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# Transient Thermal Modeling of the Nonscanning ERBE Detector

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Contract NAS1-16508
November 1983

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86N12743\*# ISSUE 3 PAGE 453 CATEGORY 43 RPT#: NASA-CR-172240 NAS 1.26:172240 CNT#: NAS1-16508 83/11/00 107 PAGES UNCLASSIFIED

DOCUMENT

UTTL: Transient thermal modeling of the nonscanning ERBE detector TLSP: Fina Report, 13 Nov. 1980 - 30 Sep. 1982

AUTH: A/MAHAN, J. R.

CORP: Virginia Polytechnic Inst. and State Univ., Blacksburg. AVAIL.NTIS

SAP: HC A06/MF A01 CIO: UNITED STATES

MAJS: /\*HEAT TRANSFER/\*MONTE CARLO METHOD/\*THERMODYNAMIC PROPERTIES/\*TRANSIENT RESPONSE

MINS: / EARTH RADIATION BUDGET EXPERIMENT/ MATHEMATICAL MODELS/ RADIOMETERS

ABA: E.A.K.

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# Transient Thermal Modeling of the Nonscanning ERBE Detector

#### 1. Introduction

This report summarizes the activity at VPI under Contract NAS1-16508 between 13 November 1980 and 30 September 1982. During this period a numerical model of the ERBE wide-field-of-view total radiometer channel was developed, and the model was applied to the TRW calibration procedure. Other activities included consulting with NASA and contractor scientists and engineers on the design and calibration of radiometric instruments and on the interpretation of data from such instruments.

Three major documents have resulted from this work: a mid-term report entitled, Application of the Monte Carlo Method to the Transient Thermal Modeling of a Diffuse-Specular Radiometer Cavity (Leo Eskin's M.S. thesis) 1, a Users' Manual for the cavity transient thermal model program (included as an appendix to reference 1, and the TRW Calibration Simulation Users' Manual, the final revised version which appears in the Appendix of this final report. In addition, a paper was presented at the Fourth AMS Conference on Atmospheric Radiation in Toronto in June of 1981<sup>2</sup>.

The final numerical model of the ERBE instrument, the "TRW Calibration Simulation Model", consists of two sub-models which interact with each other within the structure of the overall model. One of these sub-models computes the transient thermal response of the cone

cavity, and the other computes the transient thermal response of the instrument body itself. The details of the cavity sub-model are described in reference 1 and thus are not reiterated here.

In the present report we document the instrument body sub-model and its interaction with the cavity sub-model and its surroundings. We indicate in some detail how this model can be modified to take into account the "two-temperature" aspect of the TRW calibration source with its mounting collar, or baffle. We also outline how the "TRW Calibration Model" can be modified and extended to serve as a flight simulation model, including operation as a visible channel. Finally, we suggest a procedure for using the model to establish a real-time interpretation strategy.

#### 2. The TRW Calibration Simulation Model

The Users' Manual for the "TRW Calibration Simulation Model" appears in the Appendix of this report. This is the final revised version originally transmitted to NASA Langley by Dave Parekh in September, 1982. At that time Mr. Parekh also sent NASA a magnetic tape containing this model. In this section we describe the parts of this model not already documented in Reference 1. This is also where we suggest the modifications referred to in the Introduction.

#### 2.1. TRW FRONT

"TRW FRONT" is a subprogram of the TRW Calibration Simulation Model in which Monte Carlo techniques are used to compute the

end" of the instrument, between these surfaces and the blackbody calibration source, and between these surfaces and the cavity. By "front end", we mean the field-of-view (FOV) limiter and the surface in the plane of the precision aperture to which the FOV limiter is attached. The general principles of the Monte Carlo technique are described in great detail in Reference 1. Because the logic of "TRW FRONT" is very similar to that of the cavity distribution factor program documented in Reference 1, we will not reiterate the details here. However, the program listing, with notes, appears on pages 47-76 of the Appendix.

The case documented in the Appendix is for the wide field-of-view total channel. It can be easily modified for any other geometry by changing values of the data read by the program, "READ(5,97)", page 49. These data appear after the subroutines which serve "TRW FRONT", on page 75 of the Appendix. Each element of this data set is carefully labeled on page 76, described on page 47, and referred to in Fig. 1 of the February, 1982, Progress Report.

## 2.1.1. Modification to Account for a Two-Temperature Calibration Source

The present version of "TRW FRONT" assumes that the instrument faces a uniform temperature blackbody source. The blackbody calibration source is thus modeled as a perfectly absorbing sector in the plane of the FOV limiter. This is exactly equivalent to any real

as the cavity is isothermal. If, however, as in the case of the actual TRW calibration source, the calibration cavity is composed of two surfaces having different temperatures, "TRW FRONT" must be modified to take into account its actual shape. The page-by-page procedure for this modification is outlined below:

1. Page 49: change NN3 = NN2 + 3

to NN3 = NN2 + 4

change MM3 = M3 + 1

to MM3 = M3 - NTH

2. Page 50: change

GOTO(101, 102,..., 110),J

101 CONTINUE

102 CONTINUE

to

GOTO(100, 101, 102,..., 110),J 100 CONTINUE

statements similar to those presently between "101 CONTINUE" and "102 CONTINUE" which establish the x, y and z coordinants of a randomly selected source location on the controlled surface of the calibration cavity

101 CONTINUE

statements which establish the x, y and z coordinants of a randomly selected source location on the concentric collar, or baffle, which couples the calibration source to the ERBE instrument

102 CONTINUE

3. Page 55: change IF(TZ1.EQ.H2) GOTO 201

to IF (TZ1.GT.H2) GOTO 201

- 4. Page 56: change three statements between "201 CONTINUE" and "GOTO 23" to the logic necessary to compute values of UNX, UNY and UNZ, depending on whether the source point is on the blackbody surface itself or on the baffle. This logic depends on the geometry of the calibration source and its interface with the instrument, but should be easy to implement.
- depending on the geometry of the calibration source/radiometer interface. First, determine if the source point is on the calibration black-body surface or on the baffle by testing the value of TZI in a manner similar to that on page 60. Then compute TXJ and TYJ in a manner similar to that in the present version of the program. If the source point is on the black-body surface, begin the search to see if the energy packet has entered the instrument field of view by comparing the value of

DSQRT(TXJ\*TXJ + TYJ\*TYJ)

with the radius of the FOV limiter at TZJ

= H2. If this value exceeds the radius of the

FOV limiter, the ray is reabsorbed somewhere

within the calibration source, in which case

go to "RETURN". However, if this value is

less than the radius of the FOV limiter at TZJ = H2, the energy packet has entered the instrument field of view, in which case the search continues as presently organized under "11 CONTINUE". The logic is similar if the source point is on the baffle rather than the blackbody surface itself.

#### 2.1.2. Adaptation to a "Real" Source Field

If the "TRW Calibration Simulation Model" is to be modified as a flight simulator, the calibration blackbody source must be replaced with a "real" source field. This real field must, of course, be time varying. It is impractical to work in terms of distribution factors from the earth scene to the surfaces of the instrument, since these factors would all be essentially zero (the <u>fraction</u> of radiation emitted diffusely from Kansas which is absorbed by a 5-mm square surface in earth orbit is too small to compute). We also have demonstrated that an artificial "grid" placed over the FOV inlet plane to represent the earth scene is unsatisfactory because of its "focusing" effect. The most efficient way to adapt the existing model to a "real" field, such as GENDAT, is outlined below:

1. Use GENDAT (or some equivalent source field model) to compute the flux incident to the disk defined by the FOV limiter, as well as to the disk defined by the precision aperture. These calculations are both within the present capabilities of GENDAT. incident to the FOV limiter and the "floor" surrounding the precision aperture. Assume that none of this energy is specularly reflected into the cone cavity, in keeping with the design principle of the FOV limiter. The fraction of this energy which is diffusely reflected into the precision aperture is

$$\sum_{i=1}^{n} D_{i-A} \rho_{i}^{d} ,$$

where  $D_{i-A}$  is the distribution factor from the ith front end surface element to the disk defined by the precision aperture, computed in "TRW FRONT", and  $\rho_i^d$  is the diffuse component of the reflectivity of this element. Note that this calculation assumes that the energy from the scene incident to the front end optics is uniformly distributed across the FOV limiter and floor. This assumption is of minor consequence in view of the fact that it is approximately true for any realistic earth scene, and that the result in any case is a small contribution to the total energy input to the cavity.

3. The fraction of the flux entering the cavity directly from the earth scene which is absorbed by cavity surface element j is

$$D_{A-j} = \epsilon_j \frac{A_j}{A_A} D_{j-A}$$

where  $A_j$  is the surface area of cavity element j,  $\epsilon_j$  is its diffuse emissivity,  $A_A$  is the precision aperture opening area, and  $D_{j-A}$  is the distribution factor from cavity element j to the precision aperture opening computed in "TRW NODE". This calculation assumes that the flux from the earth scene incident to the precision aperture opening is diffuse. Although this assumption is not strictly valid, the error associated with it will be negligible in view of the cavity properties.

#### 2.1.3. Inclusion of a Filter Dome

The extension of the "TRW Calibration Simulation Model" to include the visible channel would be a major modification. It would involve altering the paths of the energy packets which intercepted the filter dome as follows:

- Upon the intersection of a ray with the filter dome, a random number would be generated to determine if it was reflected. If not reflected, another random number would be generated to see if it was absorbed or transmitted.
- 2. If the energy packet is absorbed, the dome element would be treated like any other surface element; that is, the counter for the distribution factor from the source element to the dome element would be incremented.
- If the energy packet is transmitted, diffraction can be included according to Snell's law.
- 4. If the energy packet is reflected, we have to generate

another random number to determine if the reflection is specular or diffuse.

- a) If specular, treat the dome element like any other surface.
- b) If diffuse, the possibility of forward scattering should be included; that is, uniformly distributed scattering in  $4\pi$  space rather than  $2\pi$  space.

After reflection, or scattering, the energy packet is traced to its next surface just as in the case without a dome.

#### 2.2. TRW TEMP

This subprogram computes the transient temperature distributions in the active cavity and body of the instrument based on the distribution factors and conductances computed in the previous subprograms (or user generated in the case of the body elements). As presently configured, this calculation is based on an assumed isothermal blackbody source. Arbitrary heat sink and mounting beam temperatures can be introduced to correspond to actual calibration conditions. A detailed annotated listing of this subprogram appears in the Appendix, pages 77-98. The data listings appear on pages 99-104.

The thermal mass of a typical instrument body node is much larger than that of a typical cavity node. For this reason the temperatures of the body nodes change much more slowly than those of the cavity nodes. Thus, we can realize a certain efficiency by using a "two-timing" technique to compute the transient temperature distributions in the instrument body and on the active cavity surfaces. The body

node temperature calculations are performed only after every tenth cavity node temperature calculation. The radiation heat transfer from the FOV limiter to the active cavity used in each cavity temperature calculation is then based on the quasi-steady temperatures of the FOV limiter nodes. The conduction heat transfer from the cavity nodes to the body nodes is also calculated based on the quasi-steady body node temperatures and stored in a buffer until needed for the next body node temperature calculation, at which time this accumulated energy is distributed to the appropriate body nodes.

#### 2.2.1. Detailed Description of the Body Node Temperature Calculation

The active cavity node temperature calculation is well documented in Reference 1. In this section the body node temperature calculation and its interaction with the cavity node temperature calculation is described.

The values of certain constants, defined on page 77, are established on page 80. In particular, the user must change "HST", "TCLAMP", "HEAT" and "TSOURC" to conform to the actual calibration conditions.

The statements on pages 80-86 establish the values of the constants, coefficients, conductances, distribution factors, and initial temperatures for the calculations. The actual calculations begin on page 86 with the "280" DO-loop, which extends to the end of the subprogram.

The first step in this sequence is to check to see if it is time to do a body temperature calculation ("IF(KOUNT.EQ.10) GOTO 678", page 87). The actual body temperature calculations lie between

"678 CONTINUE", page 87, and "679 CONTINUE", page 88. Between "219 CONTINUE", page 87, and "GOTO 219", page 88, an estimate of the new node temperatures is computed based on the old node temperatures, the input from the cavity, the input from the field (calibration source), and the inputs from the heat sink and mounting beam. This calculation loop is repeated until convergence is obtained to a cavity temperature distribution whose rms value no longer changes from one loop to the next. In this loop, "TGBODY" represents the "guess" for the temperature distribution, and "TUBODY" represents the updated temperature distribution. At the end of each cycle through this loop, the guess is updated and, if there is a significant difference between the update and the preceding guess, the loop is repeated. In the present version of the subprogram a change of one part in a million is tolerated in the rms difference between the last two iterations. This tolerance can be changed by the user in the statement "IF(TEST.LE.0.000001) GOTO 241", page 88.

The updated temperature distribution is actually computed in the "230" DO-loop, pages 87-88. The logic in this loop is "hard wired" in that, if the body shape or node configuration is changed, the logic in this loop must be changed accordingly. These changes would be made in the statements between "220 CONTINUE" and "637 CONTINUE", page 87. Between "637 CONTINUE", page 87, and "230 CONTINUE", page 88, the actual temperature calculation is made based on Newton's method. From the statement "230 CONTINUE", page 88, down to the statement "IF(TEST.LE.0.000001) GOTO 241", page 88, the relative rms change in the temperature distribution is calculated and tested.

If the temperature distribution has converged, the body node temperatures "T1BODY" are changed to their new values, "TUBODY", in the "260" DO-loop, page 88. The clock is updated and the new body temperature distributions are written out in the statements which immediately follow. Finally, the power input to the cavity by radiation from the FOV limiter, "QBOD", is computed between "246 CONTINUE" and "679 CONTINUE", page 88.

The statements after "679 CONTINUE" in which the cavity node temperatures are computed, are for the most part identical to those documented in Reference 1. The main differences are on page 91, where the heat input from the front end, "QBOD", is added ("IF (I.LE.4) Q2 = Q2 + QBOD"), and on page 93, where the heat conduction "QTOT" is buffered from the cavity to the body nodes ("IF(IJK.LE.4) QTOT(LL) = QTOT(LL) + ...").

