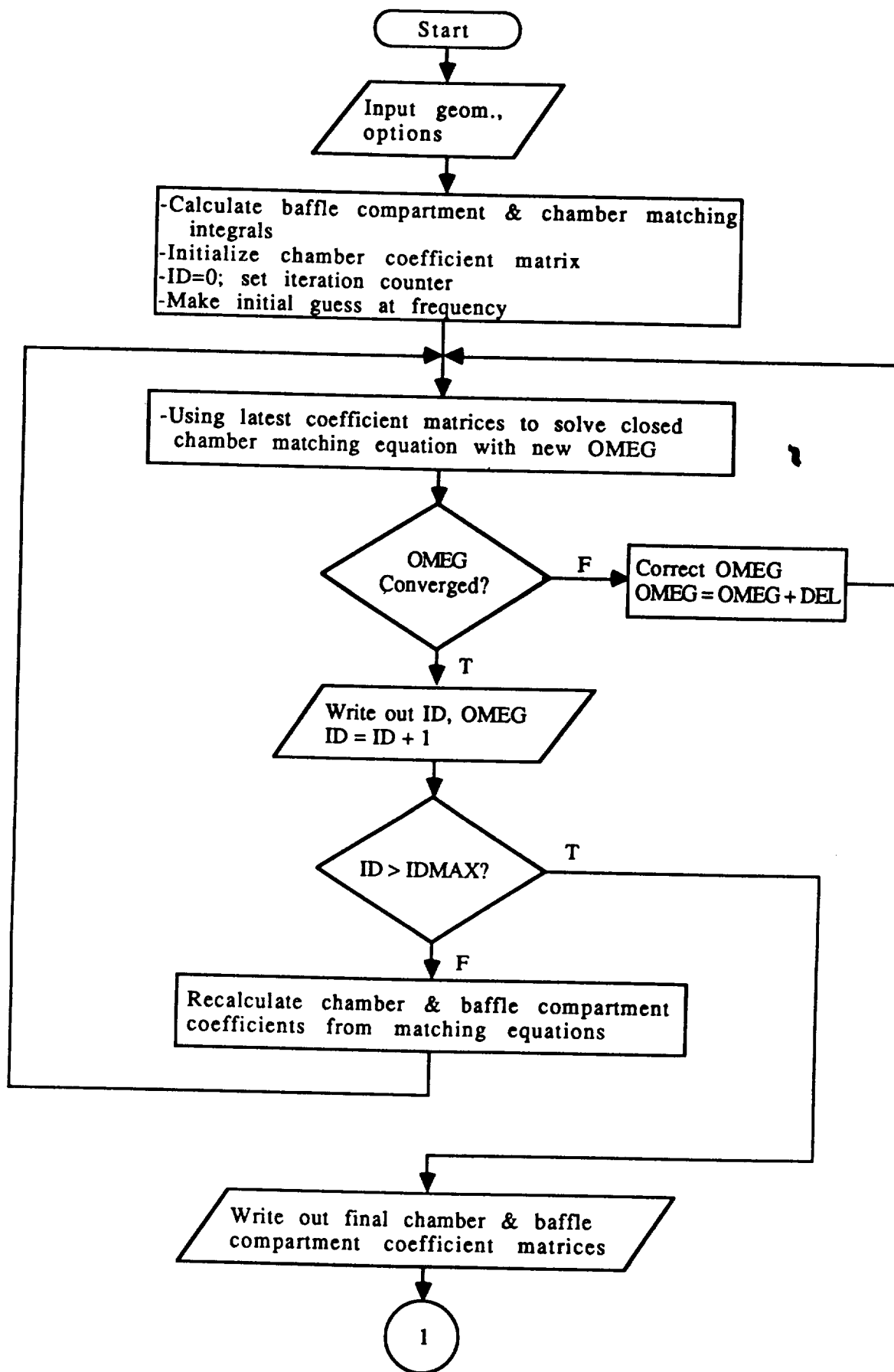


Figure 2: 2-D & 3-D program flowchart

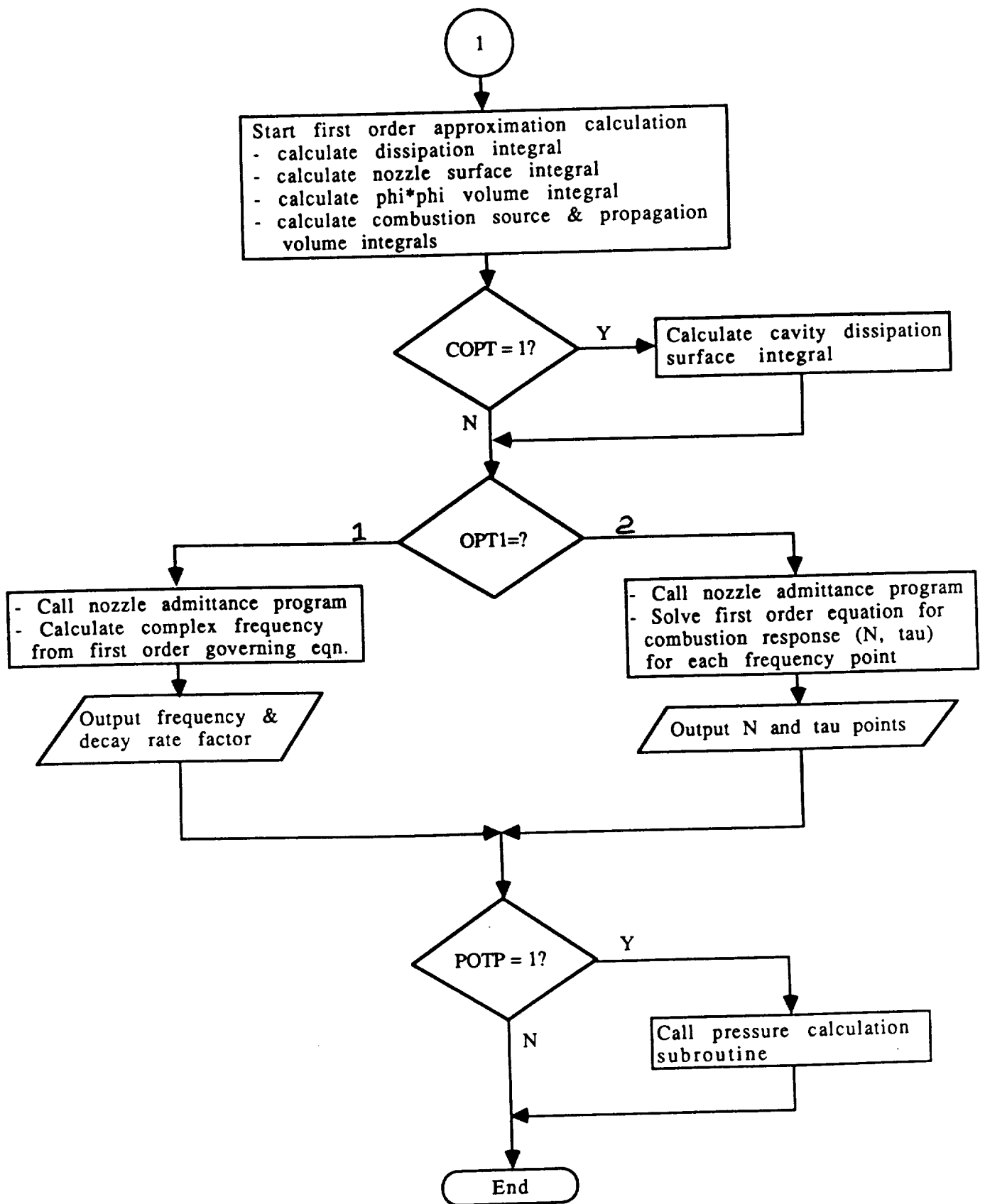


TABLE 1: subroutines and functions in DIST2D

Function CCOSH	- Performs hyperbolic cosine with a complex argument.
Subroutine NTAU1	- Calculates $n, \tau$ from combustion response.
Subroutine PRES	- Calculates pressure at a specific location in the chamber.
Function BI2	- Evaluates $\int z \phi \frac{d\phi}{dz} dz$ in the baffle compartments
Function CHI2	- Evaluates $\int z \phi \frac{d\phi}{dz} dz$ in the main chamber
Function BI3	- Evaluates $\int \phi \frac{d\phi}{dz} dz$ in the baffle compartments
Function BI1	- Evaluates $\int \phi \phi dz$ in the baffle compartments
Function CHI3	- Evaluates $\int \phi \frac{d\phi}{dz} dz$ in the main chamber
Function CHI1	- Evaluates $\int \phi \phi dz$ in the main chamber
Subroutines NOZINI, NOZADM, INTGRT, MACH, CALADM are used in the calculation of the nozzle admittance and were obtained from Refs. 4 & 5.	

iteration is complete and a check on the iteration limit is made. If the iteration limit has not been exceeded the main chamber and baffle compartment coefficients are recalculated (using Eqs. 9 & 10 Ref. 1) with the newly calculated frequency value, but if the iteration limit has been exceeded the coefficient vectors are printed out and the zero order calculation for  $\tilde{\phi}$  and  $\tilde{\omega}$  is complete.

At this point the program starts the solution computation for Equation (2) given in the Theory section above. Using the zero order velocity potential solution all the chamber surface and volume integrals are evaluated. These integrals include the combustion source and propagation volume integrals, the  $\tilde{\phi}$  squared volume integral, the nozzle surface integral, baffle blade dissipation integral, and cavity absorber surface integral if applicable. Depending on the option being run the program either solves the govern-

ing equation for complex frequency (Option 1) or solves the equation for combustion response (Option 2). Finally, the pressure calculation subroutine is called if this option is desired.

### 3.1.2 - DISTRIBUTED COMBUSTION PROGRAM INPUT

All input variables for the main program are read in from file 'DISIN2'. A list of the inputs is described in Table 2. In the file the variables appear in the same order as in Table 2 and need only be separated by commas.

The variables OPT1, COPT, POPT are the option variables. If variable (OPT1 = 1) option one, which calculates complex frequency from input values of  $n$  and  $r$ , is executed. If the variable (OPT1 = 2) option two is executed which generates a stability map. A radial cavity absorber is present if the variable (COPT = 1). The cavity admittance must be supplied along with the aperture width. The absorber is assumed to be located at the injector ( $z = 0$ ) and extends downstream from there. If the pressure option is to be run, set variable (POPT = 1). A separate input file called 'PRESPT2' must be supplied. This file described in Table 3 contains the total number of Z,Y locations to be calculated on the first line followed by each coordinate pair on successive lines.

TABLE 2: Input variables for file 'DISIN2'

<u>VARIABLE</u>	<u>VARIABLE DESCRIPTION</u>	<u>DIMENSIONS</u>
MC	: number of series terms to represent series solution in the main chamber (maximum of 50 terms. default value of 30 terms)	( none )
MB	: number of series terms to represent series solution in the baffle compartments (maximum of 50 terms. default value of 30 terms)	( none )
ALENGTH	: chamber length	( ft )
ZB	: baffle blade length	( ft )
T	: baffle blade thickness	( ft )
R	: chamber whole height	( ft )
MUB	: number of evenly spaced baffle compartments (maximum of 5 compartments)	( none )

```

HST      : nozzle throat radius ( half height )           ( ft )
RC       : radius of curvature at nozzle throat           ( ft )
RE       : radius of curvature at nozzle entrance         ( ft )
ALPHA    : nozzle convergence half angle                 ( deg )
AO       : chamber speed of sound at stagnation conditions ( ft/s )
PO       : chamber pressure at stagnation conditions      ( psf )
GAMMA    : ratio of specific heats                       ( none )
PAMP     : peak to peak pressure amplitude ( percent of main chamber )
MHAT     : dominating transverse mode number in main chamber ( none )
IDMAX    : maximum number of frequency iterations for successive approximation (default value of 5) ( none )
ZS       : z location where combustion starts            ( ft )
ZE       : z location where combustion is completed       ( ft )
OPT1     : option selection as described above ( 1 OR 2 )

IF (OPT1 - 1) THEN

VAL1     : combustion interaction index (n)                ( none )
VAL2     : combustion time lag (τ)                        ( msec )
VAL3     : not used in this option

IF (OPT1 - 2) THEN

VAL1     : starting frequency                             (hertz )
VAL2     : ending frequency                               (hertz )
VAL3     : frequency increment                            (hertz )

COPT     : cavity option ( 0 if no cavity, 1 if cavity present )
BETACR   : real part of cavity admittance                 (ft/s)/(psf)
BETACI   : imaginary part of cavity admittance            (ft/s)/(psf)
ZA       : z location where cavity ends (aperture width) ( ft )
POPT     : pressure option ( 1 if pressure points are to be calculated,
           0 if no pressure calculations are to be made)

```

TABLE 3: Input variables for file 'PRESPT2'

<u>VARIABLE</u>	<u>VARIABLE DESCRIPTION</u>	<u>DIMENSIONS</u>
N	: number of points to be calculated	( none )
Z,Y	: N number of sets of z,y locations for pressure calculations	( ft )

### 3.1.3 - DISTRIBUTED COMBUSTION PROGRAM OUTPUT

The output file generated is called 'DISOUT2'. This file starts with the chamber geometry and operating conditions, this includes calculated values for nozzle inlet Mach number, steady state pressure and sound speed. For each  $\omega$  frequency iteration the iteration number and frequency is printed out. After

the iteration limit is reached the final main chamber and baffle compartment Fourier coefficient vectors are printed out. Next, a calculation of nozzle admittance based on  $\tilde{\omega}$  is printed out. At this point if Option one is selected the chamber complex frequency is printed out ( $\omega_R + i\lambda$ ), this is also represented by a decay/growth rate factor which gives the ratio of the amplitude after one period to that of the previous period. Decay rate ( $\lambda$ ) is also represented in decibels per cycle. If Option two is selected the output consist of a list containing frequency, interaction index (n), and combustion time lag ( $\tau$ ). For Option two an additional output file called 'NTDATA2' is generated which contains  $\tau, n$  values for plotting purposes.

#### 3.1.4 - DISTRIBUTED COMBUSTION PROGRAM RESULTS AND DISCUSSION

The following section is intended to show some of the capabilities and predictions of the distributed combustion program. A series of  $n, \tau$  plots show the effects of baffle blade length, combustion zone variations, and acoustic absorbers. A pressure profile is also included which shows the effects of baffle blade length.

Run time for the two dimensional program does not present a problem. Run time for a typical test case in which 30 term vectors were kept required approximately 75 CPU seconds on the VAX 11/780 machine at Colorado State University.

The baffled chamber geometry and other input parameters used for the series of plots are listed in Table 4. The effect of baffle length on pressure amplitude is shown in Figure 3. The long baffle (33.3% of chamber length) induced a large pressure amplitude decrease from the injector face ( $z = 0$ ) to the nozzle ( $z = \text{ALENGTH}$ ) while the short (6.67%) baffle had only a minor effect. A chamber with no baffle at all would show no amplitude decrease. This result is important as an aid in understanding other stability

predictions to be discussed. The effect of moving a concentrated combustion zone down the chamber is illustrated in Figure 4 for a baffle blade length that is 6.67% of the chamber length. The same combustion zone movement is shown in Figure 5 for a baffle blade length which is 33.3% of the chamber length. In both cases (ZE - ZS) = .048ft, and ZS is varied from 0 to 1.632ft. For the 6.67% baffle moving the combustion zone downstream does not have a large effect on stability. This result is due to the fact that the pressure amplitude does not change significantly downstream. The large pressure amplitude decrease with Z for the 33.3% baffle explains the large shifts in the  $n, r$  curves as the concentrated combustion zone is moved toward the nozzle. In interpreting these results it must be remembered that combustion input is proportional to local pressure amplitude, in the  $n, r$  model. For all combustion zone locations the large baffle causes a stabilizing shift upward of the  $n, r$  curves as well as a potentially destabilizing flattening of the curves. Figure 6 shows a set of curves for a 33.3% baffle with a combustion zone starting at the injector and extending different distances downstream (i.e. ZS = 0, ZE = .408, 1.02, 2.04 (ft)). As can be readily seen distributing the combustion has a stabilizing effect. The final plot shows the clear stabilizing effect of an acoustic absorber. The absorber used had a slot width of .136ft and a pure real admittance of .04222(ft/s)/(psf).

TABLE 4: Input parameters used for 2-D plots

MC	- 30	RE	- .11 ft
MB	- 30	ALPHA	- 30 deg
ALENGTH	- 2.04 ft	AO	- 3850 ft/sec
T	- .068 ft	PO	- 43,200 psf
R (chamber height)	- 1.36 ft	GAMMA	- 1.2
MUB	- 2 compartments	PAMP	- 20
HST	- .25 ft	MHAT	- 1
RC	- .25 ft	IDMAX	- 4
MACH	- .2232		

### 3.1.5 - CONCENTRATED COMBUSTION PROGRAM DESCRIPTION

The two-dimensional concentrated combustion program is in essence the same program that appears in Reference 1. The following modifications have been made. First, The corrected concentrated combustion model described in the Theory section has been implemented. Secondly, the nozzle admittance prediction model and program appearing in References 4 & 5 (Aerojet's NOZADM program) has been added. Finally, all of the nondimensional inputs and outputs have been dimensionalized. It should be noted that the two-dimensional concentrated program is only capable of predicting frequency and decay rate from a given  $n, r$ .

### 3.1.6 - CONCENTRATED COMBUSTION PROGRAM INPUT

All input variables for the main program are read in from file 'CON2IN'. A list of the required inputs is described in Table 5. The variables appear in the same order in Table 5 as they do in the file and need only to be separated by commas.

TABLE 5: Input variables for file 'CON2IN'

<u>VARIABLE</u>	<u>VARIABLE DESCRIPTION</u>	<u>DIMENSIONS</u>
MC	: number of series terms to represent series solution in the main chamber (maximum of 50 terms. default value of 30 terms)	( none )
MB	: number of series terms to represent series solution in the baffle compartments (maximum of 50 terms. default value of 30 terms)	( none )
IDMAX	: maximum number of frequency iterations for successive approximation (default value of 5)	( none )
ALENGTH	: chamber length	( ft )
ZB	: baffle blade length	( ft )
T	: baffle blade thickness	( ft )
DR	: chamber whole height	( ft )
MUB	: number of evenly spaced baffle compartments (maximum of 5 compartments)	( none )
HST	: nozzle throat radius ( half height )	( ft )
RC	: radius of curvature at nozzle throat	( ft )
RE	: radius of curvature at nozzle entrance	( ft )
ALPHA	: nozzle convergence half angle	( deg )

AO	: chamber speed of sound at stagnation conditions	( ft/s )
PO	: chamber pressure at stagnation conditions	( psf )
GAMMA	: ratio of specific heats	( none )
PAMP	: peak to peak pressure amplitude	( percent of main chamber )
MHAT	: dominating transverse mode number in main chamber	( none )
AN	: combustion interaction index (n)	( none )
TAU	: combustion time lag ( $\tau$ )	( msec )

### 3.1.7 - CONCENTRATED COMBUSTION PROGRAM OUTPUT

The output of the concentrated combustion program is written into file 'CON2OUT'. The output starts with a listing of the geometrical inputs and operating conditions such as nozzle inlet Mach number, steady state chamber pressure and sound speed. For each frequency iteration the program outputs the complex frequency (hertz) and iteration number. After the iteration limit is reached the final chamber and baffle compartment vectors are printed out. The final output includes the decay rate with baffle dissipation included, this is also presented as decay in decibels/cycle and the decay/growth rate factor.

ALNGTH = 2.04 ft      Y = 1.36 ft (top of chamber)      ZS = 0.0, ZE = 2.04

A: 33.3% baffle      B: 6.67% baffle

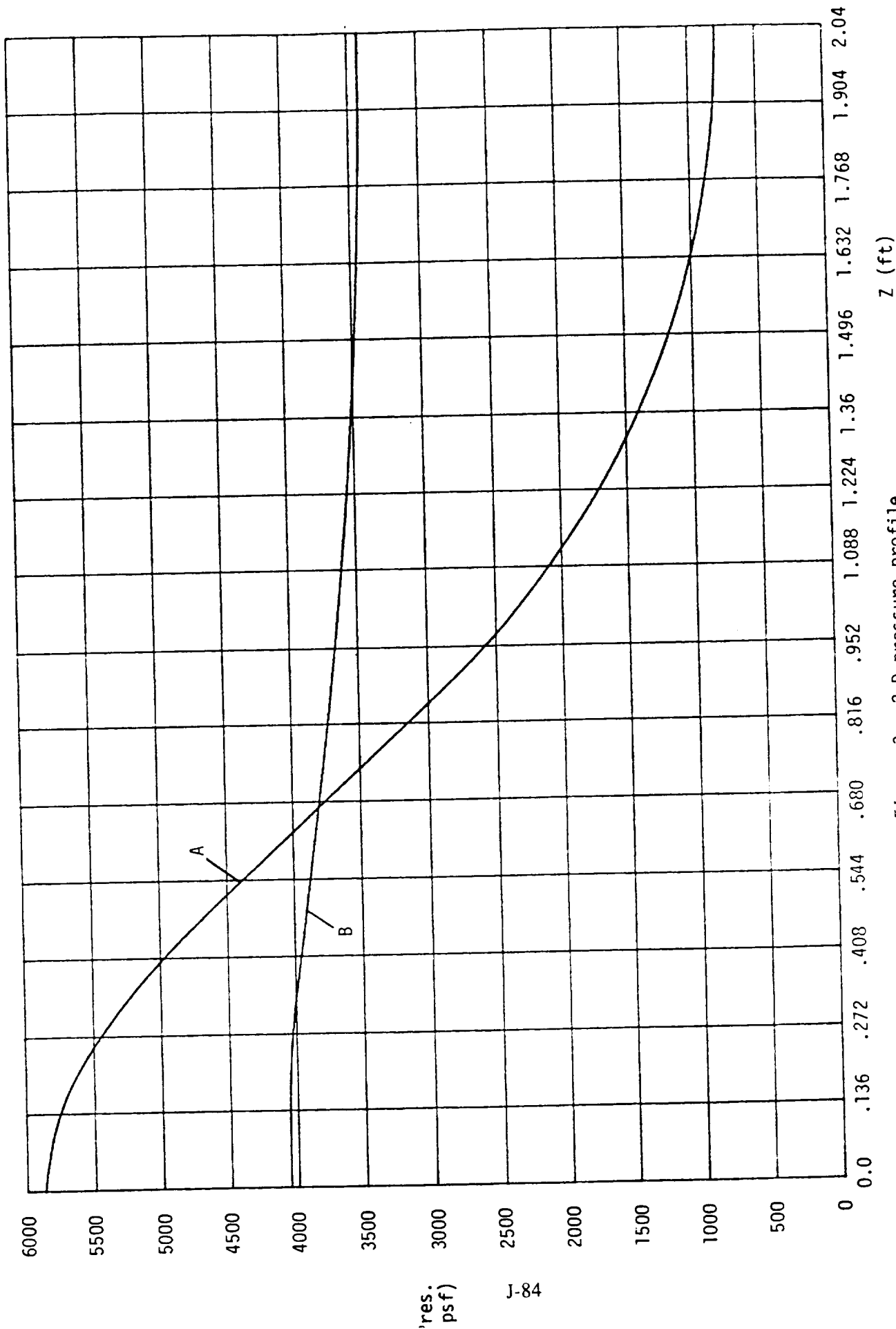


Figure 3: 2-D pressure profile

Figure 4

NEUTRAL STABILITY MAP FOR A 2-D CHAMBER, CONCENTRATED COMBUSTION ZONES

Chamber dimensions (ft): length = 2.04 whole height = 1.36  
Baffle dimensions (ft): length = 0.136 thickness = 0.068  
Steady state speed of sound = 3850.0 ft/s  
Steady state pressure = 43200.0 psf  
Ratio of specific heats = 1.2 Mode: First transverse  
Frequency range (Hz): 1200 to 1700 No. of baffles: 1

ZONES (distance in feet from injector plate):

A: ZS = 0.0 to ZE = 0.408 (no baffle present)  
B: ZS = 0.0 to ZE = 0.408  
C: ZS = 0.476 to ZE = 0.884  
D: ZS = 1.632 to ZE = 2.04

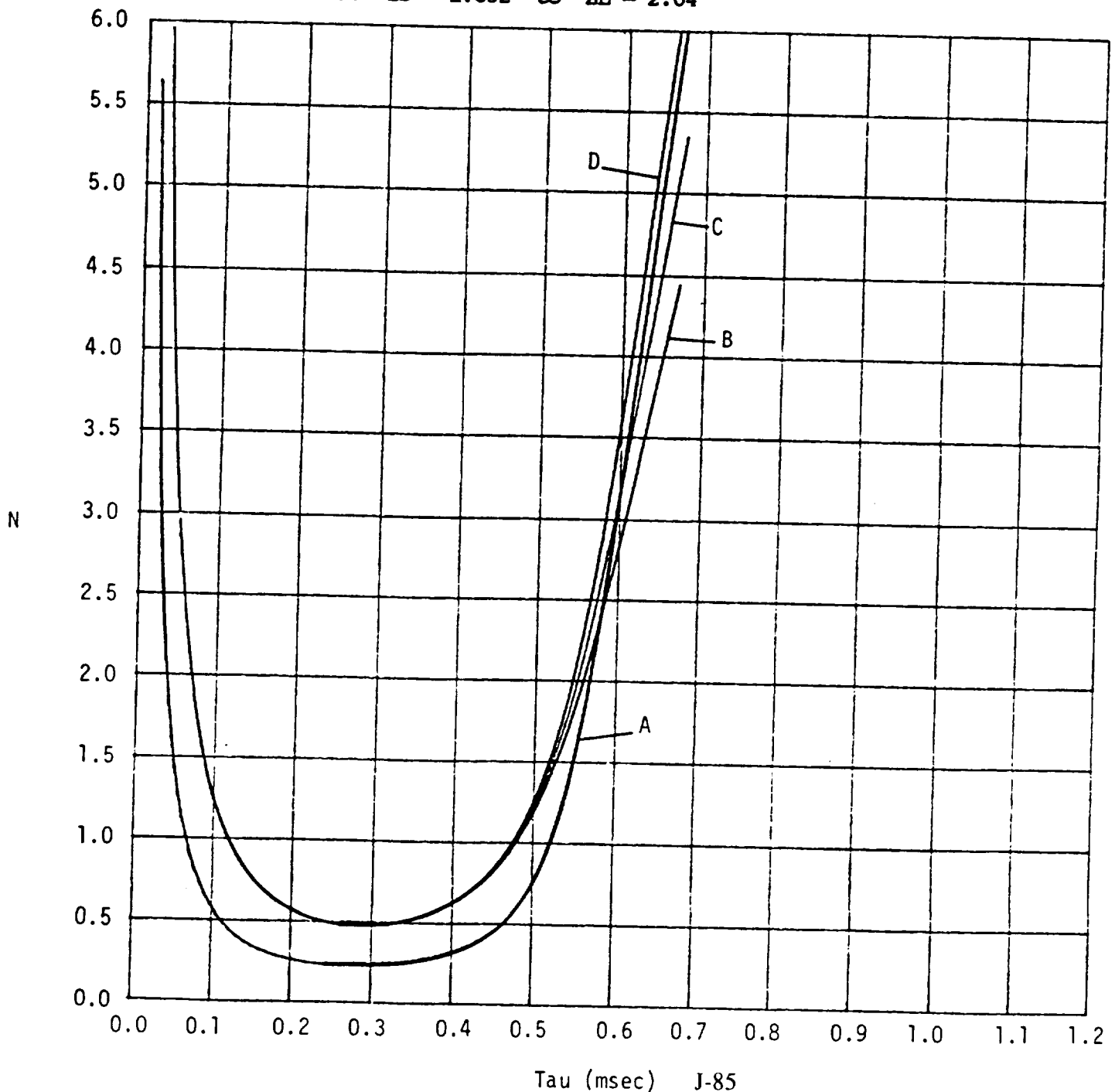


Figure 5

NEUTRAL STABILITY MAP FOR A 2-D CHAMBER, CONCENTRATED COMBUSTION ZONES

Chamber dimensions (ft): length = 2.04 whole height = 1.36  
 Baffle dimensions (ft): length = 0.68 thickness = 0.068  
 Steady state speed of sound = 3850.0 ft/s  
 Steady state pressure = 43200.0 psf  
 Ratio of specific heats = 1.2 Mode: First transverse  
 Frequency range (Hz): 700 to 1200 No. of baffles: 1

ZONES (distance in feet from injector plate):  
 A: ZS = 0.0 to ZE = 0.408  
 B: ZS = 0.476 to ZE = 0.884  
 C: ZS = 1.632 to ZE = 2.04

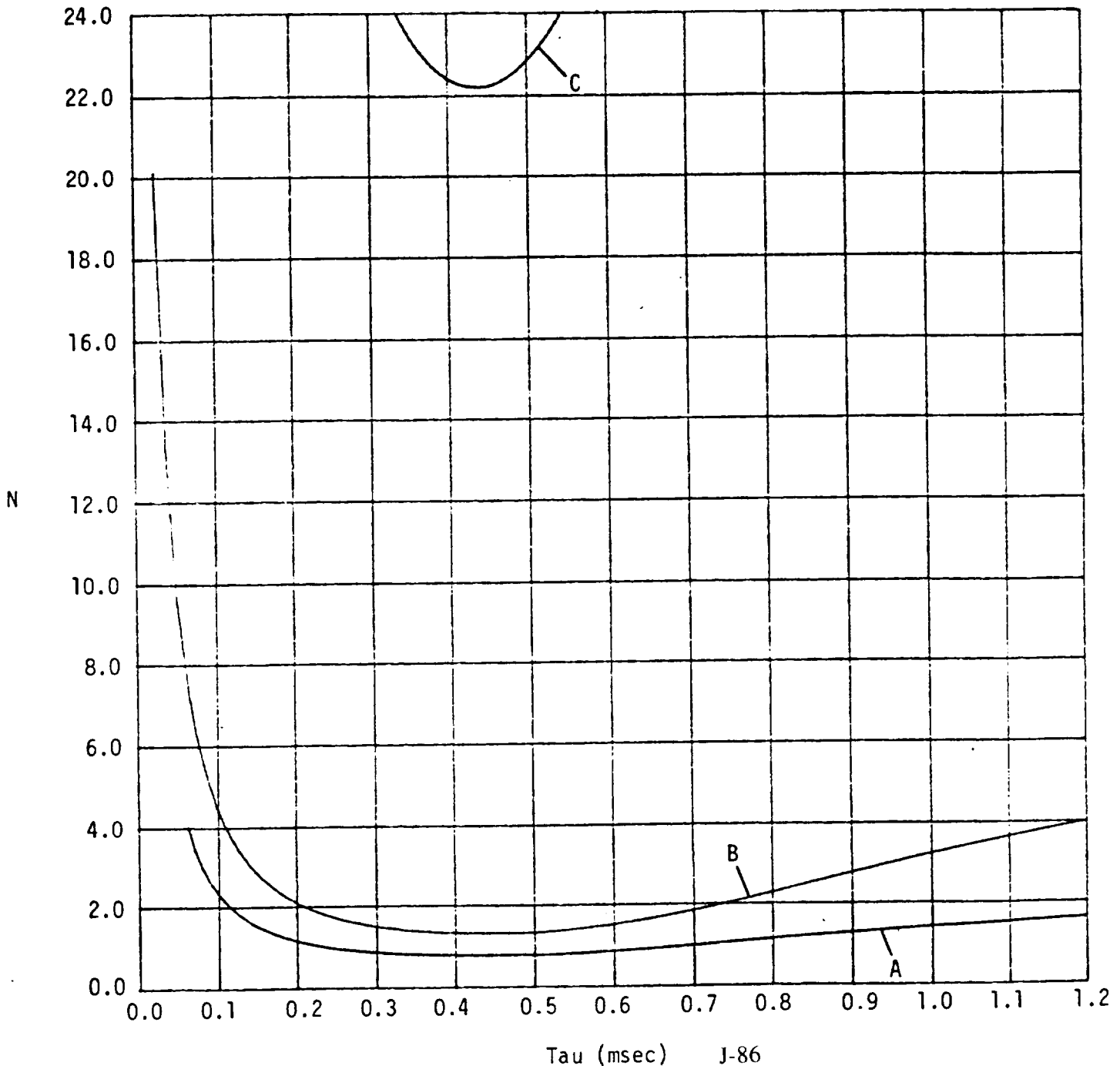


Figure 6

NEUTRAL STABILITY MAP FOR A 2-D CHAMBER, DISTRIBUTED COMBUSTION ZONES

Chamber dimensions (ft): length = 2.04 whole height = 1.36  
 Baffle dimensions (ft): length = 0.68 thickness = 0.068  
 Steady state speed of sound = 3850.0 ft/s  
 Steady state pressure = 43200.0 psf  
 Ratio of specific heats = 1.2 Mode: First transverse  
 Frequency range (Hz): 700 to 1200 No. of baffles: 1

ZONES (distance in feet from injector plate):

A: ZS = 0.0 to ZE = 2.04 (no baffle present)

B: ZS = 0.0 to ZE = 0.408

C: ZS = 0.0 to ZE = 1.02

D: ZS = 0.0 to ZE = 2.04

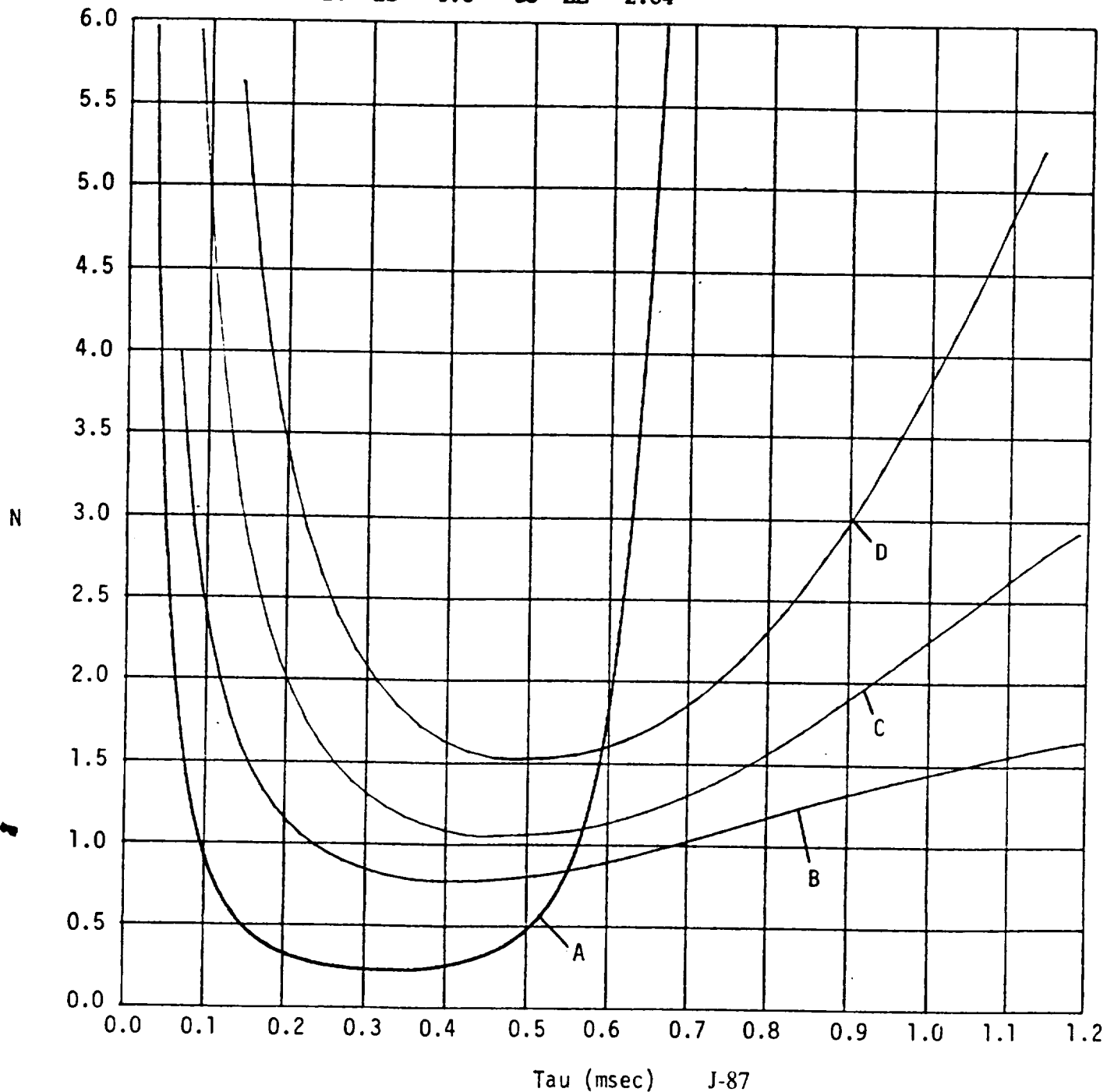


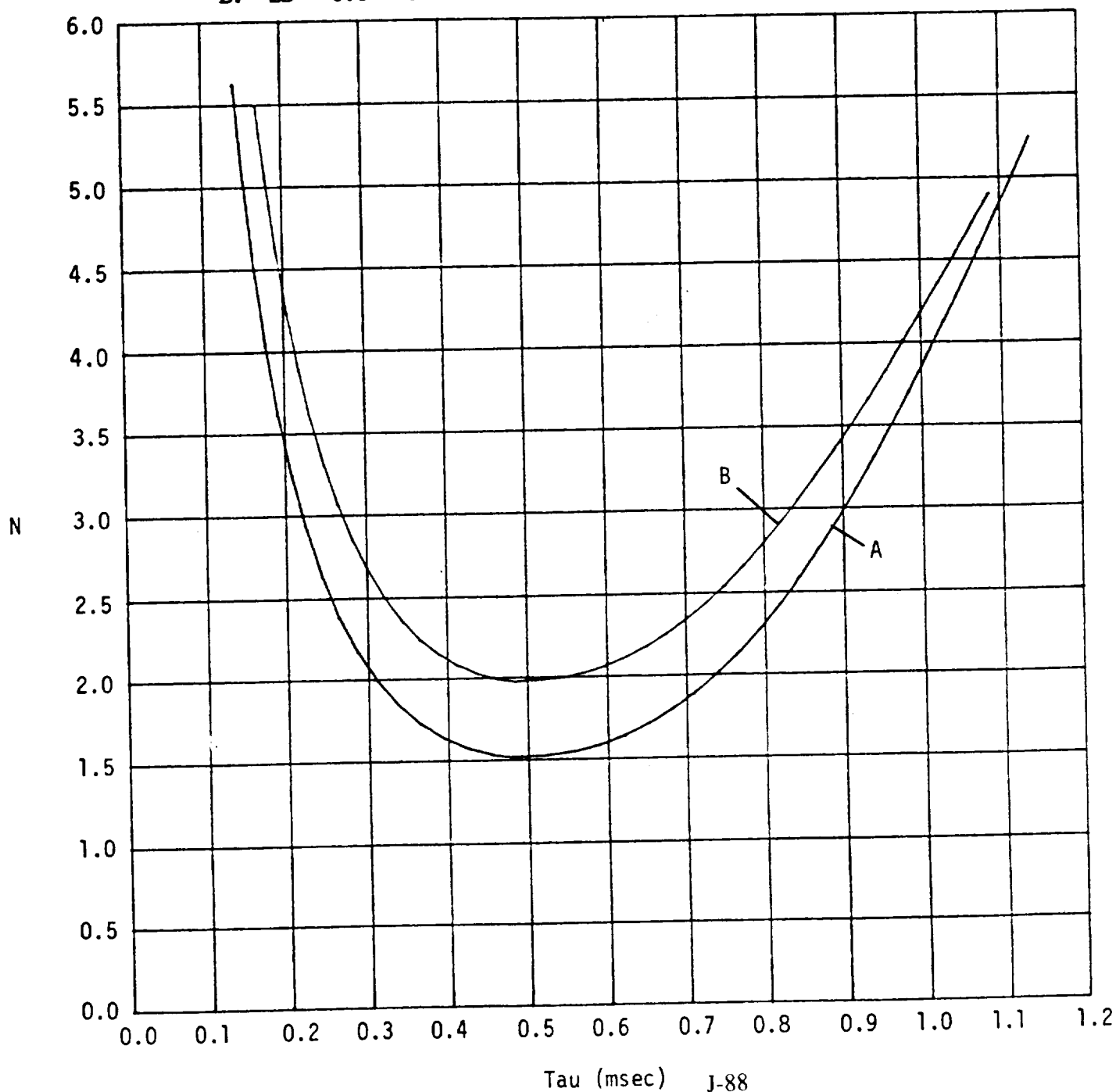
Figure 7

NEUTRAL STABILITY MAP FOR A 2-D CHAMBER, DISTRIBUTED COMBUSTION ZONES

Chamber dimensions (ft): length = 2.04      whole height = 1.36  
 Baffle dimensions (ft): length = 0.68      thickness = 0.068  
 Steady state speed of sound = 3850.0 ft/s  
 Steady state pressure = 43200.0 psf  
 Ratio of specific heats = 1.2      Mode: First transverse  
 Frequency range (Hz): 700 to 1200      No. of baffles: 1  
 Cavity admittance (ft/s)/(psf): real = 0.04222      imaginary = 0.0  
 Cavity aperture width = 0.136 ft

ZONES (distance in feet from injector plate):

A: ZS = 0.0 to ZE = 2.04 - without cavity dissipation  
 B: ZS = 0.0 to ZE = 2.04 - with cavity dissipation



### 3.2 - THREE-DIMENSIONAL PROGRAMS

This section will describe the three-dimensional distributed and concentrated combustion programs. The parallels between the two-dimensional and three-dimensional programs are strong, this fact will simplify the three-dimensional discussion somewhat. An overview of the distributed combustion program (DIST3D) will cover the program structure, input, output, and results. Three-dimensional distributed combustion program runs and listing are located in the appendix. A brief discussion of the three-dimensional concentrated combustion program (CON3D) will be presented last. This discussion will cover the program input, output, and current status. The appendix includes a sample run from (CON3D) and listing.

#### 3.2.1 - DISTRIBUTED COMBUSTION PROGRAM DESCRIPTION

The computer program DIST3D consists of a main program and seventeen subprograms which are listed in Table 6. The three-dimensional program like the two-dimensional has two main running options; Option one which calculates complex frequency from a given  $n, r$  point, and Option two which produces a  $n, r$  stability map. Additional program options include the capability of placing a radial acoustic absorber in the chamber and of making pressure amplitude and phase calculations at any location. The three-dimensional program has an additional option that determines the waveform type present in the main chamber. This waveform can either take the form of a spinning wave or standing wave.

A description of the program structure follows. (see Figure 2 for program flow chart.) After all input variables and program options are read in the program proceeds with the zero order (closed baffled chamber,  $\phi$ ) solution. The iteration counter is initialized along with the chamber coefficient matrix

TABLE 6: subroutines and functions in DIST3D

Function CCOSH	- Performs hyperbolic cosine with a complex argument.
Subroutine NTAU1	- Calculates $n, r$ from combustion response.
Subroutine PRES	- Calculates pressure at a specific location in the chamber.
Function BI2	- Evaluates $\int z \phi \frac{d\phi}{dz} dz$ in the baffle compartments
Function CHI2	- Evaluates $\int z \phi \frac{d\phi}{dz} dz$ in the main chamber
Function BI3	- Evaluates $\int \phi \frac{d\phi}{dz} dz$ in the baffle compartments
Function BI1	- Evaluates $\int \phi \phi dz$ in the baffle compartments
Function CHI3	- Evaluates $\int \phi \frac{d\phi}{dz} dz$ in the main chamber
Function CHI1	- Evaluates $\int \phi \phi dz$ in the main chamber
Function BESSCAL	- Calculates the value of the Bessel function of integer and half integer order and of arbitrary argument
Subroutine ROOT	- Calculates the root of the derivative of Bessel functions
Subroutine VDISP	- Calculated baffle blade tip dissipation
Subroutines NOZINI, NOZADM, INTGRT, MACH, CALADM are used in the calculation of the nozzle admittance and were obtained from Refs. 2 & 3.	

and an initial approximation to the frequency,  $\tilde{\omega}$  is made. The program then proceeds into a loop that iterates on frequency,  $\tilde{\omega}$  until a correct solution to the matching condition equation (Eq. 23 Ref. 1) converges. The converged frequency and iteration counter are then printed out. At this point one full iteration is complete and a check on the iteration limit is made. If the iteration limit has not been exceeded the main chamber and baffle compartment coefficients are recalculated (using Eqs. 21 & 22 Ref. 1) with the newly calculated frequency value, but if the iteration limit has been exceeded the coefficient matrices are printed out and the zero order calculation for  $\tilde{\phi}$  and  $\tilde{\omega}$  is complete.

At this point the program starts the solution computation for Equation (2) given in the Theory section above. Using the zero order velocity potential solution all the chamber surface and volume integrals are evaluated. These integrals include the combustion source and propagation volume integrals, the  $\bar{\phi}$  squared volume integral, the nozzle surface integral, baffle blade dissipation integral, and cavity absorber surface integral if applicable. Depending on the option being run the program either solves the governing equation for complex frequency (Option 1) or solves the equation for combustion response (Option 2). Finally, the pressure calculation subroutine is called if this option is desired.

### 3.2.2 - DISTRIBUTED COMBUSTION PROGRAM INPUT

All input variables for the main program are read in from file 'DISIN3'. A list of the inputs is described in Table 7. In the file the variables appear in the same order as in Table 7 and need only be separated by commas.

The variables OPT1, COPT, POPT, MX are the option variables. If variable (OPT1 = 1) program Option one is executed, if (OPT1 = 2) program Option two is executed. A radial cavity absorber is present if the variable (COPT = 1). The cavity admittance must be supplied along with the aperture width. The absorber is assumed to be located at the injector ( $z = 0$ ) and extends downstream from there. If the pressure option is to be run, set variable (POPT = 1). A separate input file called 'PRESPT3' must be supplied. This file described in Table 8 contains the total number of R, THETA, Z locations to be calculated on the first line followed by each coordinate pair on successive lines. The variable MX determines the waveform type in the main chamber, (MX = 0) for standing waves and (MX = 1) for spinning waves.

TABLE 7: Input variables for file 'DISIN3'

<u>VARIABLE</u>	<u>VARIABLE DESCRIPTION</u>	<u>DIMENSIONS</u>
MX	: variable that determines what type of main chamber solution is present. (MX = 0 for standing waves, MX = 1 for spinning waves)	( none )
MC	: number of Fourier series terms to represent the main chamber solution (maximum of 20 terms. default value of 11 terms)	( none )
LC	: number of Bessel series terms to represent solution in the main chamber (maximum of 20 terms. default value of 8 terms)	( none )
MB	: number of Fourier series terms to represent the baffle compartment solution (maximum of 20 terms. default value of 11 terms)	( none )
LC	: number of Bessel series terms to represent solution in the baffle compartments (maximum of 20 terms. default value of 8 terms)	( none )
IDMAX	: maximum number of frequency iterations for successive approximation (default value of 9)	( none )
ALENGTH	: chamber length	( ft )
ZB	: baffle blade length	( ft )
T	: baffle blade thickness	( ft )
RCHAMB	: chamber radius	( ft )
MUB	: number of evenly spaced baffle compartments (maximum of 12 compartments)	( none )
HST	: nozzle throat radius	( ft )
RC	: radius of curvature at nozzle throat	( ft )
RE	: radius of curvature at nozzle entrance	( ft )
ALPHA	: nozzle convergence half angle	( deg )
AO	: chamber speed of sound at stagnation conditions	( ft/s )
PO	: chamber pressure at stagnation conditions	( psf )
GAMMA	: ratio of specific heats	( none )
PAMP	: peak to peak pressure amplitude ( percent of main chamber )	( none )
MHAT	: dominating transverse mode number in main chamber	( none )
LHAT	: dominating radial mode number in main chamber	( none )
ZS	: z location where combustion starts	( ft )
ZE	: z location where combustion is completed	( ft )
OPT1	: option selection as described above ( 1 OR 2 )	
IF (OPT1 = 1) THEN		
VAL1	: combustion interaction index (n)	( none )
VAL2	: combustion time lag ( $\tau$ )	( msec )
VAL3	: not used in this option	
IF (OPT1 = 2) THEN		
VAL1	: starting frequency	(hertz )
VAL2	: ending frequency	(hertz )
VAL3	: frequency increment	(hertz )
COPT	: cavity option ( 0 if no cavity, 1 if cavity present )	

BETACR : real part of cavity admittance (ft/s)/(psf)  
 BETACI : imaginary part of cavity admittance (ft/s)/(psf)  
 ZA : z location where cavity ends (aperture width) ( ft )  
 POPT : pressure option ( 1 if pressure points are to be calculated,  
 0 if no pressure calculations are to be made)

TABLE 8: Input variables for file 'PRESPT3'

<u>VARIABLE</u>	<u>VARIABLE DESCRIPTION</u>	<u>DIMENSIONS</u>
N	: number of points to be calculated : N number of sets of r,theta,z locations for pressure calculations	( none )
R	: radius	( feet )
THETA	: angle counterclockwise from first baffle blade	( deg. )
Z	: axial distance downstream from injector	( feet )

### 3.2.3 - DISTRIBUTED COMBUSTION PROGRAM OUTPUT

The output file generated is called 'DISOUT3'. This file starts with the chamber geometry and operating conditions, this includes calculated values for nozzle inlet Mach number, steady state pressure and sound speed. For each  $\omega$  frequency iteration the iteration number and frequency is printed out. After the iteration limit is reached the final main chamber and baffle compartment Fourier-Bessel coefficient matrices are printed out. Next, a calculation of nozzle admittance based on  $\tilde{\omega}$  is printed out. At this point if Option one is selected the chamber complex frequency is printed out ( $\omega_R + i\lambda$ ), this is also represented by the decay/growth rate factor. If Option two is selected the output consist of a list containing frequency, interaction index (n), and combustion time lag ( $\tau$ ). For Option two an additional output file called 'NTDATA3' is generated which contains  $\tau, n$  values for plotting purposes.

### 3.2.4 - DISTRIBUTED COMBUSTION PROGRAM RESULTS AND DISCUSSION

The following section is intended to show some of the capabilities and predictions of the three-dimensional distributed combustion program. The same series of n, $\tau$  plots that were presented in the two-dimensional section are

presented here.

Run time for the three-dimensional program can present a problem if large matrices are kept. Table 9 contains run time data for the three-dimensional program run in Option two with variable matrix sizes and constant iteration limit of five. All timings were made from the VAX 11/780 machine at Colorado State University. The convergence of the results for these runs was very good, that is the smaller matrices results compared well to those obtained with the larger matrices. This result, though, is for only one specific case and should not be assumed true for all cases.

TABLE 9: run times for program 'DIST3D'

MATRIX SIZE		CPU TIME
<u>MC</u> & <u>MB</u>	<u>LC</u> & <u>LB</u>	
20	20	4.6 hours
11	8	13 min.
11	4	3.4 min.
8	4	1.93 min.

The baffled chamber geometry and other input parameters used for the series of plots are listed in Table 10. The effect of baffle length on pressure amplitude is shown in Figure 8. The long baffle (33.3% of chamber length) induced a large pressure amplitude decrease from the injector face ( $z = 0$ ) to the nozzle ( $z = \text{ALENGTH}$ ) while the short (6.67%) baffle had only a minor effect. This result is consistent with the two-dimensional predictions. The effect of moving a concentrated combustion zone down the chamber is illustrated in Figure 9 for a baffle blade length that is 6.67% of the chamber length. The same combustion zone movement is shown in Figure 10 for a baffle blade length which is 33.3% of the chamber length. In both cases ( $ZE - ZS$ ) equals .1869ft, and  $ZS$  is varied from 0 to 1.3706ft. For the 6.67% baffle moving the combustion zone downstream does not have a large effect on stability. Again this result is due to the fact that the pressure amplitude does

not change significantly downstream. The large pressure amplitude decrease with Z for the 33.3% baffle explains the large shifts in the  $n, r$  curves as the concentrated combustion zone is moved toward the nozzle. As in the two-dimensional case all the combustion zone locations for the large baffle cause a stabilizing shift upward of the  $n, r$  curves as well as a potentially destabilizing flattening of the curves. Figure 11 shows a set of curves for a 33.3% baffle with a combustion zone anchored at the injector and extending different distances downstream (i.e.  $ZS = 0$ ,  $ZE = .1869$ ,  $0.77875$ ,  $1.5575$  (ft)). As can be readily seen distributing the combustion has a stabilizing effect. The final plot shows stabilizing effect of an acoustic absorber. The absorber used had a slot width of  $.0623$  ft and a pure real admittance of  $.04467$  (ft/s)/(psf).