#### 2.2.2. Modifications for the Addition of a Filter Dome

If a filter dome is installed, additional body nodes must be provided. The conductances and distribution factors associated with these new nodes must be calculated and included in the data files. The calculation of the distribution factors would be carried out in "TRW FRONT" as discussed in Section 2.1.3. However, the conductances must be provided by the user. What follows is a page-by-page outline of the modifications required for the inclusion of the filter dome.

1. Page 80: Increase all dimensions "104" to a number which includes the filter dome nodes, and change all dimensions "106" to this new number plus 2 (the number "106" presently refers to the cavity aperture).

Increase the dimension of "SOURCE" to
41 + n, where n is the number of filter
dome nodes.

Increase the value of "NBODY" to include the additional filter dome nodes.

- 2. Page 82: Change the upper limit of the "601" DO-loop to include the filter dome nodes (that is, increase 41 to 41 + n, where n is the total number of filter dome nodes).

Change "IF (J.EQ.41) JJ = 106" to "IF (J.EQ. number1) JJ = number2", where number1 = 41 + n (n being the number of filter dome nodes) and number2 = total number of body plus filter dome nodes + 2.

After "IF (J.GE.33.AND.J.LT.41) etc.", add a similar statement which converts the dome node position indices (established in the data set) to node numbers.

Change the upper limit in the two "401" DO-loops to the total number of body plus filter dome nodes.

4. Page 87: Between "220 CONTINUE" and "636 CONTINUE", add statements similar to those for the existing body nodes, for the filter dome nodes.

- 5. Page 88: Change the upper limit of the "246" and

  "661" DO-loops to the total number of body

  plus filter dome nodes.

  In the line above "661 CONTINUE", change

  "106" to the total number of body plus

  filter dome nodes + 2.
- 6. Pages 100,101 & 104:Add thermal physical properties and conductance data (user generated) for filter dome nodes.

# 3. Use of the Model to Establish a Real-Time Flight Data Interpretation Strategy

The "TRW Calibration Simulation Model" could be used to establish a real-time flight data interpretation strategy. The first step would be to adjust the model (with the filter dome modifications for the visible channel) until it is in good agreement with the actual instrument in the calibration environment. The next step would be to implement the modifications outlined above to convert the model into a flight simulator. Finally, the "calibrated" model would be used to establish an empirical relation between the actual source energy flux at the instrument, computed independently of the instrument response using, for example, GENDAT, the active cavity heater power signal, and one or more (the more the better) key instrument temperatures. It would be hoped that these "key instrument temperatures" would play only secondary roles in this relation. A logical and yet relatively simple form, consistent with Figs. 1 and 2 of the July, 1982, Progress Report, is

suggested by the superposition theorem $^3$ :

$$Q_{\text{heater}}(t) = A \int_{0}^{t} \frac{\partial}{\partial t} Q_{\text{source}}(t-\tau)d\tau + B \int_{0}^{t} \frac{\partial}{\partial t} T_{\text{HS}}(t-\tau)d\tau + C \int_{0}^{t} \frac{\partial}{\partial t} T_{\text{G}}(t-\tau)d\tau .$$

In this relation,  $Q_{\text{heater}}(t)$  is the instantaneous heater power (or at least the "count" which represents this power),  $Q_{\text{source}}(t-\tau)$  is the flux incident to the precision aperture from the source a time delay  $\tau$  earlier,  $T_{\text{HS}}(t-\tau)$  is the heat sink temperature a time delay  $\tau$  earlier, and  $T_{\text{C}}(t-\tau)$  is the temperature of the mounting beam a time delay  $\tau$  earlier. The sensitivity coefficients A, B and C would be obtained by exercising the calibrated model. This relation is inconvenient because  $Q_{\text{source}}$  appears under an integral. However, it should be possible to write an algorithm capable of inverting this integral in real time.

It is possible that a more complicated model involving more key temperatures will be required. In any case, the numerical model, modified and calibrated as described in this report, should prove very useful in defining this relation.

Once flight data are obtained and interpreted by some data interpretation strategy, the result can be introduced into the complete model to verify that the correct count is indeed obtained.

#### References

- 1. Eskin, L.D., Application of the Monte Carlo Method to the Transient Thermal Modeling of a Diffuse-Specular Radiometer Cavity, Master's Thesis, Virginia Polytechnic Institute and State University, Blacksburg, Virginia, August 1981.
- Mahan, J. R., and L. D. Eskin, "Application of Monte Carlo Techniques to Transient Thermal Modeling of Cavity Radiometers Having Diffuse-Specular Surfaces," Fourth American Meteorological Society Conference on Atmospheric Radiation, Toronto, Ontario, Canada, June 16-18, 1981.
- 3. Eckert, E. R. G., and R. M. Drake, Jr., Analysis of Heat and Mass Transfer (McGraw-Hill, New York, 1972), p. 170.

#### 5. APPENDIX

Transient Thermal Modeling of the Nonscanning ERBE Detector

TRW Calibration Simulation

USERS MANUAL

Final Revision: March 1983

#### 1. TRW SHAPE

This program subdivides a TRW-type cavity into NN2\*NTH elements, where NN2 is the number of divisions along the cavity axis and NTH is the number of circumferential divisions. When the cavity has been subdivided into elements, the node surface areas and thermal capacities and the internode thermal conductances are computed and stored in online disk files.

This program requires two previously generated data file areas on system disk. The first file name should be named A51961.DATA10 and be of sufficient size to hold M2 + 1 card images. This file will be written to as unit 10. The second file should be named A51961.DATA15 and should be of sufficient size to hold M2\*M2 + NN2 + 2 card images. This file will be written to as unit 15.

#### 1.1 Nomenclature

- ZTB Matrix of the z-coordinate for each node level top boundary. Must be dimensioned NN2.
- DZ Matrix of the z-coordinate width of each node level. Must be dimensioned NN2.
- AA Matrix of the node surface areas (m<sup>n</sup><sup>4</sup>2). Must be dimensioned M2.
- CC Matrix of the node thermal capacities (W-s/K). Must be dimensioned M2.
- VX(I,I) X-position of the Ith node in Cartesian coordinates.
- VY(I,I) Y-position of the Ith node in Cartesian coordinates.
- VZ(I,I) Z-position of the Ith node in Cartesian coordinates.
- VX(I,J) "Half-conductance" between nodes I and J where I > J.
- VY(J,I) "Half-conductance" between nodes I and J where J > I.

(Note: VX, VY, and VZ must all be dimensioned M2 by M2.)

- A Distance from plane of precision aperture to junction of barrel and cone (mm).
- B Distance from plane of precision aperture to apex of cone (mm).
- C Radius of barrel (mm).

(Note: For the purpose of stating these dimensions, it is assumed that there is no "space ring" gap between the precision aperture and the cavity opening.)

```
Slope of cone (-).
ALPHA
         Ratio of circumference to diameter of a circle (-).
PΙ
COND
         Thermal conductivity (W/m-K).
         Mass density (kg/m**3).
DENS
         Specific heat (W-s/kg-K).
SPEC
THCK
         Thickness (mm).
         Number of columns (or rows) in arrays VX, VY,
NARSZ
         and VZ. This must be the dimensioned value.
         The number of circumferential divisions desired.
NTH
         Floating point version of "NTH".
FNTH
         Number of axial divisions of barrel.
N 1
N2
         Number of axial divisions of cavity.
FN1, FN2 Floating point versions of N1 and N2.
         Total number of axial divisions through barrel.
NN 1
NN2
         Total number of axial divisions of cavity.
         Total number of cylinder nodes.
M 1
M2
         Total number of cavity nodes.
Ι
         Node number.
         Index of node level.
J
FJ
         Real value of J.
         Temporary value of z-coordinate of node level J.
TZ
         Temporary value of z-coordinate of node level J-1.
TZ 1
         Temporary value of z-coordinate of node level J+1.
TZ2
         "Half-conductance" to node to the "left" of a
GK 1
         node on level J.
         "Half-conductance" to node "above" a node on
GK2
         level J. i.e., on level J-1.
         "Half-conductance" to node to the "right" of a
GK3
         node on level J; note that GK3 = GK1.
         "Half-conductance" to node "below" a node on
GK4
         level J. i.e., on level J+1.
         Temporary z-coordinate of the boundary between
TZB1
         nodes on level J and nodes on level J-1.
         Temporary z-coordinate of the boundary between
TZB2
         nodes on level J and nodes on level J+1.
R
          Radial location of nodes at node level J.
          Radial position of boundary between nodes on
RB1
          level J and nodes on level J-1.
          Radial position of boundary between nodes on
RB2
          level J and nodes on level J+1.
AREA
          Area occupied by a node at node level J (m**2).
          Auxillary counter defined in "10" DO-loop used
 J
          to determine which segment of the cavity is
          being considered.
FJ
          Floating point version of J.
          Temporary z-coordinate of node; later converted
 TZ
          to a permanent value, VZ(I,I).
          Radial coordinate of node; later resolved into
 R
          x- and y-coordinates, VX(I,I) and VY(I,I).
          Circumferential angular location of node.
 THETA
```

#### 1.2 Description of Program

It is necessary to specify double precision when this program is compiled on IBM equipment. This step may not be necessary with other processors.

IMPLICIT REAL\*8 (A-H,O-Z)

Dimension subscripted variables. The dimensions shown permit the cavity to be divided into 100 nodes, with 10 divisions in each direction. Of course, these can be increased by the user if desired.

DIMENSION AA(100), CC(100)
DIMENSION VX(100,100), VY(100,100), VZ(100,100)
DIMENSION ZTB(10), DZ(10)

Establish COMMON for subroutines called.

COMMON /NODE1/ FN1,FN2,N1,N2,M1,M2 COMMON /NODE2/ FNTH,PI,NTH,NN1,NN2,MM2

Read in and write out the cavity dimensions and other constants.

READ(5,97) A,B,C,N1,N2,NTH,NSHOTN,ABS,REFR,DT,ELIMIT WRITE(6,96) A,B,C,N1,N2,NTH,NSHOTN,ABS,REFR,DT,ELIMIT

Establish the values of the physical constants.

COND = 406.64D0

ALPHA = C/(B-A)

SPEC = 234.04D0

THCK = 2.0D-1

PI = 3.1415926D0

DENS = 10524.15D0

The cavity is subdivided into NTH\*NN2 elements, or nodes, in this portion of the program.

NARSZ = 100

FNTH = NTH

FN1 = N1

FN2 = N2

NN1 = N1

NN2 = NN1 + N2

M1 = NTH\*NN1

M2 = NTH\*NN2

MM2 = M2 + 1

I = 0

In the "10" DO-loop we compute temporary z-coordinates and the r-coordinate of each node. We use computed GO TO statements to direct the calculation to the proper relations depending on which segment of the cavity is being considered (as determined by the value of "J").

DO 10 J = 1, NN2

Calculation of the node coordinates, area, thermal capacity and conductances to surrounding nodes is different depending on the surface on which the node is located. Thus we must branch to the appropriate segment of the program.

If the node lies on the barrel, branch to the barrel calculation segment.

IF(J.LE.NN1) GOTO 11

If the node lies on the cone, branch to the "cone" calculation segment.

GOTO 12

The node is on the barrel; thus we compute the node coordinates, area, and half-conductances for the barrel.

```
11 CONTINUE

FJ = J

TZ = A*(2.0D0*FJ - 1.0D0)/(2.0D0*FN1)

IF(J.GT.NN2) GOTO 900

DZ(J) = A/FN1

ZTB(J) = TZ - DZ(J)/2.0D0

900 CONTINUE

R = C

AREA = 2.0D-6*PI*C*A/(FN1*FNTH)

GK1 = A*COND*THCK*1.0D-3*FNTH/(PI*R*FN1)

GK2 = COND*PI*R*THCK*4.0D-3/(FNTH*A)

GK3 = GK1

GK4 = GK2

IF(J.EQ.1) GK2 = 0.0D0

GOTO 17
```

The node is on the cone; thus we compute the coordinates, area, and half-conductances for the cone.

```
12 CONTINUE -
    FJ = J
    TZ = 2*(FN1 + FN2 - FJ) + 1.0D0
    TZ = (A-B)*DSQRT(TZ)/DSQRT(2.0D0*FN2) + B
    R = ALPHA*(B-TZ)
    AREA = PI*C*1.0D-6*DSQRT(C*C + (B-A)*(B-A))
    AREA = AREA/(FNTH*FN2)
    TZ1 = 2*(FN1 + FN2 - FJ)
    TZ1 = (A - B)*DSQRT(TZ1)/DSQRT(2.0D0*FN2)
    TZ1 = TZ1 + B
    IF(J.EQ.NN2) GOTO 800
    TZ2 = 2*(FN1 + FN2 - FJ) - 2.0D0
    TZ2 = (A - B)*DSQRT(TZ2)/DSQRT(2.0D0*FN2)
    TZ2 = TZ2 + B
1800 \text{ TZB2} = (A - B) *DSQRT((FN1+FN2-FJ)/FN2) + B
    TZB1 = (A - B)*DSQRT((FN1+FN2-FJ+1.0D0)/FN2)
    TZB1 = TZB1 + B
    IF(J.GT.NN2) GOTO 901
    ZTB(J) = TZB1
    DZ(J) = TZB2 - TZB1
1901 CONTINUE
1802 \text{ GK1} = (TZB2-TZB1)*(TZB2-TZB1)
    GK1 = GK1 + ALPHA*ALPHA*(TZB2-TZB1)*(TZB2-TZB1)
     GK1 = COND*THCK*FNTH*DSQRT(GK1)*1.0D-3/(PI*R)
     GK2 = (TZ-TZB1)*(TZ-TZB1)
     GK2 = GK2 + ALPHA*ALPHA*(TZ-TZB1)*(TZ-TZB1)
     GK2 = COND*PI*THCK*2.0D-3*ALPHA*(B-TZB1)
     GK2 = GK2/(DSQRT(GK2)*FNTH)
     GK3 = GK1
     IF(J.NE.NN2) GOTO 803
     GK4 = 0.0D0
     GOTO 17
1803 \text{ GK4} = (TZB2-TZ)*(TZB2-TZ)
     GK4 = GK4 + ALPHA*ALPHA*(TZB2-TZ)*(TZB2-TZ)
     GK4 = COND*PI*THCK*2.0D-3*ALPHA*(TZB2-TZ)
     GK4 = GK4/(DSQRT(GK4)*FNTH)
     GOTO 17
```

This segment of the program is common to both parts of the cavity, barrel and cone. In this section, the coordinates, half-conductances, and areas are stored, and the thermal capacities are computed and stored.

```
17 CONTINUE
    DO 10 K = 1. NTH
    FK = K
    I = I + 1
    VZ(I,I) = TZ
    THETA = 6.283185306D0*FK/FNTH
    VX(I,I) = R*DCOS(THETA)
    VY(I,I) = R*DSIN(THETA)
    IF(I.GT.M2) GOTO 936
    AA(I) = AREA
    CC(I) = AREA*THCK*DENS*SPEC*1.0D-3
1936 CONTINUE
    IF(J.GT.NN2) GOTO 10
    ICK = I - 1
    ICK = ICK/NTH
    ICK = I - ICK*NTH
    IF(ICK.EQ.1) GOTO 362
    VX(I,I-1) = GK1
    GOTO 363
1362 CONTINUE
    VX(I+NTH-1.I) = GK1
363 CONTINUE
    IF(I.LE.NTH) GOTO 364
    VX(I,I-NTH) = GK2
1364 CONTINUE
    IF(ICK.EQ.NTH) GOTO 365
    VY(I+1,I) = GK3
    GOTO 366
1365 CONTINUE
    VY(I,I-NTH+1) = GK3
1366 CONTINUE
    IF(I.GT.(M2-NTH)) GOTO 10
    VY(I+NTH,I) = GK4
 10 CONTINUE
```

(Note: The conductance between node I and adjacent node J is given by VX(I,J)\*VY(I,J)/(VX(I,J) + VY(I,J)), where I > J; that is, series conductances add like parallel resistances.)

Write out results.

```
DO 201 I = 1, M2
| WRITE(10,2) I, AA(I), CC(I)
|201 CONTINUE
```

Create a trip card image.

```
ISTAR = 0
AASTAR = 0.0D0
CCSTAR = 0.0D0
WRITE(10,2) ISTAR, AASTAR, CCSTAR
```

Continue writing.

```
DO 202 I = 1, M2
DO 202 J = 1, M2
WRITE(15,1) I, J, VX(I,J), VY(I,J), VZ(I,J)
202 CONTINUE
```

Create another trip card image.

```
I = 0
J = 0
VDUMMY = 0.0D0
WRITE(15,1) I, J, VDUMMY, VDUMMY
```

And write some more.

```
DO 203 J = 1, NN2
WRITE(15,2) J, ZTB(J), DZ(J)
203 CONTINUE
```

And one last trip card image.

```
J = 0.0D0
ZDUMMY = 0.0D0
WRITE(15,2) J, ZDUMMY, ZDUMMY
```

Subroutine POSMTX is a subroutine which writes out the node position matrix on the line printer so that the values can be verified by the user. Thus, it is optional and does not have to be called.

```
CALL POSMTX(VX, VY, VZ, NARSZ)
```

#### Format statements:

End of Program TRW SHAPE.

STOP END

#### 1.3. Subroutine 'POSMTX'

This subroutine rearranges the node coordinate results and writes them out with the appropriate labels so that they can be checked by the user. Use of this subroutine is optional.

```
SUBROUTINE POSMTX(VX, VY, VZ, NARSZ)
    IMPLICIT REAL*8 (A-H,O-Z)
    DIMENSION VX(NARSZ, NARSZ), VY(NARSZ, NARSZ)
    DIMENSION VZ(NARSZ, NARSZ)
    COMMON /NODE1/ FN1,FN2,N1,N2,M1,M2
    COMMON /NODE2/ FNTH, PI, NTH, NN1, NN2, MM2
    WRITE(6,1)
    DO 100 I = 1, M2
    R = DSQRT(VX(I,I)*VX(I,I) + VY(I,I)*VY(I,I))
    PHI = DARCOS(VX(I,I)/R)
    IF(VY(I,I).LT.0.0D0) PHI = 2.0D0*PI - PHI
    WRITE(6,2) I, VX(I,I), VY(I,I), VZ(I,I), R, PHI
1100 CONTINUE
     RETURN
   1 FORMAT('1',5X'NODE',T20,'X-COORDINATE',T40,
            'Y-COORDINATE', T60, 'Z-COORDINATE', T83,
            'RADIUS',T103,'ANGLE'//)
  2 FORMAT(6X, I3, T18, D14.5, T38, D14.5, T58, D14.5,
            T80, D14.5, T100, D14.5)
```

End of Subroutine 'POSMTX'.

1		i i
1		
ı	END	
1	END	
1		
1		

#### 2. TRW NODES

This program uses Monte Carlo techniques to compute the distribution factors for the radiation emitted from the cavity nodes for a TRW-type cavity radiometer. The "photon" specular reflections are computed "exactly". Program 'TRW SHAPE' must be executed before this program is submitted.

This program requires one previously generated data file area on system disk. The file should be named A51961.DATA12 and should be of sufficient size to hold (M2+1)\*NN2 + 1 card images. This file will be written to as unit 12.

#### 2.1. Nomenclature

DF	Matrix of distribution factors. Must be dimensioned M2 + 1. DF is written out for
	each source node.
XPOS	Matrix of node x-coordinates. Must be dimensioned M2.
YPOS	Matrix of node y-coordinates. Must be dimensioned M2.
ZPOS	Matrix of node z-coordinates. Must be dimensioned M2.
ZTB	Matrix of node level upper boundary z-coordinates. Must be dimensioned NN2.
DZ	Matrix of node level z-coordinate widths. Must be dimensioned NN2.
Z	Matrix of random numbers generated by 'GGUBS', a subroutine called by Subroutine 'RAND'.
A	Distance from plane of precision aperture to
В	junction of barrel and cone (mm). Distance from plane of precision aperture to
С	apex of cone (mm). Radius of barrel (mm).

(Note: For the purpose of stating these dimensions it is assumed that there is no "space ring" gap between the precision aperture and the cavity opening.)

Slope of cone. ALPHA Number of packets into which energy is divided NSHOTS for the Monte Carlo procedure. Real value of nshots. **RSHOTS** Total number of cavity nodes. M2 Number of circumferential divisions. NTH Ratio of circumference of circle to its diameter. PΙ Index of aperture. MM2 Seed number needed for random number generator. SEED Absorptivity. ABS Ratio of specular component of reflectivity to REFR diffuse + specular (= total) reflectivity. X-coordinate of the source position. TXI Y-coordinate of the source position. TYI Z-coordinate of the source position. TZI X-coordinate of the previous source position. TXIOLD Y-coordinate of the previous source position. TYIOLD Z-coordinate of the previous source position. TZIOLD X-coordinate of the receiving position. TXJ Y-coordinate of the receiving position. TYJ Z-coordinate of the receiving position. TZJ

(Note: All of the above coordinates (TXI, TYI,..., TZJ) advance with each reflection. Initially TXI, TYI, and TZI are used in a test to determine the first trip through position (TXI,TYI,TZI) by a given "photon".)

NSD Test variable whose value is set to "1" when a specular reflection occurs and to "2" when a diffuse reflection occurs.

#### 2.2. Description of Program

IBM computers require double precision to execute this program. Other compilers may not need this first step:

IMPLICIT REAL\*8 (A-H, O-Z)
REAL\*4 RN

Dimension subscripted variables. The subscript range shown below permits the cavity to be divided into up to 100 nodes, with up to 10 divisions in each direction.

DIMENSION DF(101), ZTB(10), DZ(10) DIMENSION XPOS(100), YPOS(100), ZPOS(100) Establish variables common to subroutines called.

```
COMMON /NODE1/ FN1,FN2,N1,N2,NN1
COMMON /NODE2/ FNTH,NTH,NN2,M1,M2,MM2
COMMON /GEOM/ A,B,C,ALPHA,PI
```

Read in and write out cavity dimensions and property values.

READ(5,97) A,B,C,N1,N2,NTH,NSHOTS,ABS,REFR,DT,ELIMIT WRITE(6,96) A,B,C,N1,N2,NTH,NSHOTS,ABS,REFR,DT,ELIMIT

Establish values of constants.

```
ALPHA = C/(B-A)
PI = 3.141592D0
FNTH = NTH
FN1 = N1
FN2 = N2
NN1 = N1
NN2 = NN1 + N2
M1 = NN1*NTH
M2 = NN2*NTH
MM2 = M2 + 1
RSHOTS = NSHOTS
SEED = 1234567.D0
RN = -1.0D0
```

Read in values of x-, y- and z-positions of nodes; store as XPOS, YPOS and ZPOS, respectively.

```
11 CONTINUE

READ(15,1) I, J, XX, YY, ZZ

IF(I.NE.J) GOTO 11

IF(I.EQ.O) GOTO 12

XPOS(J) = XX

YPOS(J) = YY

ZPOS(J) = ZZ

GOTO 11

12 CONTINUE
```

Read in values of ZTB and DZ.

In the "20" DO-loop we loop through the first nodes of each ring of cavity nodes, treating each as a diffuse source. The distribution factors not computed directly in this loop can be determined when needed from symmetry.

```
I = MM2
DO 20 J = 1, M2, NTH
```

Zero the 'DF' vector.

```
DO 9 I1 = 1, MM2
DF(I1) = 0.0D0
9 CONTINUE
```

In the "10" DO-loop we allow node J to emit NSHOTS volleys of energy packets and follow them through the cavity until they are either absorbed or leave the cavity through the aperture.

```
DO 10 ISHOT = 1, NSHOTS
    CALL RAND (SEED, IN, RN)
    JL = J - 1
    JL = JL/NTH + 1
    TZI = ZTB(JL) + RN*DZ(JL)
    TZIOLD = TZI
    TZJ = TZI
    CALL RAND(SEED, IN, RN)
    PHII = (2.0D0*RN + 1.0D0)*PI/FNTH
    IF(J.GT.M1) GOTO 902
    R = C
    GOTO 903
|902 R = ALPHA*(B - TZI)
1903 CONTINUE
    TXI = R*DCOS(PHII)
    TXIOLD = TXI
    TXJ = TXI
    TYI = R*DSIN(PHII)
    TYIOLD = TYI
    TYJ = TYI
```

We return to this junction ("13 CONTINUE") each time a reflection occurs.

```
13 CONTINUE
```

If this is the first time this volley has passed through J, the volley striking node J is "emitted" diffusely.

```
IF((TXI.EQ.TXIOLD).AND.(TYI.EQ.TYIOLD)
& .AND.(TZI.EQ.TZIOLD)) GOTO 137
```

Test to see if the energy packet is absorbed or reflected. If the random number RN is less than or equal to ABS, the absorptivity, the packet is absorbed; otherwise it is reflected.

```
CALL RAND(SEED, IN, RN)
IF(RN.LE.ABS) GOTO 16
```

Test to see if reflected energy packet is specularly reflected or diffusely reflected. If the random number RN is less than or equal to REFR, the reflectivity ratio, the packet is specularly reflected; otherwise it is diffusely reflected.

```
CALL RAND(SEED, IN, RN)
NSD = 1
IF(RN.LE.REFR) GOTO 17
```

Randomly select the direction of the diffuse reflection.

```
| 137 CONTINUE | CALL RAND(SEED, IN, RN) | IF(RN.EQ.1.0D0) GOTO 14 | THETA = ARSIN(SQRT(RN)) | GOTO 15 | 14 CONTINUE | THETA = 0.0D0 | 15 CONTINUE | CALL RAND(SEED, IN, RN) | PHI = 2.0D0*PI*RN | NSD = 2 | GOTO 17 | 115 CONTINUE |
```

Compute the x-, y-, and z-components of the unit vector in the direction (THETA, PHI) with respect to the central coordinate system.

CALL VECTOR (UNX, UNY, UNZ, PHI, THETA, VOX, VOY, VOZ)

Call the search subroutine which identifies the receiving point of the diffusely reflected photon.

CALL SEARCH(TXI, TYI, TZI, TXJ, TYJ, TZJ, VOX, VOY, VOZ)

If the z-coordinate of the receiving position is negative, the photon has escaped.

IF(TZJ.LT.O.ODO) GOTO 24

The photon is reflected to the next position whose coordinates are (TXJ, TYJ, TZJ).

**GOTO 25** 

The photon is absorbed. Call Subroutine NODE which determines the absorbing node.

16 CONTINUE

CALL NODE (ZTB, TXJ, TYJ, TZJ, J1, JJ1, JJ2)

DF(J1) = DF(J1) + 1.0D0

GOTO 10

17 CONTINUE

Determine the inward-directed unit normal vector of the source position. This depends on the shape of the source surface.

IF(TZI.LE.A) GOTO 19

The reflecting surface is the cone.

```
18 CONTINUE

UN = ALPHA*ALPHA*ALPHA*ALPHA

UN = UN*(B - TZI)*(B - TZI)

UN = UN + TXI*TXI + TYI*TYI

UN = DSQRT(UN)

UNX = - TXI/UN

UNY = - TYI/UN

UNZ = - ALPHA*ALPHA*(B - TZI)/UN

GOTO 23
```

The reflecting surface is the barrel.

```
19 CONTINUE

UNX = - TXI/C

UNY = - TYI/C

UNZ = 0.0D0

23 CONTINUE
```

If reflection is diffuse return to the diffuse sequence; otherwise continue in the specular sequence.

```
IF(NSD.EQ.2) GOTO 115
```

Compute the "ideal" vector of the reflected packet.

```
VDOT = (TXI-TXIOLD)*UNX

VDOT = VDOT + (TYI-TYIOLD)*UNY

VDOT = VDOT + (TZI-TZIOLD)*UNZ

VOX = TXI - TXIOLD - 2.0D0*VDOT*UNX

VOY = TYI - TYIOLD - 2.0D0*VDOT*UNY

VOZ = TZI - TZIOLD - 2.0D0*VDOT*UNZ

VMAG = DSQRT(VOX*VOX + VOY*VOY + VOZ*VOZ)

VOX = VOX/VMAG

VOY = VOY/VMAG
```

Call the search subroutine which identifies the receiving point of the "exactly" reflected photon.

```
CALL SEARCH(TXI, TYI, TZI, TXJ, TYJ, TZJ, VOX, VOY, VOZ)
```

If the z-coordinate of the receiving position is negative, the photon has escaped.