TABLE 10: Input parameters used for 3-D plots

MC	- 8	RE	- .216 ft
LC	- 2	MACH	- .06615
MB	- 8	ALPHA	- 45 deg
LC	- 2	MX	- 1 (spinning wave)
ALENGTH	- 1.5575 ft	AO	- 3850 ft/sec
T	- .03115 ft	PO	- 43,200 psf
RCHAMB	- .623 ft	GAMMA	- 1.2
MUB	- 3 compartments	PAMP	- 20
HST	- .208 ft	MHAT	- 1
IDMAX	- 8	LHAT	- 1
RC	- .216 ft		

### 3.2.5 - CONCENTRATED COMBUSTION PROGRAM DESCRIPTION

The three-dimensional concentrated combustion program is a modified version of the program that appears in Reference 1. The same modifications made to the two-dimensional program were made to the three-dimensional program and are listed here again. First, The corrected concentrated combustion model described in the Theory section has been implemented. Secondly, the axisymmetric nozzle admittance prediction model and program References 2 (Aerojet's

NOZADM program) has been added. Finally, all of the nondimensional inputs and outputs have been dimensionalized. The three-dimensional concentrated program is only capable of predicting frequency and decay rate from a given  $n, \tau$  point.

### 3.2.6 - CONCENTRATED COMBUSTION PROGRAM INPUT

All input variables for the main program are read in from file 'CON3IN'. A list of the required inputs is described in Table 11. The variables appear in the same order in Table 11 as they do in the file and need only to be separated by commas.

TABLE 11: Input variables for file 'CON3IN'

<u>VARIABLE</u>	<u>VARIABLE DESCRIPTION</u>	<u>DIMENSIONS</u>
MX	: variable that determines what type of main chamber waveform is present. (MX = 0 for standing waves, MX = 1 for spinning waves)	( none )
MC	: number of Fourier series terms to represent the main chamber solution (maximum of 20 terms. default value of 11 terms)	( none )
LC	: number of Bessel series terms to represent solution in the main chamber (maximum of 20 terms. default value of 8 terms)	( none )
MB	: number of Fourier series terms to represent the baffle compartment solution (maximum of 20 terms. default value of 11 terms)	( none )
LB	: number of Bessel series terms to represent solution in the baffle compartments (maximum of 20 terms. default value of 8 terms)	( none )
IDMAX	: maximum number of frequency iterations for successive approximation (default value of 9)	( none )
ALENGTH	: chamber length	( ft )
ZB	: baffle blade length	( ft )
T	: baffle blade thickness	( ft )
RCHAMB	: chamber radius	( ft )
MUB	: number of evenly spaced baffle compartments (maximum of 12 compartments)	( none )
HST	: nozzle throat radius	( ft )
RC	: radius of curvature at nozzle throat	( ft )
RE	: radius of curvature at nozzle entrance	( ft )
ALPHA	: nozzle convergence half angle	( deg )
AO	: chamber speed of sound at stagnation conditions	( ft/s )
PO	: chamber pressure at stagnation conditions	( psf )
GAMMA	: ratio of specific heats	( none )
PAMP	: peak to peak pressure amplitude ( percent of main chamber )	

MHAT	: dominating transverse mode number in main chamber	( none )
LHAT	: dominating radial mode number in main chamber	( none )
AN	: combustion interaction index (n)	( none )
TAU	: combustion time lag ( $\tau$ )	( msec )

### 3.2.7 - CONCENTRATED COMBUSTION PROGRAM OUTPUT

The output of the concentrated combustion program is written into file 'CON3OUT'. The output starts with a listing of the geometrical inputs and operating conditions such as nozzle inlet Mach number, steady state chamber pressure and sound speed. For each frequency iteration the program outputs the complex frequency (hertz) and iteration number. After the iteration limit is reached the final chamber and baffle compartment matrices are printed out. The final output includes the decay rate with baffle dissipation included, this is also presented as decay in decibels/cycle and the decay/growth rate factor.

AL LENGTH = 1.558 ft      R = 0.623 ft (wall of chamber)      ZS = 0.0, ZE = 1.558

A: 33.3% baffle      B: 6.67% baffle

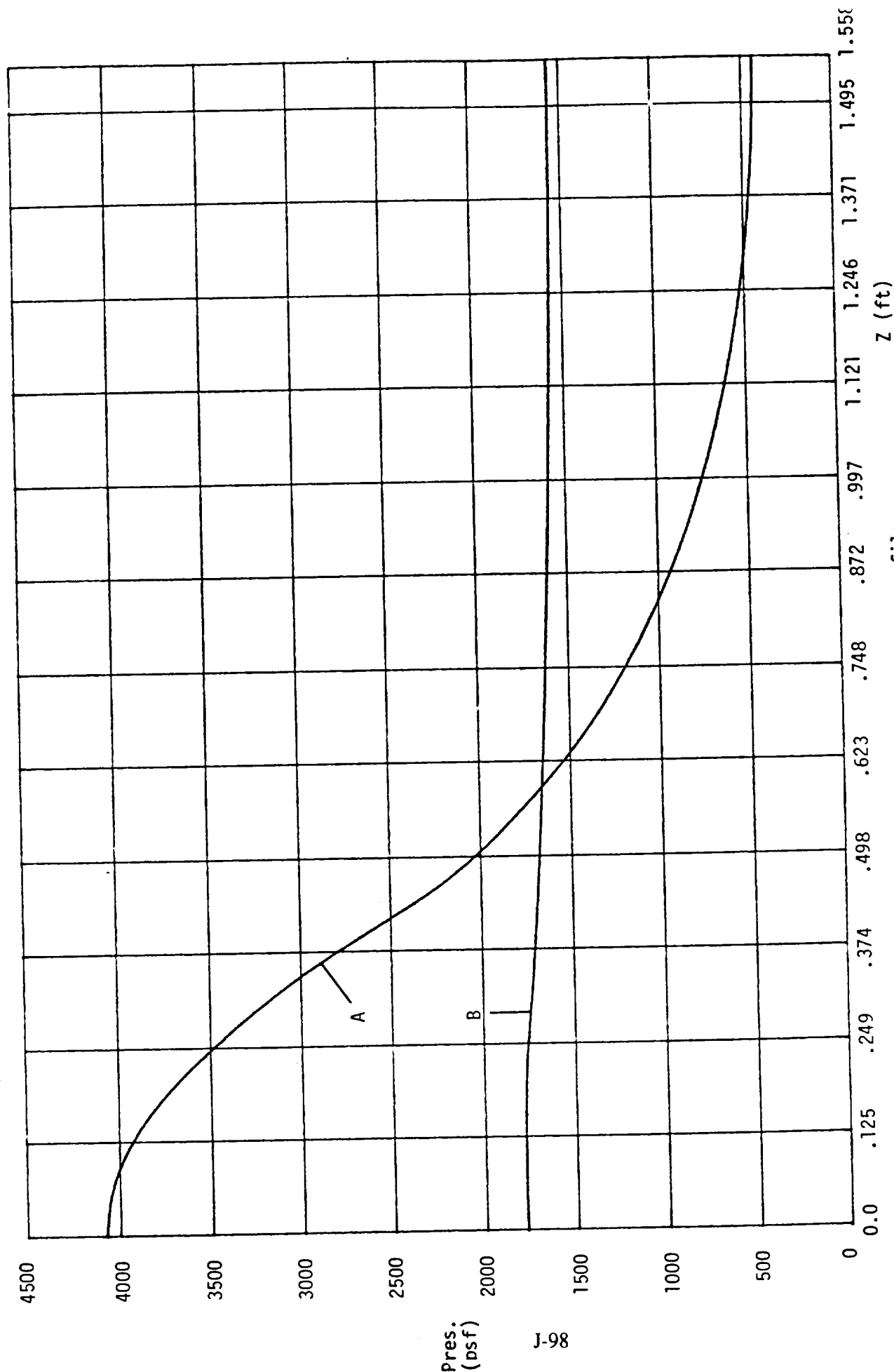


Figure 8: 3-D pressure profile

Figure 9

NEUTRAL STABILITY MAP FOR A 3-D CHAMBER, CONCENTRATED COMBUSTION ZONES

Chamber dimensions (ft): length = 1.5575 radius = 0.623  
 Baffle dimensions (ft): length = 0.1038 thickness = 0.03115  
 Steady state speed of sound = 3850.0 ft/s  
 Steady state pressure = 43200.0 psf  
 Ratio of specific heats = 1.2 Mode: First tangential  
 Frequency range (Hz): 1800 to 1950 No. of baffles: 3

ZONES (distance in feet from injector plate):

A: ZS = 0.0 to ZE = 0.1869 (no baffles present)  
 B: ZS = 0.0 to ZE = 0.1869  
 C: ZS = 0.4257 to ZE = 0.6126  
 D: ZS = 1.3706 to ZE = 1.5575

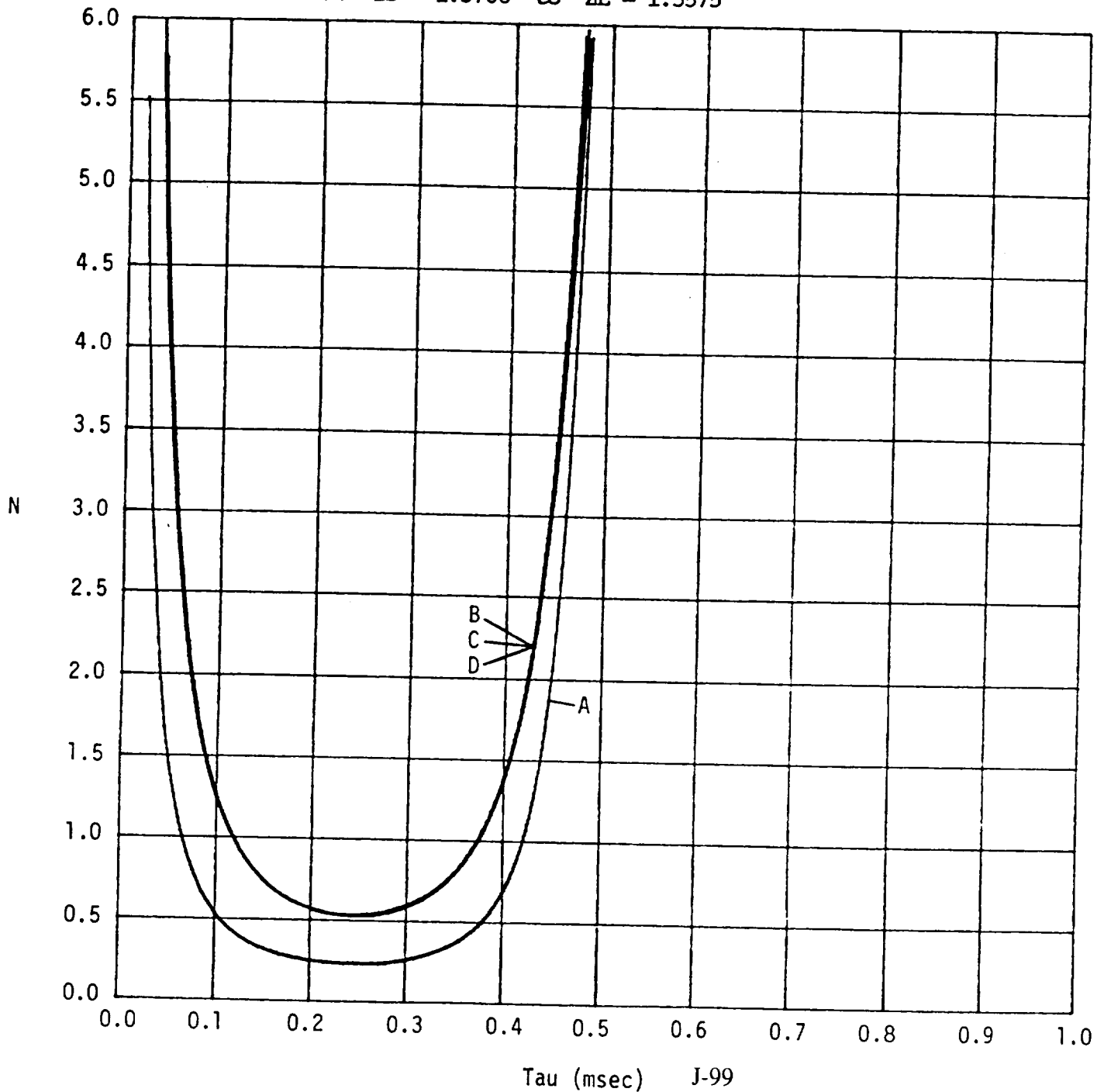


Figure 10

NEUTRAL STABILITY MAP FOR A 3-D CHAMBER, CONCENTRATED COMBUSTION ZONES

Chamber dimensions (ft): length = 1.5575 radius = 0.623  
 Baffle dimensions (ft): length = 0.51917 thickness = 0.03115  
 Steady state speed of sound = 3850.0 ft/s  
 Steady state pressure = 43200.0 psf  
 Ratio of specific heats = 1.2 Mode: First tangential  
 Frequency range (Hz): 1000 to 1500 No. of baffles: 3

ZONES (distance in feet from injector plate):

A: ZS = 0.0 to ZE = 0.1869  
 B: ZS = 0.4257 to ZE = 0.6126  
 C: ZS = 1.3706 to ZE = 1.5575

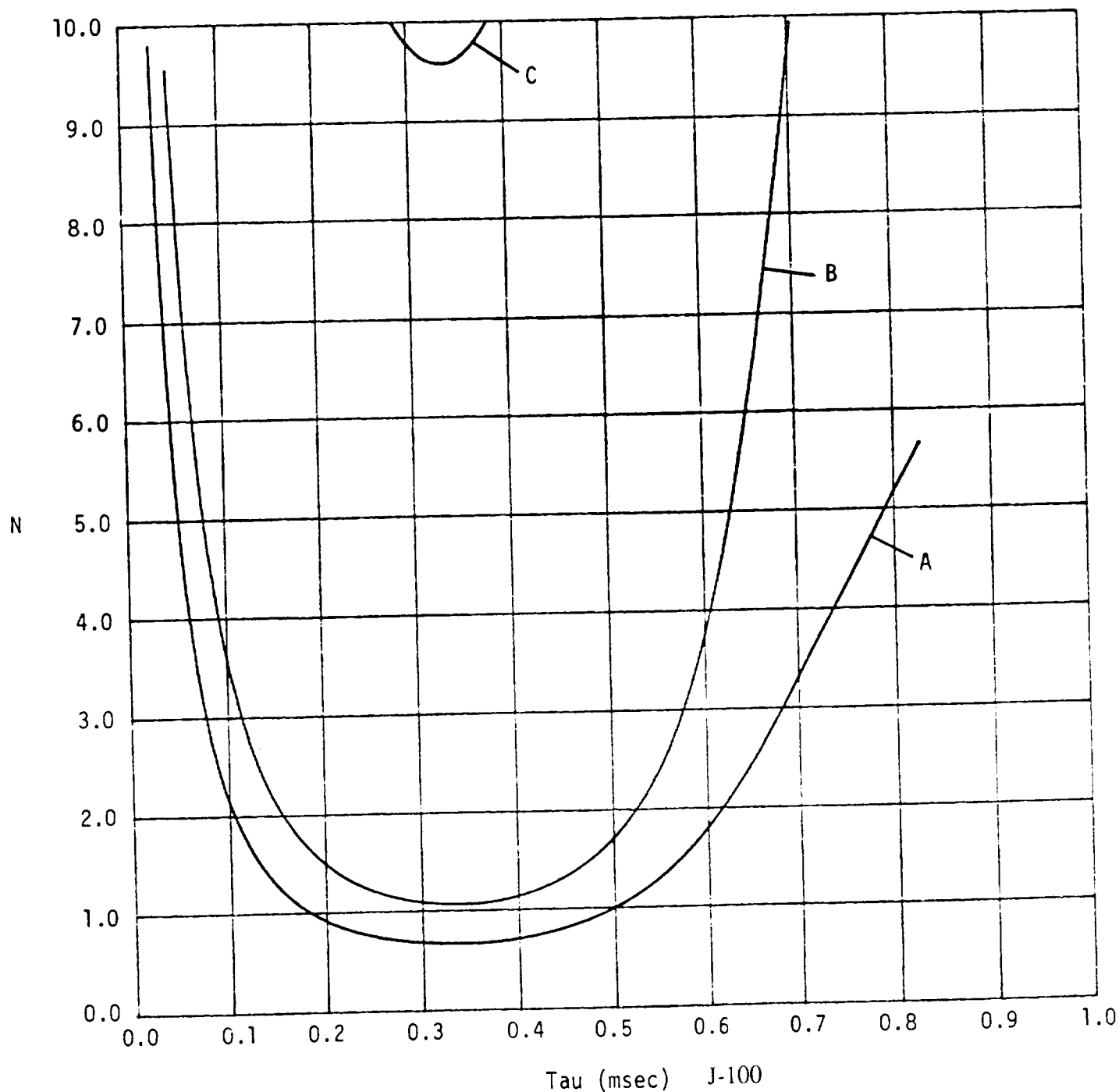


Figure 11

NEUTRAL STABILITY MAP FOR A 3-D CHAMBER, DISTRIBUTED COMBUSTION ZONES

Chamber dimensions (ft): length = 1.5575      radius = 0.623  
 Baffle dimensions (ft): length = 0.51917      thickness = 0.03115  
 Steady state speed of sound = 3850.0 ft/s  
 Steady state pressure = 43200.0 psf  
 Ratio of specific heats = 1.2      Mode: First tangential  
 Frequency range (Hz): 1000 to 1500      No. of baffles: 3

ZONES (distance in feet from injector plate):

A: ZS = 0.0 to ZE = 0.1869 (no baffles present)

B: ZS = 0.0 to ZE = 0.1869

C: ZS = 0.0 to ZE = 0.77875

D: ZS = 0.0 to ZE = 1.5575

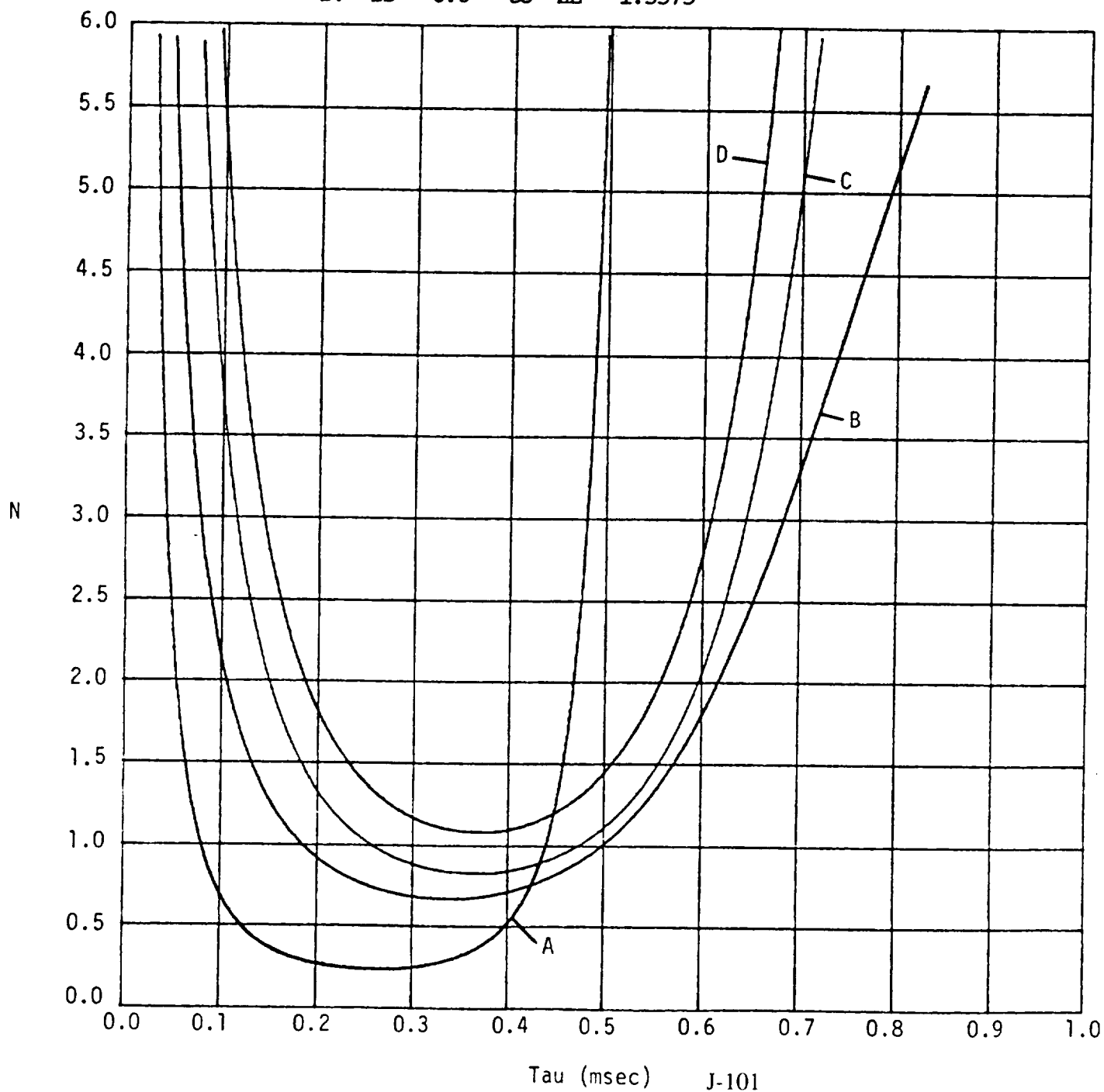


Figure 12

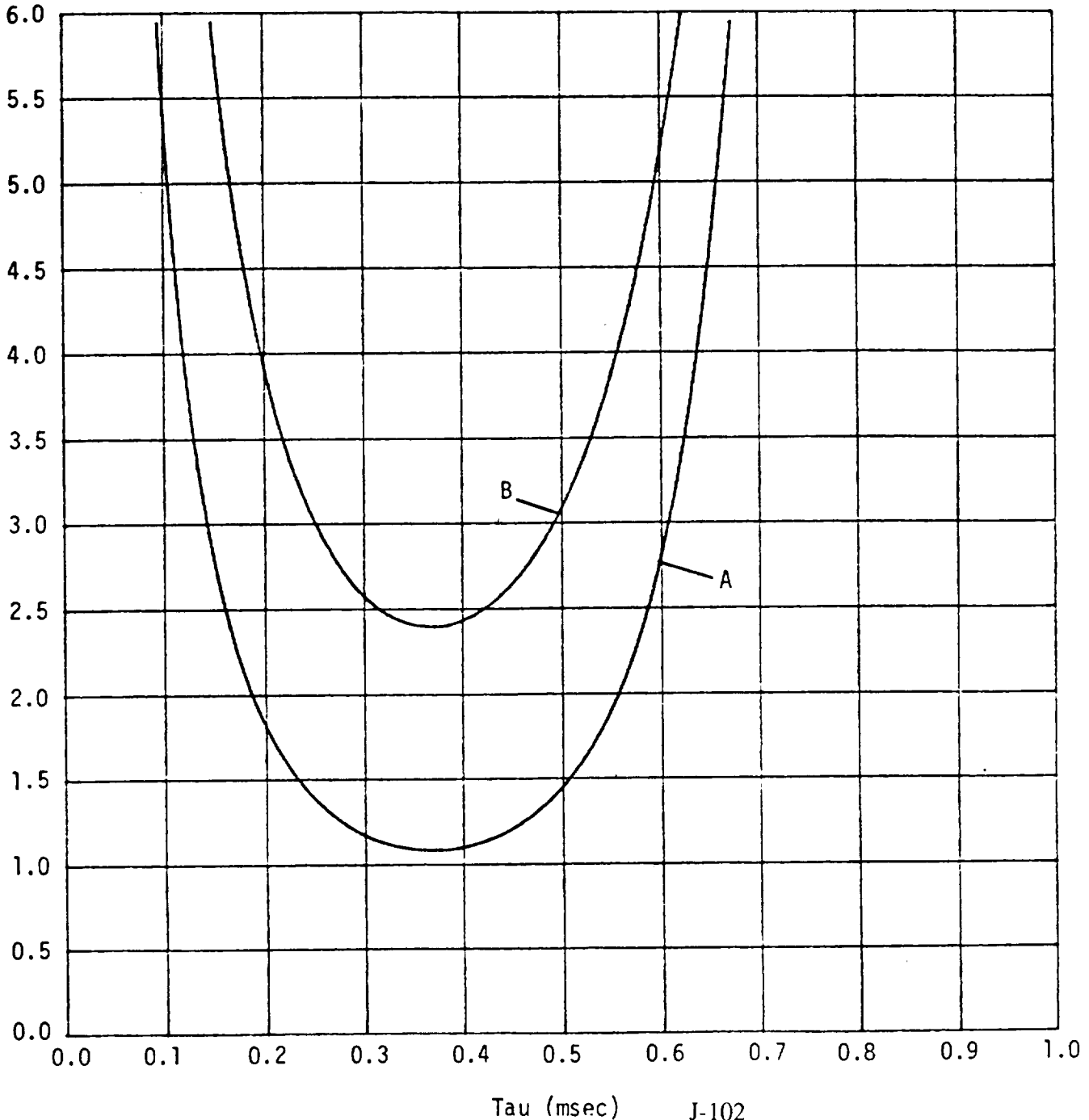
NEUTRAL STABILITY MAP FOR A 3-D CHAMBER, DISTRIBUTED COMBUSTION ZONES

Chamber dimensions (ft): length = 1.5575 radius = 0.623  
 Baffle dimensions (ft): length = 0.51917 thickness = 0.03115  
 Steady state speed of sound = 3850.0 ft/s  
 Steady state pressure = 43200.0 psf  
 Ratio of specific heats = 1.2 Mode: First tangential  
 Frequency range (Hz): 1000 to 1500 No. of baffles: 3  
 Cavity admittance (ft/s)/(psf): real = 0.04467 imaginary = 0.0  
 Cavity aperture width = 0.0623 ft

ZONES (distance in feet from injector plate):

A: ZS = 0.0 to ZE = 1.5575 - without cavity dissipation

B: ZS = 0.0 to ZE = 1.5575 - with cavity dissipation



#### 4. - REFERENCES

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4. "Computer Code For use in High Frequency Combustion Stability Analyses" T.V. Nguyen, ATC Thermodynamic Analysis Report, No. 9980:1807, Feb. 1987
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**PART C**

**COMBUSTION RESPONSE PREDICTION MODEL  
(CRP)**

ENGINEERING ANALYSIS REPORT		NUMBER: 9980:1998
		DATE: 27 JULY 1987
SUBJECT: COMBUSTION RESPONSE MODEL	PAGE 1 OF _____	
	NO. OF ENCLOSURES _____	
	NO. OF APPENDICES _____	
ADDITIONAL INFORMATION AND WORK NOTES INCLUDED IN MICROFILM FILE CDN _____		

PREPARED FOR: J. L. Pieper

As part of the development of analytical models for use in the LOX/Hydrocarbon Injector Characterization Program (contract number F04611-85-C-0100), a computer code has been developed to calculate the combustion response factors.

The code is capable of calculating the combustion response factors at sub-critical or super-critical chamber pressures. Significant effort was expended to devise a scheme to treat cases where the droplet temperature reaches the boiling temperature or the critical temperature of the propellant, and to generalize the propellant properties input. Correlations for determining the droplet and the combustion gas properties, for example the vapor pressure and the heat of evaporation of the droplet as functions of the temperature, were implemented into the code. A great amount of effort was also expended in several numerical tests to study the sensitivity of the solution to the artificial parameters, e.g. time steps, integration step sizes. The test results are also used to obtain the guidelines for specifying values of the parameters.

The attachment describes the theory, the computer code and the calculated results.

KEYWORDS: Misc (21), Chamber (52), Combustion Stability (105), LOX/HC (153), Model Development (209), Computer Program - New Develop. (210), 1987 (272), T. V. Nguyen (357)

DISTRIBUTION: R. Hewitt, J. Hulka,  
J. Hyde, J. Ito, M. Lausten, Y. Jones,  
S. Mercer, J. Muss, K. Niiya,  
D. Rousar, R. Schindler,  
J. Van Kleeck, 9980 File

PREPARED BY: *Thong Van Nguyen*

Thong Van Nguyen

REVIEWED BY:

*R. E. Walker* / *J. J. Fang*  
R. E. Walker / J. J. Fang

APPROVED BY:

*J. W. Salmon*  
J. W. SALMON, MANAGER  
ENGINEERING ANALYSIS DEPT.

W.O. NO: KAE626

J-105

# COMBUSTION RESPONSE PREDICTION

by

Thong Van Nguyen

Aerojet TechSystems Company  
Sacramento, CA 95813

July 15, 1987

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## NOMENCLATURE

### Symbols:

$A_d$	Droplet surface area.
$c$	Speed of sound.
$C_D$	Drag Coefficient.
$C_{PL}$	Specific heat of propellant liquid.
$D$	Drag force.
$h$	Convection heat transfer coefficient.
$I$	Imaginary part of the combustion response factor.
$J_\nu$	Bessel function of the $\nu^{th}$ order.
$k$	Thermal conductivity of the combustion gas.
$K_g$	Mass transfer coefficient.
$m$	Instantaneous droplet mass.
$m_i$	Initial droplet mass.
$M$	Molecular weight of the combustion gas.
$M_V$	Molecular weight of the propellant.
$n$	Pressure interaction index.
$n_d$	Number of droplets in an array.
$p$	Chamber pressure.
$p_v$	Propellant vapor pressure.
$Pr$	Combustion gas Prandtl number.
$r$	Radial coordinate normalized by the chamber radius.
$r_d$	Droplet radius.
$R$	Universal gas constant.
$Re$	Reynolds number.
$S_{\nu\eta}$	Eigenvalues correspond to the radial and the tangential resonance modes $\nu$ and $\eta$ .

$Sc$     Combustion gas Schmidt number.  
 $t$        Time.  
 $t_{qg}$     Droplet life time.  
 $T$        Temperature of the combustion gas.  
 $T_d$      Temperature of the droplet.  
 $T_f$      Film temperature.  
 $u$        Axial gas velocity component.  
 $u_f$      Final gas velocity.  
 $v$        Radial gas velocity component.  
 $V$        Gas velocity vector.  
 $V_d$      Droplet velocity vector.  
 $v_\psi$     Radial velocity of the propellant vapor leaving the droplet.  
 $w$        Tangential velocity component.  
 $W$        Total evaporation rate of an array of droplets  
 $x$        Axial coordinate.  
 $Y$        Combustion response factor  
 $\nu$        Radial resonance mode number.  
 $\eta$        Tangential resonance mode number.  
 $\Delta p$     Maximum pressure amplitude.  
 $\omega$        Angular frequency of the acoustic oscillations.  
 $\phi$        Phase angle of the acoustic oscillations.  
 $\gamma$        Specific heat ratio of the combustion gas.  
 $\rho$        Density of the combustion gas.  
 $\rho_v$       Density of the propellant vapor.  
 $\rho_l$       Density of the propellant liquid.  
 $\dot{\psi}$        Vaporization rate.  
 $\lambda$        Heat of Evaporation.

- $\mathcal{D}$  Diffusion coefficient.
- $\mu$  Viscosity of the combustion gas.
- $\mathcal{R}$  Real part of the combustion factor.
- $\tau$  Sensitive time lag.
- $\theta$  Tangential coordinate.

Subscripts:

- d Droplet.
- l Propellant liquid.
- v Propellant Vapor.
- ' Perturbation component.

Superscripts:

- Mean component.
- $\rightarrow$  Vector quantities.

## I. INTRODUCTION

Aerojet TechSystems Company is currently conducting a program (contract F04611-85-C-0100) to formulate a procedure (Ref. 1) which can accurately characterize injector designs for large thrust (0.5 to 2.0 million pounds) high pressure (500 to 3000 psia) LOX/hydrocarbon engines. As part of the development of models for use in the procedure, a computer code, Combustion Response Prediction (CRP), has been developed to calculate the combustion response factor which indicates the open-loop response of the burning rate to a specified acoustic oscillations in the combustion chamber.

### 1.1 High-Frequency Combustion Stability

Combustion instability, characterized by organized pressure oscillations in rocket combustion chamber, can cause severe vibrations on various engine system components and payloads. In addition, combustion instabilities may cause excessive mechanical stresses and heat loads on the injector and combustion chamber walls.

Combustion instabilities have been generally classified according to their frequency range: low, intermediate and high frequency. Significant efforts have been devoted to the understanding of high-frequency instability because it is the most common in new engine developments and is the most destructive. High-frequency instability results from the coupling between the combustion process and the acoustic waves in the chamber.

## 1.2 Combustion Response

Analytical models capable of characterizing combustion instability are obviously useful and valuable to engine designers during the development stage. As mentioned in the above section, high-frequency instability results from the coupling between the combustion process and the acoustic waves in the chamber. Thus, the stability of a given engine with specified operating conditions can be determined from the chamber transfer function and the burning transfer function. The chamber transfer function is defined in reference 11 as the ratio of the pressure oscillation to the burning rate oscillation normalized respectively by the mean pressure and the mean burning rate. Conversely, the burning transfer function or the combustion response is defined as the ratio of the burning rate oscillation to the pressure oscillation also normalized by the mean burning rate and the mean pressure. The burning transfer function indicates the response of the combustion process to the acoustic waves in the chamber. Previously, a computer code, HIFI (Ref. 2) has been developed to calculate the chamber transfer function. The burning transfer function was more difficult to predict analytically, therefore in the past it was expressed in term of an interaction index,  $n$  and a combustion time lag,  $\tau$ . The values of  $n$  and  $\tau$  are determined empirically.

## 1.3 Objective of the Present Study

The objective of the present study is to provide a computer code to predict the burning transfer function. Results from the

code are used together with the HIFI's prediction of the chamber transfer function to predict the high-frequency stability of rocket engines.

#### 1.4 Approach

The approach taken is to modify the Agosta and Hammer's computer model (Ref. 4) which was developed to study the vaporization response of oxygen droplets. Modifications made to the model include the following:

1. The model was extended from 1T traveling mode to mixed radial and tangential modes up to a combination of 8R and 8T. The acoustic modes can be either standing or spinning.
2. Subroutines for calculating the Bessel functions of any order were developed and implemented into the computer code. The values of the Bessel functions are calculated internally instead of being input by the users.
3. The finite-thermal-conductivity assumption was replaced by the uniform-droplet-temperature assumption. This was done since predictions of droplet evaporation rates in steady gas environments using the latter assumption agree better with the experimental data. The latter assumption is also necessary to reduce the computer time requirement to a practical level.
4. The time step used in the calculation of droplet evaporation history is determined internally by the computer code instead of being input by the users. The value of the time step is determined based on the period of oscillations and the droplet

lifetime.

5. The original model calculates the response factor for droplets injected at a specified radial and circumferential location on the injector. Because pressure and velocity oscillations vary with radial and circumferential locations, the response factor also varies with the locations. To account for this effect, the resultant response factor is obtained by averaging the response factors calculated at several different radial and circumferential locations.

6. The symmetry of the standing modes is taken into account to reduce the number of circumferential injection locations. This results in substantial saving of computer time.

7. The original computer code requires the users to input expressions for calculating the heat of vaporization and the vapor pressure as functions of temperature, and the diffusion coefficient as a function of pressure and temperature. In general, this requires the users to search literature for appropriate correlations and to compute the parameters used in the correlations for the propellants of interest. Modifications were made to the code so that the correlations are built into the code. Watson's correlation is used for calculating the heat of vaporization, Reidel's correlation is used for calculating the vapor pressure, and Mathur and Thodos' correlation is used for calculating the diffusion coefficient. The parameters used in these correlations are calculated internally by the code using user input data, e.g. molecular weight, critical temperature and pressure. These data

are existing for most propellants and can be easily found in existing literature, for example the Aerojet Handbook of Properties and Performance of Liquid Rocket Propellants (Ref. 5)

## II. THEORY

The theory in the present study follows closely references 3 and 4. In the present study, mixing and reaction is assumed to be so fast that the burning rate is assumed to be vaporization limited. Therefore, the terms burning rate and vaporization rate are interchangeable within the context of the present study.

The following sections describe in details the theory, the equations, and the calculation procedure used in the present model to calculate the response factor.

### 2.1 Theory Description

First, the equations describing the pressure and the velocity oscillations are prescribed for an acoustic mode in the chamber. The evaporation rates of a single droplet injected into the chamber is calculated assuming the heat and the mass transfer processes between the droplet and the surrounding combustion gas are at quasi-steady state. The convection heat rate and the mass vaporization rate are calculated using the Ranz-Marshall's correlations for the heat transfer and the mass transfer coefficient (Ref. 6). An energy balance is then applied to the droplet to calculate the rate at which the droplet is heated up (see equation 11 in chapter III). The theory just described is used to calculate the vaporization history of single droplets injected into the chamber. The vaporization history includes the temporal variations of the diameter, the temperature and the vaporization rate of the droplet.

The continuous injection of propellant is simulated by arrays of single droplets that are injected from the various radial and circumferential locations. Each of the arrays are comprised of droplets that are injected from the same location but at different times during an oscillation period of the acoustic fields. The previously described procedure for calculating the evaporation history of a single droplet is used to calculate the evaporation histories of each of the droplets in the array. The total evaporation rate of the array is then calculated by summing the evaporation rates of each of the individual droplets. The total evaporation rate includes the perturbation component that is induced by the acoustic oscillations. An in-phase response factor and an out-of-phase response factor are then calculated. The in-phase response factor is defined as the normalized ratio of the combustion rate perturbation component that is in phase with the pressure oscillation. Similarly, the out-of-phase response factor is defined as the normalized ratio of the combustion rate perturbation component that is out of phase with the pressure oscillation. The ratios are normalized by the mean combustion rate and the mean pressure. Because pressure and velocity oscillations vary with radial and circumferential locations, the response factor for an array of droplets injected from one radial and circumferential location may be different, in general, from the response factors calculated for arrays of droplets injected from other locations. In order to account for this effect, the resultant response factor is obtained by averaging the response factors calculated at various injection locations.

## 2.2 Equation Description

### 2.2.1 Chamber Acoustics

The expressions for the instantaneous values, which comprise the mean and the perturbation components, of the pressure and the velocity of a gas in a closed cylinder have been derived in reference 7 as:

$$P = \bar{P} \left\{ 1 + \Delta P \frac{J_\nu(s_{\nu\eta} r)}{J_\nu(s_{\nu\eta})} \sin(\nu\theta - \omega t + \phi) \right\} \quad (1)$$

$$v = \frac{\bar{c} \Delta P}{\gamma} \frac{J'_\nu(s_{\nu\eta} r)}{J_\nu(s_{\nu\eta})} \cos(\nu\theta - \omega t + \phi) \quad (2)$$

$$w = -\frac{\bar{c} \nu \Delta P}{\gamma s_{\nu\eta}} \frac{1}{r} \frac{J_\nu(s_{\nu\eta} r)}{J_\nu(s_{\nu\eta})} \sin(\nu\theta - \omega t + \phi) \quad (3)$$

for a spinning wave motion, and

$$P = \bar{P} \left\{ 1 + \Delta P \frac{J_\nu(s_{\nu\eta} r)}{J_\nu(s_{\nu\eta})} \cos(\nu\theta) \sin(\omega t - \phi) \right\} \quad (4)$$

$$v = \frac{\bar{c} \Delta P}{\gamma} \frac{J'_\nu(s_{\nu\eta} r)}{J_\nu(s_{\nu\eta})} \cos(\nu\theta) \cos(\omega t - \phi) \quad (5)$$

$$w = -\frac{\bar{c} \nu \Delta P}{\gamma s_{\nu\eta}} \frac{1}{r} \frac{J_\nu(s_{\nu\eta} r)}{J_\nu(s_{\nu\eta})} \sin(\nu\theta) \cos(\omega t - \phi) \quad (6)$$

for a standing wave motion.

The above expressions are for the transverse wave motion and have been written on the assumptions that the chamber is a closed-end cylinder. In these expressions  $p$ ,  $v$  and  $w$  are the instantaneous pressure, and radial and tangential velocities of the gas, respectively;  $\bar{p}$  is the mean pressure;  $\Delta p$  is the maximum

amplitude of the pressure oscillations for a particular mode;  $\bar{c}$  is the mean speed of sound in the chamber;  $\gamma$  is the specific heat ratio of the gas in the chamber;  $r$  is the radial coordinate normalized by the chamber radius;  $\theta$  is the tangential coordinates,  $\omega$  is the angular frequency;  $t$  is the time;  $\phi$  is the phase angle;  $\nu$  and  $\eta$  are the numbers of the radial and tangential resonance modes, respectively.  $J_\nu$  is the  $\nu^{\text{th}}$ -order Bessel function of the first kind;  $J'_\nu$  is the derivative of  $J_\nu$  with respect to  $S_{\nu\eta} r$ ; and the values of  $S_{\nu\eta}$  are given in table 2.1 for selected values of  $\nu$  and  $\eta$ . It should be noted that the maximum amplitude of the pressure oscillation has been normalized by the mean pressure, and that the means of the radial and the tangential velocity components have been assumed to be equal to zero.

The instantaneous temperature of the gas,  $T$  can be related to the pressure using the following isentropic relations:

$$\frac{T}{\bar{T}} = \left( \frac{P}{\bar{P}} \right)^{\frac{\gamma-1}{\gamma}} \quad (7)$$

where  $\bar{T}$  is the mean temperature of the gas. The instantaneous density of the gas,  $\rho$  can be related to the temperature and the pressure using the equation of state:

$$\rho = \frac{P M}{R T} \quad (8)$$

where  $R$  is the universal gas constant, and  $M$  is the molecular weight of the gas.

The mean sound speed used in equations (2), (3), (5) and (6)

$\nu \backslash \eta$	0	1	2	3	4
0	0.0000	3.8318	7.0155	10.1734	13.3238
1	1.8413	5.3313	8.5263	11.7059	14.8635
2	3.0543	6.7060	9.9695	13.1705	16.3476
3	4.2013	8.0151	11.3459	14.5858	17.7890
4	5.3175	9.2825	12.6820	15.9640	19.1961
5	6.4154	10.5199	13.9873	17.3127	20.5755
6	7.5012	11.7348	15.2681	18.6375	21.9318
7	8.5778	12.9324	16.5295	19.9419	23.2682
8	9.6475	14.1155	17.7739	21.2290	24.5874

Table 2.1: Selected Values of  $S_{\nu\eta}$  .

can be expressed in terms of the molecular weight and the mean temperature of the gas:

$$\bar{c} = \sqrt{\gamma \frac{R}{M} \bar{T}} \quad (9)$$

While the means of the radial and the circumferential components of the velocity are assumed to be zero, the mean of the axial component,  $\bar{u}$  is non-zero and it is a function of the axial coordinate,  $x$

$$\bar{u} = \bar{u}(x) \quad (10.a)$$

which is assumed to be known apriori. The axial profile of the mean velocity can be obtained, for example, from the steady-state performance analysis. For simplicity, in the present analysis, the velocity is assumed to be in the following form:

$$\bar{u} = \bar{u}_f \left( 1 - \frac{m}{m_i} \right) \quad (10.b)$$

where  $\bar{u}_f$  is the final gas velocity,  $m$  is the instantaneous value of the droplet mass, and  $m_i$  is the initial value of the droplet mass. This assumption implies the gas mean velocity varying with the axial coordinate. It simplifies the analysis since one does not have to compute and keep track of the axial location of the droplet.

Equations (1) through (10) completely describe the temporal and spatial variations of the gas properties, i.e. pressure,

density, temperature and velocities.

### 2.2.2 Vaporization of a Single Droplet

Neglecting the radiation heat transfer, an energy balance applied to a droplet which undergoes heat and mass transfer simultaneously yields:

$$hA_d(T - T_d) = \dot{\psi}\lambda + \frac{1}{2}\dot{\psi}v_v^2 + mC_{pL}\frac{dT_d}{dt} \quad (11)$$

where  $h$  is the convective heat transfer coefficient,  $A_d$  is the surface area of the droplet,  $T$  is the gas temperature,  $T_d$  is the droplet temperature,  $\dot{\psi}$  is the mass evaporation rate,  $\lambda$  is the enthalpy of evaporation,  $m$  is the instantaneous mass of the droplet,  $C_{pL}$  is the specific heat of the liquid propellant, and  $v_v$  is the velocity of the vapor leaving the droplet. Assuming the vaporizing mass leaving the droplet radially,  $v_v$  can be related to the mass evaporation rate,  $\dot{\psi}$ , the droplet surface area,  $A_d$ , and the vapor density,  $\rho_v$ :

$$v_v = \frac{\dot{\psi}}{A_d \rho_v} \quad (11.a)$$

The left-hand side of equation (11) represents the heat rate transferred to the droplet by convection; the first term on the right-hand side of the equation represents the heat required to vaporize the mass leaving the droplet; the second term on the right hand side of the equation is the kinetic energy that is imparted to the vaporizing mass; and the remaining term represents the energy required to heat up the droplet. The kinetic energy term is very

small compared to the other two terms when the droplet temperature is far below the critical point of the propellant. Therefore, it has been neglected in many of the past studies. It becomes increasingly important as the droplet temperature increases, especially as the temperature approaches the critical temperature of the propellant because at this temperature, the heat of evaporation in the first term approaches zero and the mass "evaporation" rate approaches infinity. In rocket engines with high chamber pressure such as those considered in the Lox/hydrocarbon Injector Characterization Program, the temperature of the propellant droplet is expected to be very high and it may approach or even exceed the critical temperature of the propellant. Thus, the kinetic energy term must be included. In the above equation, the temperature of the droplet has been assumed to be uniform.

The mass evaporation rate is given by:

$$\dot{m} = A_d K_g p \ln \frac{p}{p - p_v} \quad (12)$$

where  $p$  is the total gas pressure,  $p_v$  is the vapor pressure of the droplet, and  $K_g$  is the mass transfer coefficient. The heat transfer and the mass transfer coefficients are obtained from the empirical correlations of reference 6:

$$\frac{2r_d h}{k} = 2 + 0.6 (Pr)^{1/3} (Re)^{1/2} \quad (13)$$

$$\frac{2r_d R T_f K_g}{M_v g} = 2 + 0.6 (Sc)^{1/3} (Re)^{1/2} \quad (14)$$

In equations (13) and (14);  $r_d$  is the droplet radius;  $k$  is the gas thermal conductivity;  $Pr$  and  $Sc$  are the gas Prandtl and Schmidt numbers, respectively;  $R$  is the universal gas constant;  $M_v$  is the molecular weight of the propellant vapor,  $D$  is the binary diffusion coefficient of the vapor and the combustion gas,  $T_f$  is the film temperature which is defined as the arithmetic mean of the gas and the droplet temperatures:

$$T_f = \frac{1}{2} (T + T_d) \quad (15)$$

and  $Re$  is the Reynolds number which bases on the relative velocity of the gas and the droplet:

$$Re = \frac{2r_d |\vec{V} - \vec{V}_d| \rho}{\mu} \quad (16)$$

where  $\mu$  is the gas viscosity, and  $\vec{V}$  and  $\vec{V}_d$  are the velocities of the gas and the droplet, respectively. The velocity of the droplet is changing with time due to the drag exerted on the droplet by the surrounding gas. Although, the viscous effects of the gas on the droplet motion are considered, acceleration or deceleration of the gas through viscous interaction with the droplet is neglected. The drag exerted on the droplet is written as:

$$\vec{D} = C_D \pi r_d^2 \left( \frac{1}{2} \rho \Delta \vec{V} |\Delta \vec{V}| \right) \quad (17)$$

where  $C_D$  is the drag coefficient

$$C_D = 27 Re^{-0.84} \quad (18)$$

and

$$\Delta \vec{V} = \vec{V} - \vec{V}_d \quad (19)$$

is the gas velocity relative to the droplet. Newton's second law is then applied to calculate the rate of change of the droplet velocity:

$$\frac{d\vec{V}_d}{dt} = \frac{\vec{D}}{\frac{4}{3}\pi r_d^3 \rho_L} \quad (20)$$

where  $\rho_L$  is the density of the propellant liquid.

The perturbation components of the gas velocity and the density are included in the calculation of the Reynolds number, therefore the mass evaporation rates are affected by the acoustic oscillations. Nevertheless, The effects of the acoustic perturbations on the radial and the circumferential components of the droplet velocity are, neglected, i.e. the radial and the circumferential components of the droplet velocity always remain to be zero (the radial and the circumferential components of the droplet velocity at the instant of injection have also been assumed to be zero).

The time rate of change of droplet velocity is given by equation (20). The time rate of change of droplet temperature can be obtained by rearranging equation (11):

$$\frac{dT_d}{dt} = \frac{1}{m C_{pL}} \left\{ h A_d (T - T_d) - \dot{\Psi} \lambda - \frac{1}{2} \dot{\Psi} v_r^2 \right\} \quad (21)$$

and the time rate of change of droplet radius can be related to the mass evaporation rate, by assuming constant propellant liquid

density:

$$\frac{dr_d}{dt} = - \frac{\dot{\psi}}{A_d \rho_L} \quad (22)$$

The radius, temperature and velocity of the droplet at the next time step can be calculated from the known values of the functions and their derivatives at the current time using the Euler's explicit integration, which is illustrated in the following equation for a general function,  $f$ :

$$f \Big|_{t+\Delta t} = f \Big|_t + \frac{df}{dt} \Big|_t \Delta t \quad (23)$$

It should be noted that the gas transport properties such as  $k$  and  $\mathcal{D}$ , used in all of the above equations are evaluated at the film temperature.

Equations 1 through 23 are sufficient to calculate the variations of the diameter, temperature, velocity and the evaporation rate of a droplet as functions of time provided that the time and the location of injection are known and that the initial properties, i.e., diameter, temperature and velocity of the droplet are known at the instant of injection.

When the droplet temperature reaches the boiling temperature of the propellant, numerical problem arises as the mass evaporation rate calculated using equation (12) approaches infinity. Under this condition, the mass evaporation rate is calculated directly from the energy equation, equation (11), with the droplet-heat-up term being set to zero. This approach not only avoids the

numerical problem but is also more realistic since the evaporation rate is finite. The approach is extended also to the cases where the droplet temperature reaches the critical temperature of the propellant. It has been implied that when the droplet temperature reaches the boiling or the critical temperature of the propellant, no further heating of the droplet is allowed.

As previously mentioned in this section, the temperature of the droplet has been assumed to be uniform. Two other assumptions - temperature gradients exist inside the droplet (finite thermal conductivity), and uniform temperature inside the droplet with a step change at the droplet surface (zero thermal conductivity or onion skin) - are also considered but found to be unacceptable in the present study. The analysis assuming zero-thermal-conductivity is not described here since it is simple and similar to the analysis assuming uniform-temperature described in this section. The analysis assuming finite-thermal-conductivity is also not described here since it has been described in reference 4. In this reference, the temperature distribution inside the droplet is assumed to be spherically symmetric.

Although the finite-thermal-conductivity assumption appears to be more realistic than the other assumptions, the assumption of spherically symmetric temperature distribution is dubious especially when the droplet is under strong convection environments such as those in the combustion chambers.

The uniform-temperature-assumption is justifiable when the thermal conductivity of the droplet is high or strong circulating

flows exist inside the droplet. The circulating flows inside a droplet in a strong convection environment are generally believed to exist as a result of the shear force at the droplet surface. Therefore, the uniform-droplet-temperature assumption has been used commonly in the combustion or vaporization studies of droplet under convection, for example references 8 and 9. This assumption is adopted in the present study because calculations of the evaporation histories of droplets vaporizing in steady environments assuming uniform-droplet-temperature agree well with the experimental data (see section 4.1). Furthermore, the analysis using the assumption is simple and requires significantly less computer time than the analysis using the finite-thermal-conductivity assumption.

The analysis using the zero-thermal-conductivity assumption is also simple and has approximately the same computer time requirement as its uniform-temperature counterpart. Its predictions of the evaporation histories of droplets vaporizing in steady environments, however, did not agree well with the data (see section 4.1).

### 2.2.3 Evaporation of an Array of Droplets

The section 2.2.2 described equations used for the calculation of the evaporation rate of a single droplet. The evaporation rate of an array of droplets is simply the sum of the evaporation rates of each of the individual droplets that constitute the array.

The continuous injection of propellant into the chamber is simulated by the repeated injections of single droplets. Because

of the oscillations in the gas properties which directly affect the vaporization rate, the evaporation history of a droplet injected at one instant of time may be, in general, different from the evaporation histories of droplets injected at the same location but at different times. Droplets injected at times which are multiple oscillation periods apart, however, will have identical vaporization histories since they experience the identical time histories of gas thermodynamics and flow fields. The droplets will eventually disappear after they completely vaporize. If the droplets in an array are injected into the chamber at times evenly distributed over one oscillation period, then at "steady-state" conditions the number of droplets disappearing are equal to the number of droplets entering the chamber. Thus, the instantaneous total evaporation rate,  $W(t)$  is obtained by appropriately summing the evaporation rates of each of the droplets which exist at the time.