IF(TZJ.LT.0.0D0) GOTO 24

The photon is reflected to the next position whose coordinates are (TXJ, TYJ, TZJ).

GOTO 25

The packet has left the cavity through the aperture. Record and fire another volley.

24 CONTINUE DF(MM2) = DF(MM2) + 1.0D0 GOTO 10

A reflection to another surface has occurred. Thus, increment TXI, TYI, TZI, TXIOLD, TYIOLD, TZIOLD, TXJ, TYJ, and TZJ.

25 CONTINUE TXIOLD = TXI TYIOLD = TYI TZIOLD = TZI TXI = TXJ TYI = TYJ TZI = TZJ

Loop back to test sequence to determine disposition of photon when it arrives at this new surface.

GOTO 13 10 CONTINUE Write vector of distribution factors corresponding to node  $J_{\:\raisebox{1pt}{\text{\circle*{1.5}}}}$ 

```
JJ = (J + NTH - 1)/NTH
DO 30 JK = 1, MM2
DF(JK) = DF(JK)/RSHOTS
WRITE(12,3) JJ, JK, DF(JK)
30 CONTINUE
20 CONTINUE
```

Generate a "trip" card image at the end of the file.

```
JJ = 0

JK = 0

DFDUM = 0.0D0

WRITE(12,3) JJ, JK, DFDUM
```

#### Format statements:

```
1 FORMAT(5X, 215, 3D14.5)
2 FORMAT(8X, I3)
 3 FORMAT(5X, 2I5, D14.5)
4 FORMAT(5X, I5, 2D14.5)
96 FORMAT(5X, 'THE SYSTEM VARIABLES ARE AS ',
          'FOLLOWS:' ,//10X,'A = ',F5.2 ,' MM',
          /,10X,'B = ',F5.2,' MM',/,10X,'C = ',
          F5.2, 'MM', /, 10X, 'N1 = ', I2, /, 10X, 'N2 '
  &
          ,'= ',I2,/,10X,'NTH =',I2,/,10X,'NSHOTN = '
          ,16,/,10X,'ABS = ',F3.1,/,10X,'REFR = '
          .F3.1,/,10X,'DT = ',F4.1,/,10X,'ELIMIT = '
  &
          ,E8.2///)
97 FORMAT(F5.2,F5.2,F5.2,I2,I2,I2,I6,F3.1,F3.1,F4.1,
          E8.2,4I1)
```

End of Program 'TRW NODES'.

```
STOP
END
```

#### 2.3. Subroutine SEARCH

This subroutine calculates the magnitude of the vector from the source position to the receiving position of the reflected photon. Since the direction of this vector is known (VOX, VOY, VOZ), the magnitude is all that is needed to calculate the receiving position.

SUBROUTINE SEARCH(TXI, TYI, TZI, TXJ, TYJ, TZJ, VOX, VOY, VOZ)

When this subroutine is compiled on IBM equipment, double precision must be specified. This step may not be necessary on other processors.

IMPLICIT REAL\*8 (A-H,O-Z)

Establish COMMON variables.

COMMON /GEOM/ A,B,C,ALPHA,PI

If the source position is on the cone and VOZ is positive, search for the receiving position on the cone. If the z-component of the direction unit vector is zero or positive, the escape check may be skipped. Otherwise, begin the search on the aperture, proceeding to the cylinder and cone as necessary.

IF((TZI.GE.A).AND.(VOZ.GE.O.ODO)) GOTO 10

Check to see if the photon has escaped.

IF(VOZ.GE.O.ODO) GOTO 5

VMAG1 = - TZI/VOZ

TXJ = TXI + VMAG1\*VOX

TYJ = TYI + VMAG1\*VOY

RADIUS = DSQRT(TXJ\*TXJ + TYJ\*TYJ)

IF(RADIUS.GE.C) GOTO 5

TZJ = -1.0D0

RETURN

Begin the search for the receiving position on the cylinder.

```
5 A1 = VOX*VOX + VOY*VOY
B1 = 2.0D0*(TXI*VOX + TYI*VOY)
C1 = TXI*TXI + TYI*TYI - C*C
VMAG2 = (-B1 + DSQRT(B1*B1 - 4.0D0*A1*C1))
VMAG3 = VMAG2/(2.0D0*A1)
VMAG3 = (-B1 - DSQRT(B1*B1 - 4.0D0*A1*C1))
VMAG3 = VMAG3/(2.0D0*A1)
VMAGB = VMAG2
IF(VMAGB.LE.0.0D0) VMAGB = VMAG3
TXJ = TXI + VMAGB*VOX
TYJ = TYI + VMAGB*VOZ
```

Check to see if the receiving position calculated is on the barrel. If not, continue to search on the cone.

```
IF(TZJ.LE.A) RETURN
```

The receiving position is on the cone.

```
10 A2 = VOX*VOX+VOY*VOY-ALPHA*ALPHA*VOZ*VOZ

B2 = ALPHA*ALPHA*VOZ*(B-TZI)

B2 = B2 + TXI*VOX + TYI*VOY

B2 = B2*2.0D0

C2 = 2.0D0*B*TZI - TZI*TZI - B*B

C2 = C2*ALPHA*ALPHA

C2 = C2 + TXI*TXI + TYI*TYI

VMAG4 = - B2 + DSQRT(B2*B2 - 4.0D0*A2*C2)

VMAG5 = - B2 - DSQRT(B2*B2 - 4.0D0*A2*C2)

VMAG5 = VMAG5/(2.0D0*A2)

VMAG5 = VMAG5/(2.0D0*A2)

VMAGC = VMAG4

IF(VMAGC.LE.0.0D0) VMAGC = VMAG5
```

The equation for the magnitude is quadratic, giving two solutions. The zero or negative root is ignored.

Calculate the receiving position on the cone.

```
TXJ = TXI + VMAGC*VOX
TYJ = TYI + VMAGC*VOY
TZJ = TZI + VMAGC*VOZ
RETURN
END
```

## 2.4. Subroutine VECTOR

This subroutine converts the randomly selected diffuse angle with respect to the source position into the central coordinate system.

### 2.4.1. Nomenclature

UNX	X-component of inward-directed unit normal vector of source position.
UNY	Y-component of inward-directed unit normal vector of source position.
UNZ	Z-component of inward-directed unit normal vector of source position.
UTX	X-component of unit tangent vector on source position.
UTY	Y-component of unit tangent vector on source position.
UTZ	Z-component of unit tangent vector on source position.
USX	X-component of the mutually orthogonal vector.
USY	Y-component of the mutually orthogonal vector.
USZ	Z-component of the mutually orthogonal vector.
THETA	Angle of diffuse vector with respect to normal vector.
PHI	Angle of diffuse vector with respect to tangent vector.
VOX	X-component of diffuse vector with respect to central coordinate system.
YOY	Y-component of diffuse vector with respect to central coordinate system.
VOZ	Z-component of diffuse vector with respect to central coordinate system.

#### 2.4.2. Subroutine Description

SUBROUTINE VECTOR(UNX, UNY, UNZ, PHI, THETA, VOX, VOY, VOZ)

When this subroutine is compiled on IBM equipment, it is necessary to specify double precision. This step may not be necessary with other compilers.

IMPLICIT REAL\*8 (A-H, O-Z)

Compute the x-, y- and z-components of the unit tangent vector, UT, and the mutually orthogonal vector, US, on the source position.

If the surface is cylindrical, go to 11.

```
IF(UNZ.EQ.0.0D0) GOTO 11
```

Calculations for the general surface.

```
UTX = UNX*DSQRT(UNZ*UNZ/(1.0D0 - UNZ*UNZ))

UTY = UNY*DSQRT(UNZ*UNZ/(1.0D0 - UNZ*UNZ))

UTZ = DSQRT(1.0D0 - UNZ*UNZ)

GOTO 13
```

Calculations for the cylindrical surfaces.

```
| 11 CONTINUE
| UTX = 0.0D0
| UTY = 0.0D0
| UTZ = 1.0D0
| 13 CONTINUE
```

Compute the x-, y- and z-components of the unit vector on the source position which is mutually orthogonal to UN and UT.

```
USX = UNY*UTZ - UNZ*UTY
USY = UNZ*UTX - UNX*UTZ
USZ = 0.0D0
```

Compute components of the vector at angle (THETA, PHI) with respect to the UN,UT,US-coordinate system and express the result in the I,J,K-coordinate system.

```
SINTH = DSIN(THETA)

COSTH = DCOS(THETA)

SINPH = DSIN(PHI)

COSPH = DCOS(PHI)

VOX = SINTH*COSPH*UTX + SINTH*SINPH*USX

VOX = VOX + COSTH*UNX

VOY = SINTH*COSPH*UTY + SINTH*SINPH*USY

VOY = VOY + COSTH*UNY

VOZ = SINTH*COSPH*UTZ + SINTH*SINPH*USZ

VOZ = VOZ + COSTH*UNZ
```

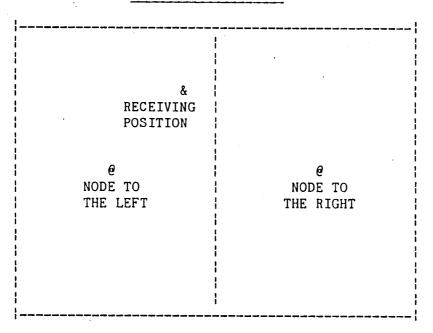
# End of Subroutine 'VECTOR'.

RETURN	
END	

#### 2.5. Subroutine NODE

This subroutine identifies the node which absorbs a given photon.  $% \left( 1\right) =\left( 1\right) \left( 1\right) +\left( 1\right) \left( 1\right) \left( 1\right) +\left( 1\right) \left( 1\right)$ 

2.5.1. Nomenclature



PHI	Circumferential angle of the receiving position.
PHILE	Circumferential angle of the left edge of the
	node to the right.
NRA	Number of the node ring "above" the receiving position.
NRB	Number of the node ring "below" the receiving position.
NLP	Position of the node to the left.
NRP	Position of the node to the right.

2.5.2. Subroutine Description

SUBROUTINE NODE(ZTB, TXJ, TYJ, TZJ, J1, JJ1, JJ2)

When this subroutine is compiled on IBM equipment, it is necessary to specify double precision. This step may not be necessary on other compilers.

IMPLICIT REAL\*8 (A-H, O-Z)

Dimension the subscripted variables.

DIMENSION ZTB(NN2)

Establish COMMON variables.

COMMON /NODE1/ FN1,FN2,FN3,N1,N2,N3,NN1
COMMON /NODE2/ FNTH,NTH,NN2,NN3,M1,M2,M3,MM2
COMMON /GEOM/ A,B,C,ALPHA,PI

Determine which surface the receiving position lies on. If the receiving position lies on the cone, jump to the cone calculation segment.

IF(TZJ.GT.A) GOTO 11

The receiving position is on the cylinder.

PHI = DAR OS(TXJ/C)IF(TYJ.LT.0.0D0) PHI = 2.0D0\*PI - PHINRA = (1.0D0 + 2.0D0\*FN1\*TZJ/A)/2.0D0NRB = NRA + 1NRP = FNTH\*PHI/(2.0DO\*PI)IF(NRP.EQ.O) NRP = NTHNLP = NRP + 1IF(NLP.EQ.(NTH+1)) NLP = 1PHILE = (2.0D0\*NRP + 1.0D0)\*PI/FNTHIF(NRP.EQ.NTH) PHILE = PHILE - 2.0D0\*PI JJ2 = NRPIF(PHI.GT.PHILE) JJ2 = NLP JJ1 = NRAIF(NRA.NE.O) GOTO 998 JJ1 = NRBGOTO 999 1998 IF(TZJ.GT.ZTB(NRB)) JJ1 = NRB1999 J1 = (JJ1 - 1.0D0)\*NTH + JJ2RETURN

The receiving position is on the cone.

```
11 R = DSQRT(TXJ*TXJ + TYJ*TYJ)
    PHI = DAR OS(TXJ/R)
    IF(TYJ.LT.0.0D0) PHI = 2.0D0*PI - PHI
    FNRA = (TZJ-B)*(DSQRT(2.0DO*FN2))/(A-B)
    FNRA = FNRA*((TZJ-B)*(DSQRT(2.0D0*FN2))/(A-B))
    FNRA = FN1 + FN2 - (FNRA-1.0D0)/2.0D0
    NRA = FNRA
    NRB = NRA + 1
    NRP = FNTH*PHI/(2.0DO*PI)
    IF(NRP.EQ.O)NRP = NTH
    NLP = NRP + 1
    IF(NLP.EQ.(NTH+1)) NLP = 1
    PHILE = (2.0D0*NRP + 1.0D0)*PI/FNTH
    IF(NRP.EQ.NTH) PHILE = PHILE - 2.0DO*PI
    JJ2 = NRP
    IF(PHI.GT.PHILE) JJ2 = NLP
    JJ1 = NRA
    IF(NRA.EQ.NN2) GOTO 997
    IF(TZJ.GT.ZTB(NRB)) JJ1 = NRB
|997 J1 = (JJ1 - 1.0D0)*NTH + JJ2
```

End of Subroutine 'NODE'.

RETURN END

#### 2.6. Subroutine RAND

This subroutine uses library Subroutine GGUBS to obtain random numbers in an efficient way. A sequence of 100 random numbers is generated each time the 100th number from the previous generation is used.

### 2.6.1. Nomenclature

SEED A seed number required by GGUBS. It must be specified in the main program as a double precision integer value.

An integer between 1 and 100 inclusive which represents which of the 100 random numbers of

a given generation is obtained.

RN . The random number passed back to MAIN.

## 2.6.2. Subroutine Description

SUBROUTINE RAND(SEED, IN, RN)

IN

When this subroutine is compiled on IBM equipment, it is necessary to specify double precision. This step may not be necessary with other processors.

IMPLICIT REAL\*8 (A-H,O-Z)
REAL\*4 RN, Z(100)

If this is the first call to RAND, RN is equal to the negative value set in MAIN. In this case we must call GGUBS to generate the first generation of 100 random numbers.

IF(RN.LT.O.ODO) GOTO 11

If this is not the first call to RAND, then increment to the next value in this generation of 100 random numbers.

IN = IN + 1

If we have not yet used up our allotment of random numbers from this generation, we simply take the next one in the sequence back to MAIN.

IF(IN.LE.100) GOTO 13

However, if we have used all the members of this generation, we must create a new generation of 100 random numbers.

11 CONTINUE IN = 1 DSEED = SEED NUMBER = 100 CALL GGUBS(DSEED, NUMBER, Z) SEED = DSEED 13 CONTINUE RN = Z(IN)

End of Subroutine ! RAND'.

RETURN END

#### 3. TRW FRONT

This program uses Monte Carlo techniques to compute the distribution factors for the radiation emitted from the 'front end' nodes for a TRW-type cavity radiometer. The specular reflections are computed "exactly".

This program requires one previously generated data file area on system disk. The file should be named A51961.DATA17 and should be of sufficient size to hold (M2+1)\*NN2 + 1 card images. This file will be written to as unit 17, and will contain the distribution factor matrix after the program has been executed.

#### 3.1. Nomenclature

DF	Matrix of distribution factors. Must be dimensioned M2 + 1. DF is written out for
	each source node.
Н1	Distance to floor, or face, of instrument (cm).
п і Н2	Distance to face of F-O-V limiter (cm).
R1	Radius which describes the inside face of the
N I	F-O-V limiter; also outside radius of the face
	of the instrument (cm).
R2	Inside radius of the face of the instrument (cm).
R3	Radius of precision aperture (cm).
ης <b>F1</b>	Radial distance to the inner edge of the first
r i	floor node (cm).
F2	Radial distance to the inner edge of the second
r 2	floor node (cm).
F3	Radial distance to the inner edge of the third
r 3	floor node (cm).
FOV1	Axial distance from the floor level (Z = H1) to
1011	the lower edge of the top F-O-V limiter
	node (cm).
FOV2	Axial distance from the floor level (Z = H1) to
	the lower edge of the central F-O-V limiter
	node (cm).
EFOV	Emissivity (= absorptivity = 1 - reflectivity)
	of the face of the F-O-V limiter (-).
RFOV	Ratio of specular to specular + diffuse
	reflectivity for the face of the F-O-V limiter (-).
EF	Emissivity (= absorptivity = 1 - reflectivity)
	of the face of the instrument (-).
RF	Ratio of specular to specular + diffuse
	reflectivity for the face of the instrument (-).
NSHOTS	Number of "photons" emitted in Monte Carlo method.
RSHOTS	Real value of "NSHOTS".

(Note: The distances H1, H2, R1, R2, R3, and R4 are all measured with respect to the origin of coordinates (0,0,0) located on the axis in the plane of the precision aperture. See Fig. 1 of the February Progress Report.)

ΡΙ	Ratio of circumference to diameter of a circle (-).
NTH	Number of circumferential divisions (set at 4).
FNTH	Floating point version of NTH.
N 1	Number of axial divisions on F-O-V limiter (-).
	(set at 3).
FN1	Floating point version of N1.
N2	Number of radial divisions on floor (set at 4).
FN2	Floating point version of N2.
NN2	Number of axial divisions on F-O-V limiter
	plus number of radial divisions on floor (-).
NN 3	Total number of "front end" nodes divided by NTH.
M1	Total number of nodes on F-O-V limiter (-).
M2	Total number of nodes on the face of the
	instrument including $F=0=V$ limiter and floor $(-)$ .
М3	Total number of nodes (= $M1 + M2 + 3*NTH$ ),
•	including the edge of the floor, the top surface
	of the precision aperture, and the imaginary pie
	shaped segment in the plane of the F-O-V limiter.
SEED	Seed number needed for the random number generator.
RN.	Uniformly distributed random number between O
	and 1 generated by Subroutine RAND.
MM3	Total number of possible receiving nodes,
	including cavity opening.
TXI	X-coordinate of the source position.
TYI	Y-coordinate of the source position.
TZI	Z-coordinate of the source position.
TXIOLD	X-coordinate of the previous source position.
TYIOLD	Y-coordinate of the previous source position.
TZIOLD	Z-coordinate of the previous source position.
TXJ	X-coordinate of the receiving position.
TYJ	Y-coordinate of the receiving position.
TZJ	Z-coordinate of the receiving position.

## 3.2. Program Description

When this program is compiled on IBM equipment, double precision must be specified. This step may not be necessary on other processors.

```
IMPLICIT REAL*8 (A-H, O-Z)
REAL*4 RN
Dimension the subscripted variable.
```

Identify the COMMON variables for the subroutines called.

COMMON H1, H2, R1, R2, R3, FOV1, FOV2, F1, F2, F3, PI

Read in and write out values of front end dimensions and optical properties.

READ(5,97) NSHOTS,H1,H2,R1,R2,R3,F1,F2,F3,F0V1,F0V2, & EFOV,RFOV,EF,RF WRITE(6,96) H1,H2,R1,R2,R3,F1,F2,F3,F0V1,F0V2, & EFOV,RFOV,EF,RF,NSHOTS

Set values of constants.

PI = 3.141592D0NTH = 4FNTH = NTHN1 = 3FN1 = N1N2 = 4FN2 = N2NN2 = N1 + N2NN3 = NN2 + 3M1 = N1\*NTHM2 = NN2\*NTHM3 = NN3\*NTHMM3 = M3 + 1RSHOTS = NSHOTS SEED = 1234567.D0RN = -1.0D0

In the "20" DO-loop we loop through the first nodes of each ring of NTH nodes, treating each as a diffuse source. The distribution factors not computed directly in this loop can be determined when needed from symmetry.

DO 20 J = 1, NN3 JJ = (J-1)\*NTH + 1

Zero the "DF" vector.

DO 9 I1 = 1, MM3 DF(I1) = 0.0D0 9 CONTINUE In the "10" DO-loop we allow node JJ to emit NSHOTS volleys of energy packets and follow them until they are either absorbed or leave the instrument.

```
DO 10 ISHOT = 1, NSHOTS
```

Locate emission point on node JJ.

```
CALL RAND (SEED, IN, RN)
RN1 = RN
CALL RAND(SEED, IN, RN)
RN2 = RN
GOTO(101, 102, 103, 104, 105, 106, 107, 108, 109, 110), J
101 CONTINUE
```

The source point is on the blackbody, which we model as a black sector in the plane of the F-O-V limiter. Note that this limits the analysis to a reasonably uniform calibration source.

```
ANGLE = RN1*PI/2.0D0

RADIUS = DSQRT(R1*R1 - H2*H2)

RADIUS = RADIUS*DSQRT(RN2)

TXI = RADIUS*DCOS(ANGLE)

TYI = RADIUS*DSIN(ANGLE)

TZI = H2

GOTO 111
```

The source point is on the upper ring of the F-O-V limiter face.

```
| 102 CONTINUE
| ANGLE1 = RN1*PI/2.0D0
| RAD1 = DSQRT(R1*R1 - H2*H2)
| A = DARSIN(RAD1/R1)
| RAD2 = DSQRT(R1*R1 - (H1+F0V1)*(H1+F0V1))
| B = DARSIN(RAD2/R1)
| ANGLE2 = DARCOS(DCOS(A) - RN2*(DCOS(A)-DCOS(B)))
| TXI = R1*DSIN(ANGLE2)*DCOS(ANGLE1)
| TYI = R1*DSIN(ANGLE2)*DSIN(ANGLE1)
| TZI = R1*DCOS(ANGLE2)
| GOTO 111
```

The source point is on the middle ring of the F-O-V limiter face.

```
103 CONTINUE

ANGLE1 = RN1*PI/2.0D0

RAD1 = DSQRT(R1*R1 - (H1+F0V1)*(H1+F0V1))

A = DARSIN(RAD1/R1)

RAD2 = DSQRT(R1*R1 - (H1+F0V2)*(H1+F0V2))

B = DARSIN(RAD2/R1)

ANGLE2 = DARCOS(DCOS(A) - RN2*(DCOS(A)-DCOS(B)))

TXI = R1*DSIN(ANGLE2)*DCOS(ANGLE1)

TYI = R1*DSIN(ANGLE2)*DSIN(ANGLE1)

TZI = R1*DCOS(ANGLE2)

GOTO 111
```

The source point is on the bottom ring of the F-O-V limiter face.

```
In the continue and the
```

The source point is on the first (outer) ring of the floor.

```
INTERPOLATION IN
```

The source point is on the second ring of the floor.

```
| 106 CONTINUE
| ANGLE = RN1*PI/2.0D0
| RADIUS = DSQRT(F2*F2 + RN2*(F1*F1-F2*F2))
| TXI = RADIUS*DCOS(ANGLE)
| TYI = RADIUS*DSIN(ANGLE)
| TZI = H1
| GOTO 111
```

The source point is on the third ring of the floor.

```
| 107 CONTINUE
| ANGLE = RN1*PI/2.0D0
| RADIUS = DSQRT(F3*F3 + RN2*(F2*F2-F3*F3))
| TXI = RADIUS*DCOS(ANGLE)
| TYI = RADIUS*DSIN(ANGLE)
| TZI = H1
| GOTO 111
```

The source point is on the fourth (inner) ring of the floor.

```
| 108 CONTINUE
| ANGLE = RN1*PI/2.0D0
| RADIUS = DSQRT(R2*R2 + RN2*(F3*F3-R2*R2))
| TXI = RADIUS*DCOS(ANGLE)
| TYI = RADIUS*DSIN(ANGLE)
| TZI = H1
| GOTO 111
```

The source point is on the inner edge of the floor.

```
| 109 CONTINUE
| ANGLE = RN1*PI/2.0D0
| ZLEVEL = RN2*H1
| TXI = R2*DCOS(ANGLE)
| TYI = R2*DSIN(ANGLE)
| TZI = ZLEVEL
| GOTO 111
```

The source point is on the precision aperture.

```
| 110 CONTINUE
| ANGLE = RN1*PI/2.0D0
| RADIUS = R3*DSQRT(RN2)
| TXI = RADIUS*DCOS(ANGLE)
| TYI = RADIUS*DSIN(ANGLE)
| TZI = 0.0D0
```

A reflection to another surface has occurred; thus, increment TXIOLD, TYIOLD, TZIOLD, TXJ, TYJ, and TZJ.

```
| 111 CONTINUE
| TXIOLD = TXI
| TYIOLD = TYI
| TZIOLD = TZI
| TXJ = TXI
| TYJ = TYI
| TZJ = TZI
```

(Note: All of the above coordinates advance with each reflection. Initially TXI, TYI, and TZI are used in a test to determine the first trip through position (TXI,TYI,TZI) by a given photon.)

We return to this junction ("13 CONTINUE") each time a reflection occurs.

```
1 13 CONTINUE
```

If this is the first time this volley has passed through J, the volley striking node J is "emitted" diffusely.

Test to see if the energy packet is absorbed or reflected. if the random number RN is less than or equal to ABS, the absorptivity, the packet is absorbed; otherwise it is reflected.

```
ABS = EF
REFR = RF
IF(TZJ.GT.H1) ABS = EFOV
IF(TZJ.GT.H1) REFR = RFOV
CALL RAND(SEED, IN, RN)
IF(RN.LE.ABS) GOTO 16
```

Test to see if the reflected energy packet is specularly reflected or diffusely reflected. If the random number RN is less than or equal to REFR, the reflectivity ratio, the packet is specularly reflected; otherwise it is diffusely reflected.

CALL RAND(SEED, IN, RN)
NSD = 1
IF(RN.LE.REFR) GOTO 17

Randomly select the direction of the diffuse reflection.

| 137 CONTINUE | CALL RAND(SEED, IN, RN) | IF(RN.EQ.1.0D0) GOTO 14 | THETA = ARSIN(SQRT(RN)) | GOTO 15 | 14 CONTINUE | THETA = 0.0D0 | 15 CONTINUE | CALL RAND(SEED, IN, RN) | PHI = 2.0DO\*PI\*RN | NSD = 2 | GOTO 17 | 115 CONTINUE

Call the vector subroutine which computes the x-, y-, and z-components of the unit vector in direction (THETA,PHI) with respect to the central coordinate system.

CALL VECTOR(UNX, UNY, UNZ, PHI, THETA, VOX, VOY, VOZ)

Call the SEARCH subroutine which identifies the receiving point of the diffusely reflected photon.

CALL SEARCH(TXI, TYI, TZI, TXJ, TYJ, TZJ, VOX, VOY, VOZ)

If the z-coordinate of the receiving position is H2, the photon has escaped.

IF(TZJ.EQ.H2) GOTO 24

If the z-coordinate of the receiving position is 0 and the r-coordinate is less than R3, the photon has entered the cavity.

```
RAD = DSQRT(TXJ*TXJ + TYJ*TYJ)
IF(TZJ.EQ.O.ODO.AND.RAD.LT.R3) GOTO 244
```

The photon is reflected to the next position whose coordinates are (TXJ, TYJ, TZJ).

GOTO 25

The photon is absorbed. Call the NODE subroutine which determines the absorbing node.

16 CONTINUE

CALL NODE(TXJ,TYJ,TZJ,J1)

DF(J1) = DF(J1) + 1.000

GOTO 10

Determine the inward-directed unit normal vector of the source position. This depends on the shape of the source surface.

17 CONTINUE
IF(TZI.EQ.H2) GOTO 201
IF(TZI.GT.H1) GOTO 202
IF(TZI.EQ.H1) GOTO 203
IF(TZI.GT.0.ODO) GOTO 204

The source point is on the precision aperture.

UNX = 0.0D0 UNY = 0.0D0 UNZ = 1.0D0 GOTO 23 The source is on the imaginary surface in the plane of F-O-V limiter.

| 201 CONTINUE | UNX = 0.0D0 | UNY = 0.0D0 | UNZ = -1.0D0 | GOTO 23

The source point is on the F-O-V limiter.

| 202 CONTINUE | UNX = -TXI/R1 | UNY = -TYI/R1 | UNZ = -TZI/R1 | GOTO 23

The source point is on the floor.

| 203 CONTINUE | UNX = 0.0D0 | UNY = 0.0D0 | UNZ = 1.0D0 | GOTO 23

The source point is on the inner edge of the floor.

| CONTINUE | UNX = -TXI/R2 | UNY = -TYI/R2 | UNZ = 0.0D0

If the reflection is diffuse, return to the diffuse sequence; otherwise continue in the specular sequence.

| 23 CONTINUE | IF(NSD.EQ.2) GOTO 115 Compute the "ideal" vector of the reflected packet.