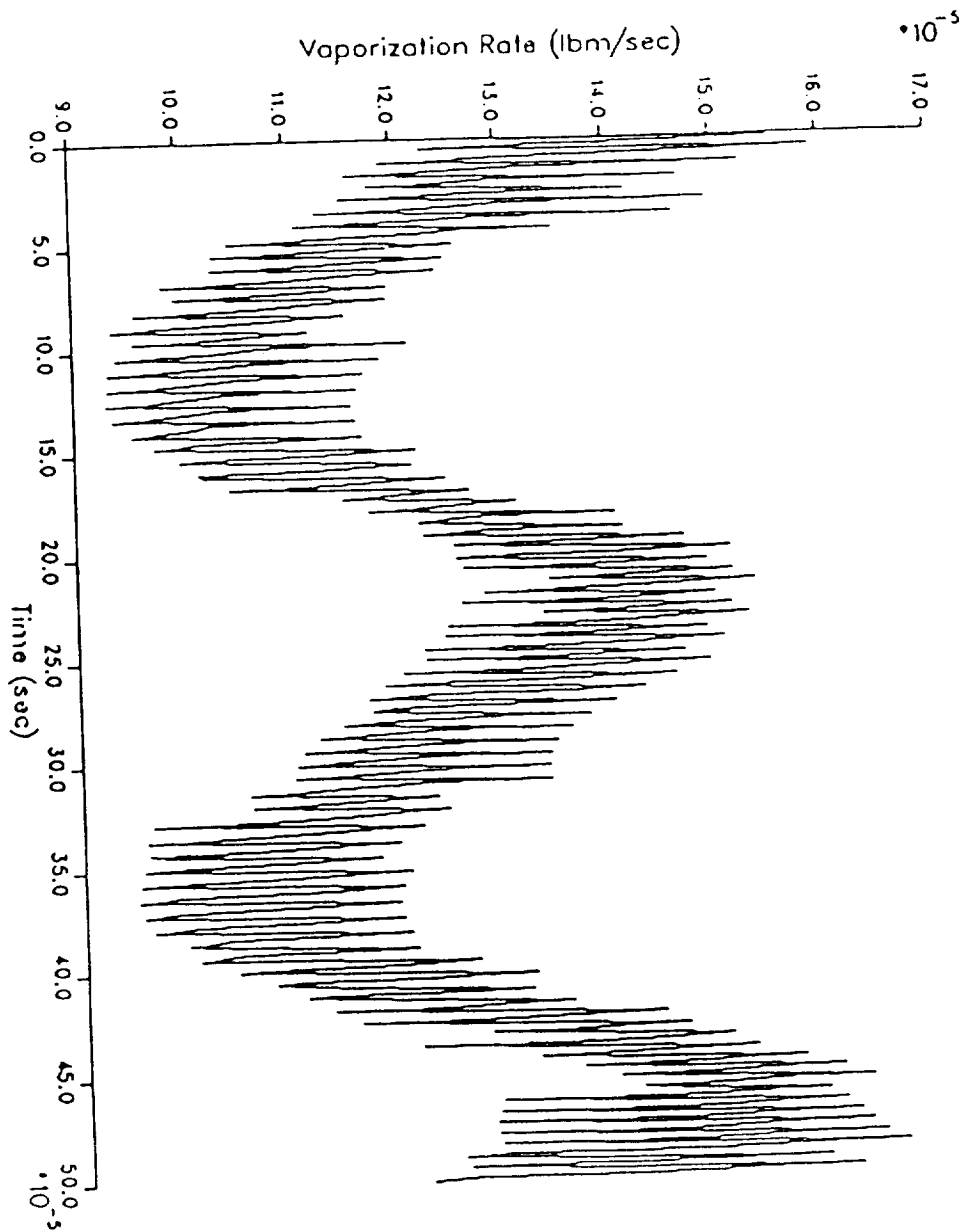
$$W(t) = \sum_{n=1}^{n_d} \dot{\psi}_n(t) \quad (24)$$

where  $\dot{\psi}_n(t)$  is the evaporation rate of the droplet  $n$ , and  $n_d$  is the number of droplets existing at time  $t$ . This total burning rate is, of course, periodic and dependent on the frequency of the acoustic fields.

#### 2.2.4 In-Phase and Out-Of-Phase Response Factors

The total burning rate (hereafter will be referred to simply as the burning rate) includes a perturbation component that is induced by the chamber acoustic fields. Figure 2.1 shows, as an

Figure 2.1 : Total Vaporization Rate of an Array of Droplets which Includes the Perturbation Component Induced by the Acoustic Oscillations.



example, the perturbation component of the burning rate which is induced by the oscillations of the pressure and the velocity and other thermodynamic properties. The pressure and velocity oscillations are shown in figures 2.2, 2.3 and 2.4, respectively. The spikes in the burning rate curve are the results of the use of a finite number of droplets to represent the continuous injection of the propellant. In this example, 80 droplets are injected at equal time intervals during an oscillation period.

The correlations between the the burning rate and the pressure oscillations describe the magnitude of the response of the burning rate to the acoustic fields in the chamber. Thus, they indicate the relative stability of the combustion.

The in-phase response factor defined as:

$$\mathcal{R} = \frac{\int_0^{2\pi} w' p' d(\omega t)}{\int_0^{2\pi} p'^2 d(\omega t)} \quad (25)$$

is a correlation of the burning rate component that is in phase with the pressure oscillation.

The out-of-phase response factor defined as:

$$\mathcal{I} = \frac{\int_0^{2\pi} w' p^* d(\omega t)}{\int_0^{2\pi} p^{*2} d(\omega t)} \quad (26)$$

is a correlation of the burning rate component that is out of phase with the pressure oscillation.

In equations (25) and (26),  $w'$  is the normalized perturbation component of the burning rate. It is defined in terms of the

Figure 2.2 : Normalized Acoustic Pressure Oscillation Component.

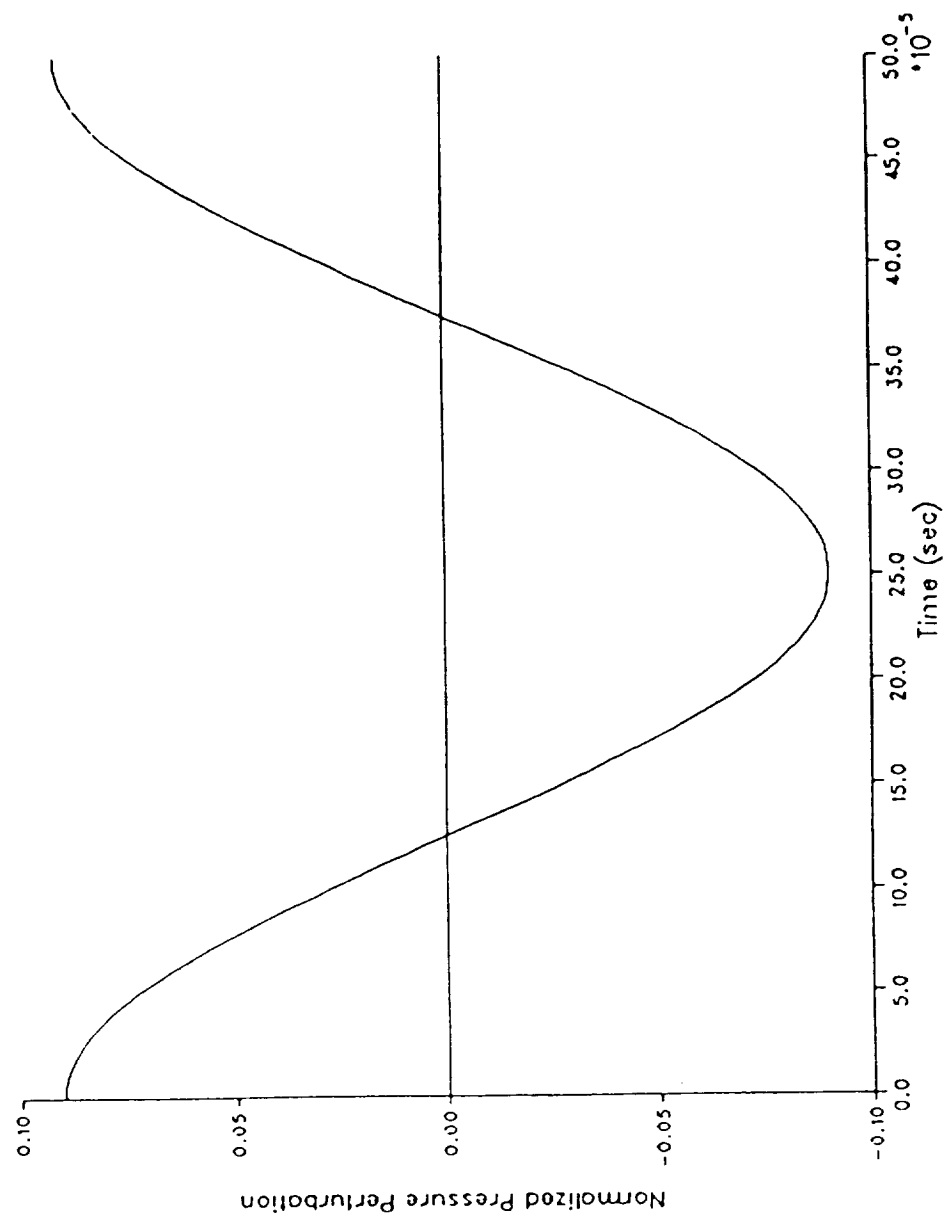


Figure 2.3 : Acoustic Radial Velocity Component.

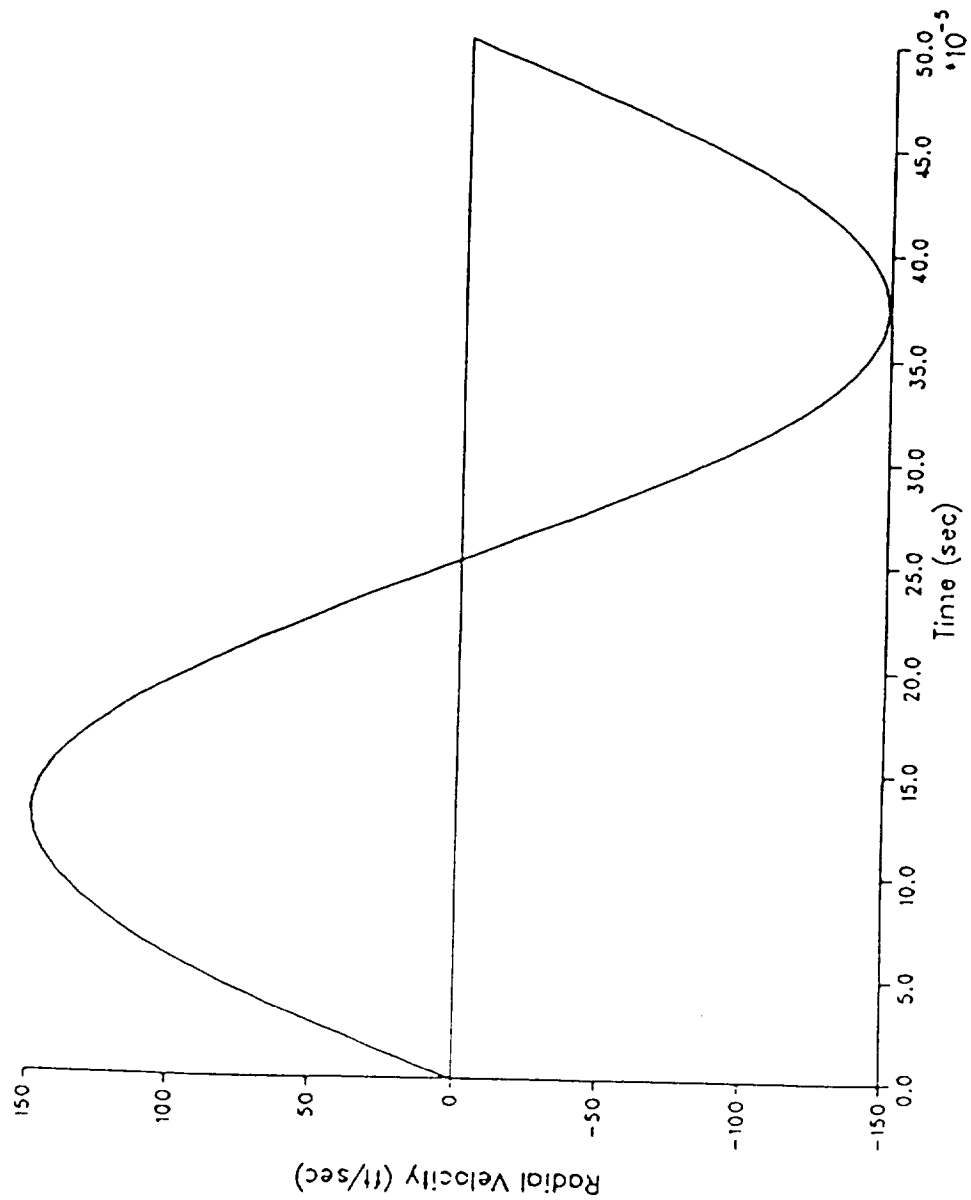
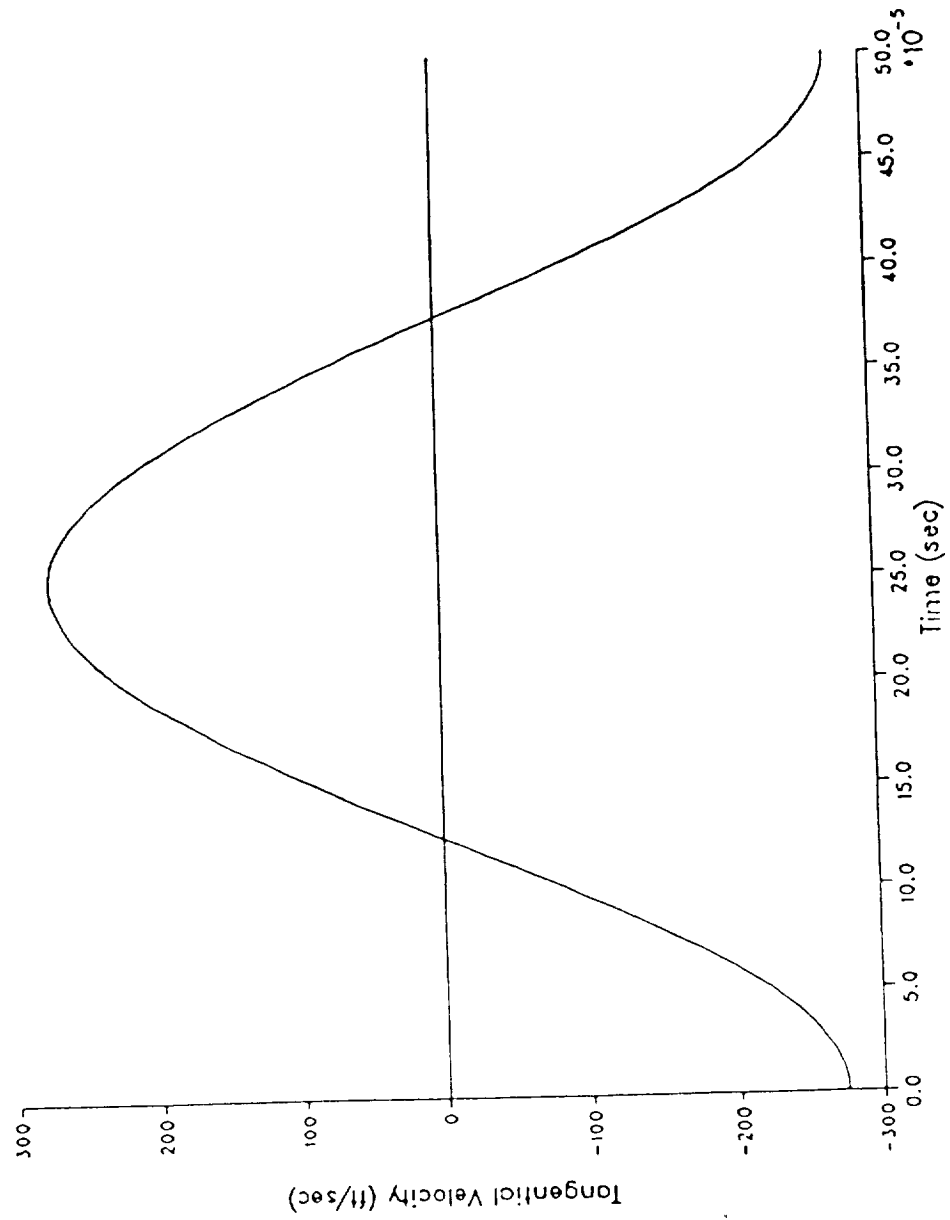


Figure 2.4 : Acoustic Tangential Velocity Component.



instantaneous burning rate,  $W$  and the mean burning rate,  $\bar{W}$  as follows:

$$w' = \frac{W - \bar{W}}{\bar{W}} \quad (27)$$

where

$$\bar{W} = \frac{1}{2\pi} \int_0^{2\pi} W d(\omega t) \quad (28)$$

$p'$  is the perturbation component of the pressure at  $r$  and  $\theta$ , and  $p^*$  is a sinusoidal function having the same amplitude and frequency but 90 degrees out of phase with  $p'$ .

The response factors defined by equations (25) and (26) vary with radial and circumferential position because of the spatial variations of the pressure and the burning rate. To account for this effect, the chamber cross-sectional area is divided into a number of equal areas in both radial and circumferential directions. The response factors are calculated at each of the centers of the areas. They are then averaged to obtain the resultant response factor.

### 2.3 Calculation Procedure

The following procedure is used to calculate the in-phase and the out-of-phase response factors as functions of the frequency:

1. Beginning with the first droplet injected at time  $t$ ; equations (1) to (9) are used to calculate the instantaneous gas pressure, velocity, temperature and density at the time  $t$ .
2. Calculate the mean gas velocity using equation (10).

3. Calculate the film temperature using equation (15). Evaluate the gas transport properties, i.e. thermal conductivity, viscosity, etc... at the film temperature.

4. Calculate the vapor pressure at the droplet temperature. This step is skipped when the droplet temperature reaches the boiling temperature or the critical temperature of the propellant.

5. Calculate the heat and the mass transfer coefficients using equations 13 and 14. The mass transfer coefficient need not be calculated when the droplet temperature reaches the boiling temperature or the critical temperature of the propellant.

6. Calculate the mass evaporation rate using equation (12). If the droplet is at the boiling temperature or the critical temperature of the propellant, the mass evaporation rate is calculated by solving equation (11) with the droplet-heat-up term being zero. Store the time and the corresponding mass evaporation rate of the droplet in an array for use at a later time to calculate the total evaporation rate.

7. Calculate the droplet drag and the droplet acceleration using equations (17) and (20). Calculate the rate of change of droplet radius using equation (22), and the rate of change of the droplet temperature using equation (21). The rate of change of the droplet temperature is set to zero when the droplet temperature reaches the boiling temperature or the critical temperature of the propellant.

8. Calculate the droplet radius, velocity and temperature at

the next time step using the Euler's explicit integration illustrated in equation (23).

9. Time is incremented and steps 1 to 8 are repeated until more than 99 percent of the droplet mass has been evaporated.

10. Repeat steps 1 to 9 for each of the droplets injected from the same radial and tangential location but at different times during a period of oscillation.

11. Calculate the total evaporation rate of the entire droplet array from the evaporation histories of each of the droplets which have been stored in step 6.

12. Calculate the in-phase and the out-of-phase response factors using equations (25) and (26).

13. Repeat steps 1 to 12 to obtain the response factors for droplets injected from different radial and circumferential locations.

14. Average the response factors calculated in step 13 in order to obtain the overall response factor.

The whole procedure is then repeated to calculate the overall response factors for other values of the frequency.

### III. PROGRAM DESCRIPTION

The Combustion Response Prediction (CRP) computer code consists of a main program, four subroutines and six function subroutines. The main program and the subroutines are described in the next section. Program input and output are described in sections 3.2 and 3.3. A listing of the computer code is provided in appendix A. Input and output for a sample case are provided in appendix B.

#### 3.1 Program Description

\* Main Program: reads all input to the code which include a problem-description title, chamber acoustic resonance modes, frequency range of interests, combustion gas properties, and droplet initial properties. It calculates the vaporization histories of each of the droplets in an array and calls subroutine OUTPUT, if the input variable DEBUG is set to TRUE, to print out the vaporization histories. Each array is a set of several droplets injected from the same radial and circumferential injection location but at different times. It then calls subroutine SUM to compute the total evaporation rate of the entire array of droplets. Next, it calculates the response factor of the droplets in the array. The procedure is then repeated to calculate the response factors for other arrays of droplets injected from other locations. Finally, it averages the response factors calculated at

various locations to obtain the overall response factor at a particular frequency. This whole procedure is then repeated for a number of frequencies. Loops over all droplets, injection locations, and frequencies are made in the main program.

- \* Subroutine OUTPUT: called by the main program to print intermediate results, e.g., time step, instantaneous radius, evaporation rate, and velocity of the droplet. It also prints out other information for debugging, e.g., Reynolds number, diffusion coefficient, vapor pressure.
- \* Subroutine MATHUR: calculates the parameter used in the Mathur and Thodos' correlation to calculate the binary diffusion coefficient as a function of pressure and temperature.
- \* Subroutine REIDEL: calculates the parameters used in the Reidel's correlation to calculate the vapor pressure as a function of temperature.
- \* Subroutine SUM: calculates the total evaporation rate of an array of droplets by summing the individual evaporation rates of each of the droplets.
- \* Function FWDOT: given the evaporation history of a single droplet, the function calculates the evaporation rate of the droplet - or of other droplets injected at a number of oscillation periods apart - at a specified

time.

- \* Function BJ: compute the second or higher-order Bessel function of the first kind.
- \* Function BJ0: compute the zeroth-order Bessel function of the first kind.
- \* Function BJ1: compute the first-order Bessel function of the first kind.
- \* Function SL: calculates the heat of vaporization as a function of temperature using Watson's correlation.
- \* Function PV: calculates the vapor pressure as a function of temperature using Reidel's correlation.

### 3.2 Input Description

All input with the exception of the problem description title are made using Fortran namelists. The problem description title can be specified using any number of lines but at least one line must be used although it can be a blank line. Following the problem description are the namelists INPUT, WAVES, CBGAS and DROPS. Variables in these namelists are described in tables 3.1 through 3.4. Input for a sample case is provided in appendix B.1.

Namelist INPUT is used to input artificial parameters such as the number of droplets injected per cycle, the number of time steps between output, and the number of time steps in a droplet life time, etc. If the variable DEBUG is set to TRUE, the vaporization histories of droplets and the intermediate results are written to a

Name	Type	Unit	Description and Remarks
DEBUG	L		=TRUE, Intermediate results are output to a debug file =FALSE, No intermediate output
NDTFQ	I		Number of time steps in one period of oscillation.
NDTLF	I		Number of time steps in the droplet life time
JA	I		Number of time steps between print-outs of intermediate results
NP	I		Number of droplets injected per cycle
NY	I		Number of integration steps used in the calculations of the response factors
NRAD	I		Number of radial injection locations
NCIRC	I		Number of circumferential injection locations

Table 3.1: Descriptions of Namelist INPUT Variables

Name	Type	Unit	Description and Remarks
FREQ	R	cps	First frequency value
DFREQ	R	cps	Frequency increment
NFREQ	I		Number of frequencies
DPF	R		Maximum pressure amplitude normalized by the mean pressure
PHIF	R	deg	Pressure oscillation phase angle
MTANG	I		Tangential resonance mode number
NRADI	I		Radial resonance mode number

Table 3.2: Descriptions of Namelist WAVES Variables

Name	Type	Unit	Description and Remarks
PO	R	psf	Mean chamber pressure
TFO	R	°R	Mean chamber temperature
GAMMA	R		Chamber gas specific heat ratio
PCB	R	psf	Chamber gas critical pressure
TCB	R	°R	Chamber gas critical temperature
TBB	R	°R	Chamber gas normal boiling point
EMB	R	lbm/lb-mole	Chamber gas molecular weight
PR	R		Chamber gas Prandtl number
AKB	R	Btu/ft-s-°R	Chamber gas thermal conductivity
VIS	R	lbm/ft-s	Chamber gas viscosity
VGAF	R	ft/s	Chamber gas final velocity

Table 3.3: Descriptions of Namelist CBGAS Variables

Name	Type	Unit	Description and Remarks
SIT	R	ft	Initial droplet radius
TO	R	°R	Initial droplet temperature
RHOL	R	lbm/ft**3	Propellant liquid density
CPL	R	Btu/lbm-R	Propellant liquid specific heat
VDI	R	ft/s	Initial droplet velocity
PCA	R	psf	Propellant critical pressure
TCA	R	°R	Propellant critical temperature
TBA	R	°R	Propellant normal boiling point
EMA	R	lbm/lb-mole	Propellant molecular weight
SLA	R	Btu/lbm	Propellant heat of vaporization at normal boiling point

Table 3.4: Descriptions of Namelist DROPS Variables

debug file.

The variables NDTFQ and NDTLF are used to calculate the time step for use in the calculation of the vaporization histories of the droplets. The time step based on the oscillation period is:

$$\Delta t_f = \frac{2\pi}{\omega * NDTFQ}$$

where  $\omega$  is the angular frequency of the oscillation, and the time step based on the droplet life time,  $t_{99}$  is:

$$\Delta t_{99} = \frac{t_{99}}{NDTLF}$$

The time step used in the calculations of the vaporization histories is the smaller of the two time steps.

The variable NY is the number of integration steps used in the numerical integration of equations (25) and (26) to obtain the in-phase and the out-of-phase response factors. For cases where the temporal variations of the droplet vaporization rates are steep, a large value must be specified for NY (finer integration step size) to avoid losing accuracy in the results.

The overall response factor is the average of the response factors calculated at various radial and tangential locations. The number of the locations are specified by the variable NRAD and NCIRC.

All of the variables described above are artificial parameters in the model. The higher the value specified for these parameters,

the more accurate the solutions will be and, of course, the more computer time is required. Recommended values for the variables are given in table 4.1 of section 4.2.

Namelist WAVES is used to input the chamber acoustic resonance modes, frequency range of interests, the maximum amplitude and the phase angle of the pressure oscillations. The variables FREQ, DFREQ, and NFREQ specify the frequency calculation domain. The amplitudes of the pressure oscillations vary spatially, the variables DPF is the maximum amplitude,  $\Delta p$ , used in equations (1) and (4).

Namelist CBGAS is used to input the combustion gas properties, for example, mean pressure and temperature, critical pressure and temperature, molecular weight, Prandtl number, etc. In general, the combustion gas is the mixture of several gas components. Therefore, the pseudo values of the critical pressure, temperature and molecular weight of the mixture must be calculated. These values can be calculated from the compositions of the mixture and the properties of the species components that constitute the mixture. The compositions, the thermodynamic properties and the transport properties can be calculated using the standard TRAN72 computer program (Ref. 12)

Namelist DROPS is used to input droplet properties such as its initial radius, velocity, and temperature. The namelist is also used to input thermodynamic properties of the propellant. These properties can be obtained from the Aerojet Handbook of Properties and Performance of Liquid Rocket Propellant (Ref. 5).

### 3.3 Output Description

Output from the code begins with the echo of input data which include the problem-description title, and the values of the namelists' variables. Although the problem description can be input using any number of lines, only the first line is output. Next, the descriptions and the values of selected input variables are output. Following the echo of the input data is the estimated droplet life time. The last section of the output is the calculated stability results which include the real part (in-phase) and the imaginary part (out-of-phase) of the response factor. The results are also output in the polar form (magnitude and phase angle). These results are output for each of the frequencies whose range is specified in the namelist WAVES. A sample output file is provided in appendix B.2.

Three additional files are also generated to be input to TELLEGRAF, a computer graphic program available at ATC, for plotting the calculated results. The first file contains the data used for plotting the real and the imaginary parts of the response factor versus the frequency. The second and the third files contain the data used for plotting the magnitude and the phase angle of the response factor, respectively, versus the frequency. Using the VAX conventions for file identifications, the file names of the three files are the same as that of the input file. The file types of the three files are RIM, MAG, and PHA, respectively.

In addition to the files described above, a debug file is also generated if the variable DEBUG in the namelist INPUT is set to

TRUE. This file contains intermediate results that are useful for debugging purposes. Quantities written to this file begin with the values of the parameters used in the vapor-pressure correlation. Next, the vaporization history of the first droplet <sup>is</sup> ~~are~~ output. The vaporization history information includes the following quantities as functions of time: the droplet radius and its time rate of change, the droplet temperature, the vapor pressure, the mass evaporation rate, and the absolute and the relative velocities of the droplet. Other quantities output along with the droplet vaporization history include the instantaneous gas pressure and temperature, the diffusion coefficient, the Reynolds number, and the heat transfer coefficient. In addition, at each of the specified frequency<sup>ies</sup>, the response factors calculated at various radial and ciurcumferential injection locations are output. The file name of this debug file is the same as that of the input file, the file type of the debug file is DBG.

#### IV. RESULTS AND DISCUSSION.

Calculations of the response factors were made for the case of n-heptane droplets vaporizing in combustion gases composed of the products of the stoichiometric reaction with oxygen. The results are discussed in section 4.3.

Before the calculations were made for the above case, several tests were performed to study the effects of the "artificial" parameters, e.g. time steps, number of droplets injected per cycle, on the solutions. The test results are discussed in section 4.2.

In addition, the model was used to calculate the evaporation history of Chinese Kerosene (RP2) and n-heptane droplets vaporizing in steady environments (no oscillations in pressure and velocity) for which experimental data are available. This provides, to some degree, the verifications of the single-droplet-vaporization model used in the present study. The following section discusses the calculated results and the comparisons of the results with the data.

##### 4.1 Calculations of Droplets Vaporizing in Steady Ambient Gases.

###### 4.1.1 Chinese Kerosene (RP2) Droplet Vaporizing in Airstream.

The model was used to calculate the evaporation history of a Chinese-Kerosene (RP2) droplet vaporizing in a steady hot crossflowing air stream, for which experimental data are available (Ref. 8). In the experiment, the initial diameter and temperature

of the droplet are 1.15 mm and 294°K. The temperature and the velocity of the air stream are 516°K and 4.85 m/s. The experiment was conducted at atmospheric pressure. Three sets of calculations were made using three different assumptions on the droplet thermal conductivity - finite thermal conductivity with spherically symmetric temperature distribution inside the droplet, zero thermal conductivity (onion skin), and uniform droplet temperature (infinite thermal conductivity or strong circulating flows inside the droplet). Figure 4.1 shows the calculated temporal variations of the droplet diameter and the experimental data. It can be seen that only the prediction using uniform-droplet-temperature assumption agrees well with the data. Based on these comparisons and the other justifications discussed in section 2.2.1, the uniform-temperature-assumption was selected for use in the present study.

#### 4.1.2 n-Heptane Droplet Vaporizing in Quiescent Nitrogen Gas.

The uniform-droplet-temperature model is then used to calculate the evaporation histories of n-heptane droplets vaporizing in hot quiescent nitrogen gas at 1 atmosphere and at 50 atmospheres. In both cases, the initial droplet diameter and temperature are 0.0354 in. and 527°R. The ambient nitrogen temperature is 1031°R. Following Baer's practice (Ref. 9), natural convection effects are accounted for in the present calculation by replacing the Reynolds number in the heat and the mass transfer correlations with the square root of the Grashof number. Figures 4.2 and 4.3 show the comparisons between the calculated results and the data taken from reference 9 for both

Figure 4.1 : Chinese RP-2 Kerosene Droplet Vaporizing in Hot Air.

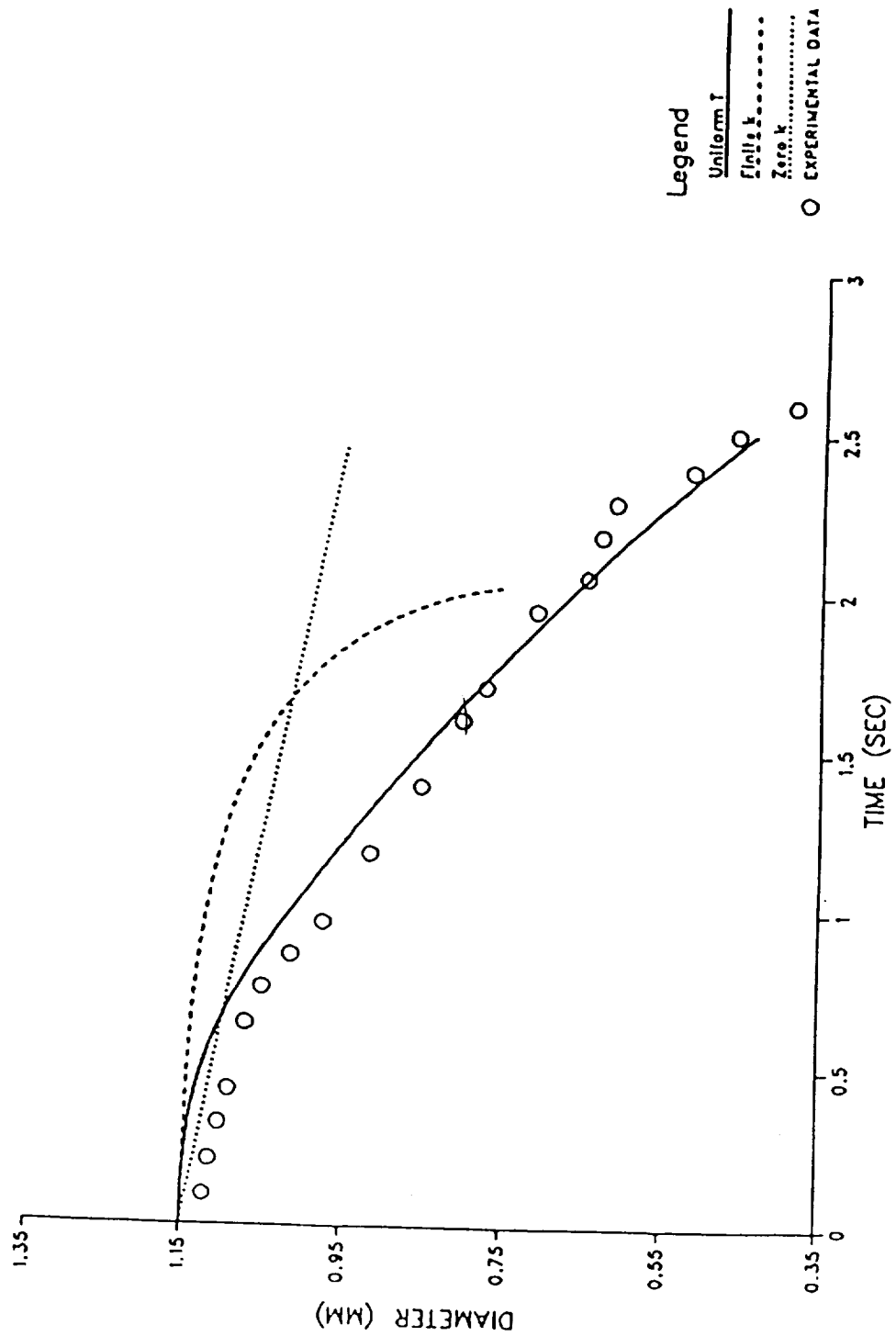


Figure 4.2 : n-Heptane Droplet Vaporizing in Nitrogen Gas at 1 atm.

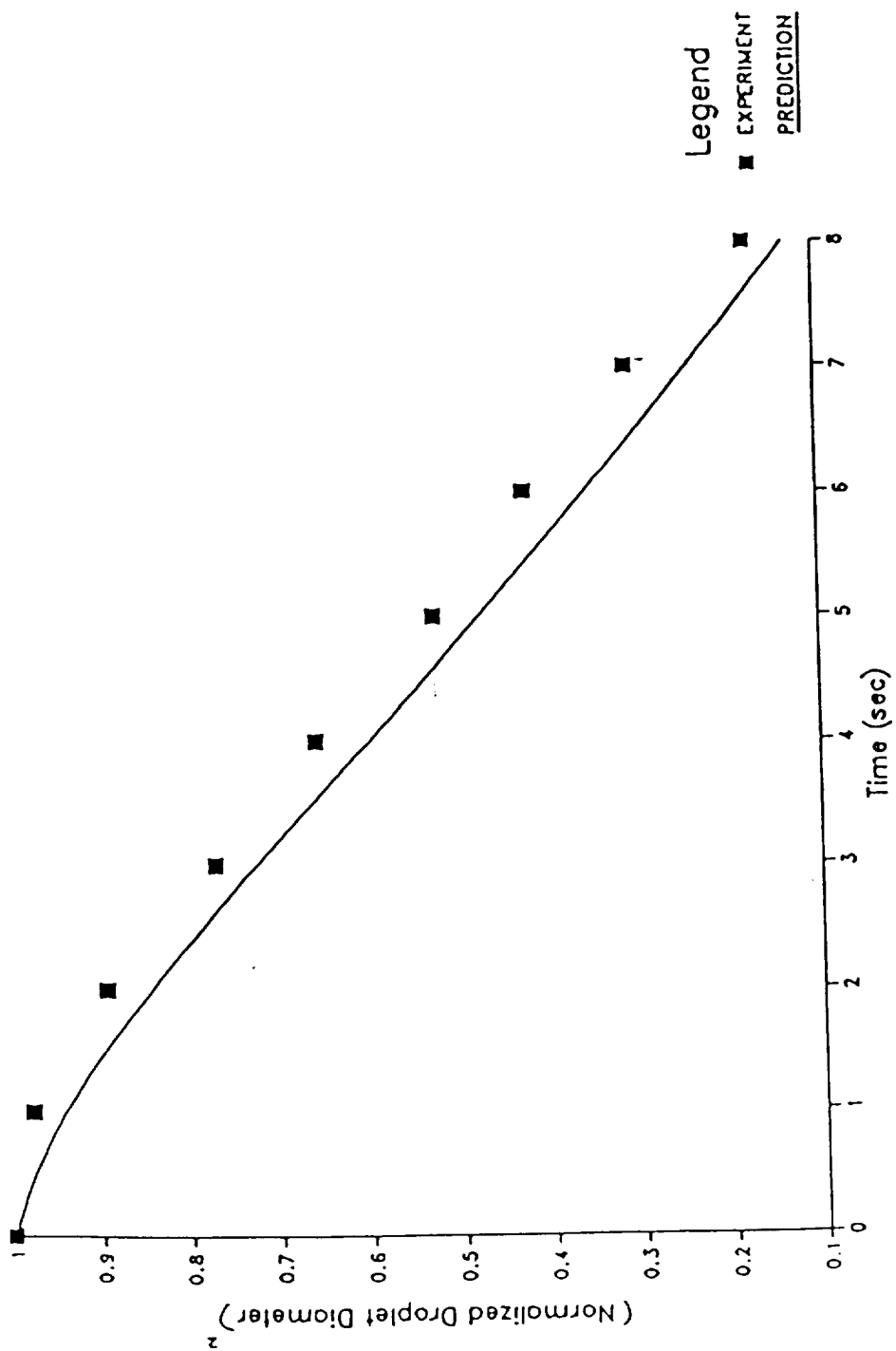
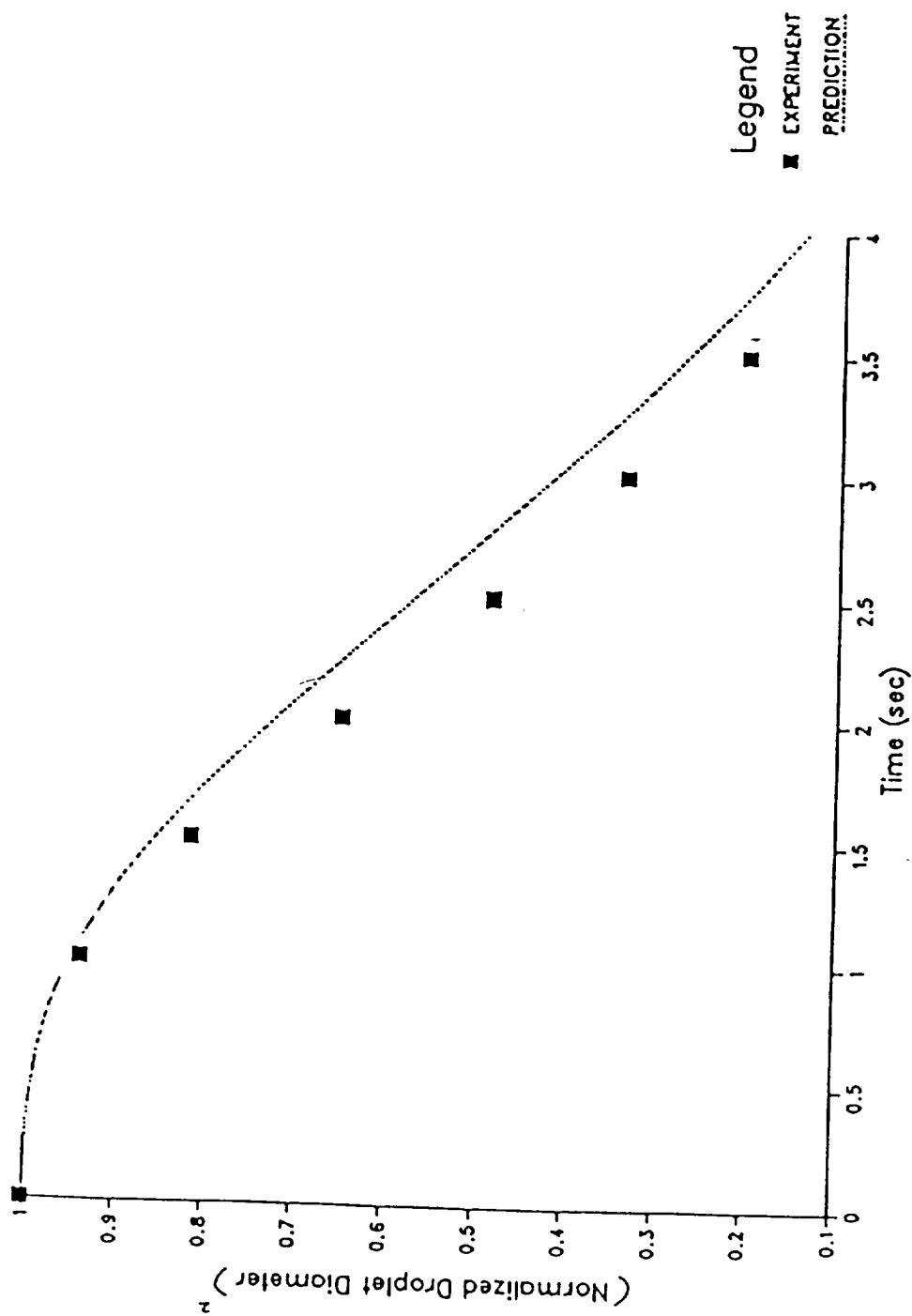


Figure 4.3 : n-Heptane Droplet Vaporizing in Nitrogen Gas at  
50 atm.

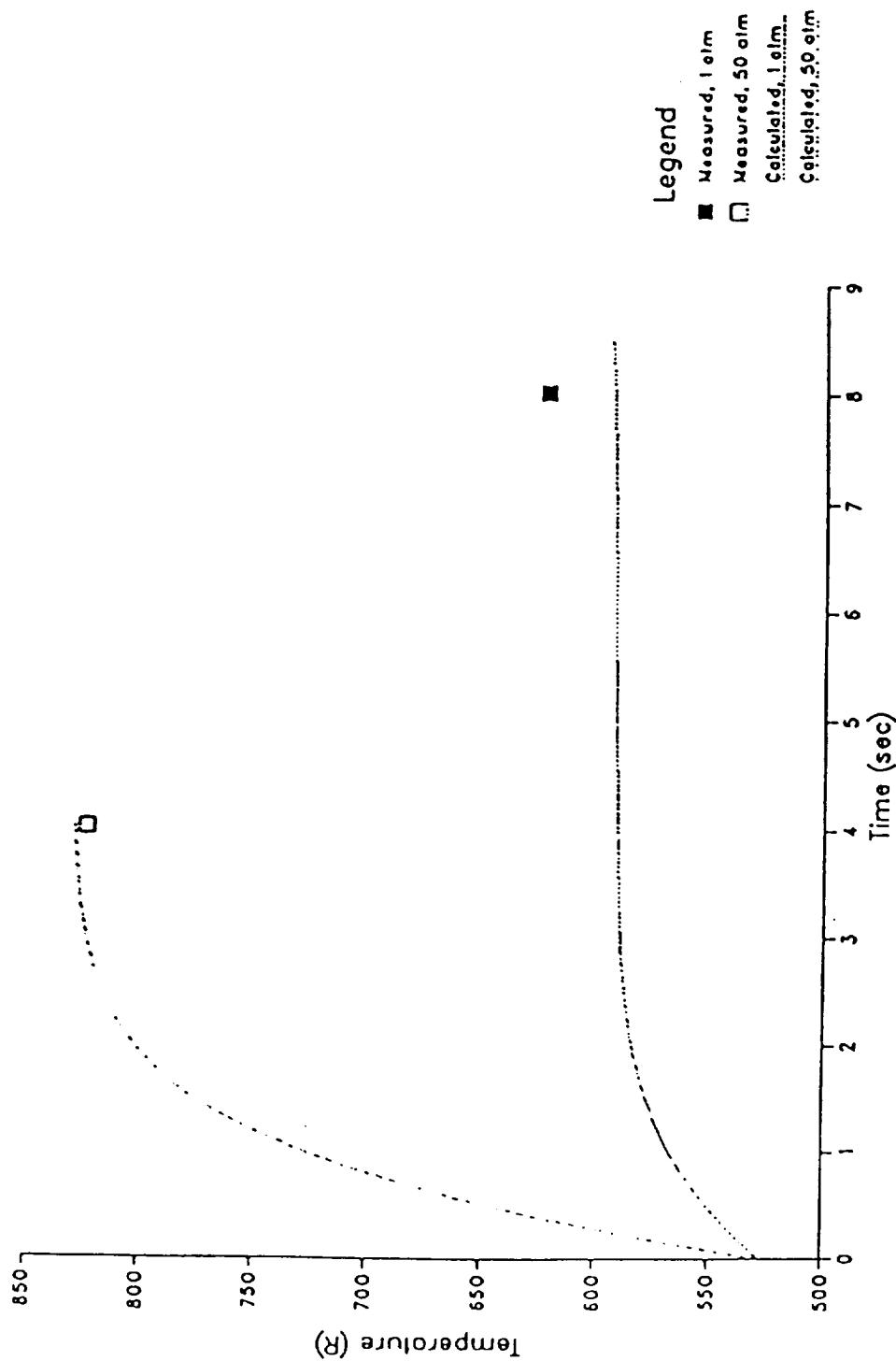


cases. The data were originally reported in reference 10. In the figures, the square of the normalized (by the initial diameter) droplet diameter is plotted versus time. It can be seen from the figure that the rate of change of the droplet diameter is overpredicted for the lower pressure case while it is underpredicted for the higher pressure case. Nevertheless, the agreements between the calculated results and the data are good. The calculated droplet temperatures for the two cases are shown as functions of time in figure 4.4. The measured wet-bulb temperatures are also shown in the figure. Comparisons between the calculated results and the data show that the wet-bulb temperature is underpredicted in the low-pressure case, and it is well predicted for the high-pressure case.

#### 4.2 Parametric Test Results

A limited number of parametric tests were performed to study the sensitivity of the solutions to the artificial parameters which have been introduced into the model. The tests were made also to obtain guidelines on what values should be specified for the parameters. While no attempts are made to describe the test results in detail, they are briefly discussed in this section. Test results show that the solutions converge as the time step decreases. The test results also show that the solutions converge as the number of droplets injected per cycle, the number of radial injection locations, the number of circumferential injection locations, and the number of integration steps used in the calculation of the response factor increase.

Figure 4.4 : Temporal Variation of the Temperatures of the n-Heptane Droplets Vaporizing in Quiescent Nitrogen Gases.



As a result of the parametric tests, the values listed in table 4.1 are recommended for the artificial parameters in order to obtain reasonable accuracy in the solutions without substantial CPU time requirements. A typical computer run using the values recommended in table 4.1 requires approximately 2 CPU minutes for each frequency on the VAX computer system at ATC. The total computer time requirement is approximately linearly proportional to the number of frequencies, NFREQ. It should be noted from the table that while the recommended value for the number of radial injection locations, NRAD, increases with the specified radial resonance mode, the recommended value for the number of tangential injection locations, NCIRC does not vary with the specified tangential mode. The reasons are that the response factors, at a given radial coordinate, do not vary with the tangential locations for the spinning waves; and that the symmetry of the standing waves have been accounted for.

#### 4.3 n-Heptane Droplet Response Factor

The computer code CRP, was used to calculate the combustion response factor of n-Heptane droplets vaporizing in the combustion gases composed of the products of the stoichiometric reaction of n-Heptane with Oxygen. The mean chamber pressure is 300 psia and the mean chamber temperature is 6280°R. The acoustic mode in the chamber is the first tangential (1T) standing mode with the maximum amplitude of the pressure oscillation normalized by the mean pressure <sup>being</sup> ~~is~~ 10 percent. The initial radius, temperature, and velocity of the droplet are 50 microns, 535°R and 50 ft/s, respectively. The number of droplets injected per cycle used in

NDTFQ	=	16
NDTLF	=	1000
JA	=	100
NP	=	16
NY	=	3200
NRAD	=	$3 * (NRADI + 1)$
NCIRC	=	3

Table 4.1: Recommended Values for the Artificial Parameters

this calculation is 16.

Figures 4.5, 4.6, and 4.7 show the temporal variations of the temperature, the radius, and the vaporization rate, respectively, of an n-Heptane droplet injected at the normalized radial distance of .707 and the tangential location is at the pressure antinode. The discontinuities in these curves are the results of the switching to a different procedure to calculate the vaporization rate when the droplet temperature reaches the boiling temperature of the propellant. In all of the figures, time equals to 0.0 is the instant the droplet is injected into the chamber, and for this particular droplet it is injected at the beginning of the oscillation period. The frequency of the acoustic fields in the chamber is 2000 Hertz.

Figure 4.5 shows that the droplet temperature reaches the boiling temperature of the propellant at approximately 62 micro-seconds. Figure 4.6 shows that the droplet completely vaporizes at approximately 82 micro-seconds. It can be seen from figure 4.7 that initially, the evaporation rate of the droplet is relative small. The rate slowly increases for approximately 62 micro-seconds. At this time the temperature of the droplet reaches the boiling point of the propellant, the vaporization rate rises sharply to a maximum then it decreases as the surface area of the droplet reduces. A small variation at the peak of the vaporization rate is the result of the oscillations in the chamber gas properties.

Figure 4.8 shows the total evaporation rate of an array of

Figure 4.5 : Temporal Variation of n-Heptane Droplet Temperature.