VDOT = (TXI-TXIOLD)\*UNX

VDOT = VDOT + (TYI-TYIOLD)\*UNY

VDOT = VDOT + (TZI-TZIOLD)\*UNZ

VOX = TXI - TXIOLD - 2.0D0\*VDOT\*UNX

VOY = TYI - TYIOLD - 2.0D0\*VDOT\*UNY

VOZ = TZI - TZIOLD - 2.0DO\*VDOT\*UNZ

VMAG = DSQRT(VOX\*VOX + VOY\*VOY + VOZ\*VOZ)

VOX = VOX/VMAG

VOY = VOY/VMAG

VOZ = VOZ/VMAG

Call the SEARCH subroutine which identifies the receiving point of the "exactly" reflected photon.

CALL SEARCH(TXI, TYI, TZI, TXJ, TYJ, TZJ, VOX, VOY, VOZ)

If the z-coordinate of the receiving position is H2, the photon has escaped.

IF(TZJ.EQ.H2) GOTO 24

If the z-coordinate of the receiving position is 0 and the r-coordinate is less than R3, the photon has entered the cavity.

RAD = DSQRT(TXJ\*TXJ + TYJ\*TYJ)
IF(TZJ.EQ.O.ODO.AND.RAD.LT.R3) GOTO 244

The photon is reflected to the next position whose coordinates are (TXJ, TYJ, TZJ).

GOTO 25

The photon has escaped through the F-O-V limiter; thus, find the pie segment through which it has escaped and increment the appropriate DF counter.

```
| 24 CONTINUE

| IF(TXJ.GE.O.ODO.AND.TYJ.GE.O.ODO) I = 1

| IF(TXJ.LT.O.ODO.AND.TYJ.GE.O.ODO) I = 2

| IF(TXJ.LT.O.ODO.AND.TYJ.LT.O.ODO) I = 3

| IF(TXJ.GE.O.ODO.AND.TYJ.LT.O.ODO) I = 4

| DF(I) = DF(I) + 1.0DO

| GOTO 10
```

The photon has entered the cavity; thus, increment the cavity DF counter.

Reflection to another surface has occurred. Thus, increment TXIOLD, TYIOLD, TZIOLD, TXJ, TYJ, and TZJ.

```
25 CONTINUE

TXIOLD = TXI

TYIOLD = TYI

TZIOLD = TZI

TXI = TXJ

TYI = TYJ

TZI = TZJ
```

Loop back to test sequence to determine disposition of the photon when it arrives at this new surface.

GOTO 13

Write the vector of distribution factors corresponding to node J.

```
10 CONTINUE

DO 30 JK = 1, MM3

DF(JK) = DF(JK)/RSHOTS

30 CONTINUE

DO 63 I = 2,MM3,4

I2 = I + 2

DF(I) = DF(I) + DF(I2)

DF(I) = DF(I)/2.0D0

DF(I2) = DF(I)

63 CONTINUE

DO 64 JK = 1,MM3

WRITE(17,3) JJ, JK, DF(JK)

64 CONTINUE
```

Generate "trip" card at end of file.

```
20 CONTINUE

    JJ = 0

    JK = 0

    DFDUM = 0.0D0

    WRITE(17,3) JJ, JK, DFDUM
```

#### Format statements:

```
1 FORMAT(5X, 215, 3D14.5)
2 FORMAT(8X, I3)
 3 FORMAT(5X, 2I5, D14.5)
4 FORMAT(5X, I5, 2D14.5)
96 FORMAT(5X, 'THE SYSTEM VARIABLES ARE AS ',
          'FOLLOWS:' ,//10X,'H1 = ',F5.2 ,' CM',
  &
          /,10X,'H2 = ',F5.2,' CM',/,10X,'R1 = ',
  Ŀ
          F5.2, CM', /, 10X, R2 = ', F5.2, /, 10X, R3
          ,'= ',F5.2,/,10X,'F1 = ',F5.2,/,10X,'F2 = ',F5.2,
          /,10X,'F3 = ',F5.2,/,10X,'FOV1 = ',F5.2,/,10X,
  å
          'FOV2 = ',F5.2,/,10X,'EFOV = ',F5.2,/,10X,
          'RFOV = ',F5.2,/,10X,'EF = ',F5.2,/,10X,'RF = ',
          F5.2./.10X.'NSHOT = '.16///)
97 FORMAT(16,14F5.2)
```

End of "TRW FRONT".

```
STOP
END
```

## 3.3. Subroutine SEARCH

This subroutine calculates the magnitude of the vector from the source position to the receiving position of the reflected photon. Since the direction of this vector is known (VOX, VOY, VOZ), the magnitude is all that is needed to calculate the receiving position.

## 3.3.1. Subroutine Description

SUBROUTINE SEARCH(TXI, TYI, TZI, TXJ, TYJ, TZJ, VOX, VOY, VOZ)

When this subroutine is compiled on IBM equipment, double precision must be specified. This step may not be necessary on other processors.

IMPLICIT REAL\*8 (A-H,O-Z)

Establish COMMON variables.

COMMON H1, H2, R1, R2, R3, FOV1, FOV2, F1, F2, F3, PI

If the source point is at the origin of coordinates,  $\ensuremath{\mathsf{RETURN}}$  .

IF(TZI.EQ.O.ODO.AND.TXI.EQ.O.ODO) GOTO 27

If the source point is on the imaginary F=0-V exit plane surface, go to 11.

IF(TZI.EQ.H2) GOTO 11

If the source point is on the F-O-V limiter, go to 12.

IF(TZI.GT.H1) GOTO 12

. If the source point is on the floor, go to 13.

IF(TZI.EQ.H1) GOTO 13

If the source point is on the inner edge of the floor, go to 14.

IF(TZI.GT.O.ODO) GOTO 14

If the source point is on the precision aperture, go to 15.

GOTO 15

The source point is on the imaginary F-O-V exit plane surface; thus begin search on the F-O-V limiter.

11 CONTINUE

VMAG = - (H2-H1)/VOZ

TXJ = TXI + VMAG\*VOX

TYJ = TYI + VMAG\*VOY

TZJ = H1

RADIUS = DSQRT(TXJ\*TXJ + TYJ\*TYJ)

If 'RADIUS' is less than 'R1', the packet has missed the F-O-V limiter, in which case we go to 16.

IF(RADIUS.LT.R1) GOTO 16

The packet has intercepted the F-O-V limiter.

20 CONTINUE

A = VOX\*TXI + VOY\*TYI + VOZ\*TZI

B = TXI\*TXI + TYI\*TYI + TZI\*TZI - R1\*R1

VMAG = - A + DSQRT(A\*A-B)

TXJ = TXI + VMAG\*VOX

TYJ = TYI + VMAG#VOY

TZJ = TZI + VMAG\*VOZ

RETURN

The packet has missed the F-O-V limiter. Check to see if it has hit the floor, in which case RETURN.

16 CONTINUE
IF(RADIUS.GT.R2) RETURN

The packet has missed the floor; check to see if it has hit the edge of the floor.

VMAG = - H2/VOZ

21 CONTINUE

TXJ = TXI + VMAG\*VOX

TYJ = TYI + VMAG\*VOY

TZJ = 0.0D0

RADIUS = DSQRT(TXJ\*TXJ + TYJ\*TYJ)

If 'RADIUS' is less than 'R2', the packet has missed the edge of the floor, in which case go to 17.

IF(RADIUS.LT.R2) GOTO 17

The packet has hit the edge of the floor.

25 CONTINUE

A = VOX\*VOX + VOY\*VOY

B = TXI\*VOX + TYI\*VOY

C = TXI\*TXI + TYI\*TYI - R2\*R2

ARG = B\*B - A\*C

VMAG = (-B + DSQRT(ARG))/A

TXJ = TXI + VMAG\*VOX

TYJ = TYI + VMAG\*VOY

TZJ = TZI + VMAG\*VOZ

RETURN

The packet has missed the edge of the floor. Check to see if it has hit the precision aperture, in which case RETURN.

17 CONTINUE IF(RADIUS.GE.R3) RETURN The packet has entered the cavity.

TXJ = 0.0D0 TYJ = 0.0D0 TXJ = 0.0D0 RETURN

The source point is on the F-O-V limiter; thus begin on the F-O-V limiter exit plane.

12 CONTINUE | IF(VOZ.GT.0.ODO) GOTO 18

The source point is not on the F-O-V exit plane; thus begin search on the F-O-V limiter.

VMAG = -(TZI-H1)/VOZ
TXJ = TXI + VMAG\*VOX
TYJ = TYI + VMAG\*VOY
TZJ = H1
RADIUS = DSQRT(TXJ\*TXJ + TYJ\*TYJ)

If 'RADIUS' is less than 'R1', the packet has missed the F-O-V limiter, in which case we continue the search on the floor.

| IF(RADIUS.LT.R1) GOTO 19

The packet has hit the F-O-V limiter. Solve for coordinates.

| GOTO 20 |

The packet has missed the F-O-V limiter, thus we continue our search on the floor. If 'RADIUS' is greater than 'R2', the packet has hit the floor, in which case we RETURN.

19 CONTINUE IF(RADIUS.GT.R2) RETURN The packet has missed the floor, in which case we continue our search on the edge of the floor.

VMAG = - TZI/VOZ GOTO 21

Search on the F-O-V limiter exit plane.

18 CONTINUE

VMAG = (H2-TZI)/VOZ

22 CONTINUE

TXJ = TXI + VMAG\*VOX

TYJ = TYI + VMAG\*VOY

TZJ = H2

RADIUS = DSQRT(TXJ\*TXJ + TYJ\*TYJ)

REXIT = DSQRT(R1\*R1 - H2\*H2)

If 'RADIUS' is less than 'REXIT', the packet has escaped, in which case we RETURN.

IF(RADIUS.LT.REXIT) RETURN

The packet has hit the F-O-V limiter.

GOTO 20

The source point is on the floor; thus begin search on the imaginary F-O-V exit plane surface.

13 CONTINUE

VMAG = (H2-H1)/VOZ

GOTO 22

The source point is on the inner edge of the floor; thus check to see if packet is emitted upward or downward.

14 CONTINUE

IF(VOZ.LT.O.ODO) GOTO 23

The packet has been emitted upward; thus resume the search on the inner edge of the floor.

VMAG = (H1-TZI)/VOZ

26 CONTINUE

TXJ = TXI + VMAG\*VOX

TYJ = TYI + VMAG\*VOY

RADIUS = DSQRT(TXJ\*TXJ + TYJ\*TYJ)

If 'RADIUS' is less than 'R2', the packet has missed the inner edge of the floor; thus resume search on the F-O-V limiter.

IF(RADIUS.LT.R2) GOTO 24

The packet has intercepted the inner edge of the floor; thus go to 25.

GOTO 25

The packet has missed the inner edge of the floor; thus resume search on the F-O-V limiter.

24 CONTINUE VMAG = (H2-TZI)/VOZ GOTO 22

The packet is emitted downward; thus begin search on the inner edge of the floor.

23 CONTINUE VMAG = -TZI/VOZ TXJ = TXI + VMAG\*VOX TYJ = TYI + VMAG\*VOY TZJ = 0.0DO RADIUS = DSQRT(TXJ\*TXJ + TYJ\*TYJ)

If 'RADIUS' is greater than 'R2', the packet has hit the inner edge of the floor, in which case go to 25; otherwise, go to 17.

IF(RADIUS.GT.R2) GOTO 25 GOTO 17 The source point is on the precision aperture; thus begin search on the inner edge of the floor.

```
| 15 CONTINUE | VMAG = H1/VOZ | GOTO 26 | 27 CONTINUE |
```

End of Subroutine 'SEARCH'.

RETURN END

## 3.4. Subroutine VECTOR

This subroutine expresses the randomly selected diffuse angle with respect to the source position in terms of the central coordinate system.

# 3.4.1. Nomenclature

X-component of inward-directed unit normal vector of source position.
vector of source position.
Y-component of inward-directed unit normal
vector of source position.
Z-component of inward-directed unit normal
vector of source position.
X-component of unit tangent vector of source
position.
Y-component of unit tangent vector of source
position.
Z-component of unit tangent vector of source
position.
X-component of the mutually orthogonal vector.
Y-component of the mutually orthogonal vector.
Z-component of the mutually orthogonal vector.
Angle of diffuse vector with respect to normal
vector.
Angle of diffuse vector with respect to tangent
vector.
X-component of diffuse vector with respect to
central coordinate system.
Y-component of diffuse vector with respect to
central coordinate system.
Z-component of diffuse vector with respect to
central coordinate system.

## 3.4.2. Subroutine Description

SUBROUTINE VECTOR(UNX, UNY, UNZ, PHI, THETA, VOX, VOY, VOZ)

When this subroutine is compiled on IBM equipment, it is necessary to specify double precision. This step may not be necessary on other processors.

IMPLICIT REAL\*8 (A-H, O-Z)

Compute the x-, y-, and z-components of the unit tangent vector, UT, and the mutually orthogonal vector, US, on the source position.

If the surface is cylindrical, go to 11.

IF(UNZ.EQ.O.ODO) GOTO 11

If the surface is plane, go to 12.

IF(UNZ.EQ.1.0D0.OR.UNZ.EQ.-1.0D0) GOTO 12

Calculations for the spherical surface:

ARG = UNZ\*UNZ/(1.0DO-UNZ\*UNZ)

UTX = UNX\*DSQRT(ARG)

UTY = UNY\*DSQRT(ARG)

UTZ = DSQRT(1.0D0 - UNZ\*UNZ)

GOTO 13

Calculations for the cylindrical surface:

11 CONTINUE

UTX = 0.0D0

UTY = 0.0D0

UTZ = 1.0D0

GOTO 13

Calculations for plane surfaces:

```
12 CONTINUE

UTX = 1.0D0

UTY = 0.0D0

UTZ = 0.0D0

13 CONTINUE
```

Compute the x-, y-, and z-components of the unit vector on the source position which is mutually orthogonal to UN and UT.

```
USX = UNY*UTZ - UNZ*UTY
USY = UNZ*UTX - UNX*UTZ
USZ = 0.0D0
```

Compute components of the vector at angle (THETA, PHI) with respect to the UN,UT,US-coordinate system and express the result in the I,J,K-coordinate system.

```
SINTH = DSIN(THETA)

COSTH = DCOS(THETA)

SINPH = DSIN(PHI)

COSPH = DCOS(PHI)

VOX = SINTH*COSPH*UTX + SINTH*SINPH*USX

VOX = VOX + COSTH*UNX

VOY = SINTH*COSPH*UTY + SINTH*SINPH*USY

VOY = VOY + COSTH*UNY

VOZ = SINTH*COSPH*UTZ + SINTH*SINPH*USZ

VOZ = VOZ + COSTH*UNZ
```

End of Subroutine 'VECTOR'.

<u> </u>		 	<del></del>	
i				
•				
	RETURN			
1	REIUNN			
	END			
	END			
:				
'				

#### 3.5. Subroutine NODE

This subroutine identifes the node which absorbs a given photon.

## 3.5.1. Subroutine Description

SUBROUTINE NODE(TXJ,TYJ,TZJ,J1)

When this subroutine is compiled on IBM equipment, it is necessary to specify double precision. This step may not be necessary on other processors.

IMPLICIT REAL\*8 (A-H, O-Z)

Identify the COMMON variables.

COMMON H1, H2, R1, R2, R3, FOV1, FOV2, F1, F2, F3, PI

Check to see if absorption occurred on the F-O-V limiter, in which case go to 11.

IF(TZJ.GT.H1) GO TO 11
RADIUS = DSQRT(TXJ\*TXJ + TYJ\*TYJ)

Check to see if absorption occurred on the floor, in which case go to 12.

IF(RADIUS.LT.R1.AND.RADIUS.GE.R2) GO TO 12

Check to see if absorption occurred on the edge of the floor, in which case go to 13.

IF(TZJ.LT.H1.AND.TZJ.GT.O.ODO) GOTO 13

Absorption occurred on the precision aperture; thus determine which node.

```
IF(TXJ.GE.O.ODO.AND.TYJ.GE.O.ODO) J1 = 37
IF(TXJ.LT.O.ODO.AND.TYJ.GE.O.ODO) J1 = 38
IF(TXJ.LT.O.ODO.AND.TYJ.LT.O.ODO) J1 = 39
IF(TXJ.GE.O.ODO.AND.TYJ.LT.O.ODO) J1 = 40
RETURN
```

Absorption occurred on the F-O-V limiter; thus check to see in which ring of nodes the absorption occurred.

```
11 CONTINUE
IF(TZJ.GT.FOV1) GOTO 14
IF(TZJ.GT.FOV2) GOTO 15
```

Absorption occurred in the bottom ring of nodes of the F=0-V limiter; thus identify the node.

```
IF(TXJ.GE.O.ODO.AND.TYJ.GE.O.ODO) J1 = 13
IF(TXJ.LT.O.ODO.AND.TYJ.GE.O.ODO) J1 = 14
IF(TXJ.LT.O.ODO.AND.TYJ.LT.O.ODO) J1 = 15
IF(TXJ.GE.O.ODO.AND.TYJ.LT.O.ODO) J1 = 16
RETURN
```

Absorption occurred in the top ring of nodes of the F-O-V limiter; thus identify the node.

```
14 CONTINUE

IF(TXJ.GE.O.ODO.AND.TYJ.GE.O.ODO) J1 = 5

IF(TXJ.LT.O.ODO.AND.TYJ.GE.O.ODO) J1 = 6

IF(TXJ.LT.O.ODO.AND.TYJ.LT.O.ODO) J1 = 7

IF(TXJ.GE.O.ODO.AND.TYJ.LT.O.ODO) J1 = 8

RETURN
```

Absorption occurred in the middle ring of nodes of the F-O-V limiter: thus identify the node.

```
| 15 CONTINUE
| IF(TXJ.GE.O.ODO.AND.TYJ.GE.O.ODO) J1 = 9
| IF(TXJ.LT.O.ODO.AND.TYJ.GE.O.ODO) J1 = 10
| IF(TXJ.LT.O.ODO.AND.TYJ.LT.O.ODO) J1 = 11
| IF(TXJ.GE.O.ODO.AND.TYJ.LT.O.ODO) J1 = 12
| RETURN
```

Absorption occurred on the floor; thus check to see in which ring of nodes the absorption occurred.

```
12 CONTINUE
IF(RADIUS.GE.F1) GOTO 16
IF(RADIUS.GT.F2) GOTO 17
IF(RADIUS.GT.F3) GOTO 18
```

Absorption occurred in the innermost ring of nodes on the floor; thus identify the node.

```
IF(TXJ.GE.O.ODO.AND.TYJ.GE.O.ODO) J1 = 29
IF(TXJ.LT.O.ODO.AND.TYJ.GE.O.ODO) J1 = 30
IF(TXJ.LT.O.ODO.AND.TYJ.LT.O.ODO) J1 = 31
IF(TXJ.GE.O.ODO.AND.TYJ.LT.O.ODO) J1 = 32
RETURN
```

Absorption occurred on the outermost ring of nodes on the floor; thus identify the node.

```
16 CONTINUE

IF(TXJ.GE.O.ODO.AND.TYJ.GE.O.ODO) J1 = 17

IF(TXJ.LT.O.ODO.AND.TYJ.GE.O.ODO) J1 = 18

IF(TXJ.LT.O.ODO.AND.TYJ.LT.O.ODO) J1 = 19

IF(TXJ.GE.O.ODO.AND.TYJ.LT.O.ODO) J1 = 20

RETURN
```

Absorption occurred on the second ring of nodes from the outer edge of the floor; thus identify the node.

Absorption occurred on the third ring of nodes from the outer edge of the floor; thus identify the node.

```
| 18 CONTINUE
| IF(TXJ.GE.O.ODO.AND.TYJ.GE.O.ODO) J1 = 25
| IF(TXJ.LT.O.ODO.AND.TYJ.GE.O.ODO) J1 = 26
| IF(TXJ.LT.O.ODO.AND.TYJ.LT.O.ODO) J1 = 27
| IF(TXJ.GE.O.ODO.AND.TYJ.LT.O.ODO) J1 = 28
| RETURN
```

Absorption occurred on the edge of the floor; thus identify the node.

| 13 CONTINUE | IF(TXJ.GE.O.ODO.AND.TYJ.GE.O.ODO) J1 = 33 | IF(TXJ.LT.O.ODO.AND.TYJ.GE.O.ODO) J1 = 34 | IF(TXJ.LT.O.ODO.AND.TYJ.LT.O.ODO) J1 = 35 | IF(TXJ.GE.O.ODO.AND.TYJ.LT.O.ODO) J1 = 36

End of Subroutine 'NODE'.

RETURN END

#### 3.6. Subroutine RAND

This subroutine uses library Subroutine GGUBS to obtain random numbers in an efficient way. A sequence of 100 random numbers is generated each time the 100th number from the previous generation is used.

# 3.6.1. Nomenclature

SEED	A seed number required by GGUBS. It must be
	specified in the main program as a double pre-
	cision integer value.
IN	An integer between 1 and 100 inclusive which
	represents which of the 100 random numbers of
	a given generation is obtained.
RN	The random number passed back to MAIN.

# 3.6.2. Subroutine Description

SUBROUTINE RAND(SEED, IN, RN)

When this subroutine is compiled on IBM equipment, it is necessary to specify double precision. This step may not be necessary with other processors.

IMPLICIT REAL\*8 (A-H,O-Z)
REAL\*4 RN, Z(100)

If this is the first call to RAND, RN is equal to the negative value set in MAIN. In this case we must call GGUBS to generate the first generation of 100 random numbers.

IF(RN.LT.O.ODO) GOTO 11

If this is not the first call to RAND, then increment to the next value in this generation of 100 random numbers.

IN = IN + 1

If we have not yet used up our allotment of random numbers from this generation, we simply take the next one in the sequence back to MAIN.

IF(IN.LE.100) GOTO 13

However, if we have used all the members of this generation, we must create a new generation of 100 random numbers.

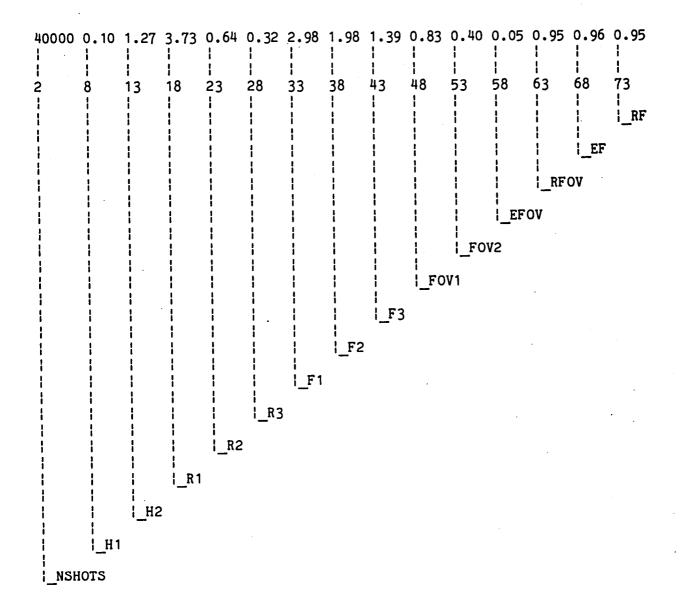
11 CONTINUE
IN = 1
DSEED = SEED
NUMBER = 100
CALL GGUBS(DSEED, NUMBER, Z)
SEED = DSEED
13 CONTINUE
RN = Z(IN)

End of Subroutine 'RAND'.

RETURN END

## 3.7. 'FRONT' DATA

These data are used only by Program 'TRW FRONT'. This is where all dimensions and properties of the body, other than the internode conductances, must be changed if another front end design is to be analyzed.



<u>Variable</u>	Type	<u>Field</u>	Description
NSHOT	16	1-6	Number of volleys fired
H1	F5.2	7-11	Distance to floor of instrument
H2	F5.2	12-16	Distance to face of F-O-V limiter
R1	F5.2	17-21	Outside radius of face
R2	F5.2	22-26	Inside radius of face
R3	F5.2	27-31	Radius of precision aperture
F1	F5.2	32-36	Radius to first floor node
F2	F5.2	37-41	Radius to second floor node
F3	F5.2	42-46	Radius to third floor node
FOV1	F5.2	47-51	Distance to top F-O-V limiter node
FOV2	F5.2	52-56	Distance to center limiter node
<b>EFOV</b>	F5.2	57-61	Emissivity of F-O-V limiter face
RFOV .	F5.2	62-66	Reflectivity ratio of F-O-V face
EF	F5.2	67-71	Emissivity of floor
RF	F5.2	72-76	Reflectivity ratio of face

Note: Please see the 'NOMENCLATURE' of this section for a more complete description of these variables. Also, see Fig. 1 of the February Progress Report.

#### 4. TRW TEMP

This program computes the transient temperature distribution in the active cavity and body of a TRW-type wide-field-of-view total spectrum radiometer subject to a uniform blackbody irradiance input and arbitrary heat sink and mounting beam temperature variations. It requires as inputs conductance matrices and radiation distribution factor matrices generated by the previous programs.

Programs TRW SHAPE, TRW NODES and TRW FRONT must be executed before this program is submitted. Also, an additional on-line disk file must be created and named A51961.DATA14 which contains the body internode conductances and the body node surface areas, emissivities, thermal capacities, volumes and heat sources. The current version of this file, which is user generated, is given in section 6. of this USER'S MANUAL.

#### 4.1. Nomenclature

TO	Matrix of "old" cavity node temperatures (K);
	i.e., temperatures at beginning of calculation
	time interval DT. Must be dimensioned M2.
TN	Matrix of "new" cavity node temperatures (K);
	i.e., temperatures at the end of calculation
	time interval DT. Must be dimensioned M2.
AA	Matrix of cavity node surface areas (m**2).
	Must be dimensioned M2.
CC	Matrix of cavity node thermal capacities
,	(W-s/K). Must be dimensioned M2.
DFC(J.I)	Matrix of cavity distribution factors for
2.0(0,2)	radiation emitted diffusely from the first node
	position on node ring J which is absorbed by
	node I (-). Must be dimensioned NN2 by (M2 + 1).
G(I,J)	Matrix of conductances for conduction path between
U(1,0)	cavity nodes I and J (W/K). Must be dimensioned
	M2 by M2.
NBODY	Number of radiometer "body" nodes (= 104).
NBODY2	NBODY + 2 (= 106).
RNBODY	Real version of NBODY.
HST	Heat sink (node "20") temperature (K).
TCLAMP	Mounting beam temperature (K).
HEAT	Proportional to cavity heater power, P (P =
HEAT	0.32*HEAT mW).
TSOURC	Source equivalent blackbody temperature (K).
GLINK	The conductance of the link between the cavity
GLINK	<u> </u>
CTC	and the thermal impedance (W/K).
SIG	Stefan-Boltzmann constant (5.6693E-08
MBODY	W/m##2-K##4).
MBODY	Body node number.

AABODY Surface area of body node 'MBODY' for purposes of calculating radiation heat transfer (cm\*\*2).

EEBODY Emissivity of body node 'MBODY' (-).

CCBODY Thermal capacity of body node 'MBODY' (W-s/cm\*\*3-K).

VVBODY Volume of body node 'MBODY' (cm##3).

QQBODY Heat source in body cavity 'MBODY'; may be due to radiation, conduction or electrical heating (W).

LIMBODY Delimiting index in DO-loop (= MBODY + 3).

ABODY(I) Surface area of body node I for purposes of calculating radiation exchange (m\*\*2).

EBODY(I) Emissivity of body node I (-).

CBODY(I) Thermal capacity of body node I  $(W-s/m^{**}3-K)$ .

VBODY(I) Volume of body node I (m##3).

QBODY(I) Heat source in body cavity I; may be due to radiation, conduction or electrical heating (W).

T1BODY(I) Initial temperature of body node I at beginning of solution; must be specified by user by specifying 'HST'. Later set equal to new temperature of node I at end of calculation interval (K).

TGBODY(I) Initial guess of body node I temperature at the end of a calculation time interval; initially set equal to 'HST', but in subsequent time steps it is updated to previous value (K).

COND Value of thermal conductance between body nodes on the same level ('MBODY') in the azimuthal direction (W/K).