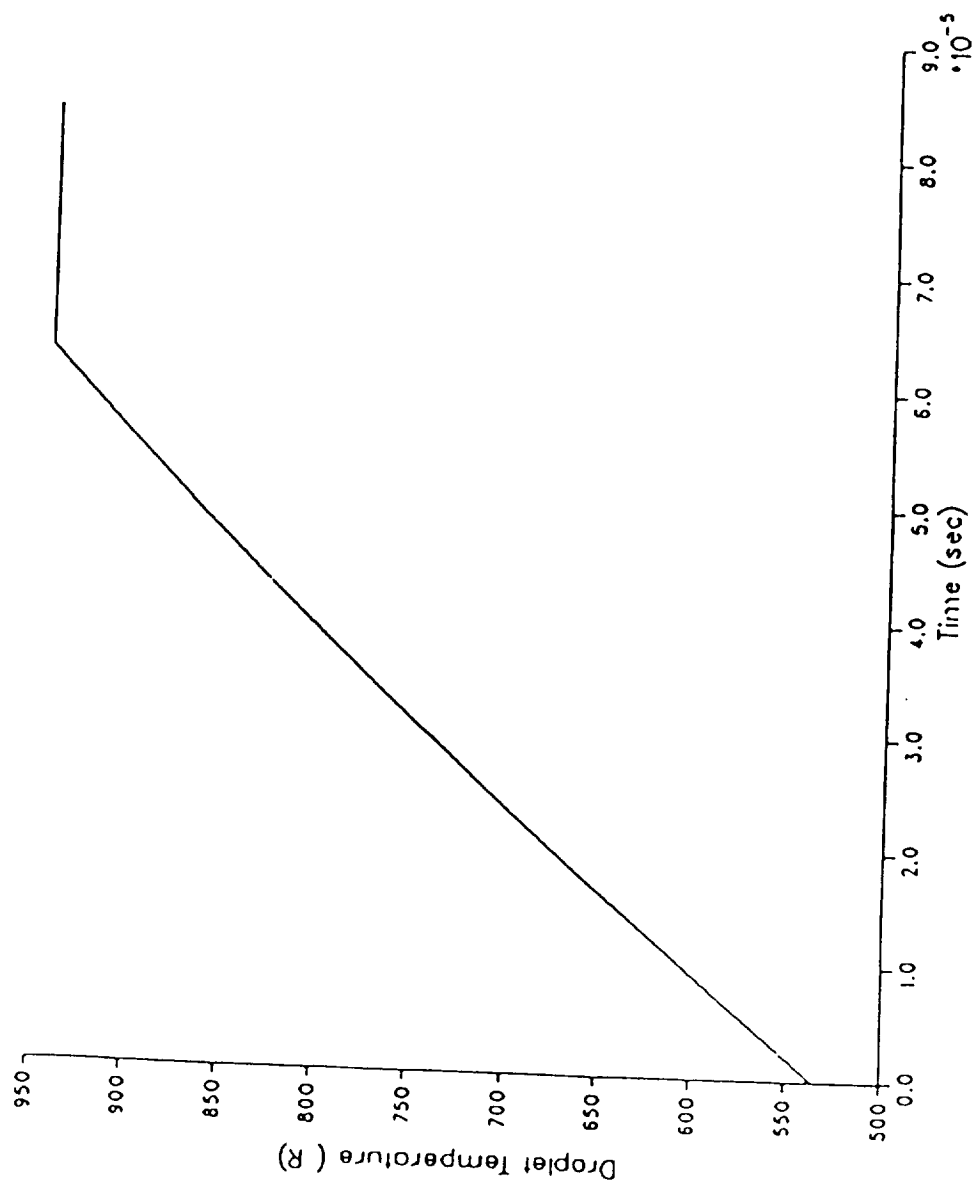


Figure 4.6 : Temporal Variation of n-Heptane Droplet Radius.

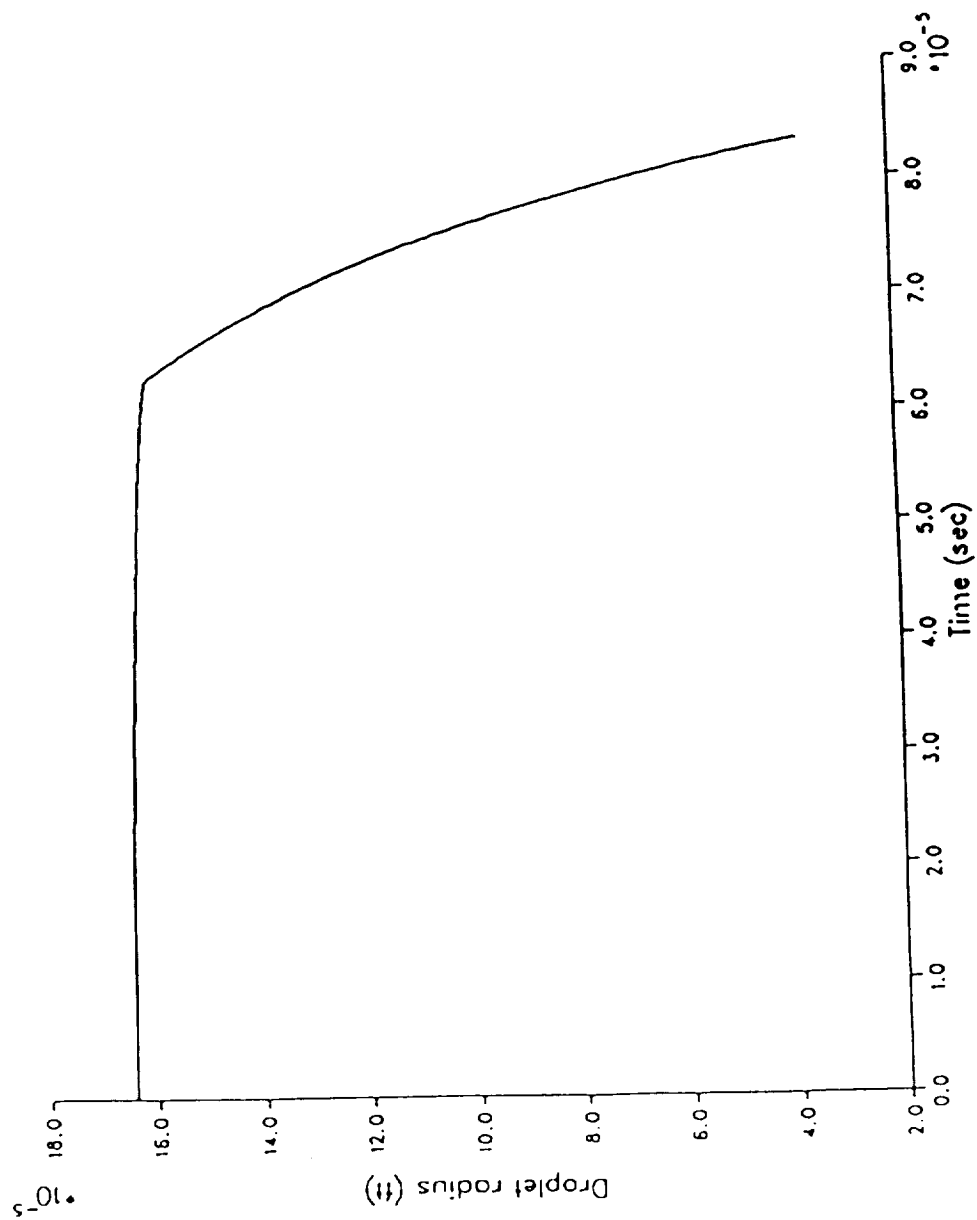


Figure 4.7 : Temporal Variation of n-Heptane Droplet Vaporization Rate.

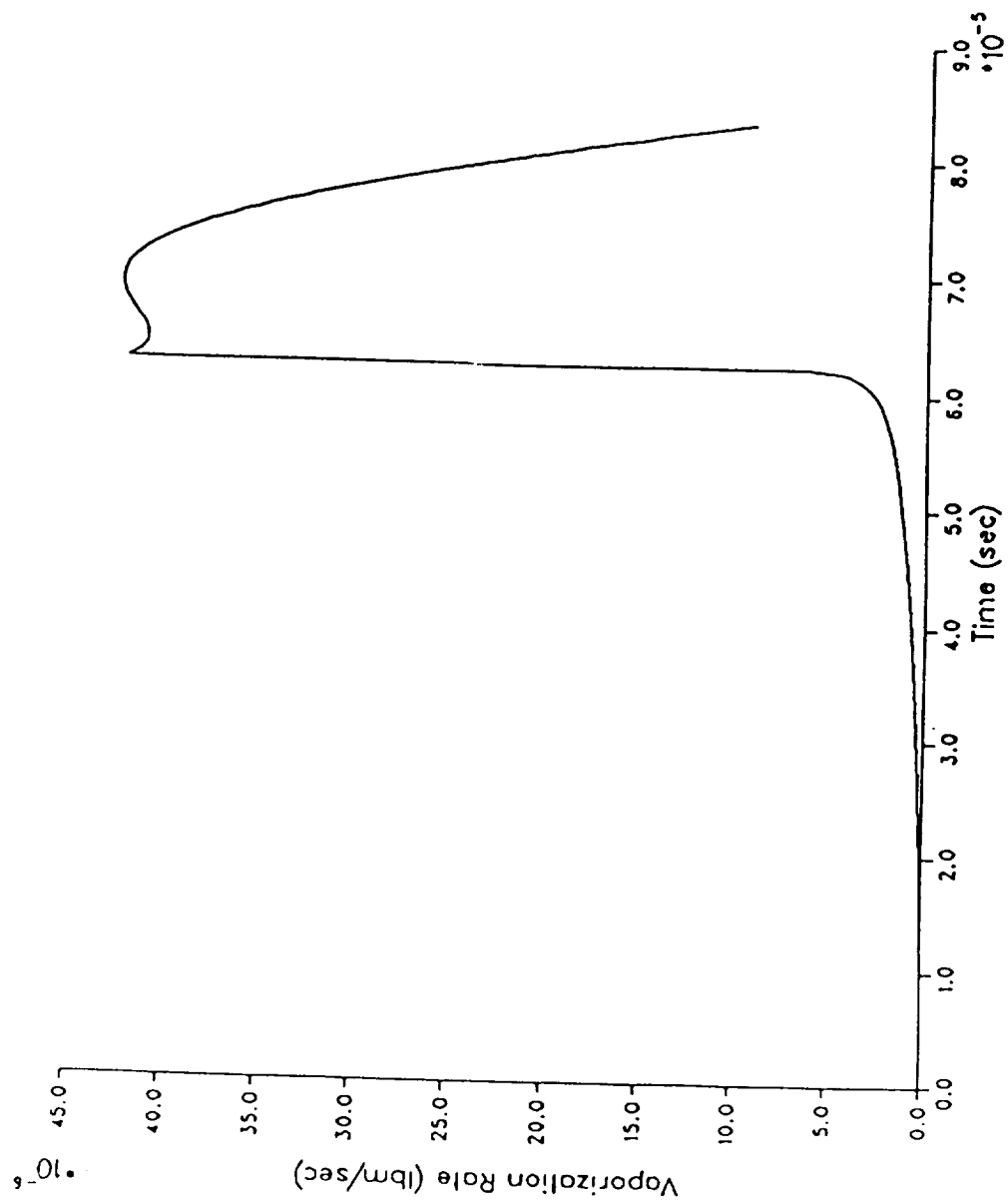
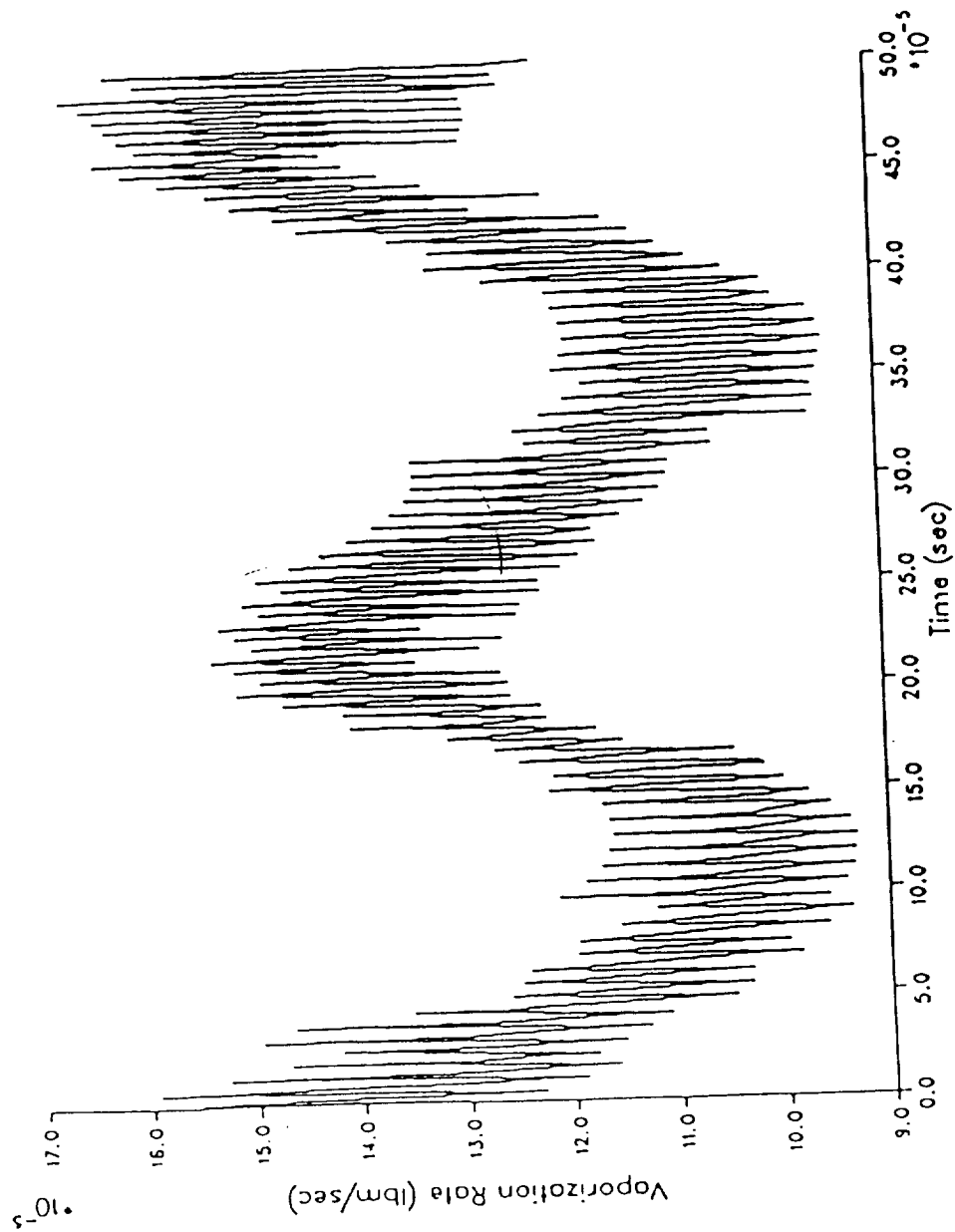


Figure 4.8 : Total Vaporization Rate of an Array of Droplets.



droplets. This plot is generated from a different computer run in which the number of droplets injected per cycle was increased to 80 so that the induced perturbation of the vaporization rate is more pronounced. The sharp spikes in the total vaporization curve are the results of the use of a finite number of droplets to represent the continuous injection of the propellant. The absolute amplitudes of the spikes are independent of the number of droplets injected per cycle. The mean burning rate is, however, approximately proportional to the number of droplets injected per cycle. Thus, as the number of droplets injected per cycle is increased, the spike amplitudes become smaller relative to the mean of the burning rate.

Finally, the calculated combustion response factors are shown in Figures 4.9, 4.10, and 4.11. Figure 4.9 is a plot of the real and the imaginary parts of the combustion response factor as functions of the frequency. Figures 4.10 and 4.11 show the combustion response factor in a different form. In these figures, the magnitude and the phase angle of the combustion response factor are plotted versus the frequency.

In the past, the combustion response is assumed to be characterized by a pressure interaction index,  $n$  and a sensitive time-lag,  $\tau$ . The values of  $n$  and  $\tau$  are determined empirically. The combustion response factor,  $Y_b$  is related to the pressure interaction index and the sensitive time lag in reference 13 as follows:

$$Y_b = n \left( 1 - e^{-i\tau\omega} \right)$$

Figure 4.9 : Calculated Real and Imaginary Parts of the Combustion Response Factors as Functions of the Frequency.

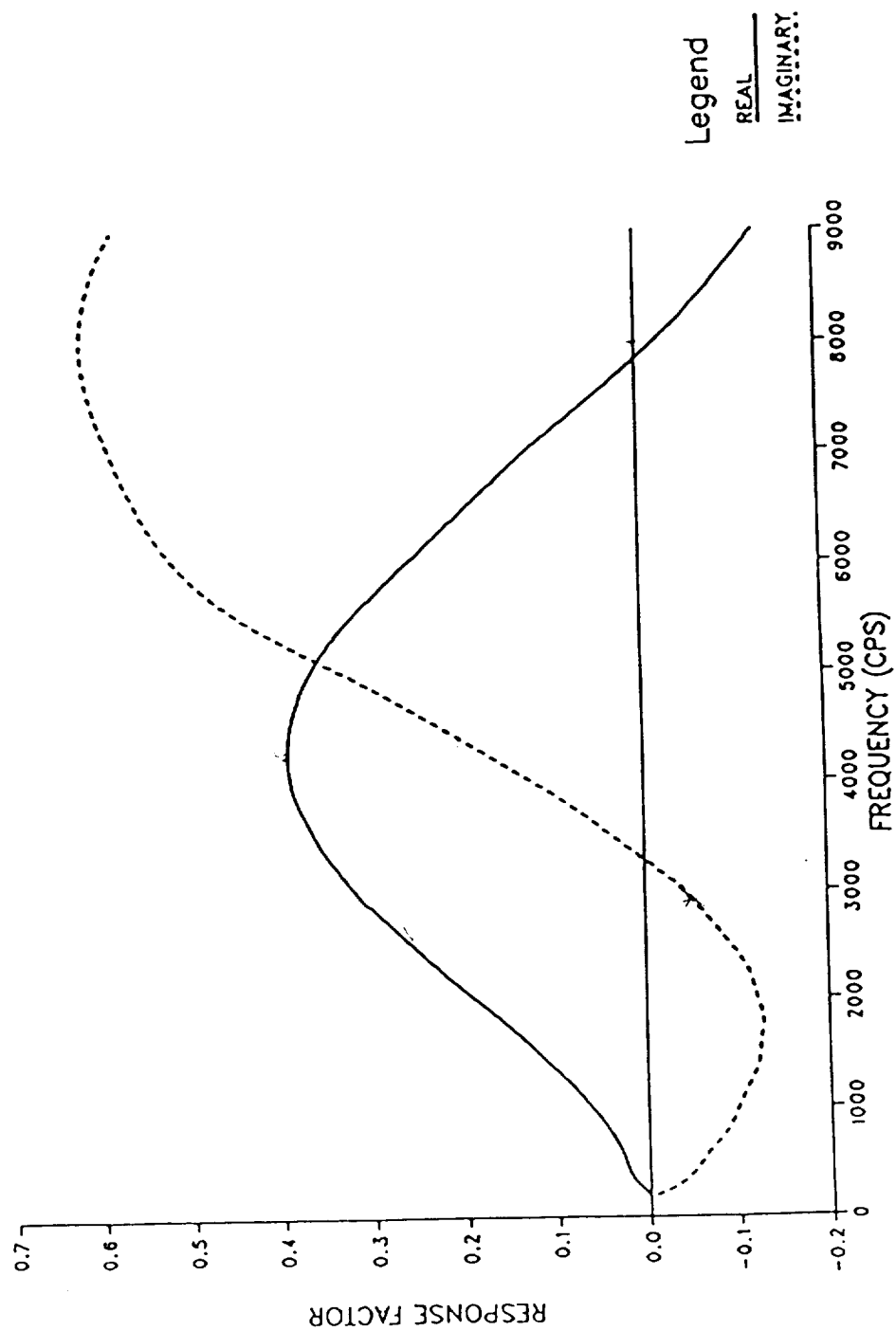


Figure 4.10: Calculated Magnitude of the Combustion Response Factors as Functions of the Frequency.

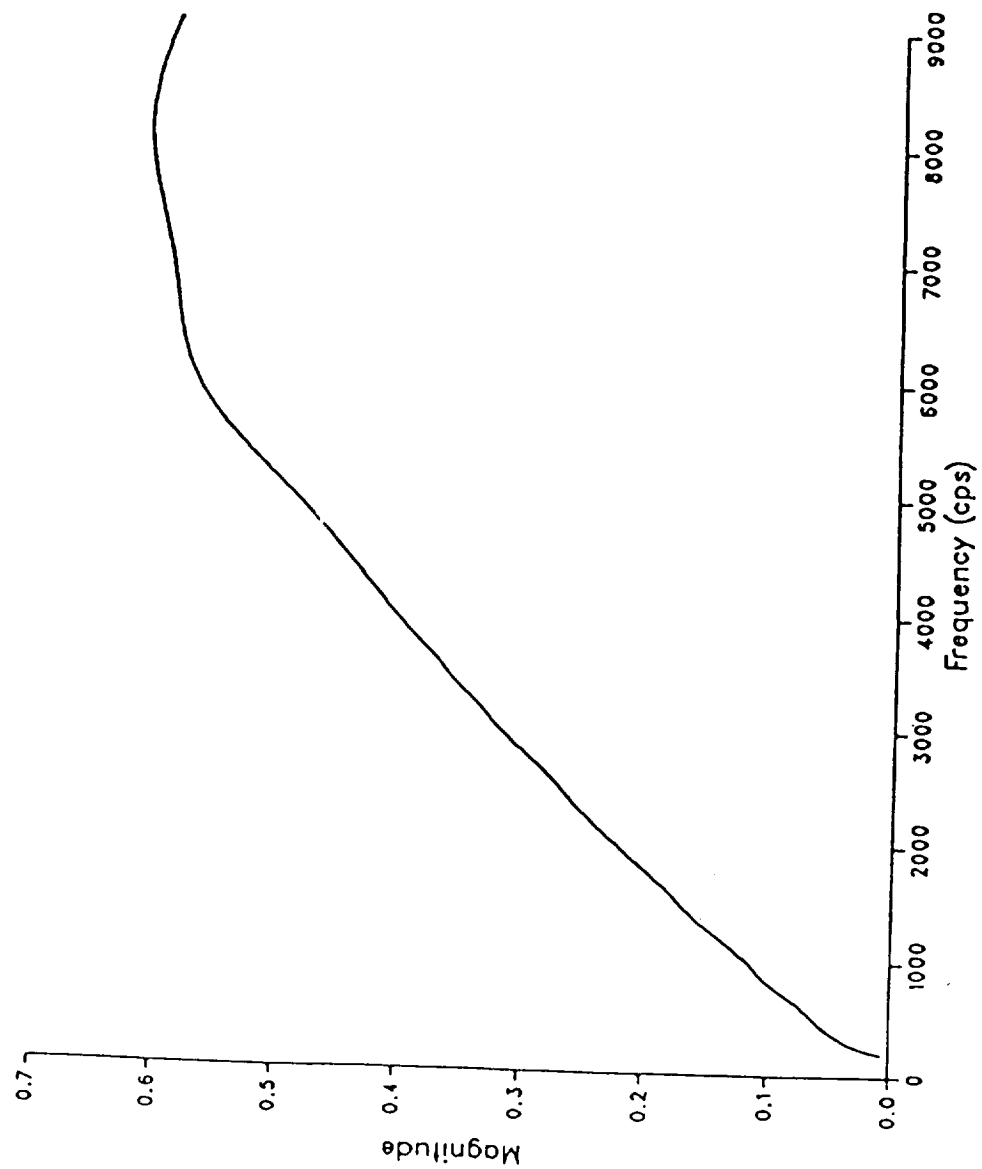
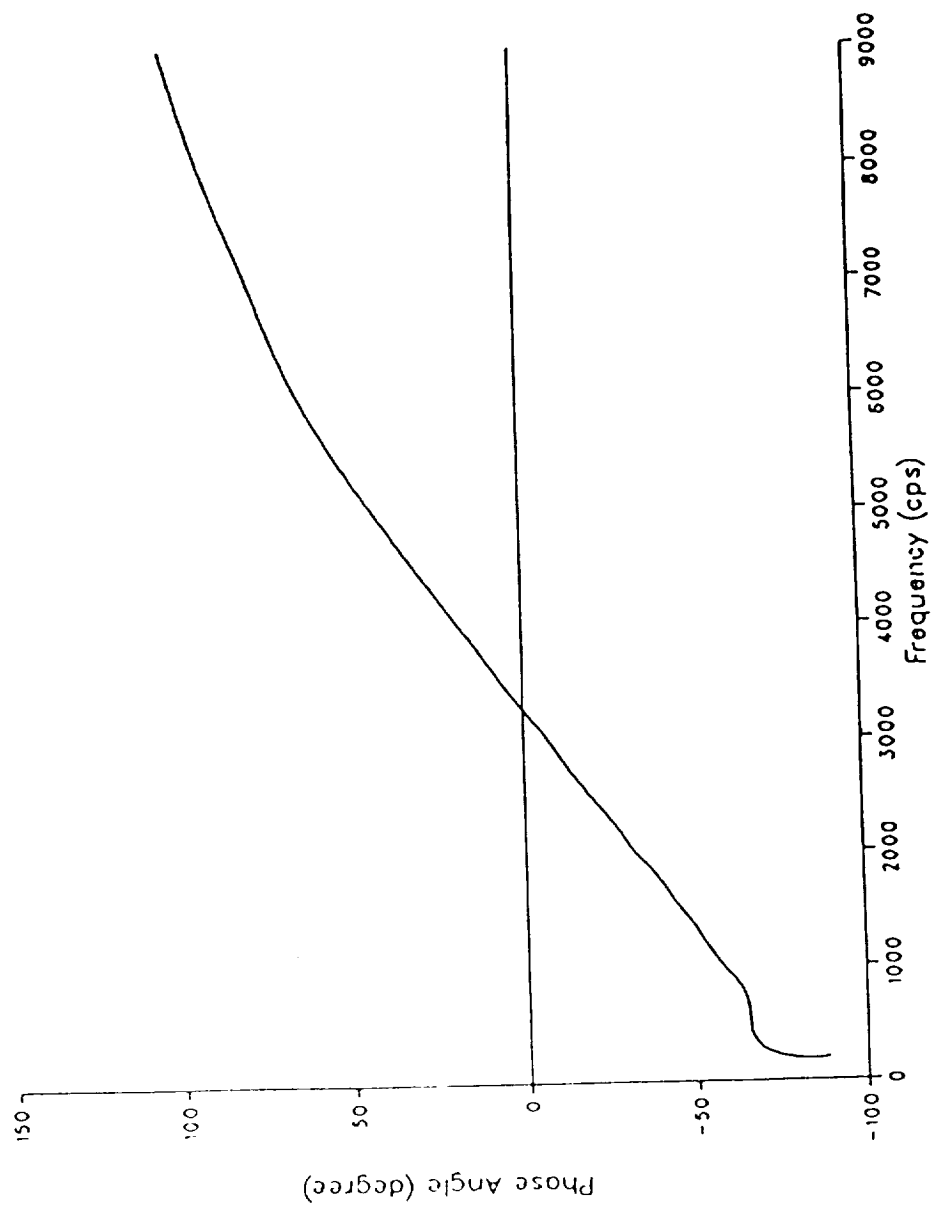


Figure 4.11: Calculated Phase Angle of the Combustion Response Factors as Functions of the Frequency.



or

$$Y_b = n \left\{ 1 - \cos(\tau\omega) + i \sin(\tau\omega) \right\}$$

The above equation shows that the real part of the response factor is a function of the angular frequency. The function has a peak value of  $2n$  at the angular frequency equal to  $\frac{\pi}{\tau}$ . Thus, the values of  $n$  and  $\tau$  can be obtained from the correlation with the plot of the combustion response factor versus the frequency. For example, in figure 4.9 the real part of the response factor has a peak value of 0.4 at 4000 Hertz, the corresponding values of  $n$  and  $\tau$  for this case are 0.2 and 0.125 milli-seconds, respectively.

## V. CONCLUSIONS AND RECOMMENDATIONS

A model has been formulated and a computer code, CRP has been developed to calculate the combustion response factors. Several statements can be made with regards to the calculated results:

1. Three different assumptions about the thermal conductivity of the droplet - zero thermal conductivity, finite thermal conductivity with spherically symmetric temperature distribution inside the droplet, and uniform droplet temperature (infinite thermal conductivity or strong recirculating flows inside the droplet) - have been tested. The latter assumption appears to be superior because calculations of the vaporization rates of droplets vaporizing in steady environments using the assumption agree well with the experimental data. Furthermore, the analysis using the uniform-droplet-temperature assumption is simple and requires less computer time.
2. Time-history calculations of n-Heptane droplets vaporizing in Nitrogen gas and of Chinese Kerosene (RP2) droplets vaporizing in air agree well with the experimental results.
3. Calculations of a 100-micron n-Heptane droplet vaporizing in combustion gases composed of the products of the stoichiometric reaction with oxygen at 300 psia show that the life time of the droplet is approximately 83 micro-seconds. The droplet temperature reaches the

boiling point of the propellant at approximately 62 micro-seconds. Only 6 percent of the droplet mass has vaporized up to this point, thus most of the vaporization takes place at the boiling temperature of the propellant and near the end of the droplet life time. These results are calculated for the case where the frequency of the acoustic oscillations is 2000 Hertz.

4. A technique was devised to calculate the vaporization rate when the droplet temperature reaches the boiling point of the propellant. The technique eliminates the numerical problem (vaporization rate blows up logarithmically) that is inherent in many of the droplet vaporization studies in the past. Calculations of droplet vaporization histories using the technique appear to be qualitatively correct. The technique is extended to the cases where the droplet temperature reaches the critical temperature of the propellant. Thus, the model can be used for both sub-critical and super-critical chamber pressures.

In addition, the following recommendations are made with regards to the future work to improve the model and/or the computer code:

1. The mean axial velocity is currently calculated using equation 10.b which is not very realistic since it does not account for the spatial distribution of the mixture

ratio. The computer code should be modified so that the axial profile of the mean velocity can be specified as input if it is available apriori, for example, from the steady-state performance analysis.

2. Modifications to the computer code to account for the dependences of the propellant density and the combustion gas viscosity on the temperature are recommended. Currently, only the dependences of the vapor pressure, heat of evaporation, and the diffusion coefficient on the temperature are accounted for.
3. A typical run using the table 4.1 - recommended values of the artificial parameters for a 1T mode requires approximately 2 CPU minutes for each frequency on the VAX computer system at ATC. Conversion of the computer code for use on the PC computer systems is possible because the memory storage required by the code is relative small. This is highly recommended because there are no charges for using the PC's. Conversion of the code for use on the CRAY-XMP at the San Diego Supercomputer Center, to which Aerojet has the access, is also recommended since it is expected that the computer run time and the cost will be significantly lower.

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**PART D**

**NASA/LeRC NON-LINEAR  
INJECTION RESPONSE MODEL  
(LEINJ)**

# NONLINEAR INJECTION ELEMENT THEORY

KEVIN BREISACHER  
NASA/LEWIS RESEARCH CENTER  
March 6, 1989

## DOME MODEL

There are three options for the model used in the dome included in the subroutine. These options are:

1. Lumped Parameter ( DomInd = 1 )
2. Longitudinal Acoustic Wave ( DomInd = 2 )
3. 3 D Acoustic Wave ( DomInd = 3 )

The lumped parameter model assumes there are no spatial derivatives in pressure or velocity.

$$P = P_d \sin \omega t \quad (1)$$

$$V_t = \frac{W}{\rho A} \quad (2)$$

$$W = \frac{\partial P}{\partial t} \frac{\partial \rho}{\partial P} Vol \quad (3)$$

$$\frac{\partial \rho}{\partial P} = \frac{g}{a^2} \quad (4)$$

where  $A$  area of tube  
 $a$  the speed of sound  
 $g$  gravitational acceleration  
 $P_d$  magnitude of pressure oscillation  
 $V_t$  velocity in the tube  
 $Vol$  volume of manifold per element  
 $W$  flowrate per element

Therefore

$$V_t = \frac{g\omega Vol}{Aa^2} P_d \cos \omega t \quad (5)$$

Note: To obtain constant pressure in the dome set "Vol" very large. For no velocity oscillations at the entrance to the tubes, set "Vol" very small (.00001).

For options 2 and 3, the manifold was modelled using the theory of Maslen and Moore for oscillations in a fluid with finite Mach number flow. Only the linear terms (small amplitudes) are included which reduces to the following wave equation:

$$-\phi'_{tt} + \nabla^2 \phi' = M^2 \frac{\partial^2 \phi'}{\partial x^2} + 2M \frac{\partial \phi'_t}{\partial x} \quad (6)$$

where  $\nabla$  gradient  
 $M$  mach number  
 $t$  time derivative  
 $u$  velocity vector  
 $x$  axial derivative

The wave is assumed to be periodic in time and separable in the  $x$ ,  $r$ , and  $\theta$  coordinates, or

$$\phi' = J_n(mr) e^{in\theta} e^{i\omega t} (e^{izB_1} + C e^{izB_2}) \quad (7)$$

For no steady state axial velocity in the dome

$$P' = -\frac{\rho a^2}{g\omega} \phi'_t \quad (8)$$

$$V' = \frac{a^2}{\omega} \nabla \phi' \quad (9)$$

where  $B_1, B_2$  complex coefficients  
 $i$  unit complex  
 $J_n$  Bessel Function of order  $n$   
 $m$  argument of Bessel function  
 $n$  number of pressure nodes in  $\theta$  direction  
 $r$  radial direction  
 $t$  time  
 $x$  axial direction  
 $\theta$  tangential direction

Assuming no steady state velocity in the dome ( $v = 0$ )

$$B_1 = -B_2 = \left( \frac{\omega^2}{a^2} - \frac{m^2}{r_d^2} \right)^{1/2} \quad (10)$$

and

$$C = 1 \quad (11)$$

where  $r_d$  is the radius of the dome. At the entrance to the tubes  $x = L$  (the effective length of the dome)

$$P' = \frac{-i\rho a^2}{g} (e^{iLB_1} + e^{-iLB_2}) \quad (12)$$

$$\frac{\partial P'}{\partial x} = \frac{\rho a^2}{g} B_1 (e^{iLB_1} - e^{-iLB_2}) \quad (13)$$

$$V' = \frac{ia^2}{\omega} B_1 (e^{iLB_1} - e^{-iLB_2}) \quad (14)$$

$$\frac{\partial V'}{\partial x} = -\frac{a^2}{\omega} B_1^2 (e^{iLB_1} - e^{-iLB_2}) \quad (15)$$

## INJECTION ELEMENT MODEL

With the time dependence of pressure, density, and velocity specified at two cells at the inlet to the element (obtained from the solution for the dome above), the time dependence for the next cell can be calculated using the equations below. This process can be repeated to march down the injection element. The dome and element calculations are repeated with a new guess for the amplitude of the oscillation in the dome until the oscillation amplitude at the exit of the element agrees with the chamber oscillation amplitude or cavitation occurs. If a calculated pressure is below the saturated vapor pressure (or below 1 psia for a gas) or exceeds the sonic velocity, the calculation procedure to match exit pressure is terminated. The calculation proceeds to match the lowest pressure to the saturation pressure and a diagnostic message is printed. The response values are adjusted to the required chamber pressure oscillations via

$$RE = RE_c \left( \frac{P_{cch}}{P_{ch}} \right)^{1/2} \quad (16)$$

where  $P_{cch}$  calculated injector face pressure amplitude  
 $P_{ch}$  chamber pressure amplitude  
 $RE_c$  calculated response

a somewhat arbitrary rule.

Continuity

$$\frac{\partial \rho}{\partial t} + \rho \frac{\partial v}{\partial x} + v \frac{\partial \rho}{\partial x} = 0 \quad (17)$$

Momentum

$$\rho \frac{\partial v}{\partial t} + \rho v \frac{\partial v}{\partial x} + g \frac{\partial p}{\partial x} + \frac{\mu v}{R^2} = 0 \quad (18)$$

State

$$\frac{\partial \rho}{\partial p} = \frac{g}{a^2} \quad (19)$$

or

$$\frac{\rho}{p} = \frac{g}{a^2} = \frac{\rho_o}{p_o} Dencon \quad (20)$$

From continuity

$$\frac{v_{n+1} - v_n}{\delta x} = \frac{\partial v}{\partial x} = \frac{-\left(\frac{\rho_o v}{p_o} \frac{\partial p}{\partial x} + \frac{\partial \rho}{\partial t}\right)}{\rho} \quad (21)$$

and from the equation of state

$$\rho = \rho_o + \frac{(p - p_o) \rho_o Dencon}{p_o} \quad (22)$$

Substituting into Eq. 17

$$\frac{p_{n+1} - p_n}{\delta x} = \frac{\partial p}{\partial x} = \frac{\left(\rho \frac{\partial v}{\partial t} - v \frac{\partial \rho}{\partial t} + \frac{\mu v}{R^2}\right)}{\frac{\rho_o a^2 Dencon}{p_o} - g} \quad (23)$$

where  $Dencon$  slope of the density/pressure relationship  
 $p_o$  the reference pressure  
 $R$  the radius of the tube  
 $\rho_o$  the reference density

At the orifice inlet, the flow variables for a fictitious cell are calculated in the orifice using the area relationship between the orifice and the tube. This fictitious cell and the previous cell in the tube are used to obtain the new conditions for calculations in the orifice via the continuity and Bernoulli equations. The same procedure is used at the orifice exit.

At the overlap points

$$W_{n+2} = W_n \quad (24)$$

$$v_{n+2} = \frac{v_n \rho_n A_n}{\rho_{n+2} A_{n+2}} \quad (25)$$

For pressure use Bernoulli's equation

$$\delta p = 1/2 \rho v^2 \quad (26)$$

$$v = C_D \left( \frac{\delta p}{\rho} \right)^{1/2} \quad (27)$$

At the upstream edge of the orifice

$$\delta p = p_n - p_{n+2} = 1/2 \rho (C D_1 v_{orf})^2 \quad (28)$$

At the downstream edge of the orifice

$$\delta p = p_{n+2} - p_n = 1/2 \rho (C D_2 v_{orf})^2 \quad (29)$$

where  $C D_1$  upstream discharge coefficient  
 $C D_2$  downstream discharge coefficient  
 $v_{orf}$  velocity in the orifice

If the velocity exceeds the speed of sound, then the velocity at that location is set equal to the speed of sound with the same direction. Momentum is ignored and continuity is used to obtain pressure via  $\rho$

$$\frac{\partial \rho}{\partial x} = - \frac{\left( \frac{\partial \rho}{\partial t} + \rho \frac{\partial v}{\partial x} \right)}{v} \quad (30)$$



**PART E**

**LUMPED PARAMETERS INJECTION RESPONSE MODEL  
(INJ)**

## Lumped Parameter Injection Response Model (INJ)

Jeffrey Muss

The lumped parameter injection responses model, INJ, is similar to the model described in NASA SP-194 (Ref. 1). The model has been limited to consider only the effects of the injection element and the propellant manifold, thereby ignoring all upstream effects. The major deviation from the SP-194 model is the extension of the model to account for element mixed patterns. This required the timelag and inertance of each element to be accounted for. This was achieved by mass weighting the individual contribution of each element. Slightly different variable normalization were applied, for computational efficiency reasons, and this results in different forms of the characteristic variables, i.e., inertance, capacitance and resistance. Expansion of these parameters will yield the traditional definitions.

The injector's resistance (R) and capacitance (C), and the element's inertance (L) are defined as:

$$R = DP_j / (n * P_c) \quad (1)$$

$$C = \frac{P_c * Vol * G_c}{A^2 * W_{tot}} \quad (2)$$

$$L = \frac{l_{orif} * W_{orif}}{A_{orif} * G_c * P_c} \quad (3)$$

where Vol is the manifold volume, in ft<sup>3</sup>, A<sub>orif</sub>, l<sub>orif</sub>, and W<sub>orif</sub> are the cross-sectional area, in ft<sup>2</sup>, length, in ft, and flowrate, in Lbm/s, of an orifice in an element, respectively, G<sub>c</sub> is the gravitational constant, A is the speed of sound in the propellant, in ft/s, W<sub>tot</sub> is the total propellant flow, in Lbm/s, and DP<sub>j</sub> and P<sub>c</sub> are in psf. L and C are in seconds, while R is nondimensional. The term "n" in the calculation of the resistance is the exponent in the equation.

$$W = k * DP_j^n \quad (4)$$

where W is mass flowrate and k is a proportionality constant. It should be noted that n=0.5 for liquid propellants, while it must be evaluated numerically for gaseous propellants.

The overall mass-weighted injection admittance (Yj) is expressed as:

$$Y_j(f) = \frac{-1}{1+MR} * \sum_{n=1, NFE} \left| \frac{FMF_n * \exp(i * w * \tau_n)}{R_f + L_{f,n} + (i * w * C_f^{-1})} \right| + \frac{-MR}{1+MR} * \sum_{m=1, NXE} \left| \frac{XMF_m * \exp(i * w * \tau_m)}{R_x + L_{x,m} + (i * w * C_x^{-1})} \right| \quad (5)$$

where f is the frequency, in hz,  $\exp(x)=e^x$ , MR is the oxidizer-to-fuel mixture ratio, Tau is the total timelag, in sec., i is the square root of -1, w is the angular frequency  $2 * \pi * f$ , NXE and NFE refer to the total number of oxidizer and fuel element types, respectively, the subscripts "x" and "f" refer to oxidizer and fuel, respectively, and FMF and XMF are the fraction of the total fuel and oxidizer mass contained in that element type, respectively.

#### References

- 1) Liquid Propellant Rocket Combustion Instability; D.T. Harrje, Ed., NASA SP-194, 1972.

**PART F**

**MCA PERFORMANCE/LIFE COMBUSTION  
MODEL DEVELOPMENT FINAL REPORT**



Aerojet  
TechSystems  
Company

## ENGINEERING AND DEVELOPMENT

<b>THERMODYNAMIC ANALYSIS REPORT</b>		NUMBER: 9980:1455
		DATE: 5 March 1986
SUBJECT:  MCA PERFORMANCE/LIFE COMBUSTION MODEL DEVELOPMENT FINAL REPORT	PAGE 1 OF _____	
	NO. OF ENCLOSURES _____	
	NO. OF APPENDICES _____	
ADDITIONAL INFORMATION AND WORK NOTES INCLUDED IN MICROFILM FILE CDN _____		



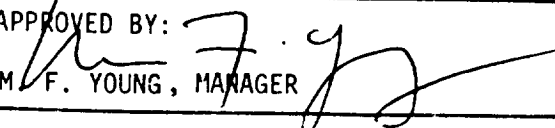
PREPARED FOR: J. A. Van Kleeck

### INTRODUCTION

The Performance/Life Combustion Model (PLC) was developed to evaluate the energy release efficiency (ERE) and the spacial combustion chamber wall mixture ratio distribution within the Space Shuttle Main Engine (SSME) for the Main Chamber Combustion and Cooling Technology Study. The output from the PLC model is to be used in conjunction with thermal, structural and performance models to assess overall thrust chamber performance and combustion chamber life.

While the model is not a rigorous CFD type of model, it does account for many of the influences in the same way a rigorous model would. PLC was developed to mechanistically account for changes in injector and chamber design parameters and variable operating conditions without reliance upon empirical user input scaling factors. The input is concise, requiring a minimum of propellant property information. The input, described in full in Appendix I, consists of propellant injection conditions, chamber wall contour information, injector pattern layout, element flowrates, and basic injector element design configuration. It can be used to evaluate any gaseous fuel-liquid oxidizer propellant combination for several types of "coaxial" injectors.

KEYWORDS: SSME (16), TCA (66), Performance (101), Compatibility (109), LOX/HC (153), Computer Program - New Development (210), 1986 (271), Muss (362)

DISTRIBUTION:  S. Brown, J. Fang, T. V. Nguyen, R. Walker, 9980 File	PREPARED BY: J. MUSS 
	REVIEWED BY: J. ITO 
	APPROVED BY:  M. F. YOUNG, MANAGER

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## MODELLING ASSUMPTIONS

The inherent strength or weakness of any model is the validity of the assumptions used in the model. This section outlines the major assumptions incorporated in PLC.

The chamber flow dynamics are modelled assuming that the flow in the near-face region is in discrete streamtubes, while flow in the remainder of the chamber is modelled as finite elements or "cells". The vapor contained within any given cell is assumed to be well mixed, and the cell possesses the physical properties characterized by the axial pressure and the cell's vaporized mixture ratio. These cells are set up by GRIDGEN so that all cells are of equal area. The number of cells is determined by the number of slices, NSLICE, and the number of radial segments, NSEQ, specified in the input, and is equal to  $NSLICE \times NSEQ$ . The number of cells is constant at any cross-section, but the cell area is a function of axial location. Figure 1 shows the grid for a chamber with 12 slices and 4 segments. Flow calculations are also based on the assumption of choked flow at the throat, but the assumption's validity is never checked by the program. Other key assumptions are that the volume occupied by the droplets is negligible, so that the vapor is assumed to occupy the entire local cross-section, and that the droplets apply no drag forces to the gases that are accelerating them.

The pressure is assumed to be constant and equal to the nominal chamber pressure throughout the cylindrical portion of the chamber. It is also assumed that the gas phase properties can be adequately estimated as a function of the local vapor mixture ratio and pressure with TDK generated one-dimensional shifting equilibrium values.

Droplets are assumed to move as a uniformly distributed ring surrounding the axis of the injection element they emanated from. The distance from the element axis (or the apparent element axis) is measured as a radius,  $R$ , and vapor generated by vaporizing droplets is distributed uniformly around the injector axis at this radius  $R$ . The radius  $R$  is changed as the droplets experience aerodynamic drag from the gases.

## PROGRAM LOGIC

This section describes the flow of information within PLC while the subroutine details are reserved for the following section. PLC models the atomization of the oxidizer, streamtube expansion of the injected gas/liquid mixture in the near-face region, vapor mixing in the cylindrical chamber section, droplet trajectories and secondary droplet shattering due to wall impingement, and 2-D isentropic flow acceleration in the convergent nozzle section.

A cylindrical coordinate system is employed to describe position, with the axial position defining the local wall radius. All internal calculations are conducted in fundamental English units, i.e. LBM-Ft-Sec. Information is transferred between subroutines by means of labelled commons which are sized at runtime based upon the problem inputs. Program output is directed to two files, "PL.HISTORY" which contains echoed input data, ERE predictions and other run information, and "MR.DIST" which contains the vapor mixture ratio as a function of axial, radial and circumferential position.

PLC is structured with a main calling program and a series of subroutines. The main program is used only to sequence the processing of information, check for error conditions, and monitor axial position. A graphical representation of this flow diagram is presented in Figure 2. The first subroutine called is SETUP. SETUP reads and echos the input file, checks it for consistency, and initializes problem specific variables. The next step is to atomize the

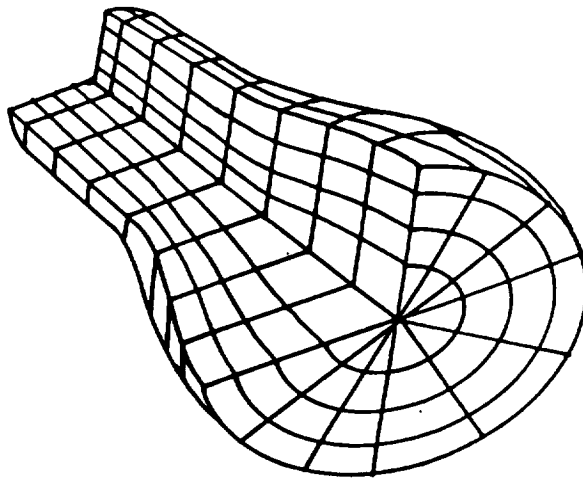


Figure 1  
Graphical Representation of  
the Computational Grid

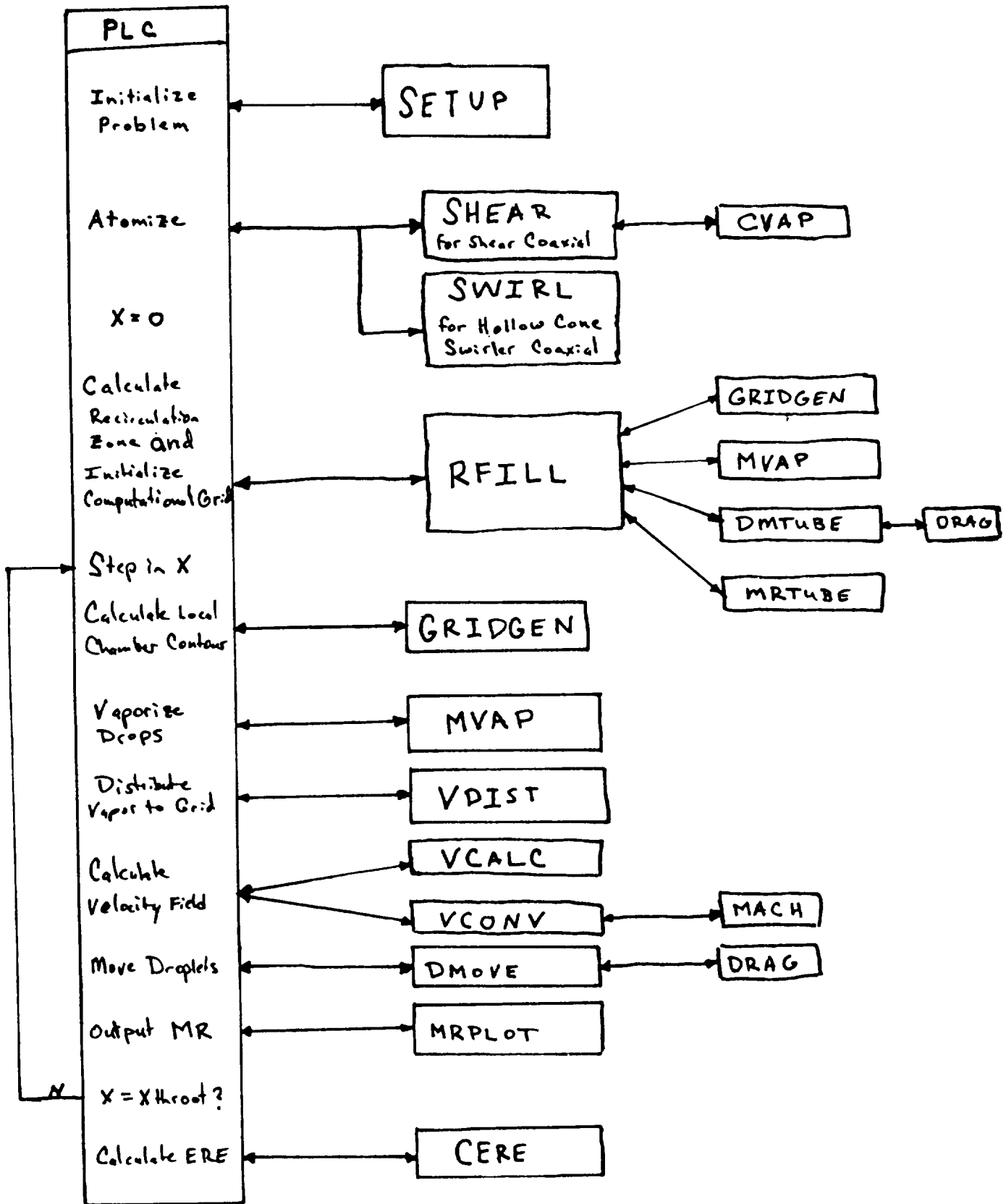


Figure 2  
PLC Program Schematic  
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oxidizer. This is done either by subroutine SHEAR for conventional shear coaxial injectors, or by SWIRL for standard and impinging hydraulically swirled coaxial injectors. These subroutines calculate the mass median dropsize, the dropsize uniformity distribution parameter,  $S_g$ , the number of drops formed per second of each size, the mean atomization length and the mean injection velocities for each element. Each injector's droplet production is characterized by three drop sizes, the mass median dropsize, which accounts for 67% of the droplets formed, and the mass median dropsize multiplied and divided by  $S_g$  which results in maximum and minimum drop sizes, respectively, of plus and minus one standard deviation. Each of these sizes represent 16.5% of the drops produced. When the oxidizer for all elements has been atomized, the droplet vaporization and chamber gas dynamics calculations begin.

The subroutine RFILL is called next to calculate the length of the near-face recirculation zone. In the near-face zone, the droplets are assumed to flow in perfectly mixed streamtubes emanating from the elements. These streamtubes expand at a rate governed primarily by the rate of oxidizer vaporization. As axial steps are taken, RFILL calls GRIDGEN to calculate the local wall radius and MVAP to calculate the oxidizer mass flow added to the gas stream due to droplet vaporization between the current and previous axial position. RFILL then calls DMTUBE to calculate the droplet acceleration resulting from aerodynamic drag on the droplet by the combustion gas and MRTUBE to output the streamtube mixture ratios. Finally RFILL calculates the cross-sectional area of the streamtubes, which is based on the conservation of linear momentum. The area of the streamtubes is summed and compared to the local cross-sectional area of the chamber. If the cross-section is not filled, another axial step is taken and the process is repeated. If the cross-section is filled, RFILL will calculate the streamtubes final location and distribute the streamtube's mass to the computational grid. When complete, the coolant flow is uniformly added to the computational grid.

With the cross-section filled, the main program calculates the chamber gas dynamics as a function of the axial location in the chamber. The main program steps in  $x$  until the throat plane is reached, calculates positional vapor mass addition, gas velocities, radial and circumferential vapor mass flux, droplet movement due to aerodynamic drag and vapor mixture ratio distribution at each axial station.

These calculations are made in the following manner. First, GRIDGEN is called to calculate local wall radius, wall angle, and to calculate the radii of the equal area computational grid cells. Next, MVAP is called to calculate droplet vaporization and dropsize reduction, and VDIST is called to distribute the vaporized mass to the computational grid. With the mass addition complete, the gas velocity components are calculated by either V CALC for the cylindrical section of the chamber, or by V CONV for the converging section. V CALC calculates an ERE based 1-D axial velocity as well as local radial and circumferential velocities based on cell-to-cell mass maldistributions. V CONV calculates a local axial and radial velocity based on isentropic acceleration and wall curvature turning of the gas. With the cross-sectional velocity profile calculated, DMOVE is called to calculate the droplets' movement resulting from aerodynamic drag. It also checks for droplet wall impingement. Finally, MRPLOT is called to report the radial and circumferential vapor mixture ratio distribution. This process is repeated until the throat plane is reached at which time CERE is called to calculate the overall vaporization, mixing and energy release efficiencies.

## SUBROUTINE DESCRIPTION

This section examines the salient features of several key subroutines. A complete listing of all the subroutines is contained in Appendix II.

SHEAR is used to calculate the mass median dropsize, distribution coefficient and total number of drops formed by a shear coaxial liquid-gas

injector. The oxidizer is assumed to be liquid and the fuel or preburner hot exhaust to be gas. The term fuel is used to represent the gas stream.

Calculations begin by estimating the oxidizer post discharge coefficient and velocity profile. These calculations consider the area ratio of the post tip to the metering diameter, the length of the final diffuser section, and the oxidizer Reynolds Number, and are based on Ito's General Hydraulic Flip Model (Ref 1). The fuel stream velocity is considered to have a flat, turbulent profile, and it is based upon a 1-D calculation. With the velocities in both streams set, the dropsize formed at any axial position is determined by equating the surface tension cohesive force to the interfacial shear force and solving for the dropsize,

$$R_m = 8 \cdot ST / (T \cdot g)$$

where  $R_m$  is the droplet radius in ft,  $ST$  is the oxidizer surface tension in lbf/ft,  $T$  is the interfacial shear stress in lbf/ft-sec<sup>2</sup>, and  $g$  is the gravitational constant. The interfacial shear stress is estimated as

$$T = m \cdot (V_f - V_x) / (T_{POST} + DR_{JET})$$

where  $m$  is the mean dynamic viscosity for the fuel and the oxidizer in lbf/ft-sec,  $V_f$  is the fuel velocity,  $V_x$  is the oxidizer free surface velocity,  $T_{POST}$  is the oxidizer post wall thickness, and  $DR_{JET}$  is the reduction in oxidizer jet radius due to mass stripping. Once the dropsize for an axial position is calculated, the oxidizer jet's radius is reduced to account for the mass of the droplets formed. First the number of drops formed is estimated by calculating how many drops will fit around the outside of the jet. The volume of these drops is removed from the jet and the resulting truncated conic section is calculated. The integrated flow in and out of the cone is then used to calculate the number of drops formed based on satisfaction of continuity.

If the injector contains a recessed post, the drops formed in the cup are vaporized between their generation location and the injector face. The percent of the droplet vaporized is based on the Priem's Generalized Length Correlation (Ref 2) which is used throughout PLC to model droplet vaporization. The oxidizer vapor is added to the fuel stream. The added mass is first used to fill the fuel injection flow area, i.e. increase the fuel  $C_d$  to 1.0, and then to accelerate the fuel velocity.

RFILL calculates the rate at which the injector element's streamtubes expand in the near-face zone. It vaporizes and accelerates the droplets at each axial step until the total cross-sectional area is filled. When the cross-section is filled, the mass in the streamtubes is distributed to the computational grid.

The area of the streamtubes is based on the conservation of linear momentum. The flow within the streamtubes is modelled as a set of rings. The inner ring is a circle of radius  $R_{CORE}$  and estimates the liquid oxidizer core radius as a function of axial location.  $R_{CORE}$  is estimated as

$$R_{CORE} = \text{MAX}(0.0, (R_{XP} - (R_{XP} / (2.0 \cdot AL)) \cdot (X + RECESS)))$$

where  $R_{XP}$  is the oxidizer post radius,  $AL$  is the element's mean atomization length,  $X$  is the current axial position, and  $RECESS$  is the oxidizer post tip recess.  $R_{CORE}$  is surrounded by an annulus of stoichiometric combustion gases. The area of this annulus,  $ACZ$ , is estimated as

$$ACZ = 1.125 \cdot W_o / RHO / VCZ \cdot ERE \quad J-192$$

where  $W_o$  is the vaporized oxidizer mass flow in the streamtube,  $RHO$  is the density of the stoichiometric zone,  $VCZ$  is the estimated combustion gas velocity, and  $ERE$  is the streamtubes characteristic velocity energy release efficiency.  $VCZ$  is estimated as

$$VCZ = (B * V_o + V_f) / 9.0$$

where  $V_o$  and  $V_f$  are the injected oxidizer and fuel velocities, respectively. The outer annulus, of outer radius equal to the streamtube's radius, is composed of the unreacted fuel.

When the cross-section is filled by the streamtubes, streamtube flow ceases and the vaporized mass is distributed to the computational flow grid. Distribution is based on final streamtube position. This is done by assuming that the streamtubes from the outer row of elements will be forced into an annular region against the wall of cross-sectional area equal to the sum of the cross-sectional area of the outer row's streamtubes. Then starting with the element of that row with a circumferential location closest to zero degrees, an angular slice of the annulus is calculated with a cross-sectional area equal to that of the element's streamtube. This process is repeated for all elements of the outer row. When complete, the vapor mass within the annular slices are transferred to the computational grid cells corresponding to each annular slices' locations. The elements' location are replaced with the coordinates of the centroid of the annular slices' area. This process is repeated for the next outermost row (the second row from the wall) assuming its annular area abutts the outer annulus' inner edge. The process is repeated for all rows. When complete, the transpiration coolant flow is uniformly added to the computational grid.

MVAP is used to vaporize the droplets in all axial flow regimes. It is based on the Priem-Heidmann Droplet Vaporization Model (Ref. 2).

VCALC calculates the gas 1-D velocity and properties at axial locations between the near-face zone and the entrance to the convergent section. It calculates radial and circumferential mass velocities due to cell-to-cell mass maldistributions resulting from differential local vaporization rates and element-to-element flow variations. These velocities are based on the tendency of mass to try to uniformly distribute itself across the cross-section rather than to remain stratified in zones of high and low mass concentration. Future development should focus on a more rigorous treatment of these phenomenae.

VCONV calculates the gas acceleration in the convergent section of the thrust chamber nozzle. It outputs axial and radial gas velocities as a function of the local wall curvature and gas properties as well as the isentropic acceleration of the gas. It is based on the Droplet Trajectory Model of Nyugen (Ref 3).

DMOVE is used to accelerate and move the droplets due to aerodynamic drag. It also checks for wall impingement of the droplets. If impingement occurs, the droplets are assumed to be vaporized at that point due to secondary droplet shattering. Both Stoke's flow and Newtonian drag expressions are used depending on the applicable Reynolds Number. The Newtonian drag equation takes the form used in SDER (Ref 4) which eliminates a stepsize constraint in order to achieve numerical stability. Rabin's correlations for drag of an accelerating liquid droplet are used to calculate drag coefficients (Ref 5). The drag is calculated for droplet assuming supercritical expansion of the droplet.

DBLINT is a general linear double interpolation routine used to calculate the various gas properties as a function of chamber pressure and local mixture

ratio. The interpolation uses UDE chamber equilibrium values at chamber pressures of 300 and 3000 psia and mixture ratios ranging from 0 to 20. The gas properties calculated are specific heat ratio, gas temperature, molecular weight, density, sonic velocity, characteristic velocity and dynamic viscosity.

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2. Priem, R. J. and Heidmann, M. F., Propellant Vaporization as a Design Criterion for Rocket-Engine Combustion Chambers, NASA Technical Report R-67, 1960
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5. Rabin, E., Schallenmueller, A. R., and Lawhead, R. B., Displacement and Shattering of Propellant Droplets, AFOSR-TR-60-75-1960

APPENDIX I  
PLC INPUT INFORMATION

CONTENTS

Input Format Information

Sample Input One: Seven Element, Single Row Shear Coaxial Injector

Sample Input Two: Thirteen Element, Single Row Hollow Cone Swirler Injector

## Input Format Information

There are a few simple steps required to run the PLC model. First an input file has to be written. When the file is complete, the user runs the CPL file PLSET to size the commons, create any necessary insert files and create an executable program. This section contains the format of PLC input files as well as two sample input files.

The input format for PLC is as follows. The file MUST be named PL.INPUT.

```

TITLE 1                      (A80)
TITLE 2                      (A80)
(BLANK LINE)
$OUTPUT
$END
(BLANK LINE)
PC HGMR                      (F7.2,F7.4)
TINJ TC RHO MU MW HOV (6E12.4) (FUEL/HOT GAS PROPERTIES)
TINJ TC RHO MU MW HOV (6E12.4) (OX PROPERTIES)
ST RHOC                      (2E12.4) (OX PROPERTIES)
(BLANK LINE)
NXP                          (I3)
XW(1) RW(1)                  (2F7.4)
.
XW(NXP) RW(NXP)              (2F7.4)
(BLANK LINE)
INPUT-TYPE NEL NROWS         (3I4)
INPUT-TYPE INPUT FORM
CFLOW CFMR                   (2F8.4)
(BLANK LINE)
INJ-TYPE                      (I2)
INJ-TYPE INPUT FORM
(BLANK LINE)
$CONT
$END

```

The inputs are further defined as follows:

NAMELIST \$OUTPUT CONTROLS PROGRAM OUTPUT  
 DEFAULT IS 0 FOR ALL EXCEPT IECHO WHERE 0-OFF

PARAMETER	VALUES	FUNCTION
IECHO	0/1	ECHO INPUT
ITRACE	0/1/2	TRACES PROGRAM PROGRESS: -1 AXIAL LOCATION -2 AND SUBROUTINE ENTRY
IAFLG	0/1/2/3	OUTPUTS DROPLET FORMATION INFORMATION: -1 MEDIAN DROPSIZE, DISTRIBUTION PARAMETER, NUMBER OF DROPS, INJECTION VELOCITY, AND ATOMIZATION LENGTH -2 AND DROPSIZES FORMED BY SHEAR COAX INJECTOR -3 AND DROPSIZES AS FORMED AND VAPORIZED BY SHEAR COAX
IRFFLG	0/1/2/3	OUTPUT RECIRCULATION ZONE CALCULATION INFORMATION: -1 PERCENT OF CROSS-SECTION FILLED AS F(X) -2 AND FINAL INJECTOR APPARENT LOCATIONS -3 AND DISTRIBUTION FOR STREAMTUBES TO GRID
IMVFLG	0/1	OUTPUTS DROPLET VAPORIZATION INFORMATION
IVDFLG	0/1	OUTPUTS DISTRIBUTION OF VAPOR
IPDFLG	0/1	OUTPUTS AXIAL AND RADIAL VELOCITY COMPONENTS
IDMFLG	0/1	OUTPUTS DROPLET MOVEMENT
IDFLG	0/1	OUTPUTS DRAG INTERACTION PARAMETERS
INRFLG	0/1/2/3	OUTPUTS MASS DISTRIBUTION INFORMATION: -0 WALL MR AS A FUNCTION OF THETA -1 AND CROSS-SECTION MR AS A F(R,THETA) -2 AND MASS FLOW AS A F(R, THETA) -3 AND FUEL AND OXIDIZER MASS FLOW AS A F(R,THETA)

## OPERATING CONDITIONS AND INJECTION PROPERTIES

PC=CHAMBER PRESSURE (PSIA)  
HGMR=FUEL/HOT GAS MIXTURE RATIO  
TINJ=INJECTION TEMPERATURE (DEG F)  
TC=CRITICAL TEMPERATURE (DEG F)  
RHO=DENSITY (LB<sub>m</sub>/CU. FT)  
MU=DYNAMIC VISCOSITY (LB<sub>m</sub>/FT-SEC)  
MW=MOLECULAR WEIGHT (LB<sub>m</sub>/LBMOLE)  
HOV=HEAT OF VAPORIZATION AT NBP (BTU/LB<sub>m</sub>)  
ST=SURFACE TENSION (LB<sub>f</sub>/FT)  
RHOC=CRITICAL DENSITY (LB<sub>m</sub>/CU. FT)

Note: the fuel/hot gas density and/or viscosity may be set to zero and the

## CHAMBER CONTOUR

NXP=NUMBER OF CHAMBER CONTOUR DESCRIPTION POINTS (MINIMUM=2)  
XW=AXIAL LOCATION OF RW (INCHES)  
RW=CHAMBER WALL RADIUS AT XW (INCHES)

## INJECTOR ELEMENT LAYOUT

INPUT-TYPE=POSITION AND FLOW INPUT TYPE:  
=1 FOR INPUT EACH ELEMENT LOCATION AND FLOWRATE  
=2 FOR INPUT # ELEMENTS/ROW, ROW POSITION AND FLOWRATE EQUATION  
NEL=TOTAL NUMBER OF ELEMENTS  
NROWS=NUMBER OF CONCENTRIC ROWS

IF INPUT-TYPE=1 INPUT-TYPE INPUT FORM IS OF THE FORMAT:

ERPOS ETPOS FMF XMF (4F8.4)

ERPOS=ELEMENT RADIAL LOCATION FROM CENTERLINE (IN)  
ETPOS=ELEMENT CIRCUMFERENTIAL POSITION FROM 0 DEGREE REF (DEG)  
FMF=HOT GAS MASS FLOWRATE (LB<sub>m</sub>/SEC)  
XMF=OX MASS FLOWRATE (LB<sub>m</sub>/SEC)

One line for each element listed by ascending radius and then by ascending circumferential location within a row

IF INPUT-TYPE=2 INPUT-TYPE INPUT FORM IS OF THE FORMAT:

NELR RROW (I4,F8.4)  
FMF\_EQ (A65)  
XMF\_EQ (A65)

NELR=NUMBER OF ELEMENTS IN THAT ROW  
RROW=RADIUS OF ELEMENTS IN ROW (IN)  
FMF\_EQ=FUEL/HOT GAS MASS FLOWRATE EQUATION (LB<sub>m</sub>/SEC)  
XMF\_EQ=OX MASS FLOWRATE EQUATION (LB<sub>m</sub>/SEC)

a function of the variable theta, the circumferential position.  
elements in a row are assumed equispaced and theta is in rads  
(0-2pi). rows should be listed in ascending radial position.  
Note: FORTRAN trig. functions use radians

CFLOW=TOTAL COOLANT FLOWRATE INTO THE THRUST CHAMBER (LB<sub>m</sub>/SEC)  
CFMR=COOLANT FLOW MIXTURE RATIO

# INJECTOR TYPE AND DIMENSIONS

ITYPE=1 FOR SHEAR COAX INJECTOR AND ITYPE INPUT IS:

RXP RFP TPOST RECESS RMS XDL FCD (7F7.4)

RXP=OX POST INNER RADIUS (IN)  
RFP=FUEL ANNULUS OUTER RADIUS (IN)  
TPOST=OX POST THICKNESS (IN)  
RECESS=OX POST RECESS (IN)  
RMS=RADIUS OF OX METERING SECTION (IN)  
XDL=OX DIFUSER SECTION LENGTH (IN)  
FCD=FUEL ANNULUS CD

ITYPE=2 FOR HOLLOW CONE SWIRL COAX AND ITYPE INPUT IS:

RXP RFP TPOST CSA FCD XCD (6F7.4)

RXP=OX POST INNER RADIUS (IN)  
RFP=FUEL ANNULUS OUTER RADIUS (IN)  
TPOST=OX POST THICKNESS (IN)  
CSA=CONE SPRAY ANGLE (DEG)  
FCD=FUEL ANNULUS CD  
XCD=OX ANNULUS CD

ITYPE=3 FOR IMPINGING HOLLOW CONE SWIRL TRIAX AND ITYPE INPUT IS:

RXP RFP TPOST PFI CSA FCD XCD (7F7.4)

RXP=OX POST INNER RADIUS (IN)  
RFP=FUEL ANNULUS OUTER RADIUS (IN)  
TPOST=OX POST THICKNESS (IN)  
CSA=CONE SPRAY ANGLE (DEG)  
PFI=PERCENT FUEL IMPINGED (XXX.XXX)  
FCD=FUEL ANNULUS CD  
XCD=OX ANNULUS CD

## NAMelist \$CONT SETS MODEL CONTROL PARAMETERS

PARAMETER	DEFAULT	FUNCTION
NSEG	NROWS+1	SETS THE NUMBER OF RADIAL GRID SEGMENTS
NSLICE	MIN(12,NEL-1)	SETS THE NUMBER OF CIRCUMFERENTIAL GRID SLICES
ASTEP	2*(-10)	ATOMIZATION STEPSIZE (FT)
XSTEP	2*(-8)	AXIAL CALCULATION STEPSIZE (FT)

The following is a sample input file for a single row, seven element shear coaxial injector. Element flows are input as formulae rather than for each element.

SAMPLE INPUT FILE ONE  
ROW TYPE INPUT FOR SINGLE ROW SEVEN ELEMENT SHEAR COAX

\$OUTPUT  
ITRACE=1, IDHFLG=0, IPDFLG=1, IVDFLG=0, IMVFLG=0, IDFLG=0,  
IMRFLG=1, IAFLG=1, IRFFLG=1,  
\$END

3006.00	0.8249				
1.0940E3	-3.999E2	0.000E0	0.000E00	2.016E0	1.953E2
-2.700E2	-1.820E2	6.535E1	6.500E-5	3.200E1	9.162E1
9.000E-6	2.7217E1				

24	
0.00	0.9604
3.00	0.9604
3.50	0.9551
4.00	0.9498
4.50	0.9445
5.00	0.9340
5.50	0.9207
6.00	0.9022
6.50	0.8810
7.00	0.8546
7.50	0.8308
8.00	0.8175
8.50	0.7779
9.00	0.7540
9.50	0.7302
10.00	0.7011
10.50	0.6773
11.00	0.6535
11.50	0.6270
12.00	0.6006
12.50	0.5794
13.00	0.5662
13.50	0.5583
14.00	0.5556

2	7	1
7	0.6791	
0.43530+0.043530*SIN(THETA)		
1	3145	
0.0934	0.00	

1								
0.0948	0.2030	0.0210	0.2550	0.0790	1.5000	0.8000		

\$CONT  
\$END

The following is an input for a single row, thirteen element hollow-cone swirl coaxial injector. In this case, each elements location and flowrate is input.

SAMPLE INPUT FILE TWO  
SEVEN ELEMENT HOLLOW CONE SWIRLER, INDIVIDUAL INPUTS

\$OUTPUT

ITRACE=1, IDMFLG=0, IPDFLG=1, IVDFLG=0, IMVFLG=0, IDFLG=0,  
IMRFLG=1, IAFLG=1, IRFFLG=1,

\$END

```
3268 84 0.8931
1.2000E3 -3.999E2 0.000E0 0.000E00 2.016E0 1.953E2
-2.650E2 -1.820E2 6.535E1 6.500E-5 3.200E1 9.162E1
9.000E-6 2.7217E1
```

```
24
0.00 1.309
3.00 1.309
3.50 1.303
4.00 1.295
4.50 1.288
5.00 1.273
5.50 1.256
6.00 1.229
6.50 1.200
7.00 1.166
7.50 1.133
8.00 1.116
8.50 1.061
9.00 1.026
9.50 0.994
10.00 0.957
10.50 0.923
11.00 0.889
11.50 0.854
12.00 0.818
12.50 0.790
13.00 0.773
13.50 0.761
14.00 0.758
```

```
1 13 1
0.926 00.000 0.48749 1.4148
0.926 27.692 0.48749 1.4148
0.926 55.385 0.48749 1.4148
0.926 83.077 0.48749 1.4148
0.926 110.769 0.48749 1.4148
0.926 138.462 0.48749 1.4148
0.926 166.154 0.48749 1.4148
0.926 193.846 0.48749 1.4148
0.926 221.538 0.48749 1.4148
0.926 249.231 0.48749 1.4148
0.926 276.923 0.48749 1.4148
0.926 304.615 0.48749 1.4148
0.926 332.308 0.48749 1.4148
0.2161 0.000
```

```
2
0.0948 0.2030 0.0210 30.00 0.800 0.5000
```

\$CONT

NSEC=3, NSLICE=5,

\$END

APPENDIX II  
PLC PROGRAMS

CONTENTS

PLC - main calling program  
SETUP - subroutine to initialize problem  
SHEAR - subroutine to atomize oxidizer for shear coaxial injectors  
CVAP - subroutine to vaporize droplets generated in the injector cup  
RFILL - subroutine to calculate the length of the recirculation zone  
DMTUBE - subroutine to accelerate droplets in streamtubes  
MRTUBE - subroutine to output streamtube MR's as a function of axial location  
GRIDGEN - subroutine to generate computational grid and calculate local contour  
MVAP - subroutine to vaporize droplets  
VDIST - subroutine to distribute vaporized oxidizer to computational grid  
VCALC - subroutine to calculate velocity field in cylindrical section of chamber  
VCONV - subroutine to calculate velocity field in convergent section of chamber  
MACH - subroutine to calculate mach number as a function of area ratio  
DMOVE - subroutine to move droplets due to aerodynamic drag  
DRAQ - subroutine to calculate aerodynamic acceleration of droplets  
MRPLOT - subroutine to output mixture ratio profile across cross-section  
CERE - subroutine to calculate TCA efficiencies  
DBLINT - subroutine to interpolate gas properties  
COMMON - common to be sized and inserted at runtime  
PLSET - CPL file to size and insert common at runtime

```

PROGRAM PLC
C
C   MAIN CALLING PROGRAM FOR COMBUSTION SECTION OF PERFORMANCE/
C   LIFE MODEL
C
C   DECLARATIONS AND OPENS
C
C *INSERT COMMON
C
C   OPEN (UNIT=5, FILE='PL. INPUT', STATUS='OLD', ERR=8000)
C   OPEN (UNIT=6, FILE='PL. HISTORY', STATUS='NEW', ERR=8100)
C   OPEN (UNIT=7, FILE='MR. DIST', STATUS='NEW', ERR=8200)
C
C   INITIALIZE PROBLEM
C
C   CALL SETUP
C   IEFLG=0
C
C   ATOMIZE LOX AND CALCULATE MASS MEDIAN DROPSIZE, DISTRIBUTION
C   COEFFICIENT, MEAN ATOMIZATION LENGTH, AND INJECTION VELOCITY
C
C   X=0.0
C   IF (ITYPE .EQ. 1) CALL SHEAR
C   IF (ITYPE .EQ. 2 .OR. ITYPE .EQ. 3) CALL SWIRL
C   IF (X .LT. 0.0) GOTO 9000
C
C   CALCULATE RECIRCULATION ZONE AND DISTRIBUTE MASS ACROSS CHAMBER
C
C   CALL RFill
C   IF (X .LE. 0.0) GOTO 9100
C   CALL MRPlot
C
C   STEP IN X
C
C 100 X=X+XSTEP
C   IF (ITRACE .GT. 0) WRITE(6,110)X
C 110 FORMAT(/1X, '***** BEGIN LOOP FOR X=', F7.4, ' FT *****', /1X)
C
C   GENERATE LOCAL GRID
C
C   CALL GRIDGEN
C   IF (RWALLX .LE. 0.0) GOTO 9200
C
C   VAPORIZE AND DISTRIBUTE MASS BETWEEN X AND X+XSTEP
C
C   CALL MVAP
C   CALL VDIST
C
C   CALCULATE VELOCITIES
C
C   IF (X .LE. X1) CALL VCalc
C   IF (X .GT. X1) CALL VCONV
C
C   MOVE DROPS AND CHECK FOR IMPINGMENT
C
C   CALL DMOVE
C   IF (X .LT. 0.0) GOTO 9300
C
C   CALCULATE VMR
C

```

```

      CALL MRPLOT
      IF (X .LT. XW(NXP)) GOTO 100
C
C      CALCULATE OVERALL ERE
C
      CALL CERE
1500  CLOSE (UNIT=5)
      CLOSE (UNIT=6)
      CLOSE (UNIT=7)
      STOP
C
C      ERROR CONDITIONS
C
8000  WRITE (1,*) 'PL INPUT DOES NOT EXIST, RUN ABORTED'
      STOP
8100  WRITE (1,*) 'PL HISTORY ALREADY EXISTS, RENAME AND RERUN'
      GOTO 1500
8200  WRITE (1,*) 'MR DIST ALREADY EXISTS, RENAME AND RERUN'
      GOTO 1500
9000  WRITE (6,*) 'ERROR IN ATOMIZATION SUBROUTINE; RUN STOPPED'
      GOTO 1500
9100  WRITE (6,*) 'ERROR IN SUBROUTINE RFILL; RUN STOPPED'
      GOTO 1500
9200  WRITE (6,*) 'ERROR IN SUBROUTINE GRIDGEN; RUN STOPPED'
      GOTO 1500
9300  WRITE (6,*) 'ERROR IN SUBROUTINE DMOVE; RUN STOPPED'
      GOTO 1500
      END

```

```

SUBROUTINE SETUP
C
C SUBROUTINE TO READ INITIAL PROBLEM VARIABLES FOR MCA PERFORMANCE/
C LIFE COMBUSTION MODEL
C CHECK RW FOR CONSISTENCY, CALCULATES WANGLE
C CONVERTS INPUT UNITS TO LBm-FT-SEC-PSI UNITS
C
$INSERT COMMON
C
C CHARACTER TITLE*80, DUMMY*80, FEQ*80, XEQ*80
C NAMEDLIST /OUTPUT/ IAFLG, IVDFLG, IMVFLG, IPDFLG, IDMFLG, IECHO,
C * ITRACE, IMRFLG, IDFLG, IRFFLG
C NAMEDLIST /CONT/ NSEG, NSLICE, XSTEP, ASTEP
C DATA IECHO, IAFLG, IVDFLG, IMVFLG, IPDFLG, IDMFLG, IMRFLG, IDFLG,
C * IRFFLG, ITRACE /1,9*0/
C
C FORMAT STATEMENTS
C
10 FORMAT(A80)
11 FORMAT(/25X, 'OUTPUT CONTROL PARAMETERS (1=ON)', /5X, 'ATOM=', I2,
C * 5X, 'MVAP=', I2, 5X, 'VDIST=', I2, 5X, 'VCALC=', I2, 5X, 'DMOVE=',
C * I2, 5X, 'DRAG=', I2, 5X, 'MRPLOT=', I2, 5X, 'RFILL=', I2, 5X,
C * 'ITRACE=', I2)
20 FORMAT(F7.2, F7.4)
21 FORMAT(6E12.4)
22 FORMAT(2E12.4)
23 FORMAT(/25X, 'FUEL AND OX PROPERTIES, PC=', F7.2, ' HOT GAS MR=',
C * F7.4, /1X, 'FUEL:', 4X,
C * 'TINJ=', F7.2, 3X, 'TCRIT=', F7.2, 3X, 'DENSITY=', F7.3, 3X,
C * 'VISCOSITY=', E12.4, 3X, 'M.W. =', F7.3, 3X, 'HEAT OF VAP=', F7.2,
C * /3X, 'OX:', 4X, 'TINJ=', F7.2, 3X, 'TCRIT=', F7.2, 3X, 'DENSITY=',
C * F7.3, 3X, 'VISCOSITY=', E12.4, 3X, 'M.W. =', F7.3, 3X,
C * 'HEAT OF VAP=', F7.2, /10X, 'SURFACE TENSION=', E12.6, 3X,
C * 'CRITICAL DENSITY=', F7.3)
30 FORMAT(I3)
31 FORMAT(/25X, 'CHAMBER CONTOUR', /)
32 FORMAT(2F7.4)
33 FORMAT(5X, 'POINT=', I2, 5X, 'X=', F7.4, 5X, 'RWALL=', F7.4)
34 FORMAT(/5X, 'X1(FT)=' , F7.4, 5X, 'X2(FT)=' , F7.4, 5X, 'CR=' , F6.3)
40 FORMAT(3I4)
41 FORMAT(/25X, 'INJECTOR ELEMENT POSITIONS', /)
42 FORMAT(2F8.4, 2F8.5)
43 FORMAT(1X, 'ROW ', I2, ' ELEMENTS')
44 FORMAT(5X, 'INJECTOR=', I3, 5X, 'RADIUS=', F8.4, 5X, 'THETA=', F8.4, 5X,
C * 'FUEL MASS FLOW=', F8.5, 5X, 'OX MASS FLOW=', F8.5)
45 FORMAT(I4, F8.4)
46 FORMAT(1X, 'ROW ', I3, ' IS CENTERED AT R=', F8.4, ' in. AND CONTAINS '
C * , I3, ' ELEMENTS')
47 FORMAT(A80)
48 FORMAT(5X, 'FUEL FLOW EQ: ', A80, /7X, 'OX FLOW EQ: ', A80, /5X,
C * 'ROW FUEL FLOW=', F8.3, 5X, 'ROW OX FLOW=', F8.3)
49 FORMAT(2F8.4)
50 FORMAT(/5X, 'COOLANT MASS FLOW=', F8.4, 5X, 'COOLANT MR=', F8.4)
51 FORMAT(/5X, 'TOTAL FUEL MASS FLOW=', F8.3, 5X, 'TOTAL FUEL VAPOR MASS'
C * , ' FLOW=', F8.3, /5X, 'TOTAL OX MASS FLOW =', F8.3, 5X, 'TOTAL'
C * ' OX VAPOR MASS FLOW =', F8.3)
60 FORMAT(I2)
61 FORMAT(7F7.4)
62 FORMAT(/20X, 'SHEAR COAXIAL INJECTOR CONFIGURATION', /5X,
C * 'OX POST RADIUS=', F7.4, 8X, 'OX POST THICKNESS=', F7.4, 4X,

```

```

*      'OX POST RECESS=',F7.4,/5X,'OX METERING RADIUS=',F7.4,4X,
*      'OX POST DIFFUSER LENGTH=',F7.4,/5X,'FUEL POST RADIUS=',
*      F7.4,6X,'FUEL CD=',F7.4)
63 FORMAT(6F7.4)
64 FORMAT(/20X,'SWIRL COAXIAL INJECTOR CONFIGURATION',//5X,
*      'OX POST RADIUS=',F7.4,8X,'OX POST THICKNESS=',F7.4,/5X,
*      'SPRAY FAN ANGLE=',F7.4,7X,'OX CD=',F7.4,/5X,
*      'FUEL POST RADIUS=',F7.4,6X,'FUEL CD=',F7.4)
65 FORMAT(/20X,'IMPINGING TRIAXIAL INJECTOR CONFIGURATION',//5X,
*      'OX POST RADIUS=',F7.4,8X,'OX POST THICKNESS=',F7.4,/5X,
*      'SPRAY FAN ANGLE=',F7.4,7X,'OX CD=',F7.4,/5X,
*      'FUEL POST RADIUS=',F7.4,5X,'FUEL CD=',F7.4,/5X,
*      'PERCENT FUEL INPINGING=',F7.4)
70 FORMAT(/25X,'CALCULATION CONTROL PARAMETERS',//5X,'# SEGMENTS=',
*      I2,5X,'# SLICES=',I3,/5X,'ATOM STEPSIZE=',E12.4,5X,
*      'VAPOR STEPSIZE=',E12.4,/)

C      IF (ITRACE .EQ. 2) WRITE(6,1)
1  FORMAT(/1X,'ENTER SETUP')
   DUMMY=''

C
C      READ AND WRITE TITLE
C
   READ(5,10)TITLE
   WRITE(6,10)TITLE
   WRITE(7,10)TITLE
   READ(5,10)TITLE
   WRITE(6,10)TITLE
   WRITE(7,10)TITLE

C
C      READ OUTPUT FLAGS
C
   READ(5,10)TITLE
   IF(TITLE .NE. DUMMY) GOTO 9000
   READ(5,OUTPUT)
   IF (IECHO .EQ. 1) WRITE(6,11) IAFLG, IMVFLG, IVDFLG, IPDFLG,
*      IDMFLG, IDFLG, IMRFLG, IRFFLG, ITRACE

C
C      READ OPERATING CONDITION AND PROPERTIES
C
   READ(5,10)TITLE
   IF(TITLE .NE. DUMMY) GOTO 9000
   READ(5,20)PC, HGMRO
   READ(5,21)FTJ, FTC, FRHO, FMU, FMW, FHV
   READ(5,21)XTJ, XTC, XRHO, XMU, XMW, XHV
   READ(5,22)XST, XRHOC
   IC=2
   IF (FMW .LE. 0.0) CALL DBLINT (PC, HGMRO, IC, FMW)
   IC=5
   IF (FRHO .LE. 0.0) CALL DBLINT (PC, HGMRO, IC, FRHO)
   IC=7
   IF (GMU .LE. 0.0) CALL DBLINT (PC, HGMRO, IC, FMU)
   IF (IECHO .EQ. 1) WRITE(6,23)PC, HGMRO, FTJ, FTC, FRHO, FMU, FMW,
*      FHV, XTJ, XTC, XRHO, XMU, XMW, XHV, XST, XRHOC

C
C      READ CHAMBER DESCRIPTION
C      CHECK FOR CONSISTENCY, CALCULATE X1, X2, CR, WANGLE
C
100 READ(5,10)TITLE
   IF(TITLE .NE. DUMMY) GOTO 9000

```

```

      READ(5,30)NXP
      IF (IECHO .EQ. 1) WRITE(6,31)
      DO 200 I=1,NXP
        READ(5,32,ERR=9110)XW(I),RW(I)
        IF (IECHO .EQ. 1) WRITE(6,33)I,XW(I),RW(I)
        XW(I)=XW(I)/12.0
        RW(I)=RW(I)/12.0
200  CONTINUE
      IF (XW(1) .GT. 0.0) GOTO 9120
      DO 250 I=1,NXP-1
        IF (XW(I) .GE. XW(I+1)) GOTO 9130
        IF (RW(I+1) .GT. RW(I)) GOTO 9140
        WANGLE(I)=ATAN((RW(I+1)-RW(I))/(XW(I+1)-XW(I)))
250  CONTINUE
      CR=(RW(1)/RW(NXP))**2
      X1=0.0
      X2=XW(NXP)
      DO 275 I=1,NXP-1
        IF (WANGLE(I) .LT. 0.0) GOTO 280
275  CONTINUE
      GOTO 9150
280  X1=XW(I)
      X2=XW(NXP)-X1
      IF (IECHO .EQ. 1) WRITE(6,34) X1, X2, CR
C
C      READ INJECTOR FACE LAYOUT, CHECK FOR CONSISTENCY, CALCULATE RROW,
C      NELR, AND TOTAL MASS FLOWS.  DECOMPOSE HOT GAS AND LOAD TOTAL VAPOR
C
300  READ(5,10)TITLE
      IF(TITLE .NE. DUMMY) GOTO 9000
      READ(5,40) INTYPE, NEL, NROWS
      IF (IECHO .EQ. 1) WRITE(6,41)
      TFF=0.0
      TXF=0.0
      THVM=0.0
      TXVM=0.0
      IF (INTYPE .EQ. 2) GOTO 400
C
C      TYPE 1 INPUT; R, THETA, FLOW FOR ALL ELEMENTS
C
      IROW=0
      R1=-1.0
      DO 310 I=1,NROWS
        NELR(I)=0
310  CONTINUE
      DO 350 I=1,NEL
        READ(5,42,ERR=9210) ERPOS(I), ETPOS(I), FMF(I), XMF(I)
        IF (ERPOS(I) .LT. R1) GOTO 9220
        IF (ERPOS(I) .GT. R1) THEN
          IROW=IROW+1
          R1=ERPOS(I)
          TC=-1.0
          RROW(IROW)=R1/12.0
          IF (IECHO .EQ. 1) WRITE(6,43) IROW
        END IF
        IF (IECHO .EQ. 1) WRITE(6,44) I, ERPOS(I), ETPOS(I), FMF(I),
*          XMF(I)
        ERPOS(I)=ERPOS(I)/12.0
        IF (ERPOS(I) .GE. RW(1) .OR. ETPOS(I) .GT. 360.0) GOTO 9230

```

```

        IF (ETPOS(I) .LE. TC) GOTO 9240
        TC=ERPOS(I)
        NELR(IROW)=NELR(IROW)+1
        FMF(I)=FMF(I)/(1.0+HGMRO)
        XMF(I)=XMF(I)+FMF(I)*HGMRO
        TFF=TFF+FMF(I)
        TXF=TXF+XMF(I)
        THVM=THVM+FMF(I)
        TXVM=TXVM+FMF(I)*HGMRO
350 CONTINUE
        IF (IROW .NE. NROWS) GOTO 9250
        GOTO 499
C
C      TYPE 2 INPUTS; R, NELR, FLOWRATES AS F(R,THETA)
C
400 ICNT=0
      DO 460 IROW=1,NROWS
        READ(5,45,ERR=9280) NELR(IROW), RROW(IROW)
        IF (IECHO .EQ. 1) WRITE(6,46) IROW, RROW(IROW), NELR(IROW)
        READ(5,47,ERR=9290) FEQ
        READ(5,47,ERR=9295) XEQ
        IEL1=ICNT+1
        IELL=IEL1+NELR(IROW)-1
        TI=6.28319/NELR(IROW)
        RFSUM=0.0
        RXSUM=0.0
        DO 450 INJ=IEL1, IELL
          ICNT=ICNT+1
          THETA=TI*(INJ-IEL1)
$INSERT FMF.EQ
$INSERT XMF.EQ
          FMF(INJ)=FMF(INJ)/(1.0+HGMRO)
          XMF(INJ)=XMF(INJ)+FMF(INJ)*HGMRO
          RFSUM=RFSUM+FMF(INJ)
          RXSUM=RXSUM+XMF(INJ)
          TFF=TFF+FMF(INJ)
          TXF=TXF+XMF(INJ)
          THVM=THVM+FMF(INJ)
          TXVM=TXVM+FMF(INJ)*HGMRO
450 CONTINUE
          IF (IECHO .EQ. 1) WRITE(6,48) FEQ, XEQ, RFSUM, RXSUM
460 CONTINUE
499 READ(5,49) CFLOW, CFMR
        IF (IECHO .EQ. 1) WRITE(6,50) CFLOW, CFMR
        TXVM=TXVM+CFLOW*CFMR/(1.0+CFMR)
        THVM=THVM+CFLOW/(1.0+CFMR)
        TXF=TXF+CFLOW*CFMR/(1.0+CFMR)
        TFF=TFF+CFLOW/(1.0+CFMR)
        IF (IECHO .EQ. 1) WRITE(6,51) TFF, THVM, TXF, TXVM
C
C      READ INJECTOR
C
      READ(5,10)TITLE
      IF(TITLE .NE. DUMMY) GOTO 9000
      READ(5,60,ERR=9300)ITYPE
      IF (ITYPE .LT. 1 .OR. ITYPE .GT. 3) GOTO 9300
      IF (ITYPE .EQ. 2) GOTO 520
      IF (ITYPE .EQ. 3) GOTO 530
C
C      SHEAR COAX INJECTOR

```

```

C
  READ(5,61)RXP, RFP, TPOST, RECESS, RMS, XDL, FCD, XCD
  IF (IECHO .EQ. 1) WRITE(6,62)RXP, TPOST, RECESS, RMS, XDL,
  *
  RECESS=RECESS/12.0
  RMS=RMS/12.0
  XDL=XDL/12.0
  GOTO 550

C
C   SWIRL COAX INJECTOR
C
520 READ(5,63)RXP, RFP, TPOST, CSA, FCD, XCD
  IF (IECHO .EQ. 1) WRITE(6,64) RXP, TPOST, CSA, XCD, RFP, FCD
  GOTO 540

C
C   INPINGING TRIAX INJECTOR
C
530 READ(5,61)RXP, RFP, TPOST, CSA, PFI, FCD, XCD
  IF (IECHO .EQ. 1) WRITE(6,65) RXP, TPOST, CSA, XCD, RFP, FCD, PFI
540 CSA=CSA*3.1416/180.0
550 RFP=RFP/12.0
  RXP=RXP/12.0
  TPOST=TPOST/12.0

C
C   READ PROBLEM CONTROL PARAMETERS
C   **** IF DEFAULTS ARE CHANGED, THEY MUST ALSO BE UPDATED IN PLSET.CPL ***
C
  NSEG=NROWS+1
  NSLICE=12
  IF (NROWS .EQ. 1 .AND. NEL .LT. 13 .AND. NEL .GT. 1) NSLICE=NEL-1
  XSTEP=2.0**(-8)
  ASTEP=2.0**(-10)
  READ(5,CONT)
  IF (IECHO .EQ. 1) WRITE(6,70)NSEG, NSLICE, ASTEP,XSTEP
  DO 600 I=1,NSLICE
    DO 600 J=1,NSEG
      XGRID(I,J)=0.0
      HGRID(I,J)=0.0
600 CONTINUE
  RETURN

C
C   ERROR CONDITION
C
9000 WRITE(6,*) 'BLANK LINE GROUP SEPERATOR MISSING, RUN STOPPED'
  GOTO 9999
9110 WRITE(6,*) 'ERROR READING XW,RW FOR POSITION=',I,', RUN STOPPED'
  GOTO 9999
9120 WRITE(6,*) 'FIRST XW NOT AT INJECTOR FACE, RUN STOPPED'
  GOTO 9999
9130 WRITE(6,*) 'XW(I) .GE. XW(I+1) FOR I=',I,', RUN STOPPED'
  GOTO 9999
9140 WRITE(6,*) 'INCONSISTENCY WITH RW FOR XW=',XW(I),RW(I),RW(I+1)
  GOTO 9999
9150 WRITE(6,*) 'NO THROAT CALCULATED, RUN STOPPED'
  GOTO 9999
9210 WRITE(6,*) 'COORDINATES FOR EL#=',I,', MISSING, RUN STOPPED'
  GOTO 9999
9220 WRITE(6,*) 'ELEMENT ',I,', NOT IN ASCENDING RADIUS ORDER, ',
  *
  'RUN STOPPED'
  GOTO 9999

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9230 WRITE(6,*) 'INJECTOR OUTSIDE CHAMBER FOR I=',I,' ', RUN STOPPED'
      GOTO 9999
9240 WRITE(6,*) 'INJECTORS NOT IN ASCENDING THETA ORDER, RUN STOPPED'
      GOTO 9999
9250 WRITE(6,*) NROWS, ' ROWS NOT INPUT, LAST ROW OF RADIUS=',RROW(IROW)
      *      , ' FEET, RUN STOPPED'
      GOTO 9999
9260 WRITE(6,*) 'ERROR OPENING FMF.EG, RUN STOPPED'
      GOTO 9998
9270 WRITE(6,*) 'ERROR OPENING XMF.EG, RUN STOPPED'
      GOTO 9998
9280 WRITE(6,*) 'NO ROW INFORMATION FOUND FOR ROW=',IROW,' RUN STOPPED'
      GOTO 9998
9290 WRITE(6,*) 'NO FUEL EQUATION FOUND FOR ROW=',IROW,' RUN STOPPED'
      GOTO 9998
9295 WRITE(6,*) 'NO OX EQUATION FOUND FOR ROW=',IROW,' RUN STOPPED'
      GOTO 9998
9300 WRITE(6,*) 'ERROR WITH INJECTOR TYPE, ITYPE=',ITYPE,' RUN STOPPED'
      GOTO 9999
9998 CLOSE(UNIT=8)
      CLOSE(UNIT=9)
9999 CLOSE(UNIT=5)
      CLOSE(UNIT=6)
      CLOSE(UNIT=7)
      STOP
      END

```

SUBROUTINE SHEAR

PROGRAM TO CALCULATE DROPLET FORMATION FOR SHEAR COAXIAL INJECTOR  
 RETURNS MASS MEDIAN DROPSIZE, NUMBER OF DROPS/SEC AND DISTRIBUTION  
 PARAMETER, SG AS WELL AS MEAN ATOMIZATION LENGTH AND INJECTION VELOCITY

OUTPUTS:

RMM=MASS MEDIAN DROPLET RADIUS, FT  
 TND=TOTAL NUMBER OF DROPS OF GENERATED BY ELEMENT  
 SG=DISTRIBUTION PARAMETER  
 AL=MEAN ATOMIZATION LENGTH, FT  
 VJ=MEAN INJECTION VELOCITY, FT/SEC IN THE FORM OF  
 VFACT, THE VAPORIZATION FACTOR IN LGEN

ERROR CONDITIONS:

X=-1 ITERATION NOT CONVERGED AFTER 20 LOOPS  
 X=-2 AX BEYOND THROAT  
 X=-3 R AND XD NOT DIMENSIONED LARGE ENOUGH, I.E. A  
 GREATER THAN ANTICIPATED NUMBER OF DROPSIZES GENERATED  
 X=-4 XND LT 1 OR RJET LT 0

\*INSERT COMMON

DIMENSION R(200), XD(200)

IF (ITRACE .EQ. 2) WRITE(6,1)  
 1 FORMAT(/1X, 'ENTER SHEAR')  
 XCD=1.0  
 QFACT=4.0/3.0\*3.1416  
 VAPC=PC\*\*0.66/(1.0-(XTJ+460.)/(XTC+460.))\*\*0.4/XHV\*\*0.8/XMW\*\*0.35  
 DO 1000 INJ=1,NEL  
 IF (XMF(INJ) .EQ. XMF(INJ-1) .AND. FMF(INJ) .EQ. FMF(INJ-1))  
 \* GOTO 1100

INITIALIZE SIZES

AX=-RECESS  
 RJETO=RXF  
 A=RF  
 B=RXF+TPOST  
 EFF=FMF(INJ)  
 ELXF=XMF(INJ)-EFF\*HGMR0  
 EVXF=EFF\*HGMR0  
 HGMR(INJ)=HGMR0  
 HGCD=FCD  
 DO 2 I=1,200  
 R(I)=0.0  
 XD(I)=0.0  
 2 CONTINUE  
 IDC=0  
 TND(INJ)=0.0  
 VTOT=0.0  
 RTOT=0.0  
 IC=4  
 CALL DBLINT(PC, (TXF/TFF), IC, VSONIC)  
 XFILL=X1  
 VT=0.62/CR\*VSONIC  
 IF (X1 .GT. 0.0) GOTO 50  
 XFILL=X2  
 VT=VSONIC

```

C      CALCULATE OX VELOCITY PROFILE
C
50      CCO=(RMS/RXP)**2
        VMEAN=ELXF/XRHO/3.1416/RMS**2
        RE=2.0*XRHO*VMEAN*RMS/XMU
        IF (RE .GE. 4000.0) RCRIT=((1.0-SQRT(CCO))/0.75)**1.25*RE**0.25
        IF (RE .LT. 4000.0) RCRIT=((1.0-SQRT(CCO))/11.28)**2*RE
        AR=(XDL/RMS/2.0)
        IF (RCRIT .LE. AR) GOTO 150

```

```

C      CALCULATE XCD
C
        RE1=4.0
        CE1=CCO+0.214*AR
        IF (RE .LT. 1.0E5) GOTO 100
        RE1=5.0
        CE1=CCO+0.095*AR
        IF (RE .LT. 1.0E6) GOTO 100
        RE1=6.0
        CE1=0.054*AR
100      RE2=5.0
        CE2=CCO+0.095*AR
        IF (RE .LT. 1.0E5 .OR. RE .GE. 1.0E6) GOTO 120
        RE2=6.0
        CE2=CCO+0.054*AR
120      CCE=CE1+(CE2-CE1)/(RE2-RE1)*(LOG10(RE)-RE1)
        IF (CCE .GT. 1.0) CCE=1.0
        XCD=(CCO-0.5)+0.31789*EXP(0.89192*(CCE-CCO+0.5))
        RJETO=SQRT(RJET*XCD)

```

```

C      CALCULATE POTENTIAL CORE SIZE AND BL PROFILE
C
150      RJET=RJETO
        REM=AIN(T(LOG10(RE)))
        FR=RE/(10.0**REM)
        DEL=RMS*EXP(LOG(0.048)+(6.0-REM-FR)*0.44)*AR**0.8
        RP=RMS-DEL
        IF (RP .LT. 0.0) RP=0.0
        ICNT=0.0
        POW=0.7
        CFACT=(ELXF-XRHO*VMEAN*3.1416*RP**2)/3.1416/XRHO/VMEAN
200      F1=(RJET**2+RJET*RP*POW+(1.0+POW)*RP**2)
        F2=POW**2+3.0*POW+2.0
        F3=2.0*POW+3.0
        FX=F1/F2-CFACT
        FP=(F2*(RJET*RP+RP**2)-F1*F3)/F2**2
        POWN=POW-FX/FP
        IF (ABS(POW-POWN) .LE. 1.0E-4) GOTO 210
        ICNT=ICNT+1
        POW=POWN
        IF (ICNT .LE. 10) GOTO 200
        GOTO 9000

```

```

C      CALCULATE FUEL AND OX FREE SURFACE VELOCITIES
C
210      VF=(EFF+EVXF)/(FCD*FRHO*3.1416*(A**2-B**2))
        VFO=VF
220      VO=VMEAN
        IF (RJET .GT. RP) VO=VMEAN*(RJETO-RJET)/(RJETO-RP)
        IF (AX .GE. 0.0 .AND. AX .LE. VFILL) VF=VFO-(VFO-VT)*AX/XFILL

```

```

      IF (VF .GT. 5500.0) WRITE(6,230) VF, INJ
230  FORMAT(/1X, '***** WARNING *****',
*      /10X, 'FUEL INJECTION VELOCITY APPROACHING SONIC VELOCITY, VF='
*      ,E11.4, ' FOR INJ=', I4)
      IF (VO .GT. 500.0) WRITE(6,240) VO, INJ
240  FORMAT(/1X, '***** WARNING *****',
*      /10X, 'OX INJECTION VELOCITY EXCESSIVELY HIGH, VX='
*      ,E11.4, ' FOR INJ=', I4)

C
C      CALCULATE INTERFACIAL SHEAR STRESS AND DROPSIZE (SCALING FACTOR OF 1500)
C
      TAU1=(XMU+FMU)/2.0*(VF-VO)/(B-RJET)/1500.0
      RM=8.0*XST/TAUI/32.1739

C
C      CALCULATE NEW X POSITION
C
      AX=AX+ASTEP
      IF (AX .GT. XW(NXP)) GOTO 9100
      IDC=IDC+1
      IF (IDC .GT. 200) GOTO 9200

C
C      CALCULATE NUMBER OF DROPS USING RJET
C
      XMI=(RJET/RJET0)**2*ELXF
      XND=MAX(1.0, AINT(3.1416*RJET/2.0/RM))
      VDROPS=XND*QFACT*RM**3.0
      VJETM=3.1416*RJET**2*ASTEP
      IF (VJETM .GT. (2.0*VDROPS)) GOTO 300

C
C      LAST DROPLETS FORMED, JET DISAPPEARS
C
      XND=MAX(1.0, AINT(VJETM/VDROPS*XND))
      RM=(VJETM/XND/QFACT)**(1./3.)
      RJET=0.0
      GOTO 325

C
C      ITERATE WITH TRUNCATED CONE FORMULA
C      IF NONCONVERGENT, RETURN INJ=-1
C
300  ICNT=0
      RN=RJET
310  FX=3.1416/3.0*ASTEP*(RJET**2.0+RJET*RN+RN**2)-VJETM+VDROPS
      FP=3.1416/3.0*ASTEP*(RJET+2.0*RN)
      RNN=RN-FX/FP
      IF (ABS(RN-RNN) .LT. 0.0005*RN) GOTO 320
      ICNT=ICNT+1
      RN=RNN
      IF (ICNT .LT. 10) GOTO 310
      GOTO 9300
320  RJET=RNN

C
C      CALCULATE NDROPS/SEC BASED ON CONTINUITY
C
325  XMO=(RJET/RJET0)**2*ELXF
      DVF=(XMI-XMO)/XRHO
      XND=DVF/(QFACT*RM**3)
      IF (IAFLO .GT. 2) WRITE(6,326) AX, VF, VO, TAU1, RM, XND
326  FORMAT(1X, 'AX, VF, VO, T, RM, XND=', 4(F9.4, 3X), 2(E12.4, 3X))
      IF (XND .LT. 1.0 .OR. RJET .LT. 0.0) GOTO 9400

```

```

C      VAPORIZE DROPLET IF STILL IN CUP
C
      IF (AX .GE. 0.0) GOTO 400
      VJ=(4.0*VD+VF)/5.0
      G1=VAPC/RM**1.45/VJ**0.75
      D1=-AX
      CALL CVAP (G1, D1, PV)
      VP=PV*XND*XRHO*QFACT*RM**3
      TXVM=TXVM+VP
      EVXF=EVXF+VP
      RM=((1.0-PV)*RM**3)**(1.0/3.0)
      IF (IAFLG .GT. 2) WRITE(6,330) (100.0*PV), RM
330    FORMAT(10X,F6.2,'% OF DROP VAPORIZED, NEW RM=',E12.5)
      HGMR(INJ)=EVXF/EFF
      IC=5
      CALL DBLINT(PC, HGMR(INJ), IC, RHO)
      IF (HGCD .GE. 1.0) GOTO 350
      HGCD=(EFF+EVXF)/(3.1416*(A**2-RJET**2)*RHO*VF)
      GOTO 400
350    VF=(EFF+EVXF)/(RHO*3.1416*(A**2-RJET**2))
      VFO=VF

C
C      INSERT INTO LIST IN ASCENDING ORDER
C
400    IF (RM .GT. 0.0) GOTO 410
      IDC=IDC-1
      GOTO 775
410    IF (IDC .GT. 1) GOTO 420
      I=1
      GOTO 700
420    DO 450 I=1, (IDC-1)
      IF (ABS(R(I)-RM) .LE. 0.005*RM) GOTO 500
      IF (R(I) .GE. RM) GOTO 600
450    CONTINUE
      I=IDC
      GOTO 700

C
C      SIMILAR SIZE ALREADY EXISTS
C
500    VI=QFACT*R(I)**3*XD(I)
      VN=QFACT*RM**3*XND
      XD(I)=XD(I)+XND
      R(I)=((VI+VN)/QFACT/XD(I))**((1.0/3.0))
      RM=R(I)
      IDC=IDC-1
      GOTO 750

C
C      MOVE LARGER DROPS DOWN
C
600    DO 650 K=IDC, (I+1), -1
      R(K)=R(K-1)
      XD(K)=XD(K-1)
650    CONTINUE
700    R(I)=RM
      XD(I)=XND

C
C      SUM FOR MEAN VALUE CALCULATIONS, THEN CONTINUE
C
750    TND(INJ)=TND(INJ)+XND
      WV=XND*QFACT*RM**3

```

```

      VTOT=VTOT+VD*VV
      RTOT=RTOT+VV
775   IF (RJET .GT. 0.0) GOTO 220
C
C   CALCULATE MEAN QUANTITIES
C
      VMEAN=(RTOT/TND(INJ))
      RMEAN=(VMEAN/QFACT)**(1./3.)
      AL(INJ)=(AX-RECESS)/2.0
      VJ=VTOT/TND(INJ)/VMEAN
      VFACT(INJ)=VAPC/VJ**0.75
C
C   CALCULATE MASS MEDIAN
C
      AM=0.0
      CN=TND(INJ)/2.0
      DO 800 I=1, IDC
        AM=AM+XD(I)
        IF (AM .LT. CN) GOTO 800
        RMM=R(I)
        GOTO 810
800   CONTINUE
810   RMO(INJ)=RMM
C
C   CALCULATE DISTRIBUTION (STANDARD DEVIATION)
C
      SUM=0.0
      VMM=QFACT*RMM**3
      DO 900 I=1, IDC
        IF (IAFLG .GT. 1) WRITE(6,820)I,R(I),XD(I)
820   FORMAT(5X,'INDEX=',I3,5X,'RM=',E10.4,5X,'NDROPS/SEC=',E11.5)
        VDI=QFACT*R(I)**3
        SUM=SUM+(VDI-VMM)**2*XD(I)
900   CONTINUE
      SD=(SQRT(SUM/TND(INJ)))/QFACT**(.1/3.)
      SG(INJ)=(SD+RMM)/RMM
C
C   CORRECT TND TO CONSERVE MASS
C
      TND(INJ)=(ELXF-EFF*(HGMR(INJ)-HGMR0))/(QFACT*XRHO*RMM**3)/
*      ((0.165/SG(INJ)**3)+0.67+(0.165*SG(INJ)**3))
C
C   FILL INITIAL DROPSIZES
C
910   RMX(1,INJ)=RMM/SG(INJ)
      RMX(2,INJ)=RMM
      RMX(3,INJ)=RMM*SG(INJ)
      DRP(1,INJ)=RJETO/2.0
      DRP(2,INJ)=RJETO/2.0
      DRP(3,INJ)=RJETO/2.0
      DXV(1,INJ)=VJ
      DXV(2,INJ)=VJ
      DXV(3,INJ)=VJ
      DRV(1,INJ)=0.0
      DRV(2,INJ)=0.0
      DRV(3,INJ)=0.0
      IF (IAFLG .GT. 0) WRITE(6,920) INJ, (RMO(INJ)*304800.),
*      (RMEAN*304800.), SG(INJ), TND(INJ), (AL(INJ)*12.0), VJ
920   FORMAT(5X,'INJECTOR NUMBER=',I3,/9X,'MASS MEDIAN (MICRONS)=',
*      F7.2,5X,'MASS MEAN (MICRONS)=',F7.2,8X,'SIGMAG=',F6.3,/9X,

```

```

      *          'NUMBER OF DROPS=' E11.5,7X, 'ATOMIZATION LENGTH (IN)=' ,
      *          F6.3,5X, 'INJECTION VELOCITY (F/S)=' ,F6.2)
1000 CONTINUE
      FCD=MIN(HGCD,1.0)
      RETURN
C
C      INJECTOR SAME AS PREVIOUS
C
1100 RMO(INJ)=RMO(INJ-1)
      SG(INJ)=SG(INJ-1)
      TND(INJ)=TND(INJ-1)
      AL(INJ)=AL(INJ-1)
      VFACT(INJ)=VFACT(INJ-1)
      HGMR(INJ)=HGMR(INJ-1)
      TXVM=TXVM+(HGMR(INJ)-HGMR0)*FMF(INJ)
      GOTO 910
C
C      ERROR CONDITION, ABORT RUN
C
9000 WRITE(6,*) 'POW NOT CONVERGENT IN SHEAR, POW=',POW,' FOR INJ=',INJ
      X=-1.0
      RETURN
9100 WRITE(6,*) 'AX BEYOND THROAT IN SHEAR FOR INJ=',INJ,' RUN STOPPED'
      X=-2.0
      RETURN
9200 WRITE(1,*) 'MORE DROPS PRODUCED THAN DIMENSIONED FOR, RUN STOPPED'
      X=-3.0
      RETURN
9300 WRITE(6,*) 'RJET NOT CONVERGENT IN SHEAR, RJET=',RN,' FOR INJ=',INJ
      X=-2.0
      RETURN
9400 WRITE(6,*) 'LESS THAN 1 DROP FORMED OF NEG RJET, XND,RJET=',XND,
      *          ', ',RJET
      X=-4.0
      RETURN
      END

```

```

SUBROUTINE CVAP (D1, G1, V)
C
C   CALCULATES CHANGE IN DROP RADIUS DUE TO EVAPORATION IN
C   INJECTOR CUP. BASED ON PRIEM-HEIDMANN GENERALIZED LENGTH CORRELATION
C
*INSERT COMMON
C
C   LOCAL BLOCK COMMON ONLY
C
COMMON /VFRAC/ GL(41), F(41)
DATA GL /0.0, .01, .02, .04, .06, .1, .2, .4, .6, 1., 1.5, 2., 3., 4., 5., 6., 7.,
*      8., 9., 10., 13., 15., 17., 20., 23., 25., 28., 30., 33., 35., 38., 40.,
*      45., 50., 55., 60., 70., 80., 90., 100., 110. /
DATA F /0.0, .0008, .003, .009, .016, .031, .055, .123, .173, .26, .343,
*      .418, .522, .60, .66, .706, .746, .779, .804, .828, .879, .902,
*      .92, .94, .955, .963, .972, .976, .982, .985, .9885, .9905,
*      .994, .996, .9973, .9982, .9992, .9996, .9998, .9999, 1.0 /
C
IF (ITRACE .EQ. 2) WRITE(6,1)
1 FORMAT(/1X, 'ENTER CVAP')
C
C   CALCULATE GENERALIZED LENGTH, GLEN
C
GLEN=0.0137466*D1/CR**0.44*G1
C
C   VAPORIZATION INTERPOLATION CALCULATION
C
IF (GL(1) .GE. GLEN) GOTO 150
IF (GLEN .GE. GL(41)) GOTO 175
DO 125 I=2,41
  IF (GLEN .GT. GL(I)) GOTO 125
  V=F(I-1)+(GLEN-GL(I-1))/(GL(I)-GL(I-1))*(F(I)-F(I-1))
RETURN
125 CONTINUE
150 V=F(1)
RETURN
175 V=F(41)
RETURN
END

```

# SUBROUTINE SWIRL

PROGRAM TO CALCULATE DROPLET FORMATION FOR HOLLOW CONE SWIRLER COAXIAL  
AND IMPINGING TRIAXIAL INJECTORS. RETURNS MASS MEDIAN DROPSIZE, NUMBER  
OF DROPS/SEC. USES A DISTRIBUTION PARAMETER OF 2.3

## OUTPUTS:

RMM=MASS MEDIAN DROPLET RADIUS, FT  
TND=TOTAL NUMBER OF DROPS OF GENERATED BY ELEMENT  
SG=DISTRIBUTION PARAMETER  
AL=MEAN ATOMIZATION LENGTH, FT  
VJ=MEAN INJECTION VELOCITY, FT/SEC IN THE FORM OF  
VFACT, THE VAPORIZATION FACTOR IN LGEN

## ERROR CONDITIONS:

X=-1 ITERATION NOT CONVERGED AFTER 20 LOOPS

\$INSERT COMMON

IF (ITRACE .EQ. 2) WRITE(6,1)  
1 FORMAT(/1X, 'ENTER SWIRL')  
VAPC=PC\*\*0.66/(1.0-(XTJ+460.)/(XTC+460.))\*\*0.4/XHV\*\*0.8/XMW\*\*0.35  
FFACT=1.0  
IF (ITYPE .EQ. 3) FFACT=(1.0-PFI/100.0)

LOOP THROUGH ALL INJECTORS

DO 500 INJ=1,NEL  
IF (XMF(INJ) .EQ. XMF(INJ-1) .AND. FMF(INJ) .EQ. FMF(INJ-1))  
\* GOTO 600

INITIALIZE SIZES

A=RFP  
B=RXPT+TPOST  
EFF=FMF(INJ)\*FFACT  
ELXF=XMF(INJ)-FMF(INJ)\*HGMRO  
EVXF=EFF\*HGMRO  
HGMR(INJ)=HGMRO

CALCULATE FUEL AND OX VELOCITIES

VF=(EFF+EVXF)/(FCD\*FRHO\*3.1416\*(A\*\*2-B\*\*2))  
VO=ELXF/(XCD\*XRHO\*3.1416\*RXPT\*\*2)  
IF (VF .GT. 5500.0) WRITE(6,50) VF  
50 FORMAT(/1X, '\*\*\*\*\* WARNING \*\*\*\*\*',  
\* /10X, 'FUEL INJECTION VELOCITY APPROACHING SONIC VELOCITY, VF=',  
\* ,E11.4, ' FOR INJ=',I4)  
IF (VO .GT. 500.0) WRITE(6,60) VO  
60 FORMAT(/1X, '\*\*\*\*\* WARNING \*\*\*\*\*',  
\* /10X, 'OX INJECTION VELOCITY EXCESSIVELY HIGH, VX=',E11.4  
\* , ' FOR INJ=',I4)

CALCULATE MOMENTUM ANGLE

PSI=ATAN((VO\*ELXF\*SIN(CSA))/((EFF+EVXF)\*VF+ELXF\*VO\*COS(CSA)))

CALCULATE ATOMIZATION LENGTH

ALEN=1.0

```

      ICNT=0
      F1=SIN(PSI)/12.0/RXP
      F2=50.6970*RXP*XCD*(XRHO*VO/XMU)**0.2
100   FX=ALEN**0.8+F1*ALEN**1.8-F2
      FP=0.8*ALEN**(-0.2)+1.8*F1*ALEN**0.8
      ALENN=ALEN-FX/FP
      IF (ABS(ALEN-ALENN) .LE. 0.001*ALEN) GOTO 200
      ICNT=ICNT + 1
      ALEN=ALENN
      IF (ICNT .LT. 10) GOTO 100
      GOTO 9000

```

C  
C  
C     CALCULATE TCRIT AND RMM

```

200   TCRIT=6.0*XCD*RXP/(1.0+SIN(PSI)*ALEN/12.0/RXP)
      RMM=0.62035*TCRIT/12.0
      RMO(INJ)=RMM
      AL(INJ)=ALEN
      VJ=VO*COS(PSI)
      VFACT(INJ)=VAPC/VJ**0.75
      SG(INJ)=2.3

```

C  
C  
C     CALCULATE TND TO CONSERVE MASS

```

      TND(INJ)=(ELXF-EFF*(HGMR(INJ)-HGMR0))/((4.1888*XRHO*RMM**3)/
*      ((0.165/SG(INJ)**3)+0.67+(0.165*SG(INJ)**3))

```

C  
C  
C     FILL INITIAL DROPSIZES

```

300   RMX(1,INJ)=RMM/SG(INJ)
      RMX(2,INJ)=RMM
      RMX(3,INJ)=RMM*SG(INJ)
      DRP(1,INJ)=ALEN*SIN(PSI)
      DRP(2,INJ)=ALEN*SIN(PSI)
      DRP(3,INJ)=ALEN*SIN(PSI)
      DXV(1,INJ)=VJ
      DXV(2,INJ)=VJ
      DXV(3,INJ)=VJ
      DRV(1,INJ)=VO*SIN(PSI)
      DRV(2,INJ)=VO*SIN(PSI)
      DRV(3,INJ)=VO*SIN(PSI)
      IF (IAFLG .GT. 0) WRITE(6,400) INJ, (RMO(INJ)*304800.),
*      SG(INJ), TND(INJ), (AL(INJ)*12.0), VJ
400   FORMAT(5X, 'INJECTOR NUMBER=', I3, /9X, 'MASS MEDIAN (MICRONS)=',
*      F7.2, 6X, 'SIGMAQ=', F6.3, 5X, 'NUMBER OF DROPS=', E11.5,
*      /9X, 'ATOMIZATION LENGTH (IN)=', F6.3, 5X,
*      'INJECTION VELOCITY (F/S)=', F6.2)
500   CONTINUE
      RETURN

```

C  
C  
C     INJECTOR SAME AS PREVIOUS

```

600   RMO(INJ)=RMO(INJ-1)
      SG(INJ)=SG(INJ-1)
      TND(INJ)=TND(INJ-1)
      AL(INJ)=AL(INJ-1)
      VFACT(INJ)=VFACT(INJ-1)
      HGMR(INJ)=HGMR(INJ-1)
      TXVM=TXVM+(HGMR(INJ)-HGMR0)*FMF(INJ)
      GOTO 300

```

```
C
C      ERROR, ABORT RUN
C
9000 WRITE(6,*) 'ALEN NOT CONVERGENT IN SWIRL, ALEN=',ALEN,
*              ' FOR INJ=',INJ
      X=-1.0
      RETURN
      END
```

```

      SUBROUTINE RFILL
C
C      SUBROUTINE TO CALCULATE RECIRCULATION ZONE FEATURES
C
*INSERT COMMON
      IF (ITRACE .EQ. 2) WRITE(6,1)
1  FORMAT(1X, 'ENTER RFILL')
C
C      CALCULATE STREAMTUBE 100% ERE C* AND COMBUSTION ZONE
C      GAS VELOCITY. STORE IN COM(1,NEL) AND COM(2,NEL)
C      INITIALIZE PREVAPORIZED OX FROM HOT GAS, STORE IN COM(3,NEL)
C
      IC=5
      CALL DBLINT(PC, HGMR(1), IC, FRHO)
      IC=6
      DO 100 INJ=1,NEL
          VMR=XMF(INJ)/FMF(INJ)
          CALL DBLINT(PC, VMR, IC, COM(1,INJ))
          VF=FMF(INJ)*(1.0+HGMR(INJ))/FRHO/
*          (FCD*3.1416*(RFP**2-RXP**2))
          VX=(XMF(INJ)-FMF(INJ)*HGMR(INJ))/XRHO/(XCD*3.1416*RXP**2)
          IF (VF .GT. 5500.0) WRITE(6,50) VF
50      FORMAT(/1X, '***** WARNING *****',
*          /10X, 'FUEL INJECTION VELOCITY APPROACHING SONIC VELOCITY, VF=',
*          ,F8.2, ' FOR INJ=', I4)
          IF (VX .GT. 500.0) WRITE(6,60) VX
60      FORMAT(/1X, '***** WARNING *****',
*          /10X, 'OX INJECTION VELOCITY EXCESSIVELY HIGH, VX=', F8.2,
*          ' FOR INJ=', I4)
          COM(2,INJ)=(VF+8.0*VX)/9.0
          COM(3,INJ)=HGMR(INJ)*FMF(INJ)
100 CONTINUE
C
C      STEP IN X UNTIL CHAMBER CROSS-SECTION IS FILLED
C
      X=0.0
200 X=X+XSTEP
      CALL GRIDGEN
      IF (RWALLX .LE. 0.0) GOTO 8000
      AXC=3.1416*RWALLX**2*(1.0-CFLOW/(TFF+TXF))
      CALL MVAP
C
C      CALCULATE RFUEL FOR EACH ELEMENT, STORE IN ERPOS AND SUM.
C      STORE OX VAPOR IN COM(3,INJ)
C
      AFILL=0.0
      DO 300 INJ=1,NEL
          COM(3,INJ)=COM(3,INJ)+XVM(1,INJ)+XVM(2,INJ)+XVM(3,INJ)
          VMR=COM(3,INJ)/FMF(INJ)
          smr=8.0
          IC=6
          CALL DBLINT(PC, VMR, IC, CS)
          ERE=CS/COM(1,INJ)*(FMF(INJ)+COM(3,INJ))/(XMF(INJ)+FMF(INJ))
          RHO=0.0
          IC=5
          CALL DBLINT(PC, SMR, IC, RHO)
          DA=1.125*COM(3,INJ)/RHO/COM(2,INJ)
          RCORE=0.0
          IF (ITYPE .EQ. 1) RCORE=MAX(0.0, (RXP-RXP/2.0/AL(INJ)*
*          (X+RECESS)))

```

```

      RCG=SQRT(RCORE**2+DA*ERE/3.1416)
      RF=SQRT(RCG**2+(FMF(INJ)*(1.0+HGMR(INJ))-0.125*COM(3,INJ))*
*      (RFP**2-(RXP+TPOST)**2)/(FMF(INJ)*(1.0+HGMR(INJ))))
      ERPOS(INJ)=3.1416*RF**2
      AFILL=AFILL+ERPOS(INJ)
300 CONTINUE
C
C   ACCELERATE PARTICLES IN STREAMWISE DIRECTION
C
      CALL DMTUBE
      IF (X .LE. 0.0) GOTO B100
      PFILL=AFILL/AXC*100.0
C
C   CHECK FOR FILLED CHAMBER
C
      IF (IRFFLG .GT. 0) WRITE(6,310) X,PFILL
      IF (IRFFLG .GT. 0) WRITE(7,310) X,PFILL
310 FORMAT(/10X,'AT AXIAL POSITION X=',F7.4,F7.2,'% OF CROSS-SECTION'
*      , ' IS FILLED')
      IF (IRFFLG .GT. 0) CALL MRTUBE
      IF (PFILL .LT. 100.0) GOTO 200
C
C   CROSS SECTION FILLED, CALCULATE APPEARENT INJECTOR LOCATION
C   DISTRIBUTE MASS INTO GRID
C
      INJF=NEL+1
      REDGE=RWALLX
      FFACT=100.0/PFILL/(1.0-CFLOW/(TFF+TXF))
      DO 1000 IROW=NROWS,1,-1
        INJL=INJF-1
        INJF=INJL-NELR(IROW)+1
        ASUM=0.0
        DO 320 INJ=INJF,INJL
          ERPOS(INJ)=ERPOS(INJ)*FFACT
          ASUM=ASUM+ERPOS(INJ)
320      CONTINUE
        RO=SQRT(MAX(0.0,(REDGE**2-ASUM/3.1416)))
        RPOS=RO+(REDGE-RO)/2.0**0.5
        IF (NELR(IROW) .EQ. 1 .AND. RO .LT. 1.0E-5) RPOS=0.0
C
C   LOCATE SEG(S) TO ADD MASS TO
C
      R1=0.0
      DO 330 J=1,NSEG
        R2=RGRID(J)
        IF (RO .GE. R1 .AND. RO .LT. R2) GOTO 340
        R1=R2
330      CONTINUE
      GOTO 9000
340      IR1=J
      DO 350 J=IR1,NSEG
        R2=RGRID(J)
        IF (REDGE .GT. R1 .AND. REDGE .LE. R2) GOTO 360
        R1=R2
350      CONTINUE
      GOTO 9100
360      IR2=J
      IF (IRFFLG .GT. 1) WRITE(6,370) IROW, RO, REDGE, IR1, IR2
370      FORMAT(/5X,'INJECTOR ROW ',I3,' CONFINED BETWEEN R='F7.4,
*      ' AND 'F7.4,/10X,'THIS CORRESPONDS TO SEGS ',I3,' THRU ',I3)

```

```

C
C   LOCATE STREAMTUBE SLICE BOUNDARY AND CORRESPONDING SLICE(S)
C
      T1=0.0
      IT2=1
      DO 500 INJ=INJF,INJL
        STXVM=COM(3,INJ)
        T2=T1+ERPOS(INJ)/ASUM*360.0
        DO 400 J=IT2,NSLICE
          IF ((THETA*J) .GT. T1) GOTO 410
400      CONTINUE
          GOTO 9200
410      IT1=J
          J=NSLICE
          IF (T2 .GT. 360.0 .AND. (T2-360.0) .LT. 1.0E-4) GOTO 430
          DO 420 J=IT1,NSLICE
            IF ((THETA*J) .GE. T2) GOTO 430
420      CONTINUE
          GOTO 9300
430      IT2=J
          IF (IRFFLG .GT. 1) WRITE(6,440) INJ, T1, T2, IT1, IT2
440      FORMAT(/10X, 'INJECTOR ', I4, ' CONFINED BETWEEN THETA=', F6.2,
*           ' AND ', F6.2, /10X, 'THIS CORRESPONDS TO SLICES ', I3,
*           ' THRU ', I3)
C
C   DISTRIBUTE MASS TO GRID
C
      IF (IR1 .NE. IR2 .OR. IT1 .NE. IT2) GOTO 450
      XGRID(IT1,IR1)=XGRID(IT1,IR1)+STXVM
      HGRID(IT1,IR1)=HGRID(IT1,IR1)+FMF(INJ)
      GOTO 490
450      RS=REDGE**2-R0**2
      TS=T2-T1
      RB=R0
      RFACT=1.0
      DO 470 ISEG=IR1,IR2
        RG=RGRID(ISEG)
        IF (ISEG .EQ. IR2) RG=REDGE
        RFACT=(RG**2-RB**2)/RS
        RB=RGRID(ISEG)
        TB=T1
        TFACT=1.0
        DO 470 ISL=IT1,IT2
          ANG=ISL*THETA
          IF (ISL .EQ. IT2) ANG=T2
          SFACT=(ANG-TB)/TS
          TB=ISL*THETA
          FADD=SFACT*RFACT*FMF(INJ)
          XADD=SFACT*RFACT*STXVM
          HGRID(ISL,ISEG)=HGRID(ISL,ISEG)+FADD
          XGRID(ISL,ISEG)=XGRID(ISL,ISEG)+XADD
          IF (IRFFLG .GT. 2) WRITE(6,460) INJ,FADD,XADD,ISL,ISEG
460      FORMAT(10X, 'INJECTOR ', I3, ' ADDS ', E11.4,
*           ' LB/S OF FUEL AND ', E11.4, ' LB/S OF OX TO ',
*           'SLICE, SEG=', I3, ', ', I2)
470      CONTINUE
C
C   CALCULATE APPEARENT INJECTOR LOCATION AND DROPLET CHARACTERISTICS
C
490      ERPOS(INJ)=RPOS

```

```

      ETPOS(INJ)=(T2+T1)/2.0
      IF (IRFFLG .GT. 1) WRITE(6,495)INJ,ERPOS(INJ),ETPOS(INJ)
495  FORMAT(10X,'INJECTOR ',I3,' HAS AN APPEARENT LOCATION OF
      *      'R, THETA=',F7.4,F7.2)
      T1=T2
      DR1=DRP(1,INJ)/RFP*(REDGE-RO)/2.0
      DR2=DRP(2,INJ)/RFP*(REDGE-RO)/2.0
      DR3=DRP(3,INJ)/RFP*(REDGE-RO)/2.0
      DRV(1,INJ)=(DR1-DRP(1,INJ))/(X-AL(INJ))
      DRV(2,INJ)=(DR2-DRP(2,INJ))/(X-AL(INJ))
      DRV(3,INJ)=(DR3-DRP(3,INJ))/(X-AL(INJ))
      DRP(1,INJ)=DR1
      DRP(2,INJ)=DR2
      DRP(3,INJ)=DR3
500  CONTINUE
      REDGE=RO
1000 CONTINUE
C
C  DISTRIBUTE COOLANT FLOW UNIFORMLY
C
      IF (CFLOW .LE. 0.0) GOTO 1100
      CHPS=CFLOW/(1.0+CFMR)/NSEG/NSLICE
      CXPS=CFLOW*CFMR/(1.0+CFMR)/NSEG/NSLICE
      DO 1050 I=1,NSLICE
        DO 1050 J=1,NSEG
          HGRID(I,J)=HGRID(I,J)+CHPS
          XGRID(I,J)=XGRID(I,J)+CXPS
1050 CONTINUE
C
C  CALCULATE MEAN GAS PROPERTIES AND PRINT FULL MR DISTRIBUTION
C
1100 OVMR=TXVM/THVM
      IC=1
      CALL DBLINT (PC, OVMR, IC, CT)
      IC=2
      CALL DBLINT (PC, OVMR, IC, GMW)
      IC=5
      CALL DBLINT (PC, OVMR, IC, GRHO)
      IC=7
      CALL DBLINT (PC, OVMR, IC, GMU)
      GXV(1)=(TXVM+THVM)/GRHO/AXC
      RETURN
C
C  ERROR MESSAGES
C
8000 WRITE(6,*) 'ERROR FROM GRIDGEN, RUN STOPPED'
      X=-1.0
      RETURN
8100 WRITE(6,*) 'ERROR FROM DMTUBE, RUN STOPPED'
      RETURN
9000 WRITE(6,*) 'ERROR IN RFILL, INNER EDGE NOT FOUND FOR ROW=',IROW
      WRITE(6,*) 'RO,RGRID=',RO,RGRID
      X=-1.0
      RETURN
9100 WRITE(6,*) 'ERROR IN RFILL, OUTER EDGE NOT FOUND FOR ROW=',IROW
      WRITE(6,*) 'REDGE,RGRID=',REDGE,RGRID
      X=-1.0
      RETURN
9200 WRITE(6,*) 'ERROR IN RFILL, FIRST SLICE NOT FOUND FOR ROW=',IROW,
      *      'ELEMENT=',INJ,'THETA=',T1

```

```
      X=-1.0  
      RETURN  
9300 WRITE(6,*) 'ERROR IN RFILL, END SLICE NOT FOUND FOR ROW=', IROW,  
      *          'ELEMENT=', INJ, 'THETA=', T2  
      X=-1.0  
      RETURN  
      END
```

```

SUBROUTINE DMTUBE
C
C SUBROUTINE TO ACCELERATE DROPS DUE TO DRAG ACCELERATIONS
C USED IN STREAMTUBE ONLY
C
*INSERT COMMON
  IF (ITRACE .EQ. 2) WRITE(6,1)
  1 FORMAT(/1X,'ENTER DMTUBE')
  IF (IDMFLG .GT. 0) WRITE(6,2)
  2 FORMAT(/1X)
C
C LOOP THROUGH FOR ALL INJECTORS AND DROPSIZES
C
  TVMR=-1.0
  DO 1000 INJ=1,NEL
C
C CALCULATE STREAMTUBE PROPERTIES
C
    TVMRN=COM(3,INJ)/FMF(INJ)
    IF (ABS(TVMRN-TVMR) .LE. 0.1) GOTO 100
    TVMR=TVMRN
    IC=5
    CALL DBLINT (PC, TVMR, IC, GRHO)
    IC=7
    CALL DBLINT (PC, TVMR, IC, GMU)
  100 DO 1000 IRM=1,3
      IF (RMX(IRM,INJ) .LE. 0.0) GOTO 1000
C
C "EXPAND" DROPLET TO SUPERCRITICAL PROPORTIONS
C
    RM=RMX(IRM,INJ)*(XRHO/XRHOC)**(1./3.)
C
C CALCULATE RELATIVE GAS VELOCITY, DRAG AND ACCELERATION
C COMBUSTION ZONE GAS VELOCITY STORED IN COM(2,NEL)
C
    RP=DRP(IRM,INJ)
    RGV=SQRT((COM(2,INJ)-DXV(IRM,INJ))**2+DRV(IRM,INJ)**2)
    DVM=SQRT(DXV(IRM,INJ)**2+DRV(IRM,INJ)**2)
    IF (RGV .LT. 1.0) GOTO 1000
    DZ=XSTEP*SQRT(1.0+((DRV(IRM,INJ)/DXV(IRM,INJ))**2)
    ICNT=0
  175 CALL DRAG (RM, COM(2,INJ), DVM, RGV, DZ, ACC)
C
C MOVE DROPLET
C
    DXVN=COM(2,INJ)-(COM(2,INJ)-DXV(IRM,INJ))*ACC
    DRVN=DRV(IRM,INJ)*ACC
    DZN=XSTEP*SQRT(1.0+((DRV(IRM,INJ)+DRVN)/(DXV(IRM,INJ)+DXVN))
    *
    **2)
    IF (ABS(DZN-DZ) .LE. 0.01*DZ) GOTO 180
    DZ=DZN
    ICNT=ICNT+1
    IF (ICNT .GT. 10) GOTO 9000
    GOTO 175
  180 DRP(IRM,INJ)=RP+(DRV(IRM,INJ)+DRVN)/(DXV(IRM,INJ)+DXVN)
    *
    *XSTEP
    DXV(IRM,INJ)=DXVN
    DRV(IRM,INJ)=DRVN
    IF (DXV(IRM,INJ) .LT. 0.0) GOTO 9100
    IF (DXV(IRM,INJ) .GT. COM(2,INJ)) GOTO 9200

```

```

      IF (IDMFLG .GT. 0) WRITE(6,200) IRM, INJ, RP, DRP(IRM,INJ),
      *                               DXV(IRM,INJ), DRV(IRM,INJ),COM(2,INJ)
200  FORMAT(1X,'DROP #',I1,' FROM INJ=',I3,' MOVED RADIALLY ',
      *      'FROM ',F7.3,' TO ',F7.3,' WITH NEW AXIAL AND ',
      *      'RADIAL VEL=',F7.2,' ',F7.2,'; GAS VELOCITY=',F7.2)
1000 CONTINUE
      RETURN
C
C      ERROR CONDITION
C
9000 WRITE(6,*) 'DZ NOT CONVERGING IN DMTUBE FOR INJ,IRM=',INJ,', ',IRM,
      *          ' AT X=',X,' RUN STOPPED'
      X=-1.0
      RETURN
9100 WRITE(6,*) 'DROP ACCELERATED TO NEGATIVE VELOCITY AT X=',X,
      *          ' INJ,IRM=',INJ,IRM,' ACC=',ACC
      X=-1.0
      RETURN
9200 WRITE(6,*) 'DROP ACCELERATED BEYOND GAS VELOCITY AT X=',X,
      *          ' INJ,IRM=',INJ,IRM,' ACC=',ACC
      X=-1.0
      RETURN
      END

```

```

      SUBROUTINE MRTUBE
C
C      SUBROUTINE TO CALCULATE VMR FOR STREAMTUBES AS A FUNCTION OF X
C
C      *INSERT COMMON
      DIMENSION OUT(10)
C
      IF (ITRACE .EQ. 2) WRITE(6,1)
1  FORMAT(/1X,'ENTER MRTUBE')
      WRITE(6,2)
2  FORMAT(/1X)
C
C      CALCULATE VMR FOR STREAMTUBES 10 AT A TIME
C
      NL1=1
100  NLL=NL1+9
      IF (NLL .GT. NEL) NLL=NEL
      ICNT=0
      DO 150 INJ=NL1,NLL
          ICNT=ICNT+1
          OUT(ICNT)=-1.0
          IF (COM(3,INJ) .GT. 0.0) OUT(ICNT)=99.99
          IF (FMF(INJ) .GT. 0.0) OUT(ICNT)=COM(3,INJ)/FMF(INJ)
150  CONTINUE
      WRITE(6,175) NL1, NLL, (OUT(I),I=1,ICNT)
175  FORMAT(5X,'TUBE MR FOR INJ=',I3,'-',I3,4X,10(F5.2,4X))
      IF (NLL .EQ. NEL) RETURN
      NL1=NLL+1
      GOTO 100
      END

```

```

      SUBROUTINE GRIDGEN
C
C      SUBROUTINE TO GENERATE COMPUTATION CELLS AT X AND RETURNS RWALLX
C
$INSERT COMMON
      IF (ITRACE .EQ. 2) WRITE(6,1)
1    FORMAT(/1X,'ENTER GRIDGEN')
C
C      CALCULATE RWALL AT X
C
      DO 100 I=2,NXP
        IF (XW(I) .GE. X) GO TO 110
100   CONTINUE
      IF (IEFLG .EQ. 0) GOTO 300
      GOTO 990
110   RWALLX=RW(I-1)+(RW(I)-RW(I-1))/(XW(I)-XW(I-1))*(X-XW(I-1))
      ANGLEX=WANGLE(I-1)
C
C      CALCULATE SLICE ANGLE IN DEGREES
C      NB: TRIG FUNCTIONS REQUIRE RADIAN INPUTS
C
120   THETA=360.0/NSLICE
C
C      CALCULATE NSEG EQUAL AREA RADIAL AREAS
C
      ASEG=RWALLX**2/NSEG
      RGRID(NSEG)=RWALLX
      DO 200 I=1,(NSEG-1)
        RGRID(I)=SQRT(FLOAT(I)/FLOAT(NSEG))*RWALLX
200   CONTINUE
      RETURN
C
C      NEXT STEP MOVES PAST THROAT
C
300   XP=X-XSTEP
      XSTEP1=XW(NXP)-XP
      IF (XSTEP1 .GE. XSTEP .OR. XSTEP1 .LE. 0) GOTO 990
      IEFLG=1
      X=XW(NXP)
      XSTEP=XSTEP1
      RWALLX=RW(NXP)
      ANGLEX=WANGLE(NXP-1)
      GOTO 120
C
C      ERROR
C
990   WRITE (6,*)'*** X BEYOND THROAT IN GRIDGEN, X,XW(NXP)=',X,XW(NXP)
      RWALLX=0.0
      RETURN
      END

```

# SUBROUTINE MVAP

CALCULATES CHANGE IN DROP RADIUS DUE TO EVAPORATION  
BASED ON PRIEM-HEIDMANN GENERALIZED LENGTH CORRELATION

\*INSERT COMMON

LOCAL BLOCK COMMON ONLY

```
COMMON /VFRAC/ GL(41), F(41)
DATA GL /0.0, .01, .02, .04, .06, .1, .2, .4, .6, 1., 1.5, 2., 3., 4., 5., 6., 7.,
*      8., 9., 10., 13., 15., 17., 20., 23., 25., 28., 30., 33., 35., 38., 40.,
*      45., 50., 55., 60., 70., 80., 90., 100., 110. /
DATA F /0.0, .0008, .003, .009, .016, .031, .055, .123, .173, .26, .343,
*      .418, .522, .60, .66, .706, .746, .779, .804, .828, .879, .902,
*      .92, .94, .955, .963, .972, .976, .982, .985, .9885, .9905,
*      .994, .996, .9973, .9982, .9992, .9996, .9998, .9999, 1.0 /
```

```
IF (ITRACE .EQ. 2) WRITE(6,1)
1 FORMAT(/1X, 'ENTER MVAP')
```

LOOP THROUGH ALL ELEMENTS AND DROPSIZES

```
DO 500 INJ=1,NEL
  DO 500 IRM=1,3
    XVM(IRM,INJ)=0.0
    IF (X .LE. AL(INJ)) GOTO 500
    IF (RMX(IRM,INJ) .LE. 0.0) GOTO 500
```

FIX CANT ANGLE, A, TO 0

A=0.0

CALCULATE RMO

```
RMI=RMO(INJ)
IF (IRM .EQ. 1) RMI=RMI/SG(INJ)
IF (IRM .EQ. 3) RMI=RMI*SG(INJ)
XND=0.67*TND(INJ)
IF (IRM .NE. 2) XND=0.165*TND(INJ)
X0=AL(INJ)
```

CALCULATE GENERALIZED LENGTH, GLEN

```
IF (X .GT. (X0+X1)) GOTO 50
D1=X-X0
D2=0.0
GOTO 60
50 D1=X1-X0
D2=X-X1
60 G1=VFACT(INJ)/RMI**1.45/COS(A)**0.75
G2=0.0137466*(D1/CR**0.44+0.83*D2/(CR**0.22*
* ((1.0+1/SQRT(CR)+1/CR)/3.0)**0.33))
GLEN=G1*G2
```

VAPORIZATION INTERPOLATION CALCULATION

```
IF (GL(1) .GE. GLEN) GOTO 150
IF (GLEN .GE. GL(41)) GOTO 175
DO 125 I=2,41
```

```

        IF (GLEN GT. GL(1)) GOTO 125
        V=F(I-1)+(GLEN-GL(I-1))/(GL(I)-GL(I-1))*(F(I)-F(I-1))
        GOTO 200
125      CONTINUE
150      V=F(1)
        GOTO 300
175      V=F(41)
        XVM(IRM, INJ)=XRHO*(4.18879*RMX(IRM, INJ)**3)*XND
        RMX(IRM, INJ)=0.0
        GOTO 300
C
C      CALCULATE MASS VAPORIZED AND CURRENT DROPLET RADIUS
C
200      VNEW=(1.0-V)*4.18879*RFI**3
        XVM(IRM, INJ)=XRHO*((4.18879*RMX(IRM, INJ)**3)-VNEW)*XND
        RMX(IRM, INJ)=(VNEW/4.18879)**(1./3.)
300      TXVM=TXVM+XVM(IRM, INJ)
        IF (IMVFLG GT. 0) WRITE(6,310) INJ, IRM, GLEN,RFI,
*
*          RMX(IRM, INJ), V, XVM(IRM, INJ)
310      FORMAT(1X, 'FROM MVAP FOR INJ, IRM=', I4, ', ', I1, 9X, 'GLEN,RFI, ',
*
*          'RMX, %VAP, MVAP: ', /1X, 5E12.5)
500 CONTINUE
        RETURN
        END

```

```

SUBROUTINE VDIST
C
C SUBROUTINE TO DISTRIBUTE VAPORIZED MASS TO VARIOUS CELLS
C
C $INSERT COMMON
C   IF (ITRACE .EQ. 2) WRITE(6,1)
C   1 FORMAT(/1X, 'ENTER VDIST')
C
C LOOP THROUGH FOR EACH ELEMENT AND DROPSIZE
C
C DO 500 INJ=1,NEL
C   DO 500 IRM=1,3
C     IF (XVM(IRM, INJ) .LE. 0.0) GOTO 500
C     IF (IVDFLG .GT. 0) WRITE(6,10) IRM, INJ
10    FORMAT(/1X, 'BEGIN DISTRIBUTING DROP# ', I1, ' FOR INJ=', I3)
C
C     DTR=(3.1416/180.)
C
C RETRIEVE INJECTOR LOCATION
C
C     RI=ERPOS(INJ)
C     TI=ETPOS(INJ)
C
C RETRIEVE DROP RELATIVE POSITION AND CALCULATE NDROPS/SLICE
C
C     DPOS=DRP(IRM, INJ)
C     DANG=THETA
C     NDIST=NSLICE-1
C     IF (ERPOS(INJ) .GT. 0.0) DANG=45.0
C     IF (ERPOS(INJ) .GT. 0.0) NDIST=7
C     VPS=XVM(IRM, INJ)/(NDIST+1)
C
C ADD MASS PROPORTIONATELY TO GRID
C
C DO 200 I=0,NDIST
C   XP=RI*COS(TI*DTR)+DPOS*COS((DANG*I+DANG/2.0)*DTR)
C   YP=RI*SIN(TI*DTR)+DPOS*SIN((DANG*I+DANG/2.0)*DTR)
C   RP=SQRT(XP**2+YP**2)
C   TP=ATAN2(YP, XP)/DTR
C   IF (TP .LT. 0.0) TP=360.0+TP
C
C CORRESPONDING SLICE
C
C   T1=0.0
C   DO 100 J=1,NSLICE
C     T2=T1+THETA
C     IF (TP .GE. T1 .AND. TP .LT. T2) GOTO 110
C     T1=T2
100    CONTINUE
110    ISLICE=J
C     R1=0.0
C     DO 120 J=1,NSEG
C       R2=RGRID(J)
C       IF (RP .GE. R1 .AND. RP .LT. R2) GOTO 130
C       R1=R2
120    CONTINUE
130    IRAD=J
C     XGRID(ISLICE, IRAD)=XGRID(ISLICE, IRAD)+VPS
C     IF (IVDFLG .GT. 0) WRITE(6,140) VPS, ISLICE, IRAD
140    FORMAT(1X, E12.5, ' LB/S OF OX VAPOR ADDED TO SLICE, SEG=',

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```
      *                213)
200      CONTINUE
500 CONTINUE
      RETURN
      END
```

```

      SUBROUTINE VCALC
C
C      SUBROUTINE TO CALCULATE RADIAL AND AXIAL GAS VELOCITIES
C      AND DISTRIBUTES MASS ACROSS GRID TO ACHIEVE UNIFORM
C      PRESSURE DISTRIBUTION ACROSS CROSS-SECTION
C
C      $INSERT COMMON
C      IF (ITRACE .EQ. 2) WRITE(6,1)
C      1 FORMAT(/1X,'ENTER VCALC')
C      AT=3.1416*RW(NXP)**2
C
C      CALCULATE OVERALL VAPOR MR
C
C      OVMR=TXVM/THVM
C      CPC=PC
C
C      CALCULATE MEAN GAS PROPERTIES AND 1-D GAS VELOCITY, GXV(1)
C
C      IC=1
C      CALL DBLINT (CPC, OVMR, IC, CT)
C      IC=2
C      CALL DBLINT (CPC, OVMR, IC, GMW)
C      IC=4
C      CALL DBLINT (CPC, OVMR, IC, VSONIC)
C      IC=5
C      CALL DBLINT (CPC, OVMR, IC, GRHO)
C      IC=7
C      CALL DBLINT (CPC, OVMR, IC, GMU)
C      IC=6
C      CALL DBLINT (PC, OVMR, IC, CSTAR)
C      TIMR=TXF/TFF
C      CALL DBLINT (PC, TIMR, IC, PCSTAR)
C      ERE=CSTAR*(THVM+TXVM)/PCSTAR/(TXF+TFF)
C      GXV(1)=0.62/CR*VSONIC*ERE
C
C      CALCULATE RADIAL VELOCITIES IN EACH SLICE DUE TO MALDISTRIBUTION
C
C      IF (NSEG .EQ. 1) GOTO 501
C      DO 500 I=1,NSLICE
C          SMEAN=0.0
C          DO 200 J=1,NSEG
C              SMEAN=SMEAN+HGRID(I,J)+XGRID(I,J)
C              GRV(I,J)=0.0
C      200 CONTINUE
C          SMEAN=SMEAN/NSEG
C
C      REDISTRIBUTE MASS STARTING AT WALL
C
C      DO 300 J=NSEG,2,-1
C          DM=SMEAN-HGRID(I,J)-XGRID(I,J)
C          IF (DM .EQ. 0.0) GOTO 300
C          SGN=1.0
C          JF=J
C          JT=J-1
C          IF (DM .GT. 0.0) JF=J-1
C          IF (DM .GT. 0.0) JT=J
C          IF (DM .GT. 0.0) SGN=-1.0
C          AMASS=HGRID(I,JF)+XGRID(I,JF)
C          IF (DM .GT. AMASS) DM=AMASS
C          DF=(HGRID(I,JF)+XGRID(I,JF))/DM*SGN

```

```

      HL=HGRID(I,JF)/DF
      XL=XGRID(I,JF)/DF
      XGRID(I,JT)=XGRID(I,JT)-XL
      HGRID(I,JT)=HGRID(I,JT)-HL
      XGRID(I,JF)=XGRID(I,JF)+XL
      HGRID(I,JF)=HGRID(I,JF)+HL
C
C      SUM MOMENTUM IN/OUT OF CELL
C
      AC=6.2832*RGRID(J-1)*THETA/360.0*XSTEP
      GRV(I,JT)=GRV(I,JT)+DM/AC
      GRV(I,JF)=GRV(I,JF)+DM/AC
300    CONTINUE
C
C      CALCULATE RADIAL VELOCITY
C
      DO 400 J=1,NSEG
        IF (HGRID(I,J) .LE. 0.0 .AND. XGRID(I,J) .LE. 0.0) THEN
          GRV(I,J)=0.0
          GOTO 400
        ELSE IF (HGRID(I,J) .LE. 0.0) THEN
          VMR=20.0
          GOTO 350
        END IF
        VMR=XGRID(I,J)/HGRID(I,J)
350      CRHO=0.0
          IC=5
          CALL DBLINT (CPC, VMR, IC, CRHO)
          GRV(I,J)=GRV(I,J)/CRHO
400    CONTINUE
      IF (IPDFLG .GT. 0) WRITE(6,450) I, (GRV(I,J), J=1,NSEG)
450    FORMAT(1X, 'FOR SLICE ', I3, ' GRV(SEG)=', 10(F7.2, 3X))
500  CONTINUE
501  CONTINUE
C
C      DISTRIBUTE MASS CIRCUMFRENTIALLY
C
      IF (NSLICE .EQ. 1) GOTO 601
      DO 600 I=1,NSLICE
        IASLICE=I+1
        IF (IASLICE .GT. NSLICE) IASLICE=1
        DO 600 J=1,NSEG
          DM=(XGRID(I,J)+HGRID(I,J)+XGRID(IASLICE,J)+HGRID(IASLICE,J))
          *      /2.0-XGRID(I,J)-HGRID(I,J)
          IF (DM .EQ. 0.0) GOTO 600
          SGN=1.0
          IFR=I
          IT=IASLICE
          IF (DM .GT. 0.0) IFR=IASLICE
          IF (DM .GT. 0.0) IT=I
          IF (DM .GT. 0.0) SGN=-1.0
          AMASS=HGRID(IFR,J)+XGRID(IFR,J)
          IF (DM .GT. AMASS) DM=AMASS
          DF=(HGRID(IFR,J)+XGRID(IFR,J))/DM*SGN
          HL=HGRID(IFR,J)/DF
          XL=XGRID(IFR,J)/DF
          XGRID(IT,J)=XGRID(IT,J)-XL
          HGRID(IT,J)=HGRID(IT,J)-HL
          XGRID(IT,J)=XGRID(IT,J)+XL
          HGRID(IT,J)=HGRID(IT,J)+HL

```

```
600 CONTINUE
601 CONTINUE
    IF (IPDFLG .GT. 0) WRITE(6,700) CPC, GXV(1), ERE
700 FORMAT(1X, 'PC=', F7.2, 5X, '1-D VELOCITY=', F7.2, 5X, 'ERE=', F7.5)
RETURN
END
```

```

      SUBROUTINE VCONV
C
C      ROUTINE TO CALCULATE GAS VELOCITY IN A CONVERGING NOZZLE
C
      *INSERT COMMON
        IF (ITRACE .EQ. 2) WRITE(6,1)
        1 FORMAT(/IX, 'ENTER VCONV')
C
C      CALCULATE GAS MW AND STAGNATION DENSITY
C
        IC=2
        OVMR=TXVM/THVM
        CALL DBLINT (CPC, OVMR, IC, GMW)
        IC=3
        CALL DBLINT (CPC, OVMR, IC, GAMMA)
        GRHO=CPC*GMW/CT/10.73
C
C      CALCULATE AREA RATIO AND MACH NUMBER
C
        ARATIO=(RWALLX/RW(NXP))**2
        CALL MACH (ARATIO, ZMACH)
        IF (ZMACH .LE. 0.0) GOTO 9000
        IF (ZMACH .GT. 1.0) GOTO 9100
C
C      CALCULATE ISENTROPIC PROPERTIES CGT, GRHO
C
        TRPG=1.0+(GAMMA-1.0)/2.0*ZMACH**2
        CGT=CT/TRPG
        GRHO=GRHO/(TRPG**(1.0/(GAMMA-1.0)))
C
C      1-D GAS VELOCITY
C
        VAVG=ZMACH*SQRT(GAMMA/GMW*CGT*49712.69)
C
C      CALCULATE AXIAL AND RADIAL GAS VELOCITY FROM LOCAL VELOCITY MAGNITUDE
C
        RO=0.0
        DO 200 J=1,NSEG
            RJ=RO+(RGRID(J)-RO)/2.0
            RO=RGRID(J)
            GANGLE=RJ/RWALLX*ANGLEX
            VGAS=VAVG
            IF (GANGL .NE. 0.0) VGAS=VAVG*ANGLEX**2/2.0/
            * (ANGLEX*SIN(ANGLEX)+COS(ANGLEX)-1.0)
            GRV(1,J)=VGAS*SIN(GANGLE)
            GXV(J)=VGAS*COS(GANGLE)
        200 CONTINUE
        IF (IPDFLG .GT. 0) WRITE(6,210) ZMACH, VAVG, (GXV(I), I=1,NSEG)
        210 FORMAT(IX, 'MACH#=', F5.3, 3X, '1-D VELOCITY=', F7.2, /2X, 'AXIAL VEL: ',
            * 10(F7.2, 2X))
        IF (IPDFLG .GT. 0) WRITE(6,211) (GRV(1,K), K=1,NSEG)
        211 FORMAT(IX, 'RADIAL VEL: ', 10(F7.2, 2X))
        RETURN
C
C      ERROR CONDITIONS
C
        9000 WRITE(6,*) 'ERROR IN VCONV, NEGATIVE MACH NO FOR ARATIO=', ARATIO
            RWALLX=0.0
            RETURN
        9100 WRITE(6,*) 'ERROR IN VCONV, SUPERSONIC MACH NO FOR ARATIO='

```

```
RWALLX=0 0  
RETURN  
END
```

```

      SUBROUTINE MACH (AR, ZM)
C
C      SUBROUTINE TO CALCULATE MACH NUMBER FOR A GIVEN AREA RATIO
C      USES NEWTON-RAPHSON ITERATION
C
*INSERT COMMON
      IF (ITRACE .EQ. 2) WRITE(6,1)
1  FORMAT(/1X, 'ENTER MACH')
C
      ZM=0.01
      IF (AR .GT. 57.0) RETURN
      IF (AR .LT. 2.2) ZM=0.4
      ICNT=0
      F1=2.0/(1.0+GAMMA)
      F2=(GAMMA-1.0)/(GAMMA+1.0)
      F3=(GAMMA+1.0)/2.0/(GAMMA-1.0)
10  ICNT=ICNT+1
      Q=(F1+F2*ZM**2)
      FX=Q**F3/ZM-AR
      FXP=(2.0*F3*F2*Q** (F3-1.0))-(Q**F3/ZM**2)
      ZMN=ZM-FX/FXP
      IF (ABS(ZM-ZMN) .LT. 0.001*ZM) RETURN
      ZM=ZMN
      IF (ICNT .LT. 21) GOTO 10
C
C      ERROR CONDITION
C
      WRITE(6,*) 'ERROR IN MACH, ROUTINE NON-CONVERGANT'
      ZM=-1.0
      RETURN
      END

```

```

      SUBROUTINE DMOVE
C
C      SUBROUTINE TO MOVE DROPLETS DUE TO DRAG ACCELERATIONS
C
C      $INSERT COMMON
      IF (ITRACE .EQ. 2) WRITE(6,1)
1  FORMAT(/1X,'ENTER DMOVE')
      IF (IDMFLG .GT. 0) WRITE(6,2)
2  FORMAT(/1X)
C
C      LOOP THROUGH FOR ALL INJECTORS AND DROPSIZES
C
      DO 1000 INJ=1,NEL
      DO 1000 IRM=1,3
      IF (RMX(IRM,INJ) .LE. 0.0) GOTO 1000
C
      DTR=(3.1416/180.)
C
C      "EXPAND" DROPLET TO SUPERCRITICAL PROPORTIONS
C
      RM=RMX(IRM,INJ)*(XRHO/XRHOC)**(1./3.)
C
C      CALCULATE CURRENT DROPLET POSITION
C
      RI=ERPOS(INJ)
      TI=ETPOS(INJ)
      DPOS=DRP(IRM,INJ)
      XP=(RI+DPOS)*COS(TI*DTR)
      YP=(RI+DPOS)*SIN(TI*DTR)
      RP=SQRT(XP**2+YP**2)
      TP=ATAN2(YP,XP)/DTR
      IF (TP .LT. 0.0) TP=360.0+TP
C
C      CALCULATE CORRESPONDING GRID LOCATION
C
      T1=0.0
      DO 100 J=1,NSLICE
      T2=T1+THETA
      IF (TP .GE. T1 .AND. TP .LT. T2) GOTO 110
      T1=T2
100  CONTINUE
110  ISLICE=J
      JSLICE=J
      R1=0.0
      DO 120 J=1,NSEG
      R2=RGRID(J)
      IF (RP .GE. R1 .AND. RP .LT. R2) GOTO 130
      R1=R2
120  CONTINUE
130  IRAD=J
      IRAD1=IRAD
      IF (X .LE. X1) IRAD1=1
C
C      CALCULATE RELATIVE GAS VELOCITY, DRAG AND ACCELERATION
C
      IF (X .GT. X1) JSLICE=1
      DRVN=DRV(IRM,INJ)
      DXVN=DXV(IRM,INJ)
      GV=SQRT(GXV(IRAD1)**2+GRV(JSLICE,IRAD)**2)
      DV=SQRT(DXV(IRM,INJ)**2+DRV(IRM,INJ)**2)

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```

      XRV=GXV(IRAD1)-DXV(IRM, INJ)
      RRV=GRV(JSLICE, IRAD)-DRV(IRM, INJ)
      RGV=SQRT(XRV**2+RRV**2)
      IF (RGV .LE. 1.0) GOTO 180
      DZ=XSTEP*SQRT(1.0+((DRV(IRM, INJ)/DXV(IRM, INJ))**2)
      ICNT=0
175    CALL DRAG (RM, GV, DV, RGV, DZ, ACC)
C
C      MOVE DROPLET
C
      DXVN=GXV(IRAD1)-XRV*ACC
      DRVN=GRV(JSLICE, IRAD)-RRV*ACC
      DZN=XSTEP*SQRT(1.0+((DRV(IRM, INJ)+DRVN)/(DXV(IRM, INJ)+DXVN))
*
*          **2)
      IF (ABS(DZN-DZ) .LE. 0.01*DZ) GOTO 180
      DZ=DZN
      ICNT=ICNT+1
      IF (ICNT .GT. 10) GOTO 9000
      GOTO 175
180    RPNEW=RP+(DRV(IRM, INJ)+DRVN)/(DXV(IRM, INJ)+DXVN)*XSTEP
C
C      CHECK FOR IMPINGEMENT
C
      IF (RPNEW .GE. RWALLX) GOTO 500
C
C      NO IMPINGEMENT, UPDATE POSITION
C
      DRP(IRM, INJ)=RPNEW-RI
      DXV(IRM, INJ)=DXVN
      DRV(IRM, INJ)=DRVN
      IF (DXV(IRM, INJ) .LT. 0.0) GOTO 9100
      IF (DXV(IRM, INJ) .GT. GXV(IRAD1)) GOTO 9200
      IF (IDMFLG .GT. 0) WRITE(6,200) IRM, INJ, RP, RPNEW,
*
*          DXV(IRM, INJ), DRV(IRM, INJ)
200    FORMAT(1X, 'DROP #', I1, ' FROM INJ=', I3, ' MOVED RADIALLY ',
*
*          'FROM ', F7.3, ' TO ', F7.3, ' WITH NEW AXIAL AND ',
*
*          'RADIAL VEL=', F7.2, ' ', F7.2)
      GOTO 1000
C
C      IMPINGEMENT OCCURS, ADD OXIDIZER TO WALL VAPOR
C
500    PFACT=0.165*TND(INJ)*XRHO*4.1888
      IF (IRM .EQ. 2) PFACT=0.67*TND(INJ)*XRHO*4.1888
      XVAPN=RMX(IRM, INJ)**3*PFACT
      TXVM=TXVM+XVAPN
      IF (ERPOS(INJ) .GT. 0.0) GOTO 550
      XVAPN=XVAPN/NSLICE
      DO 525 IK=1, NSLICE
          XGRID(IK, NSEG)=XGRID(IK, NSEG)+XVAPN
525    CONTINUE
      GOTO 575
550    XGRID(ISLICE, NSEG)=XGRID(ISLICE, NSEG)+XVAPN
575    RMX(IRM, INJ)=0.0
      WRITE(6,600) X, TP, IRM, INJ, RMO(INJ), SG(INJ)
600    FORMAT(1X, 'IMPINGEMENT OCCURS AT X, THETA=', F7.3, ' ', F6.2,
*
*          /5X, 'DROP NUMBER ', I1, ' FROM INJ=', I4, ' RMEAN, SG=',
*
*          E11.4, ' ', F6.3)
1000 CONTINUE
      RETURN

```

```

C      ERROR CONDITION
C
9000 WRITE(6,*) 'DZ NOT CONVERGING FOR INJ,IRM=',INJ,' ',IRM,' AT X=',
*      X,' RUN STOPPED'
      X=-1.0
      RETURN
9100 WRITE(6,*) 'DROP ACCELERATED TO NEGATIVE VELOCITY AT X=',X,
*      ' INJ,IRM=',INJ,IRM,' ACC=',ACC
      X=-1.0
      RETURN
9200 WRITE(6,*) 'DROP ACCELERATED TO GREATER THAN GAS VELOCITY AT X=',X,
*      ' INJ,IRM=',INJ,IRM,' ACC=',ACC
      X=-1.0
      RETURN
      END

```

```

SUBROUTINE DRAG (RM, GVM, DVM, VELR, DZ, A)
C
C SUBROUTINE TO CALCULATE THE DRAG ON A DROPLET OF RADIUS RM IN
C A GAS FIELD WITH A RELATIVE VELOCITY VELR. THE RESULTANT
C ACCELERATION IS A. CALCULATIONS, IN PART, FROM TPP AND SDER.
C
*INSERT COMMON
  IF (ITRACE .EQ. 2) WRITE(6,1)
  1 FORMAT(/1X,'ENTER DRAG')
C
C CALCULATE FLOW REGIEME
C
  RE=(2.0*GRHO*VELR*RM)/GMU
  IF (RE .GE. 0.5) GOTO 100
C
C STOKES FLOW
C
  Q=4.5*GMU/XRHOC/RM**2*DZ
  A=EXP(-(Q+DVM)/GVM)*EXP(DVM/GVM)
  IF (IDFLG .GT. 0) WRITE(6,90) RM, VELR, RE, A
  90 FORMAT(1X,'RM, VELR, RE, A=',4E12.5)
  RETURN
C
C NEWTONIAN FLOW
C
  100 IF (RE .GT. 0.5 .AND. RE .LT. 70.0) CD=27.0*RE**(-0.84)
  IF (RE .GE. 70.0 .AND. RE .LT. 59200.0) CD=0.414*RE**(0.1433)
  IF (RE .GE. 59200.0) CD=2.0
C*
  Q=0.375*CD*GRHO/XRHOC/RM*DZ
C*
  A=EXP(Q-(GVM/VELR))*EXP(GVM/VELR)
  A=EXP(-6.0*CD*GRHO*VELR*DZ/XRHOC/RM/DVM)
  IF (IDFLG .GT. 0) WRITE(6,110) RM, VELR, RE, CD, A
  110 FORMAT(1X,'RM, VELR, RE, CD, A=',5E12.5)
  RETURN
END

```

```

SUBROUTINE MRPLOT

C
C SUBROUTINE TO CALCULATE VMR(R, THETA)
C IF IMRFLG=1, ONLY WALL MR IS PRINTED; IMRFLG=1, CROSS-SECTION MR
C IS PRINTED; IMRFLG=2 PRINTS MASS DISTRIBUTION IN PL HISTORY;
C IMRFLG=3 PRINTS GRID CONTENTS IN PL HISTORY
C
$INSERT COMMON
  IF (ITRACE .EQ. 2) WRITE(6,1)
1  FORMAT(/1X, 'ENTER MRPLOT')
  IROW=NSEG
  IF (IMRFLG .GT. 0) IROW=1

C
C WRITE HEADER
C
  XV=GXV(1)
  IF (X .GT. X1) XV=GXV(NSEG)
  WRITE (7,10) X, XV, (RGRID(I), I=IROW,NSEG)
10 FORMAT(/10X, 'AT AXIAL POSITION=', F7.4, ' FT. FROM INJECTOR FACE',
*       ' WALL AXIAL VELOCITY=', F7.2, ' FT/SEC', /5X, 'RSEG (FT)', 7X,
*       10(F6.3, 4X), /10X, 10(F6.3, 4X), /10X, 10(F6.3, 4X))

C
C CALCULATE VMR IN SLICE
C
  DO 200 I=1, NSLICE
    DO 150 J=IROW, NSEG
      GRV(I, J)=-1.0
      IF (XGRID(I, J) .GT. 0.0) GRV(I, J)=99.99
      IF (HGRID(I, J) .GT. 0.0) GRV(I, J)=XGRID(I, J)/HGRID(I, J)
150   CONTINUE
    WRITE(7,160) (I*THETA), (GRV(I, J), J=IROW, NSEG)
160   FORMAT(5X, 'THETA=', F6.2, 4X, 10(F6.3, 4X), /10X, 10(F6.3, 4X), /10X,
*         5(F6.3, 4X))
    IF (IMRFLG .GT. 1) WRITE(6,170) I, ((HGRID(I, J)+XGRID(I, J)),
*                                     J=1, NSEG)
170   FORMAT(/1X, 'FOR SLICE=', I3, /1X, 'MASS: ', 10(E11.5, 2X))
    IF (IMRFLG .GT. 2) WRITE(6,180) (HGRID(I, J), J=1, NSEG)
    IF (IMRFLG .GT. 2) WRITE(6,190) (XGRID(I, J), J=1, NSEG)
180   FORMAT(1X, 'FUEL: ', 10(E11.5, 2X))
190   FORMAT(1X, 'OX: ', 10(E11.5, 2X))
200  CONTINUE
    RETURN
  END

```

```

      SUBROUTINE CERE
C
C      SUBROUTINE TO CALCULATE ENERGY RELEASE EFFICIENCY
C
      *INSERT COMMON
      IF (ITRACE .EQ. 2) WRITE(6,1)
      1 FORMAT(/1X, 'ENTER CERE')
C
      CMR=TXF/TFF
C
C      SUM C* FOR EACH CELL INCLUDING MASS WEIGHTING FACTOR
C
      CSUM=0.0
      IC=6
      CI=0.0
      DO 100 I=1, NSLICE
      DO 100 J=1, NSEG
      VMR=XGRID(I, J)/HGRID(I, J)
      CALL DBLINT (PC, VMR, IC, CI)
      WI=(XGRID(I, J)+HGRID(I, J))/(TXVM+THVM)
      CSUM=CSUM+WI*CI
100 CONTINUE
C
C      CORRECT C* FOR MASS DEFECT
C
      CSUM=CSUM*(TXVM+THVM)/(TFF+TXF)
C
C      CALCULATE C* COMBUSTION EFFICIENCY
C
      CSTAR=0.0
      CALL DBLINT (PC, CMR, IC, CSTAR)
      ERE=CSUM/CSTAR*100.0
      EVAP=MIN(100.0, (TXVM/TXF*100.0))
      EMIX=ERE/EVAP*100.0
      WRITE(6,200) THVM, TFF, TXVM, TXF, EVAP, EMIX, ERE
200 FORMAT(/5X, F8.3, ' LB/S OF FUEL VAPORIZED OUT OF ', F8.3,
*        ' LB/S TOTAL', /5X, F8.3, ' LB/S OF OX VAPORIZED OUT OF ', F8.3,
*        ' LB/S TOTAL', /5X, 'OX VAPORIZATION EFFICIENCY=', F7.3, '%',
*        /5X, 'MIXING EFFICIENCY=', F7.3, '%', /5X,
*        'ENERGY RELEASE EFFICIENCY=', F7.3, '%', //)
      RETURN
      END

```

SUBROUTINE DBLINT (PR, XMR, IVFLG, Q)

SUBROUTINE TO PERFORM DOUBLE INTERPOLATION TO CALCULATE T,  
MW, GAMMA, VSONIC, RHO C\* AND MU AS A FUNCTION OF PC AND MR

INPUTS:

PR=CHAMBER PRESSURE, PSIA

XMR=MIXTURE RATIO

IVFLG=VALUE FLAG: 1=TGAS, DEGREES R

2=MW

3=SPECIFIC HEAT RATIO

4=VSONIC, FT/SEC

5=RHO, LB/CU. FT

6=CSTAR, FT/SEC

7=MU, LB/FT-SEC

OUTPUT:

Q=VALUE OF IVFLG

IMPLICIT REAL\*8 (A-H,O-Z)

COMMON /IPROPS/ FTJ, FRHO, FMU, FTC, FHV, FMW, XRHOC,  
\* XST, XTJ, XRHQ, XMU, XTC, XHV, XMW

LOCAL COMMON ONLY

COMMON /DBLDAT/ XM(13), T(2,13), WM(2,13), CS(2,13), UM(2,13),  
\* GM(2,13)

DATA XM / 0., 1., 2., 3., 4., 5., 6., 7., 8., 9., 10., 15., 20. /  
DATA ((T(I,J),J=1,13),I=1,2) / 37.0, 1759., 3232., 4404., 5237.,

\* 5756., 6093., 6159., 6177., 6141., 6078., 5650., 5209.,

\* 37.0, 1759.,

\* 3236., 4430., 5375., 6064., 6504., 6717., 6756., 6695.,

\* 6592., 5974., 5401. /

DATA ((WM(I,J),J=1,13),I=1,2) / 2.016, 4.032, 6.048, 8.048, 9.957,

\* 11.703, 13.263, 14.135, 15.835, 16.889, 17.825, 21.291,

\* 23.501, 2.016, 4.032, 6.048, 8.059, 10.029, 11.897, 13.603,

\* 15.094, 16.355, 17.421, 18.341, 21.653, 23.722 /

DATA ((CS(I,J),J=1,13),I=1,2) / 1287., 6864., 7751., 7959., 7922.,

\* 7742., 7493., 7220., 6956., 6715., 6500., 5715., 5194.,

\* 1287., 6864., 7751., 7959., 7950., 7829., 7637., 7396.,

\* 7133., 6878., 6647., 5800., 5237. /

DATA ((UM(I,J),J=1,13),I=1,2) / 0.2875, 0.5139, 0.9856, 1.363,

\* 1.646, 1.835, 1.951, 2.014, 2.045, 2.057, 2.058, 2.006,

\* 1.920, 1.305, 0.5319, 0.9857, 1.369, 1.679, 1.908, 2.062,

\* 2.148, 2.183, 2.190, 2.182, 2.088, 1.971 /

DATA ((GM(I,J),J=1,13),I=1,2) / 1.400, 1.359, 1.283, 1.227, 1.178,

\* 1.148, 1.133, 1.126, 1.125, 1.125, 1.126, 1.135, 1.149, 1.400,

\* 1.359, 1.284, 1.237, 1.199, 1.168, 1.148, 1.138, 1.136, 1.136,

\* 1.138, 1.149, 1.165 /

R=10.73

FIND MR BOUNDS

IF(XMR .GT. XM(1)) GOTO 20

I=2

GOTO 60

20 IF(XMR .LT. XM(13)) GOTO 30

I=13

GOTO 60

30 DO 50 I=2, 13

```

        IF (XM(I) LT. XMR) GOTO 50
        GOTO 60
50 CONTINUE
60 CONTINUE
    IF (IVFLG EQ. 2) GOTO 200
    IF (IVFLG EQ. 3) GOTO 300
    IF (IVFLG EQ. 6) GOTO 600
    IF (IVFLG EQ. 7) GOTO 700
C
C    CALCULATE T
C
    T(1,1)=FTJ+460.0
    T(2,1)=FTJ+460.0
    TP1=T(1,I-1)+(XMR-XM(I-1))/(XM(I)-XM(I-1))*(T(1,I)-T(1,(I-1)))
    TP2=T(2,I-1)+(XMR-XM(I-1))/(XM(I)-XM(I-1))*(T(2,I)-T(2,(I-1)))
    Q=TP1+(PR-300.)/2700. *(TP2-TP1)
    IF (IVFLG EQ. 1) RETURN
C
C    CALCULATE MW
C
200 CMW1=WM(1,I-1)+(XMR-XM(I-1))/(XM(I)-XM(I-1))*(WM(1,I)-WM(1,(I-1)))
    CMW2=WM(2,I-1)+(XMR-XM(I-1))/(XM(I)-XM(I-1))*(WM(2,I)-WM(2,(I-1)))
    CMW=CMW1+(PR-300.)/2700. *(CMW2-CMW1)
    IF (IVFLG EQ. 4) GOTO 300
    IF (IVFLG EQ. 5) GOTO 500
    Q=CMW
    RETURN
C
C    CALCULATE GAMMA
C
300 GM1=GM(1,I-1)+(XMR-XM(I-1))/(XM(I)-XM(I-1))*(GM(1,I)-GM(1,(I-1)))
    GM2=GM(2,I-1)+(XMR-XM(I-1))/(XM(I)-XM(I-1))*(GM(2,I)-GM(2,(I-1)))
    GAMMA=GM1+(PR-300.)/2700. *(GM2-GM1)
    IF (IVFLG NE. 3) GOTO 400
    Q=GAMMA
    RETURN
C
C    CALCULATE VSONIC
C
400 IF (IVFLG NE. 4) GOTO 500
    Q=SQRT(GAMMA*R*Q/CMW*4633.056)
    RETURN
C
C    CALCULATE DENSITY
C
500 Q=PR*CMW/R/Q
    RETURN
C
C    CALCULATE CSTAR
C
600 CP1=CS(1,I-1)+(XMR-XM(I-1))/(XM(I)-XM(I-1))*(CS(1,I)-CS(1,(I-1)))
    CP2=CS(2,I-1)+(XMR-XM(I-1))/(XM(I)-XM(I-1))*(CS(2,I)-CS(2,(I-1)))
    Q=CP1+(PR-300.)/2700. *(CP2-CP1)
    RETURN
C
C    CALCULATE MU
C
700 UP1=UM(1,I-1)+(XMR-XM(I-1))/(XM(I)-XM(I-1))*(UM(1,I)-UM(1,(I-1)))
    UP2=UM(2,I-1)+(XMR-XM(I-1))/(XM(I)-XM(I-1))*(UM(2,I)-UM(2,(I-1)))
    Q=(UP1+(PR-300.)/2700. *(UP2-UP1))*3.2174E-5

```

RETURN  
END

C  
C  
C  
COMMON INSERT FOR PERFORMANCE/LIFE COMBUSTION MODEL, PLC

IMPLICIT REAL\*8 (A-H,O-Z)

COMMON /OPCOND/ PC, TFF, TXF, HGMRO

COMMON /WALL/ NXP, XW(NXP), RW(NXP), WANGLE(NXP), CR, X1, X2,  
\* X, RWALLX, ANGLEX

COMMON /FACE/ NEL, NROWS, NELR(NROWS), RROW(NROWS),

\* ERPOS(NEL), ETPOS(NEL), CFLOW, CFMR

COMMON /INJECT/ ITYPE, RXP, RFP, TPOST, RECESS, RMS, XDL, CSA,

\* PFI, FCD, XCD, FMF(NEL), XMF(NEL), HGMR(NEL)

COMMON /GRID/ NSLICE, NSEG, THETA, RGRID(NSEG), XGRID(NSLICE,NSEG),

\* HGRID(NSLICE,NSEG)

COMMON /IPROPS/ FTJ, FRHO, FMU, FTC, FHV, FMW, XRHOC,

\* XST, XTJ, XRHQ, XMU, XTC, XHV, XMW

COMMON /VAPOR/ TXVM, THVM, XVM(3,NEL)

COMMON /GAS/ CPC, CT, GAMMA, GRHO, GMU, GMW, GXV(NSEG), GRV(NSLICE,NSEG)

COMMON /DROP/ RMO(NEL), SG(NEL), VFACT(NEL), AL(NEL), TND(NEL),

\* RMX(3,NEL), DRP(3,NEL), DXV(3,NEL), DRV(3,NEL)

COMMON /FILL/ COM(3,NEL)

COMMON /MSG/ IAFLG, IVDFLG, IMVFLG, IPDFLG, IDMFLG, IMRFLG, IDFLG,

\* IRFFLG, ITRACE

COMMON /STEP/ ASTEP, XSTEP, IEFLG

```

/* CPL PROGRAM TO CREATE A SEG FOR PERFORMANCE/LIFE COMBUSTION MODEL
/* WITH COMMONS DIMENSIONED TO THE PROBLEM
/*
/* ***** NSUBS=NUMBER OF PROGRAMS WITH INSERT COMMONS *****
/*
/* RETRIEVE BASE FILES
/*
&IF [EXISTS PLC.F77 -FILE -BRIEF] &THEN DELETE PLC.F77
&IF [EXISTS COMMON -FILE -BRIEF] &THEN DELETE COMMON
&IF [EXISTS FMF.EQ -FILE -BRIEF] &THEN DELETE FMF.EQ
&IF [EXISTS XMF.EQ -FILE -BRIEF] &THEN DELETE XMF.EQ
COPY <AMFD27>E23846>PL.UFD>PLC.F77 PLC.F77
COPY <AMFD27>E23846>PL.UFD>COMMON COMMON
&S UNIT1 := [OPEN_FILE FMF.EQ -MODE W THERE]
&S UNIT2 := [OPEN_FILE XMF.EQ -MODE W THERE]
&S END := $END
/*
/* DEFAULT SIZES AND CHECK FOR NEW VARIABLES
/*
&S TYPE := 1
&S NEL := 1
&S NROWS := 1
&S NXP := 2
&S NSLICE := 12
/*
/* FIND NEW VALUES IF THEY EXIST
/*
&S UNIT := [OPEN_FILE PL.INPUT -MODE R THERE]
&IF %THERE% ^= 0 &THEN &STOP &MESSAGE PL.INPUT NOT FOUND
/*
/* LOOK FOR END OF NAMELIST OUTPUT
/*
&DO I := 1 &TO 10
  &S LINE := [READ_FILE %UNIT% OK]
  &IF %LINE% = %END% &THEN &GOTO OUT
&END
&STOP &MESSAGE NO END FOR NAMELIST OUTPUT
/*
/* LOOK FOR NXP AFTER PROPERTIES
/*
&LABEL OUT
&S LINE := [READ_FILE %UNIT% OK]
&S LINE := [READ_FILE %UNIT% OK]
&S LINE := [READ_FILE %UNIT% OK]
&S LINE := [READ_FILE %UNIT% OK]
&S LINE := [READ_FILE %UNIT% OK]
&S LINE := [READ_FILE %UNIT% OK]
&S NXP := [READ_FILE %UNIT% OK]
/*
/* LOOK FOR NEL AND NROWS AFTER WALL DATA
/*
&DO I := 1 &TO %NXP%
  &S LINE := [READ_FILE %UNIT% OK]
&END
&S LINE := [READ_FILE %UNIT% OK]
&S LINE := [READ_FILE %UNIT% OK]
&S TYPE := [TRIM [SUBSTR %LINE% 1 4] -BOTH]
&S NEL := [TRIM [SUBSTR %LINE% 5 4] -BOTH]
&S NROWS := [TRIM [SUBSTR %LINE% 9 4] -BOTH]

```

```

&S NSEG := %NROWS% + 1
/*
/* LOOK FOR NAMELIST CONT AFTER ELEMENTS AND INJECTOR CONFIGURATION
/* SETUP FMF.EQ AND XMF.EQ IF TYPE 2 INPUT
/*
&IF %TYPE% = 2 &THEN &GOTO TYPE2
&DO I := 1 &TO %NEL%
  &S LINE := [READ_FILE %UNIT% OK]
&END
&GOTO END_2
&LABEL TYPE2
&S IFS := '      IF (IROW.EQ. '
&S IFS2 := ' ) FMF(INJ) = '
&S IFS3 := ' ) XMF(INJ) = '
&S ICN := '      * '
&DO I := 1 &TO %NROWS%
  &S LINE := [READ_FILE %UNIT% OK]
  &S FMF := [READ_FILE %UNIT% OK]
  &IF %OK% ^= 0 &THEN &STOP &MESSAGE ERROR READING FMF.EQ FOR ROW %I%
  &S FMFC := %IFS%%I%%IFS2%
  &S FMF := %ICN%%FMF%
  &S J := [WRITE_FILE %UNIT1% %FMFC%]
  &S J := [WRITE_FILE %UNIT1% %FMF%]
  &S XMF := [READ_FILE %UNIT% OK]
  &IF %OK% ^= 0 &THEN &STOP &MESSAGE ERROR READING XMF.EQ FOR ROW %I%
  &S XMFC := %IFS%%I%%IFS3%
  &S XMF := %ICN%%XMF%
  &S J := [WRITE_FILE %UNIT1% %XMFC%]
  &S J := [WRITE_FILE %UNIT2% %XMF%]
&END
&LABEL END_2
&DO &UNTIL %LINE% = $CONT
  &S LINE := [READ_FILE %UNIT% OK]
&END
/*
/* READ ALL LINES AND PARSE UNTIL $END IS FOUND
/*
&LABEL NEXTLINE
&S LINE := [READ_FILE %UNIT% OK]
&IF %LINE% = %END% &THEN &GOTO NOMORE
&S LINE := [TRIM %LINE% -RIGHT ]
&DO &UNTIL [LENGTH %LINE%] = 0
  &S LINE := [TRIM %LINE% -LEFT ]
  &S NAME := [BEFORE %LINE% =]
  &S LINE := [AFTER %LINE% =]
  &S %NAME% := [BEFORE %LINE% ', ' ]
  &S LINE := [AFTER %LINE% ', ' ]
&END
&GOTO NEXTLINE
/*
/* CHANGE DIMENSIONS IN COMMON
/*
&LABEL NOMORE
&DATA SED COMMON
C/(NXP)/( %NXP% )/50 G
TOP
C/(NEL)/( %NEL% )/50 G
TOP
C/(NROWS)/( %NROWS% )/50 G
TOP

```

```
C/(NSLICE/(%NSLICE%/50 G
TOP
C/NSEG)/%NSEG%)/50 G
FILE
&END
CLOSE FMF.EQ
CLOSE XMF.EQ
CLOSE PL.INPUT
/*
/* COMPILE, SEG AND RUN
/*
F77 PLC -DD1 -SILENT 1 -OPT 0
&DATA SEG -LOAD
  LO PLC
  LI
  MAP 3
  SA
  QU
&END
DELETE PLC.BIN
&RETURN
```

**PART G**

**ADVANCED OXYGEN-HYDROGEN ROCKET  
ENGINE STUDY CHAMBER GEOMETRY DEFINITION**

AEROJET LIQUID ROCKET COMPANY  
INTER-OFFICE MEMORANDUM

9 January 1980  
RAH:sm:9751:0389  
CDN:9751:4399

TO: C. J. O'Brien

FROM: R. A. Hewitt

SUBJECT: Advanced Oxygen-Hydrocarbon Rocket Engine Study  
Chamber Geometry Definition

COPIES TO: K. Christiansen, D. Culver, O. D. Goodman, D. Kors,  
D. Lemke, J. Mellish, H. Mueggenburg, J. L. Pieper,  
J. Salmon, R. Schwantes, C. E. Taylor, 9751 Personnel,  
9751 File

ENCLOSURE: (1) Typical Rocket Engine Parameters  
(2) Chamber Pressure vs Thrust  
(3) Liquid/Liquid Engine Contraction Ratio vs Thrust  
(4) Liquid/Liquid Engine Chamber Length vs Thrust  
(5) Liquid/Hot-Gas Engine Chamber Length vs Thrust

REFERENCE: (a) C. J. O'Brien, "Advanced Oxygen-Hydrocarbon Rocket  
Engine Study", Program Plan 33542 PP, 29 Oct. 1979  
(b) ALRC Rocket Design Presentation by J. I. Ito  
(c) A. J. Pavli, NASA-Cleveland, Ohio, "Design of Injectors  
for H/C".  
(d) J. A. Mellish, "Advanced Engine Study for Mixed Mode  
OTV's", NASA CR 159491, Dec. 1978  
(e) Empirical Design Curves in C. J. O'Brien Possession  
(f) Rocketdyne Monthly Reports No. 1 and 2 in  
R. J. LaBotz Possession  
(g) R. A. Hewitt to J. A. Mellish, "COTV Geometry",  
28 September 1979, ALRC Memorandum 9751:0348

APPENDIX: (A) Universal Geometric Guidelines

INTRODUCTION

The "Advanced Oxygen-Hydrocarbon Rocket Engine Study" parametric analysis requires chamber geometry ( $D_t$ ,  $L'$ , CR) guidelines to provide

9 January 1980

## Introduction (cont.)

reasonable and typical design values. This preliminary study (See Ref. a) is directed at a thrust range of from 200,000 to 1,500,000 lbf with emphasis on the 600,000 to 1,000,000 lbf level. The chamber pressure range is from 1000 to 5000 psia. The propellant combinations are LOX/RP-1 and LOX/CH<sub>4</sub>, with LH<sub>2</sub> considered only as an additional coolant possibility. The mixture ratio range to be studied is 2.0 to 3.5 for LOX/RP-1, and 3.0 to 4.5 for the LOX/CH<sub>4</sub> propellant combination. The nozzle exit area ratio will range from about 15:1 to 100:1 as determined by attachment limit as a function of chamber pressure and optimum trajectory trade-offs.

RECOMMENDED GEOMETRY GUIDELINES

Based on the information presented in the body of this memorandum the following rocket geometry relationships are recommended for the "Advanced LOX/HC Engine Study":

Injection State	Liquid/Liquid	Liquid/Gas
Contraction Ratio CR	$\log_{10} CR = -.0715 \log_{10} F + 0.689$	3.00
Chamber Length L' (in)	$\log_{10} L' = .23 \log_{10} (F/P_c) + .85$	$\log_{10} L' = .23 \log_{10} (F/P_c) + .621$

The above contraction ratio relationship will yield a liquid/liquid value of about 1.85(@ 800,000 lbf) which can be assumed constant for purposes of this study. The estimated chamber lengths will be as follows:

## Recommended Geometry Guidelines (cont.)

<u>Injection State</u>	<u>Liquid/Liquid</u>		<u>Liquid/Gas</u>	
	<u>1000</u>	<u>5000</u>	<u>1000</u>	<u>5000</u>
<u>Thrust, lbf/Pressure, psia</u>				
200,000	23.9"	16.5"	14.1"	9.8 "
600,000	30.8	21.3	18.2	12.6
1,000,000	34.7	23.9	20.5	14.1
1,500,000	38.1	26.3	22.5	15.5

The estimation of the chamber throat diameter depends on the details of the engine design being considered. However, a diameter estimate for the propellant combinations used here are shown in the appendix. The following information is of more general interest and referred to as "Universal Guidelines" since so little data may exist to substantiate design trends in any given narrow spectrum of thrust or chamber pressure.

The designation of what propellants and engine cycle qualify for the designation liquid/liquid or liquid/gas are estimated as follows:

	<u>Liquid/Liquid</u>	<u>Average*</u>	<u>Liquid/Gas</u>
Lox/RP-1 Ambient	X		
LOX/RP-1 Regenerative	X		
LOX/RP-1 Staged Combustion			X
LOX/CH <sub>4</sub> Ambient		X	
LOX/CH <sub>4</sub> Regenerative		X	
LOX/CH <sub>4</sub> Staged Combustion			X

\*Average: Arithmetic average of the liquid/liquid and liquid/gas values.

9 January 1980

UNIVERSAL GEOMETRIC PREMISES

The "Advanced Oxygen-Hydrocarbon Rocket Engine Study" chamber geometry definition relies heavily on two premises: (1) that the average trends of existing rocket engine geometries as a function of chamber pressure and thrust will continue to be valid for future designs, and (2) that a given existing rocket geometry could theoretically have its chamber pressure and thrust increased and decreased over about a factor of 3 without significantly altering its performance, stability, or compatibility characteristics. Design information used to substantiate the assumptions made in this memorandum are largely contained in Enclosure (1) and reference (b) through (f).

EXISTING ROCKET GEOMETRIES

A table of typical rocket engine parameters are shown in Enclosure (1). The rocket geometric characteristics are defined by three dimensions; (1) throat diameter, (2) chamber diameter, and (3) chamber length from injector to throat. The non-dimensional ratios such as contraction ratio and chamber length to diameter ratios are shown for convenience. The rocket engine operational characteristics are defined by: (1) propellants, (2) mixture ratio, (3) thrust, and (4) chamber pressure.

EXTRAPOLATION OF THRUST AND CHAMBER PRESSURE

It is a premise of this study that the chamber pressure and thrust of any existing rocket engine can be theoretically increased and decreased within one order of magnitude peak-to-peak without significantly altering its value for indicating trends in future rocket engine designs. An example of this would be to assume a Titan I, second stage engine rated at 80,000 lbf thrust

## Extrapolation of Thrust and Chamber Pressure (cont.)

at 682 psia can be uprated to 250,000 lbf at 2150 psia or downrated to 25,000 lbf at 215 psia. The only geometric alteration required would be that the injector orifice size be increased from an assumed 0.100 in. dia. to 0.133 or decreased to 0.0750 in. This orifice size change is relatively insignificant due to the fact that the injection pressure drop is assumed to increase and decrease in order to maintain a nearly constant combustion time lag and constant chug stability margin by slightly varying the liquid injection pressure drop to chamber pressure ratio of about 0.15 to 0.20.

THRUST AND CHAMBER PRESSURE TRENDS

A comparison of existing rocket engine chamber pressures over three-orders-of-magnitude is made with thrust over eight orders-of-magnitude in Enclosure (2). The trend indicated is that logarithm of the rocket engine chamber pressure tends to be proportional to the logarithm of the thrust. There are no engines with very high chamber pressures used to obtain a very low thrust or vice-versa. Although low chamber pressure-large thrust studies (e.g., Big Dumb Booster & APS) have been made, no production engines have resulted. The lowest chamber pressure and highest thrust engine shown is the pressure fed Apollo SPS engine. The highest chamber pressure and lowest thrust engines include the 0.5 lbf at 125 psia, and the recent high pressure LOX/HC engines in the 4500 to 55000 lbf thrust range. The logic for the fact that the LOX/HC engines were designed for higher pressures than the trend line indicated for their thrust, lies in the fact that many are "subscale" engines. Or, in the case of the 0.5 lbf engine, it was designed to "blowdown" with tank pressure thereby lowering its thrust to near the "design diagonal" trend line. Note the "design diagonal" line bandwidth is defined by empirical data.

CHAMBER CONTRACTION RATIO TRENDS

The contraction ratio for present day liquid/liquid rocket engines is shown in Enclosure (3) over the eight orders-of-magnitude of thrust. The trend indicated narrows with increasing thrust. A weakly increasing contraction ratio with increasing thrust trend is indicated although a constant value of as low as 1.8 is also indicated. In either case in the thrust range of from 600,000 to 1,500,000 lbf a contraction ratio of less than 2.0 is indicated for a liquid/liquid LOX/HC engine. The use of engine cycles that pass hot-gas through the injector require a greater injector diameter to allow the lower density gases through with a minimum of pressure drop. This trend results in contraction ratio's of about 3.0, which tends to force liquid/hot-gas chamber pressures to higher values in order to bring chamber diameters down to comparable liquid-liquid engines operating at lower chamber pressures.

CHAMBER LENGTH TRENDS

The liquid-liquid engine chamber length from injector-to-throat is shown in Enclosure (4) for eight orders-of-magnitude of thrust. The chamber lengths shown are limited to their reasonable range of applicability as defined in Enclosure (2). That is, low pressure applies to low thrust only, and high pressure applies to high thrust only. This results in a reasonable "diagonal design" line extending from the longest, low thrust, engine and the shortest, high thrust engine, as shown below:

<u>F</u> <u>(lbf)</u>	<u>Pc</u> <u>(psia)</u>	<u>Chamber Length L' (in.)</u>	
		<u>Liquid/Gas</u>	<u>Liquid/Liquid</u>
1.0	30	2.0	3.
1.K	250	6.0	10.
1.M	1500	18.0	33.0

## Chamber Length Trends (cont.)

Note that the thrust "band-width" of the "design diagonal" is at least an order-of-magnitude wide, and results in about a  $\pm 40\%$  chamber length "band-width". For example if the 33 in. long 1 MLBF liquid/liquid engine shown above is raised in pressure to about 4000 psia, the estimated chamber length drops to about 25 in.

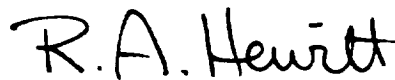
The chamber length of a liquid/hot-gas cycle engine is shown in Enclosure (5). Note that the net effect of the hot-gas cycle is to shorten the chamber relative to the liquid/liquid engines as is shown by comparing the chamber length equations:

LIQUID/LIQUID:

$$\text{LOG}_{10} L' = .23 \text{ LOG}_{10} (F/P_c) + .850$$

LIQUID/HOT-GAS:

$$\text{LOG}_{10} L' = .23 \text{ LOG}_{10} (F/P_c) + .621$$



R. A. Hewitt  
Thermodynamic Analysis  
Rocket Design Analysis

Approved by:



J. I. Ito  
Thermodynamic Analysis  
Rocket Design Analysis



J. L. Pieper, Manager  
Thermodynamic Analysis  
Rocket Design Analysis

## TYPICAL ROCKET ENGINE PARAMETERS

Symbol	Chamber Cooling Method	Mixture Ratio (O/F)	Chamber Length (in)	Throat Diameter (in)	Injector Diameter (in)	Contract. ratio (N.D.)	Number Primary Elements (N.D.)	Propellants (O/F)	Thrust (lbf)	Chamber Pressure (psia)	Chamber L/D
Titan II 1st Stage	Regen.	1.93	24	15.2	21.8	2.0	516	N <sub>2</sub> O <sub>4</sub> /A-50	215,000	785	1.1
Titan III 1st Stage	Regen.	1.93	24	15.3	21.65	2.0	504	N <sub>2</sub> O <sub>4</sub> /A-50	220,000	807	1.1
Titan II 2nd Stage	Regen.	1.80	17	9.1	14.2	2.4	818	N <sub>2</sub> O <sub>4</sub> /A-50	100,000	827	1.2
Titan III 2nd Stage	Regen.	1.80	17	9.2	14.2	2.4	500	N <sub>2</sub> O <sub>4</sub> /A-50	100,000	830	1.2
ITIP	Abla.	2.0	18	7.5	11.9	2.5	600	N <sub>2</sub> O <sub>4</sub> /A-50	8,000	105	1.5
Transtage	Abla.	2.0	18	7.5	11.9	2.5	336	N <sub>2</sub> O <sub>4</sub> /A-50	8,000	105	1.5
N-II, Delta FJ	Abla.	1.9	18	7.5	11.9	2.5	336	N <sub>2</sub> O <sub>4</sub> /A-50	9,850	125	1.5
LCAE 4K	Film Cooled	1.65	7	3.3	5.84	3.1	450	N <sub>2</sub> O <sub>4</sub> /MMH	4,000	260	1.2
LCAE 2.5K	Film Cooled	1.65	7	3.3	5.84	3.1	450	N <sub>2</sub> O <sub>4</sub> /MMH	2,500	165	1.2
OMS	Regen.	1.65	16	5.9	8.11	1.9	272	N <sub>2</sub> O <sub>4</sub> /MMH	6,000	125	2.0
Apollo/I.O.S.	Abla.	1.6	24	12.4	17	2.5	575/900	N <sub>2</sub> O <sub>4</sub> /A-50	20,000	97	1.2
M-1 Coax	Regen.	5.5	29	32	40	1.8	3248	LOX/LH <sub>2</sub>	1,500,000	1000	0.8
F-1	Regen.	2.35	48	35	39.2	1.3	702	LOX/RP-1	1,522,000	1128	1.2
ITA	Regen./F.C. 5.		7	1.92	3.5	3.3	72	O <sub>2</sub> /H <sub>2</sub>	1,500	340	2.0
ETR LOL	5.		5		2.2		48	LOX/LH <sub>2</sub>	1,200	500	2.2
ETR Coax	5.		5		2.2		36	LOX/LH <sub>2</sub>	1,200	500	2.2
100 lbf	Film	1.6	4	.69	0.94	1.86	33	N <sub>2</sub> O <sub>4</sub> /MMH	100	150	4.3
870	Film	1.6	3.9	2.04	3.0		270	N <sub>2</sub> O <sub>4</sub> /MMH	870	150	1.3
LM-A	Abla.	1.6	14	4.5	7.8	3.0	177	N <sub>2</sub> O <sub>4</sub> /A-50	3,500	120	1.8
IFAR/DELTAV	Abla.	1.1	9		4.		124/32	N <sub>2</sub> O <sub>4</sub> /N <sub>2</sub> H <sub>4</sub>	2,800	300	2.3
Fluorine T/S Fine/Coarse	Abla.	1.9	19		9.45		344/69	LF <sub>2</sub> /N <sub>2</sub> H <sub>4</sub>	7,000	100	2.0
Scaleable Fine/Coarse	Film	1.6	5		1.5		108/39	N <sub>2</sub> O <sub>4</sub> /MMH	200-300	75-130	3.3
NASA (Pavli)	Uncooled	2.8	8.5	2.6	5.32	4.2	97	LOX/CH <sub>4</sub>	4,500	600	1.6
NASA (Pavli)	Uncooled	2.8	22	2.6	5.32	4.2	37	LOX/RP-1	4,500	600	4.2
APS	Regen/F.C. 5.		6		16		200	O <sub>2</sub> /H <sub>2</sub>	1,500	15	0.4
APS	" "	5.	16		16		200	O <sub>2</sub> /H <sub>2</sub>	1,500	15	1.0
Titan I 1st Stage	Regen.	2.25	24	15.2	21.6	2.0	560	LOX/RP-1	150,000	587	1.6
Titan I 2nd Stage	Regen.	2.25	17	9.2	14.2	2.5	328	LOX/RP-1	80,000	682	1.9
ALRC H/C	Regen.	2.8	14	2.46	4.8	3.8	120	LOX/RP-1	12,000	1200	2.9
ALRC H/C	Regen.	2.8	14	2.46	4.8	3.8	120	LOX/RP-1	20,000	2000	2.9
SSME	Regen.	5.			17.8		600	LOX/RP-1	509,000	3250	

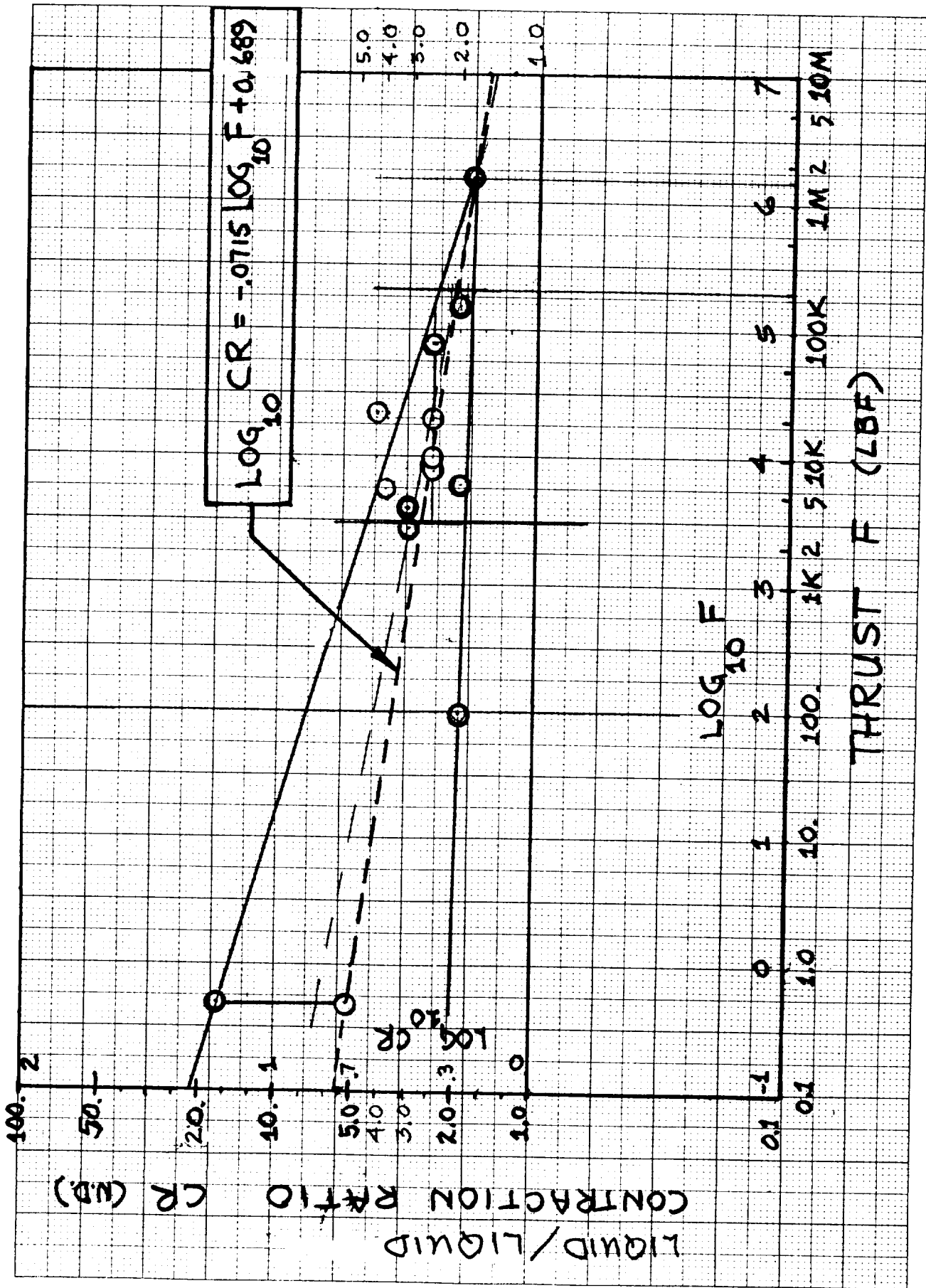
Enclosure 2



NOTE :  $F = P_c A_c \frac{I_s}{C^*}$

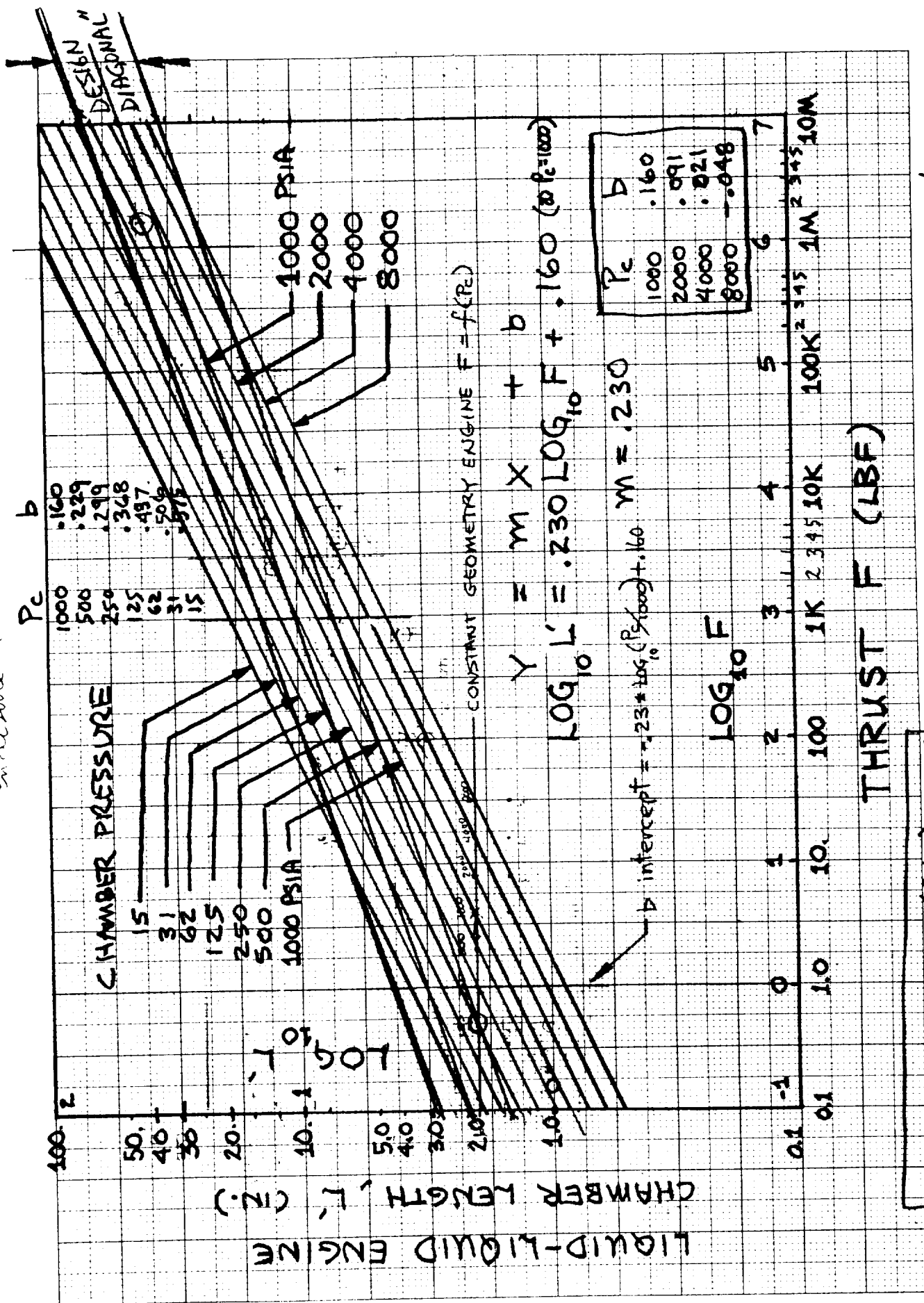
RAH 2 JAN 80

Enclosure 3



RAH 2 JAN '80

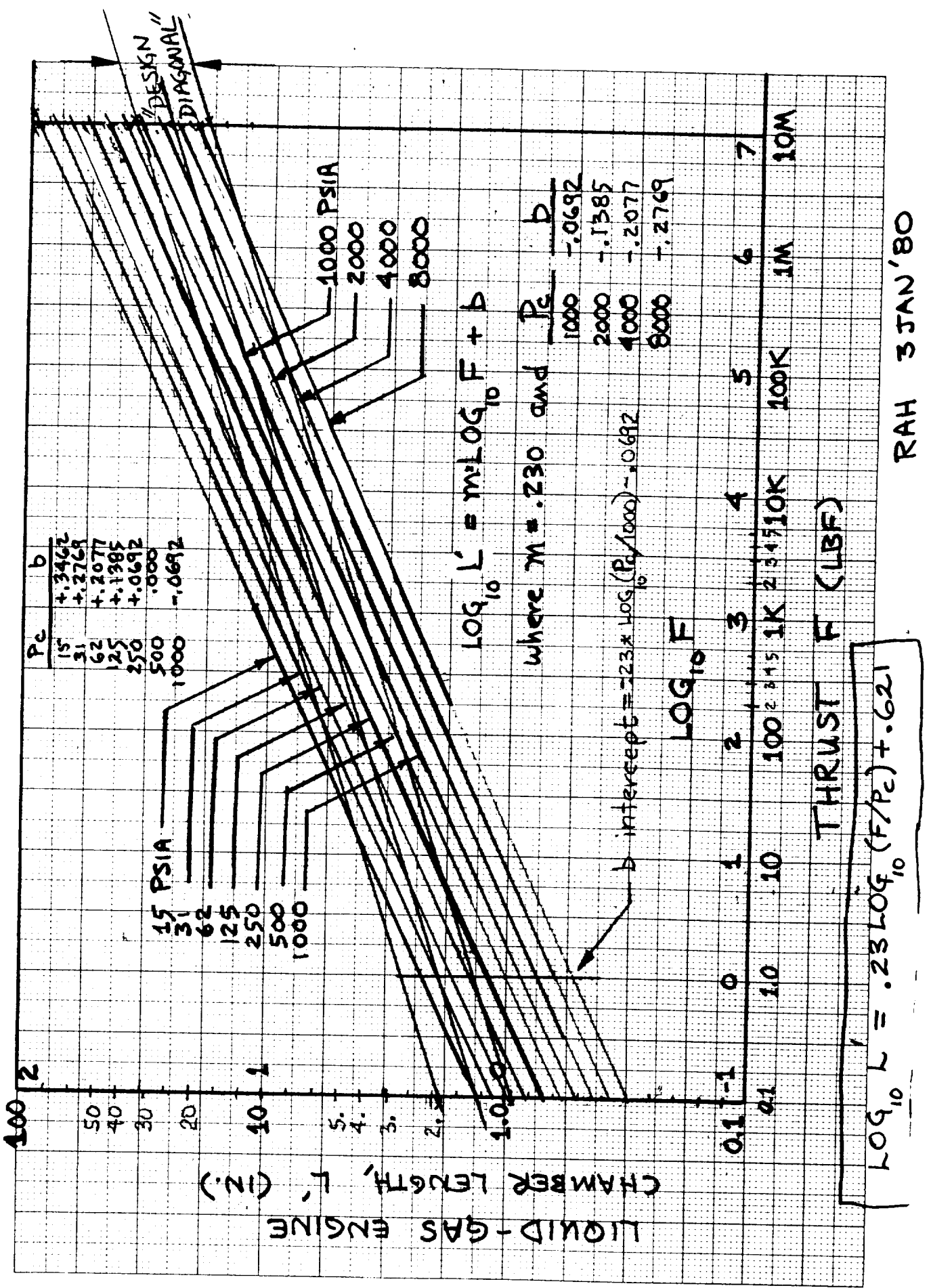
Envelope 4



$$\text{LOG}_{10} L' = .23 \text{ LOG}_{10} (F/P_c) + .85$$

RAH 2 JAN '80

Envelope 5



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## APPENDIX A

- Figure A-1 Throat Diameter vs Thrust
- Figure A-2 Liquid/Liquid Chamber Diameter
- Figure A-3 Liquid/Liquid Chamber Length to Diameter Ratio  
vs Thrust

## APPENDIX

### THROAT DIAMETER TRENDS

The exact determination of the throat diameter depends on many design details that will not be examined here. However, relatively accurate trends can be seen by using the specific impulse and characteristic velocity relationships and corrected for estimates of efficiencies and losses as follows:

$$C^* = P_c A_t g/W_T$$

and

$$I_{sp} = F/W_T$$

such that

$$D_t = \sqrt{\left[ \frac{4}{\pi g \eta_{DEL}} \right] \left[ \frac{C^* F}{P_c I_{sp}} \right]}$$

where  $\eta_{DEL} = (.95)$

$$D_t = .204 \sqrt{\frac{C^*_{ODK}}{I_{sp_{ODE}}} \left( \frac{F}{P_c} \right)}$$

$$D_{t_{Typical}} = .204 \sqrt{\frac{6000}{360} \left( \frac{F}{P_c} \right)}$$

$$D_t = .833 \sqrt{F/P_c}$$

This relationship for throat diameter is shown in Enclosure (6). Note that the "design diagonal" limits the allowable throat diameters for a given thrust and chamber pressure to the "band-width" shown.

### CHAMBER DIAMETER TRENDS

Once the estimated throat diameter and contraction ratio are known from Enclosures (3) and (6) the chamber diameter can be calculated as follows and as shown in Enclosure (7):

#### LIQUID/LIQUID (AND LIQUID/HOT-GAS IF $CR \geq 3.0$ ):

$$\text{LOG}_{10} D_c = .4643 \text{ LOG}_{10} F - 0.5 \text{ LOG}_{10} P_c + .2651$$

LIQUID/HOT-GAS (Assume:  $CR = 3.0$ , @  $P_c > 1000$  psia,  $F > 200,000$  lbf):

$$\text{LOG}_{10} D_c = 0.5 \text{ LOG}_{10} (F/P_c) + .159$$

Note that the liquid/hot-gas chamber diameter is calculated assuming a constant contraction ratio of 3.0 for the thrust and chamber pressure range considered in this study. If lower values of thrust (and at  $P_c > 1000$ ) were being considered and a liquid/liquid contraction ratio of greater than 3.0 were calculated it should be used in place of the constant 3.0 value.

### CHAMBER LENGTH TO DIAMETER RATIO TRENDS

Using the relationships for chamber length and chamber diameter shown above and in Enclosures (4), (5), and (7) the chamber  $L/D_c$  versus thrust and chamber pressure can be defined as follows:

#### LIQUID/LIQUID:

$$\frac{\text{LOG}_{10} L'}{\text{LOG}_{10} D_c} = \frac{.23 \text{ LOG}_{10} (F/P_c) + .850}{.4643 \text{ LOG}_{10} F - .5 \text{ LOG}_{10} P_c + .2651}$$

$$\text{LOG}_{10} (L'/D_c) = .27 \text{ LOG}_{10} P_c - .2343 \text{ LOG}_{10} F + .585$$

LIQUID/HOT-GAS: (Assuming:  $CR \geq 3$ ):

$$\begin{aligned}\frac{\text{LOG}_{10} L'}{\text{LOG}_{10} D_c} &= \frac{.23 \text{ LOG}_{10} (F/P_c) + .621}{.4643 \text{ LOG}_{10} F - .5 \text{ LOG}_{10} P_c + .2651} \\ \text{LOG}_{10} (L'/D_c) &= .27 \text{ LOG}_{10} P_c - .2343 \text{ LOG}_{10} F + .356\end{aligned}$$

LIQUID/HOT-GAS: (Assume:  $CR = 3$ , @  $P_c > 1000$  psi, &  $F \geq 200,000$  lbf):

$$\begin{aligned}\frac{\text{LOG}_{10} L'}{\text{LOG}_{10} D_c} &= \frac{.23 \text{ LOG}_{10} (F/P_c) + .621}{.5 \text{ LOG}_{10} (F/P_c) + .159} \\ \text{LOG}_{10} (L'/D_c) &= .462 - .27 \text{ LOG}_{10} (F/P_c)\end{aligned}$$

Note that the middle formula does not apply to this study since all the liquid/liquid contraction ratios will be less than 3.0 and all the liquid/hot-gas contraction ratios are set equal to 3.0 due to the high thrust.

Note also that the "design diagonal" band-width does not allow excessively large or small  $L'/D_c$  values to be encountered (i.e., no larger than 22 @  $F = 0.1$  lbf, or no smaller than 0.5 @  $F = 10\text{MLBF}$ ).



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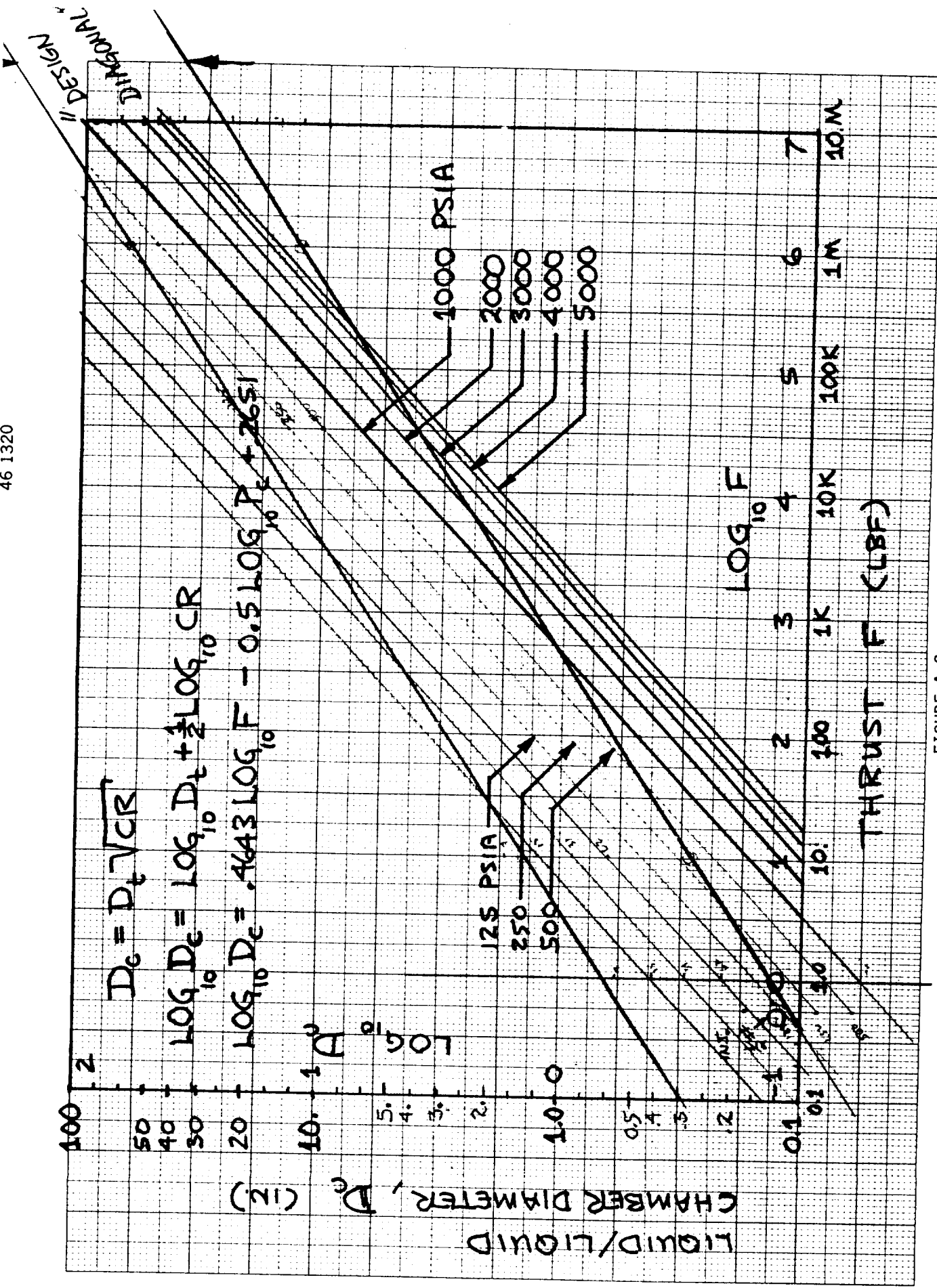


FIGURE A-2

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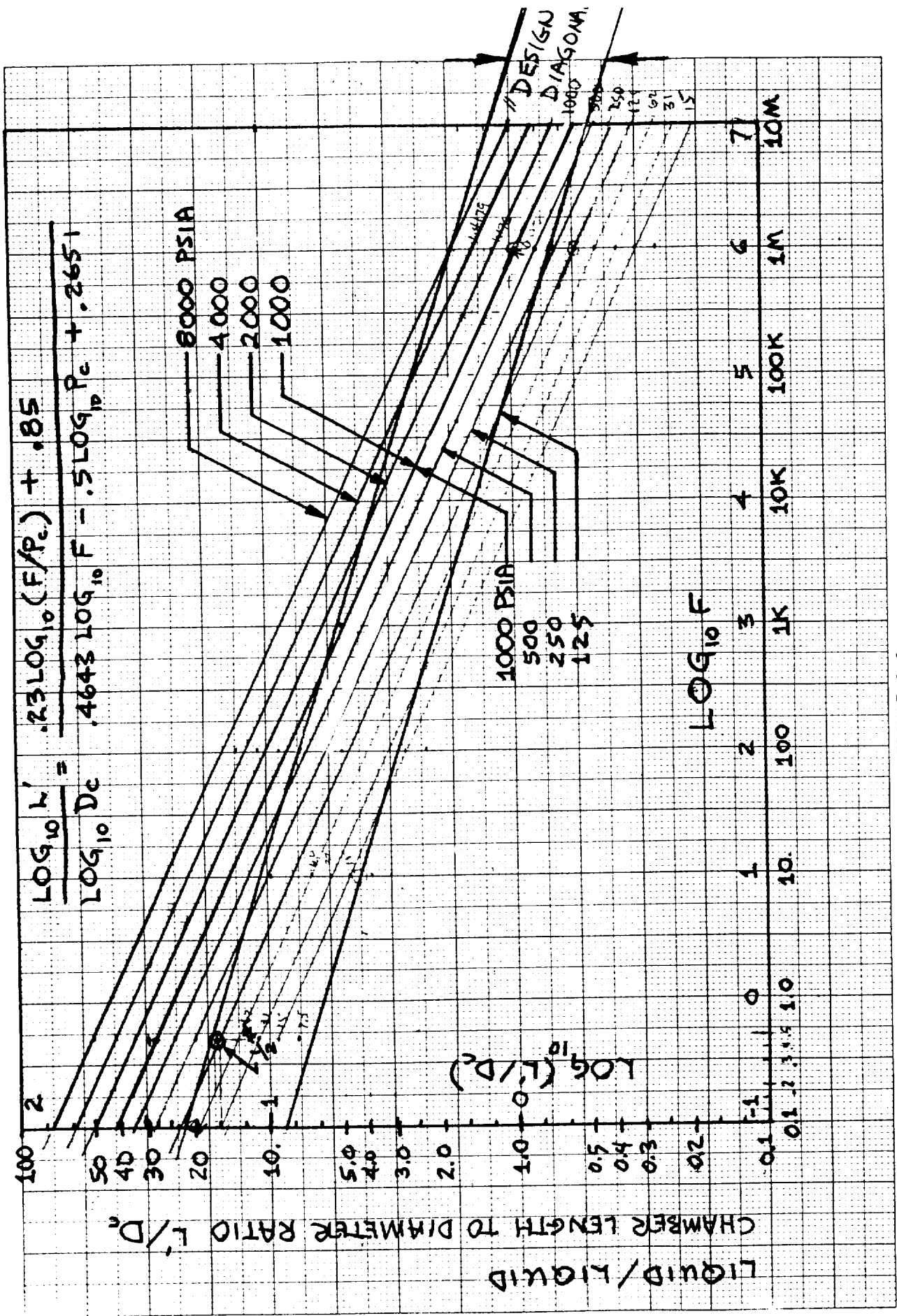
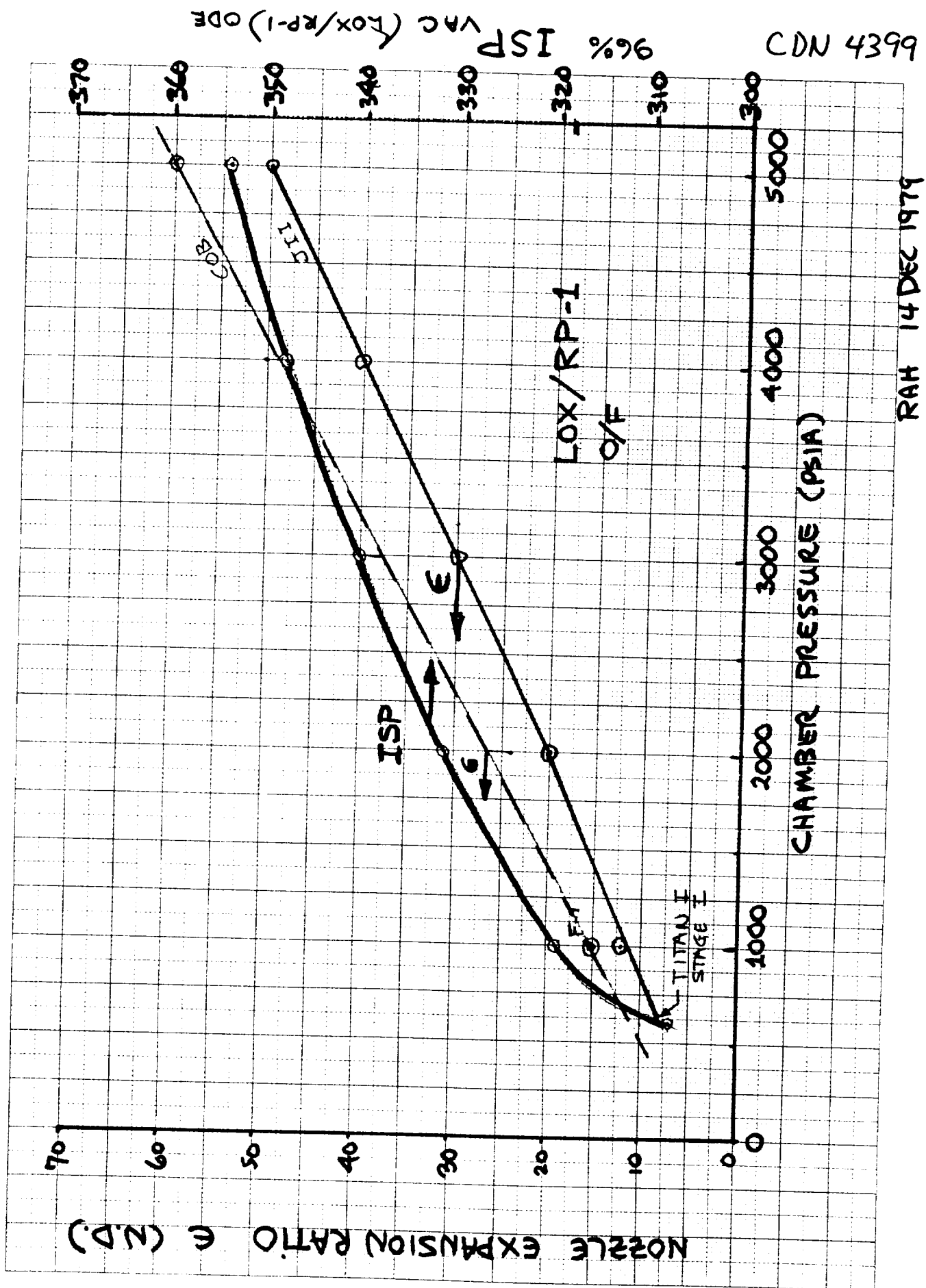
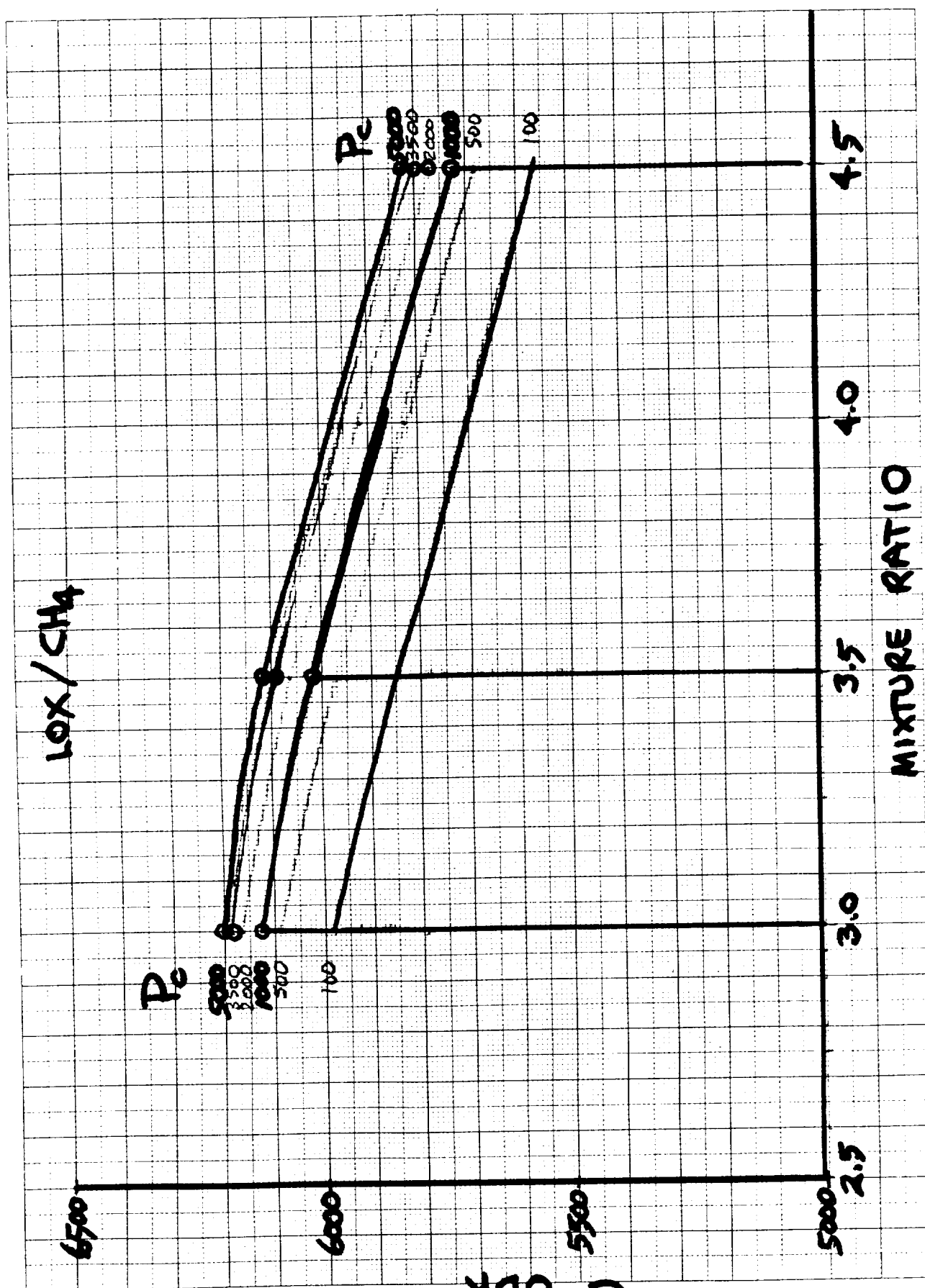
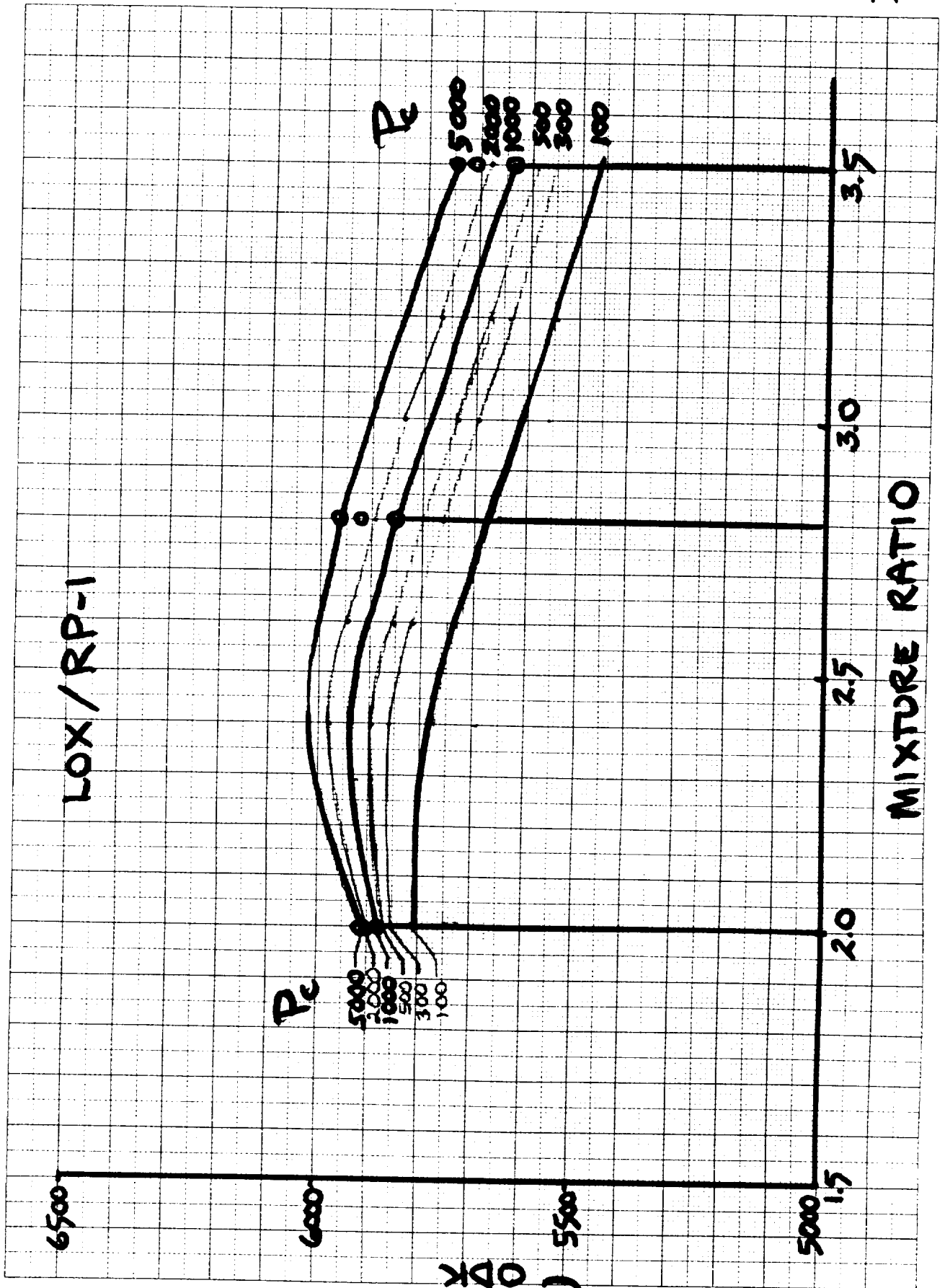


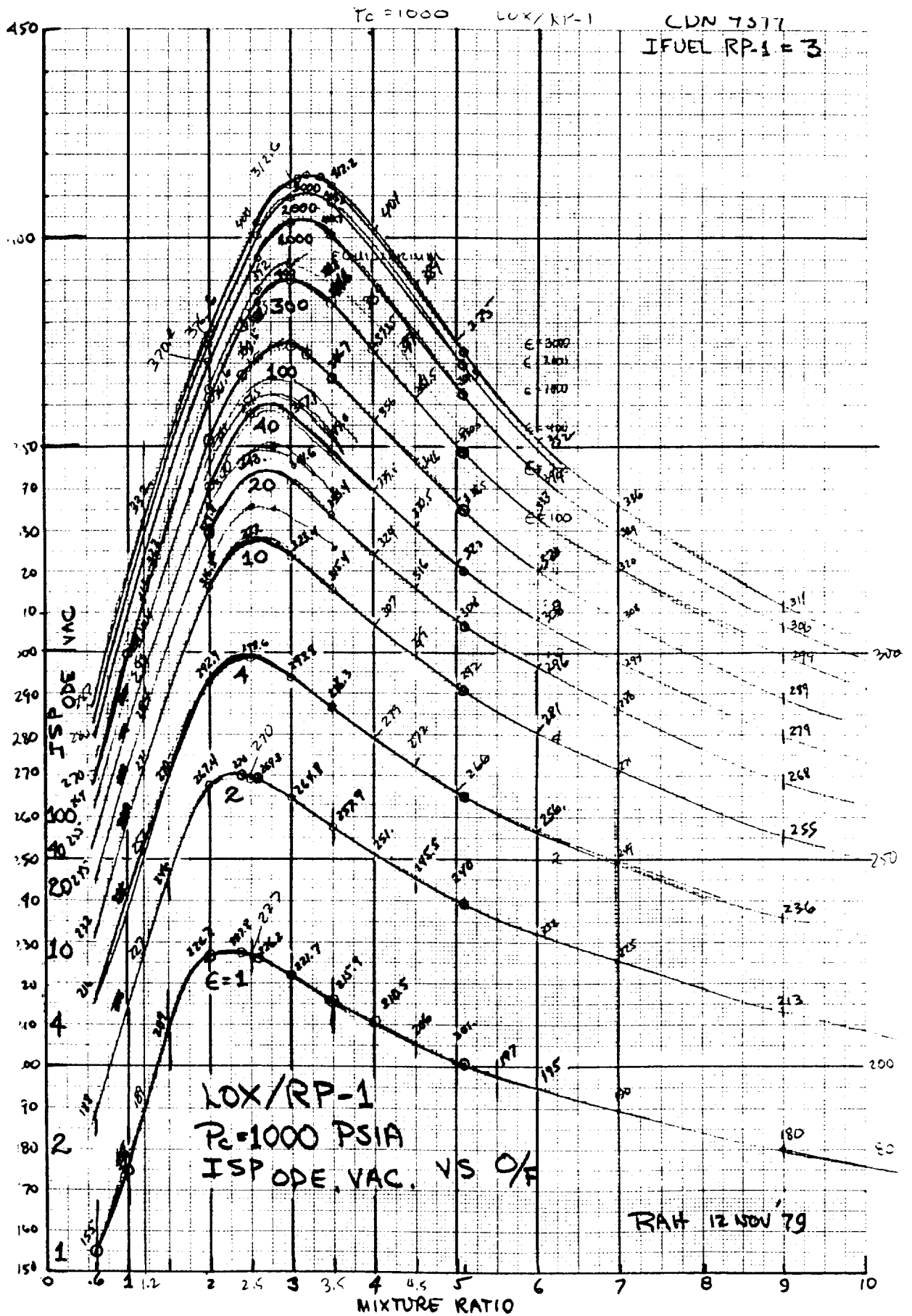
FIGURE A-3

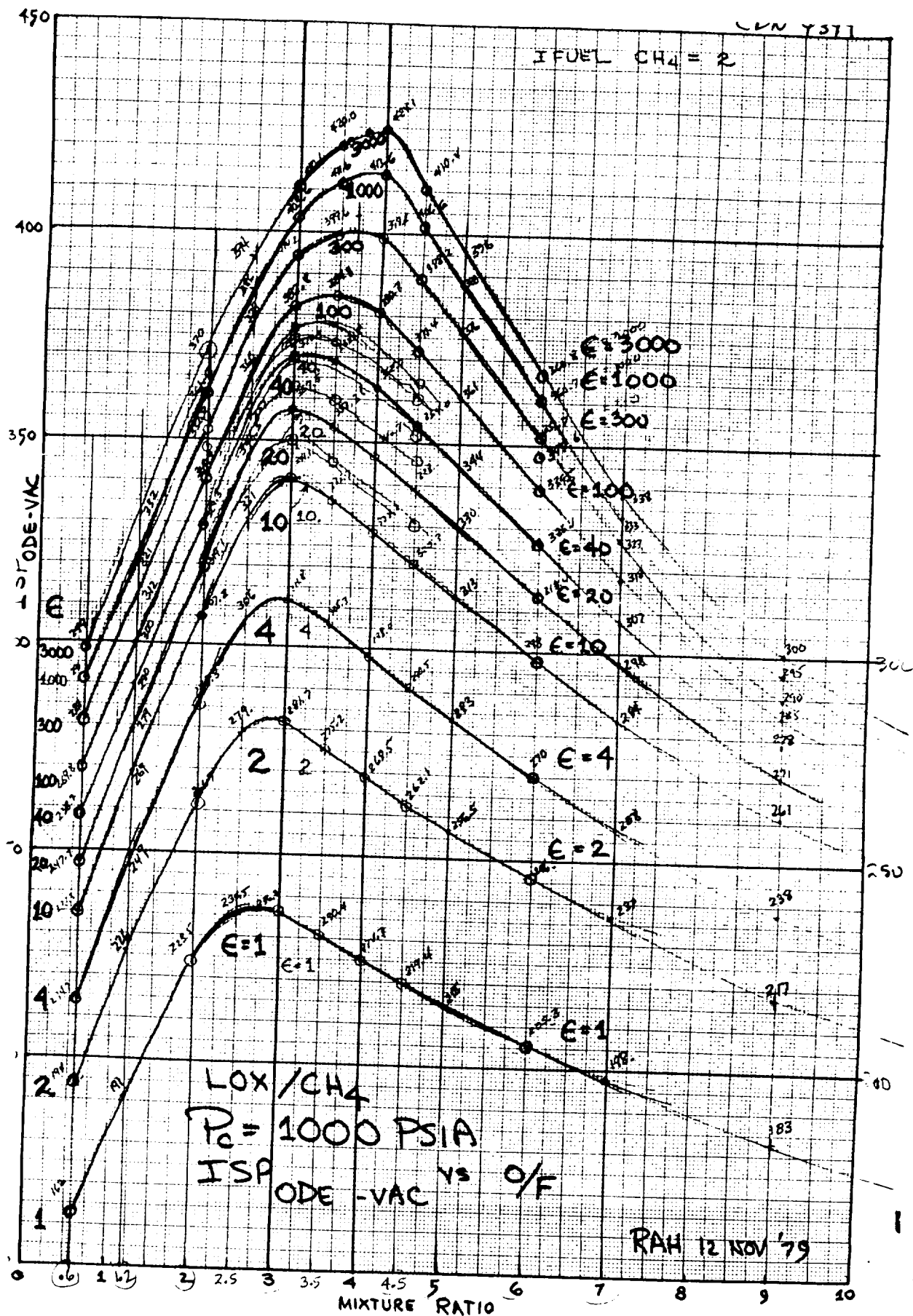




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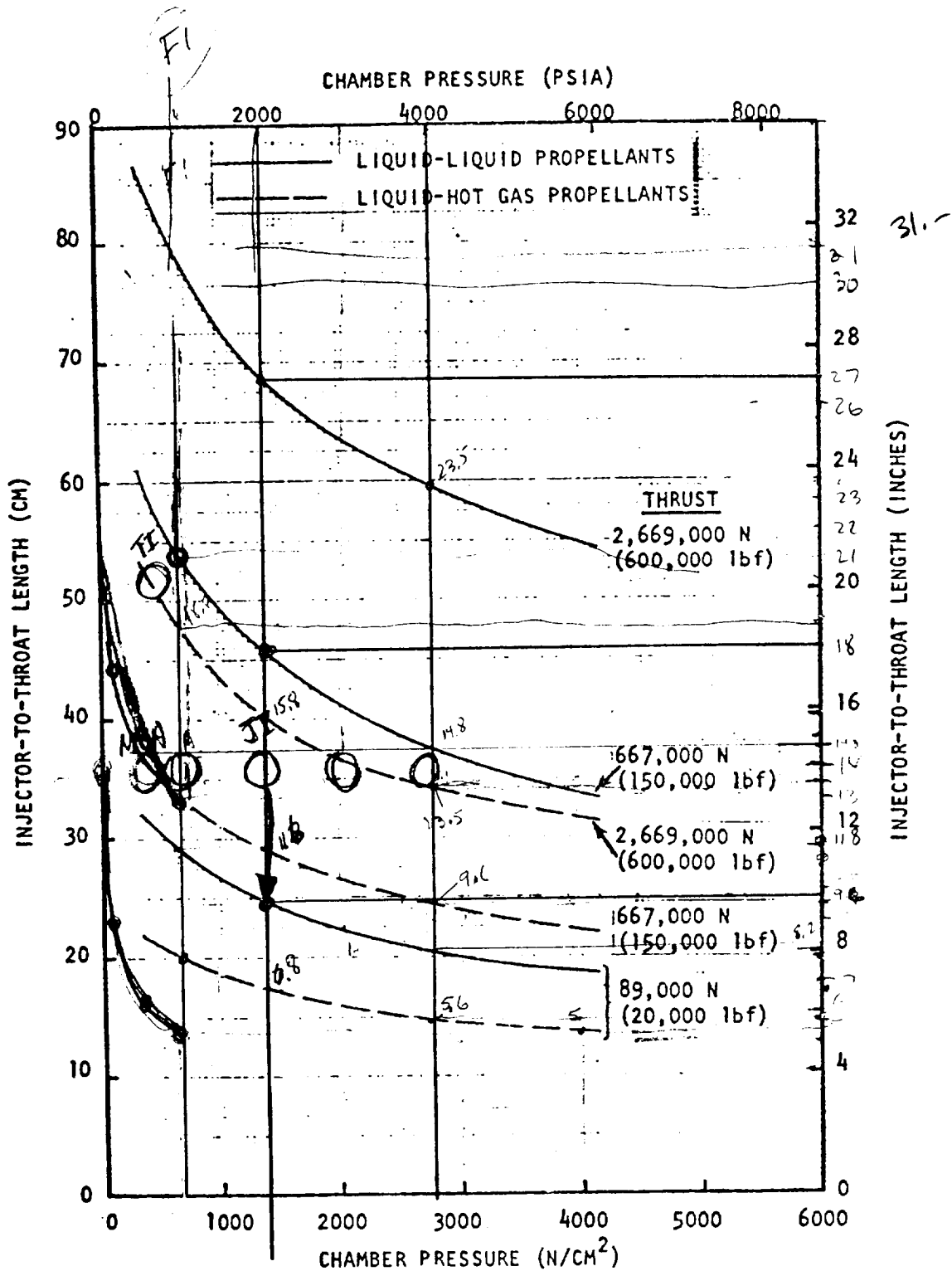


Figure 3. Injector-to-Throat Length Parametric Data

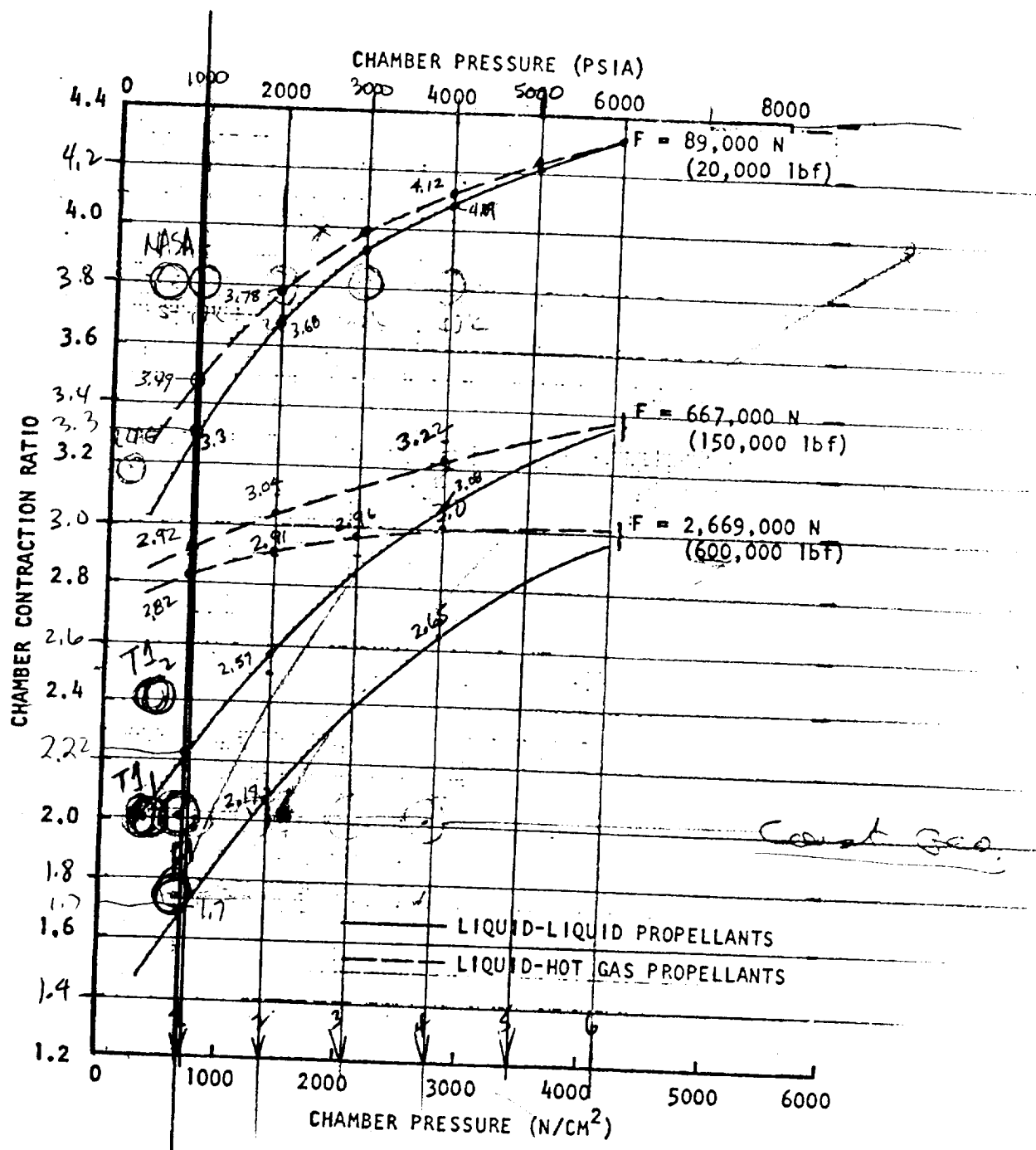


Figure 4. Chamber Contraction Ratio Parametric Data.

# INJECTION ELEMENT QUANTITY REMAINS CONSTANT FOR A GIVEN CHAMBER WHILE DESIGN PC AND ENGINE THRUST ARE VARIED

NASA Lewis Design

● ASSUMED CHAMBER DESIGN-CONSTANT

○  $D_c = 5.39$  IN. ✓

○  $D_T = 2.60$  IN. ✓

○  $L' = 14.0$  IN. (VARIABLE)

● INJECTION ELEMENT QUANTITY-CONSTANT

○  $N_E \sim 180$  PAIRS, TRANSVERSE (PLATELET) L-O-L DOUBLET VS 37 LOL ???

● VARIABLES ( $600 \text{ PSIA} \leq PC \leq 4000 \text{ PSIA}$ )

○ INJECTION VELOCITY AND  $\Delta P$   $120 \rightarrow 600 \text{ psid}$

○ ORIFICE DIAMETER  $.035 \rightarrow .06$

○ UNLIKE MIXING DISTANCE Little  $\rightarrow$  Bigger

○ BOOSTER NOZZLE Little  $\rightarrow$  Bigger

○ SENSITIVE TIME LAG large  $\rightarrow$  smaller

○ CHAMBER HEAT FLUX little  $\rightarrow$  Bigger

○  $\Delta P/PC$  RATIO-SLIGHT  $.20 \rightarrow .15$  (Decrease)

○ ATOMIZED DROP SIZE Increase

○ SPRAY ATOMIZATION LENGTH Increase

○ OPTIMUM CORE MIXTURE RATIO Increase

○ HI FREQ COMBUSTION GAIN Increase

● CONSTANTS

○ COMBUSTION EFFICIENCY

98 TO 99% DES'N GOAL

○ TOTAL COMB. TIME LAGS

$\tau_{ox} = 0.4 \text{ msec}$

$\tau_f = 0.25 \text{ msec}$

○ CHUG STABILITY MARGIN

○ AXIAL COMBUSTION PROFILE

% VAP.  $L' \text{ (IN.)}$

99 - OX 3.5

100 - OX 5

99 - F 12

99.8-TOT 14

Aerofast Liquid Rocket Company

# PLATELET COMBUSTION TECHNOLOGY

THE PROPOSED PLATELET INJECTOR CONCEPT WILL PROVIDE ESSENTIAL LOX/HC COMBUSTION TECHNOLOGY DATA OVER THE ENTIRE PC OPERATING RANGE.

LO <sub>2</sub> /RP-1 (EX0 - THDCP)		D <sub>CHAM</sub> = 5.39
N <sub>E</sub> = 180 T-LOL PAIRS		D <sub>T</sub> = 2.60
η <sub>C*</sub> = 98%		L' = 14.0

P <sub>C</sub> ΔP/P <sub>C</sub> ΔP <sub>INJ</sub> (PSID) ε (BOOSTER) O/F OPT. 96% I <sub>sp</sub> (SEC) F <sub>VAC</sub> (LB <sub>F</sub> )	600 .20 120 8 2.6 307 5.5K	1000 .185 185 12 2.65 319 9.5K	2000 .17 340 20 2.75 331 20K	3000 .16 480 30 2.8 340 30K	4000 .15 600 40 2.9 348 40K
W <sub>i</sub> (LB <sub>m</sub> /SEC) V <sub>INJ</sub> (FT/SEC) D <sub>ORIF</sub> (IN.)	OX F 13.0 5.0 125 150 .035 .023	OX F 21.5 8.1 155 185 .041 .027	OX F 43.3 15.7 210 250 .050 .033	OX F 65.3 23.3 250 300 .056 .036	OX F 87.0 30.0 280 335 .061 .039

Aerojet Liquid Rocket Company

$$W_T = \frac{F}{I_s} = \frac{5,300}{301} = 18$$

# LeRC IN-HOUSE COMBUSTION TESTING WILL PROVIDE VALID HI FREQ COMBUSTION STABILITY DATA FOR 600K ADVANCED HYDROCARBON ENGINE DEVELOPMENT

- ALRC HAS DEMONSTRATED ABILITY TO ACHIEVE DYNAMIC COMBUSTION STABILITY PER CPIA 247 WITH ACOUSTIC CAVITIES WHEN TRANSVERSE RESONANT COMBUSTION MODES ARE LIMITED  $\leq 3T$ .

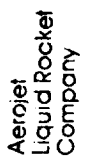
- OMS
- ITIP
- MX-AXIAL

- INJECTOR COMBUSTION CAPABILITY TO SUPPORT 1T MODE IN PROPOSED CHAMBER WILL BE INDICATIVE OF 3T COMBUSTION GAIN FOR FULL SCALE AHCE.

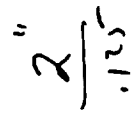
PROGRAM	LeRC IN-HOUSE	LOX/HC TECHNOLOGY	AHCE
Pc, (PSIA)	600	~2000	4250
F, (LBf)	5K	25K	600K
D <sub>CHAM.</sub> , (IN.)	5.4	5.4	14.
F <sub>1T</sub> , (HZ)	5100	5100	-
F <sub>3T</sub> , (HZ)	-	-	4500

- ELEMENT TYPES WHICH SUPPORT  $\geq 2T$  IN SUB-SCALE CHAMBERS SHOULD BE ELIMINATED FROM FURTHER CONSIDERATION FOR FULL SCALE AHCE ENGINE APPLICATION.

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**CHAMBER CONTOUR IS DESIGNED TO MINIMIZE HEAD END HEAT FLUX WHILE MAINTAINING CHAMBER COMPATIBILITY**



2

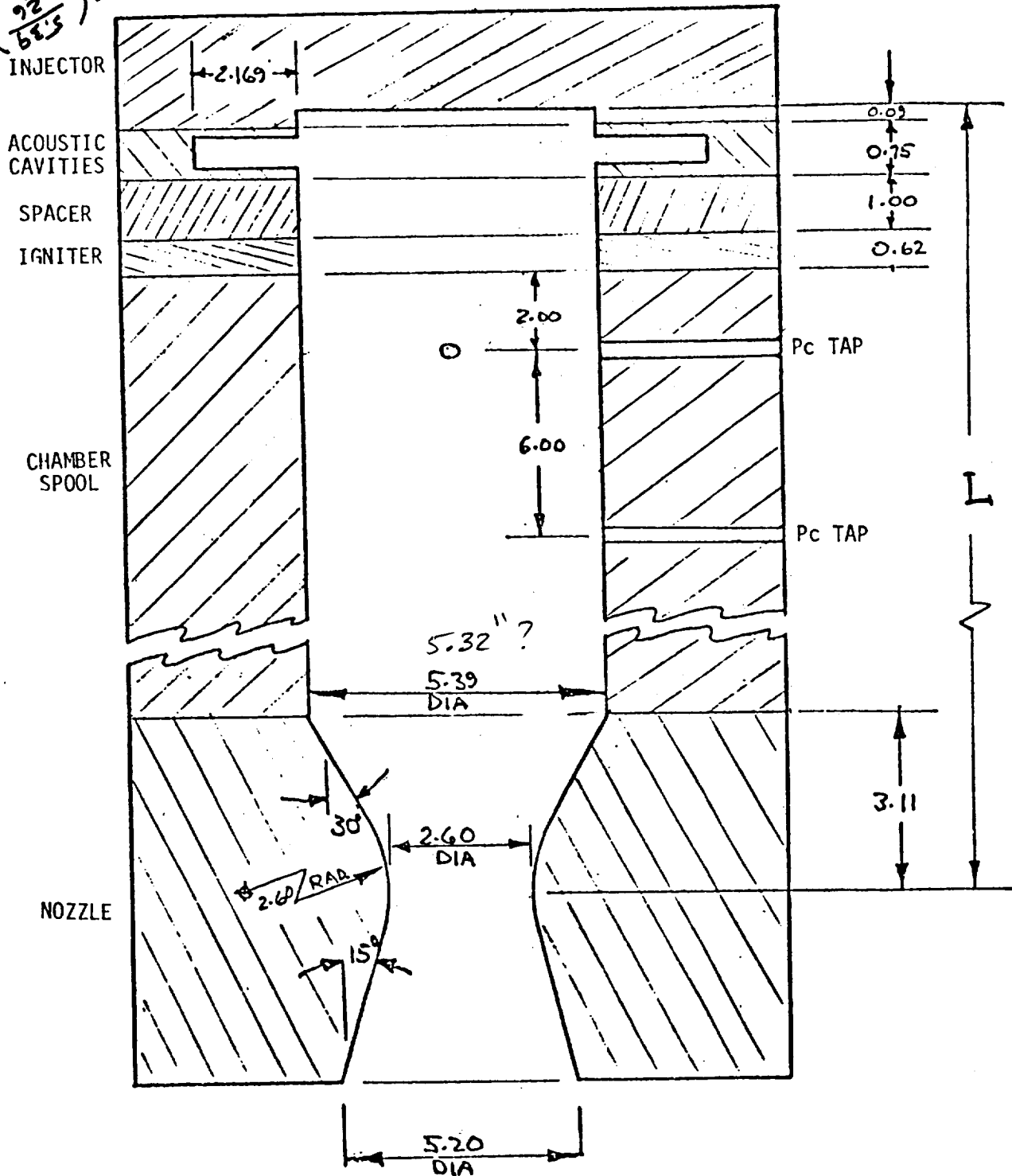
88% FVE

NASA Pauli

CDN 4399

COMBUSTION CHAMBER

SL981.6  
 $9262.4 = 277$   
 $\left(\frac{92}{68.5}\right) =$   
 2 INJECTOR



**APPENDIX K**  
**ON FILE AT NASA/LeRC**

**Contact**

**Mark Klem**  
**National Aeronautics and Space Administration**  
**Lewis Research Center**  
**Cleveland, Ohio 44135**  
**(216) 433-2450**



National Aeronautics and  
Space Administration

## Report Documentation Page

1. Report No. NASA CR - 187110	2. Government Accession No.	3. Recipient's Catalog No.	
4. Title and Subtitle User's Manual for Rocket Combustor Interactive Design (ROCCID) and Analysis Computer Program  Volume II—Appendixes A-K		5. Report Date May 1991	
		6. Performing Organization Code	
7. Author(s) J.A. Muss, T.V. Nguyen, and C.W. Johnson		8. Performing Organization Report No. None	
		10. Work Unit No. 582 - 01 - 21	
9. Performing Organization Name and Address Gencorp, Aerojet Propulsion Division Sacramento, California 95813 - 6000 and Software and Engineering Associates Carson City, Nevada 89701		11. Contract or Grant No. NAS3 - 25556	
		13. Type of Report and Period Covered Contractor Report Final	
12. Sponsoring Agency Name and Address National Aeronautics and Space Administration Lewis Research Center Cleveland, Ohio 44135 - 3191		14. Sponsoring Agency Code	
15. Supplementary Notes Project Manager, Mark D. Klem, Space Propulsion Technology Division, NASA Lewis Research Center, (216) 433 - 2450. J.A. Muss and T.V. Nguyen, Gencorp Aerojet Propulsion Division; C.W. Johnson, Software and Engineering Associates.			
16. Abstract <p>This report is the Appendixes A-K to the User's manual for the Rocket Combustor Interactive Design (ROCCID) computer program. This includes installation instructions, flow charts, subroutines model documentation and sample output files. The ROCCID program, written in FORTRAN 77, provides a standardized methodology using state-of-the-art codes and procedures for the analysis of a liquid rocket engine combustor's steady state combustion performance and combustion stability. ROCCID is currently capable of analyzing mixed element injector patterns containing impinging like doublet or unlike triplet, showerhead, shear coaxial and swirl coaxial elements as long as only one element type exists in each injector core, baffle or barrier zone. Real propellant properties of oxygen, hydrogen, methane, propane and RP - 1 are included in ROCCID. The properties of other propellants can be easily added. The analysis models in ROCCID can account for the influences of acoustic cavities, helmholtz resonators and radial thrust chamber baffles on combustion stability. ROCCID also contains the logic to interactively create a combustor design which will meet input performance and stability goals. A preliminary design results from the application of historical correlations to the input design requirements. The steady state performance and combustion stability of this design is evaluated using the analysis models, and ROCCID guides the user as to the design changes required to satisfy the user's performance and stability goals, including the design of stability aids. Output from ROCCID includes a formatted input file for the standardized JANNAF engine performance prediction procedure.</p>			
17. Key Words (Suggested by Author(s)) Combustion stability; Rocket engine design; Liquid propellant rocket engines; Launch vehicles; Booster rocket engines; Propellant combustion; Engine design; Dynamic stability; Acoustic velocity; Velocity coupling		18. Distribution Statement Unclassified - Unlimited Subject Categories 5 and 20	
19. Security Classif. (of the report) Unclassified	20. Security Classif. (of this page) Unclassified	21. No. of pages 577	22. Price* A25