GBODY(I,J) Value of thermal conductance between body nodes I and J (W/K).

DBODY(I,J) Value of radiation distribution factor from body node I to body node J (-).

MIBODY Index number of body node "level", or ring (-).

MJBODY Index number of body node "level", or ring (-).

SOURCE(I) Blackbody source-to-body node I distribution factor (-).

BODYDF Value of distribution factor, either body nodeto-body node or source-to-body node, read from unit 17 (-).

FTIME Total calculation time;, i.e., duration of the problem (s).

DT Length of cavity calculation time increment; i.e., integration step size (s).

NDT Number of time increments used in cavity calculation (-).

DTBODY Length of body calculation time increment (s).

PI Ratio of circumference to diameter of a circle (-).

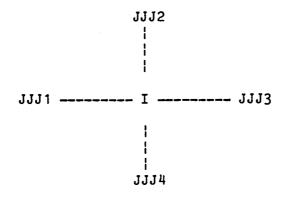
NARSZ Index used in Subroutine GAUSS; represents upper limit of size of matrix that can be inverted by the subroutine; i.e., used as argument of dimension statement in the subroutine. (-).

M2 Number of cavity nodes (-).

ELIMIT RMS error limit (%).

ITRMAX Maximum number of iterations permitted in the cavity temperature distribution calculation loop.

ITRMIN	Minimum number of iterations permitted in the cavity temperature distribution calculation loop.
AO	Aperture area (mm**2).
NTH	Number of circumferential divisions of cavity (-).
EC	Emissivity (=absorptivity) of the cavity (-)
EB	Emissivity of the back surface of the cavity (-).
ETIME	Elapsed time counter (s).
QTOT(I)	Average rate of energy transfer from the cavity
	to Node '13' of the radiometer body during ten
	successive cavity calculation time increments (W).
I1	Index indicating level of Node I.
12	Index indicating position of Node I in Level I1.
JJJ1	Index of node "to left" of Node I.
<b>JJJ2</b>	Index of node "above" Node I.
<b>J</b> JJ3	Index of Node "to right" of Node I.
<b>JJJ</b> 4	Index of node "below" Node I.



- Q1 Energy transfer rate into Node I from cavity nodes other than the receiving node itself during calculation time interval DT (W).
- J1 Index indicating level of Node J.
- J2 Index indicating position of node J in Level J1.

## 4.2. Program Description

Double precision must be specified when this program is compiled on IBM equipment. When other processors are used, the following statement may not be necessary.

IMPLICIT REAL\*8 (A-H,O-Z)

Dimension subscripted variables.

DIMENSION ABODY(104), EBODY(104), CBODY(104), VBODY(104)
DIMENSION QBODY(104), T1BODY(104), TGBODY(104), TUBODY(104)
DIMENSION QTOT(4), SOURCE(45), IANS(100)
DIMENSION DBODY(104, 106), GBODY(104, 104)
DIMENSION TO(100), TN(100), AA(100), CC(100), RHS(100)
DIMENSION Z(100, 100), G(100, 100)
DIMENSION DFC(10, 101), DFA(4, 100, 101)

Establish COMMON variables.

COMMON /EMSGC/ IER

Set constants. (Note: This is where each of these numbers must be changed when different cases are to be run.)

NBODY = 104
RNBODY = NBODY
HST = 313.0D0
TCLAMP = 326.0D0
HEAT = 400.0D0
TSOURC = 299.36D0
GLINK = 0.04D0
SIG = 5.6693D-08
NARSZ = 100
PI = 3.141592D0

Read in and write out system input variables.

READ(5,97) A,B,C,N1,N2,NTH,NSHOTN,ABS,REFR,DT,ELIMIT,MAT WRITE(6,96) A,B,C,N1,N2,NTH,NSHOTN,ABS,REFR,DT,ELIMIT,MAT

The following 'WRITE' statement causes a dimensioned sketch of the cavity to be "drawn" by the line printer.

WRITE(6,94)

Initialize radiometer body internode conductances and distribution factors to zero. This is necessary because the solution strategy considers possible interaction between all possible combinations of any two radiometer body nodes. Later we will read in the nonzero values of these two matrices.

```
NBODY2 = NBODY + 2
DO 111 I = 1,NBODY
DO 111 J = I,NBODY2
IF(J.LE.NBODY) GBODY(I,J) = 0.0D0
IF(J.LE.NBODY) GBODY(J,I) = GBODY(I,J)
DBODY(I,J) = 0.0D0
IF(J.LE.NBODY) DBODY(J,I) = DBODY(I,J)
111 CONTINUE
```

Read in values of constants and properties; initialize the node temperatures for radiometer body.

```
| 123 CONTINUE | READ(14,81) MBODY, AABODY, EEBODY, CCBODY, VVBODY, QQBODY | IF(MBODY.EQ.O) GOTO 124 | LIMBOD = MBODY + 3 | DO 223 I = MBODY, LIMBOD | ABODY(I) = AABODY*1.0D-4 | EBODY(I) = EEBODY | CBODY(I) = CCBODY*1.0D6 | VBODY(I) = VVBODY*1.0D-6 | QBODY(I) = QQBODY | T1BODY(I) = HST | TGBODY(I) = HST | CONTINUE | GOTO 123 | 124 CONTINUE | GOTO 123 | 124 CONTINUE | CONT
```

Read in nonzero elements of 'GBODY' matrix.

```
1211 CONTINUE
    READ(14,82) MBODY, COND
    IF(MBODY.EQ.O) GOTO 212
    LIMBOD = MBODY + 3
    DO 215 I = MBODY, LIMBOD
     J = I + 1
    IF(I.EQ.LIMBOD) J = MBODY
    GBODY(I,J) = COND
    GBODY(J,I) = COND
1215 CONTINUE
    GOTO 211
212 CONTINUE
     READ(14,83) MIBODY, MJBODY, GBODY(MIBODY, MJBODY)
     IF(MIBODY.EQ.O) GOTO 213
     GBODY(MJBODY, MIBODY) = GBODY(MIBODY, MJBODY)
     GOTO 212
1213 CONTINUE
```

Initialize source-to-instrument distribution factors to zero.

```
DO 601 I = 1,41
SOURCE(I) = 0.0D0
601 CONTINUE
```

Read in nonzero values of 'DBODY' matrix.

```
INN = 0
    INK = 1
1403 CONTINUE
    READ(17,92) I, J, BODYDF
    IF(I.EQ.0) GOTO 402
    INN = INN + 1
    IF(I.EQ.1) GOTO 501
    IF(I.LE.13) II = I - 4
     IF(I.EQ.17) II = 17
     IF(I.GE.21.AND.I.LT.33) II = I + 8
     IF(I.GE.33) II = I + 4
     IF(J.LT.5) JJ = 105
     IF(J.GE.5.AND.J.LT.17) JJ = J - 4
     IF(J.GE.17.AND.J.LT.21) JJ = J
     IF(J.GE.21.AND.J.LT.33) JJ = J + 8
     IF(J.GE.33.AND.J.LT.41) JJ = J + 4
     IF(J.EQ.41) JJ = 106
     DBODY(II,JJ) = DBODY(II,JJ) + BODYDF
     GOTO 403
1501 CONTINUE
     IF(INN.EQ.5) INK = INK + 4
     IF(INN.EQ.5) INN = 1
     SOURCE(INK) = SOURCE(INK) + BODYDF
     GOTO 403
1402 CONTINUE
```

Use symmetry to establish the values of the elements of the 'DBODY' matrix not computed directly.

```
DO 401 I = 1,104,4

DO 401 K = 1,104

I1 = I + 1

I2 = I + 2

I3 = I + 3

L1 = K + 3

IF(L1.GT.4) L1 = K - 1

L2 = K + 2

IF(L2.GT.4) L2 = K - 2

L3 = K + 1

IF(L3.GT.4) L3 = K - 3

DBODY(I1,K) = DBODY(I,L1)

DBODY(I2,K) = DBODY(I,L2)

DBODY(I3,K) = DBODY(I,L3)

401 CONTINUE
```

Set upper limit on calculation time; i.e., the duration of the heating or cooling problem being studied. The value used below, 400 s, permits the instrument to come into equilibrium. Also, set number of calculation time intervals.

FTIME = 400.0D0 NDT = FTIME/DT

Set calculation step size for body temperature calculation. Note that this is ten times the calculation time increment for the cavity calculation.

DTBODY = 10.0D0\*DT

Establish constants and indices.

AO = PI\*C\*C/100.0D0

FNTH = NTH

NN1 = N1

NN2 = NN1 + N2

M1 = NN1\*NTH

M2 = NN2\*NTH

FM2 = M2

MM2 = M2 + 1

EC = ABS

ITRMAX = 100

ITRMIN = 2

EB = 1.0D0

ETIME = 0.0D0

IER = 0

Write out node grid and geometry information.

WRITE(6,98) M2,NN2,NTH,NA,NTH,A,B,C IF(MAT.EQ.0) WRITE(6,93) ELIMIT

Read in values of dimensioned constants. Begin by zeroing the G(I,J), Z(I,J) and RHS(I) matrices.

DO 19 I = 1, M2

RHS(I) = 0.0D0

DO 19 J = 1, M2

G(I,J) = 0.0D0

Z(I,J) = 0.0D0

19 CONTINUE

Note: The node half-conductances must be retrieved from the lower left-hand triangle of the 'V' matrix created by 'TRW SHAPE'.

```
10 CONTINUE
READ(15,7) I, J, XX, YY, ZZ
IF((I.GT.M2).OR.(I.EQ.O)) GOTO 11
IF(I.LE.J) GOTO 10
```

If I or J are either one greater than M2, the values of XX and YY will be zero; that is, there are no conductances stored at those locations. We skip the following two statements in that case.

```
IF(XX.LE.1.0D-10) GOTO 10
   IF(YY.LE.1.0D-10) GOTO 10
   G(I,J) = 1.0D0/XX + 1.0D0/YY
   G(I,J) = 1.0D0/G(I,J)
   G(J,I) = G(I,J)
   GOTO 10
11 CONTINUE
   READ(10,1) I, AI, CI
   IF(I.EQ.0) GOTO 12
   AA(I) = AI
   CC(I) = CI
   GOTO 11
12 CONTINUE
   READ(11,2) L, M, I, DFADUM
   IF(L.EQ.O) GOTO 13
   DFA(L,M,I) = DFADUM
   GOTO 12
13 CONTINUE
   READ(12,3) I, J, DFCIJ
   IF(I.EQ.0) GOTO 14
   DFC(I,J) = DFCIJ
   GOTO 13
14 CONTINUE
```

Initialize 'old' node temperatures. This establishes the initial conditions for the solution.

```
DO 15 I = 1, M2
TO(I) = HST
TN(I) = TO(I)
15 CONTINUE
```

Write out initial cavity temperature distribution.

```
WRITE(6,8) ETIME
DO 301 I5 = NTH, M2 , NTH
I9 = I5 - NTH + 1
WRITE(6,6) (TN(I3), I3 = I9, I5)
WRITE(16,6) (TN(I3), I3 = I9, I5)
301 CONTINUE
```

Zero the 'QTOT' accumulator.

```
DO 987 I = 1,4
QTOT(I) = 0.0D0
1987 CONTINUE
```

We "march in time" by letting (if desired) the heat sink and calibration source temperatures change with each time increment DT. We assume that DT is sufficiently small that each incremental change in these temperatures is small.

The source and heat sink temperatures are changed in the 'K' DO-loop (if desired; this is an option not actually used in this example).

```
KOUNT = 0
DO 280 K = 1, NDT
```

Compute the new radiometer body temperature distribution every tenth time increment.

```
KOUNT = KOUNT + 1
    IF(KOUNT.EQ.10) GOTO 678
    GOTO 679
1678 CONTINUE
    KOUNT = 0
1219 CONTINUE
    DO 230 J = 1, NBODY
    CJ = CBODY(J)*VBODY(J)
    CJ = DTBODY/CJ
    SUM1 = 0.0D0
    SUM2 = 0.0D0
    SIGMA = SIG*AO/10000.0D0
    DO 220 I = 1.NBODY
    IF(I.EQ.J) GOTO 220
    FACTOR = EBODY(I)*SIG*ABODY(I)*DBODY(I,J)
    FACTOR = FACTOR*TGBODY(I)*TGBODY(I)*TGBODY(I)*TGBODY(I)
    FACTOR = FACTOR + GBODY(I,J)*TGBODY(I)
    SUM1 = SUM1 + FACTOR
     SUM2 = SUM2 + GBODY(I,J)
1220 CONTINUE
     IF(J.GE.1.AND.J.LT.5)
            QBODY(J) = SIGMA*SOURCE(5)*TSOURC**4
     IF(J.GE.5.AND.J.LT.9)
            QBODY(J) = SIGMA*SOURCE(9)*TSOURC**4
     IF(J.GE.9.AND.J.LT.13)
            QBODY(J) = SIGMA*SOURCE(13)*TSOURC**4
     IF(J.GE.17.AND.J.LT.21)
            QBODY(J) = SIGMA*SOURCE(17)*TSOURC**4
     IF(J.GE.29.AND.J.LT.33)
            QBODY(J) = SIGMA*SOURCE(21)*TSOURC**4
     IF(J.GE.33.AND.J.LT.37)
            QBODY(J) = SIGMA*SOURCE(25)*TSOURC**4
     IF(J.GE.37.AND.J.LT.41)
            QBODY(J) = SIGMA*(SOURCE(29)+SOURCE(33))*TSOURC**4
     IF(J.GE.41.AND.J.LT.45)
            QBODY(J) = SIGMA*SOURCE(37)*TSOURC**4
     IF(J.GT.20.AND.J.LT.25) QBODY(J) = 3.46DO*TCLAMP
     IF(J.GT.24.AND.J.LT.29) QBODY(J) = 0.22DO*TCLAMP
     IF(J.GT.48.AND.J.LT.53) GOTO 636
     GOTO 637
1636 CONTINUE
     ILL = J - 48
     QBODY(J) = QTOT(ILL)/10.0D0 + GLINK*T1BODY(J)
1637 CONTINUE
     TOP = CJ*(SUM1 + QBODY(J)) + T1BODY(J)
     BOTTOM = (DBODY(J,J) - 1.0DO)*EBODY(J)*SIG*ABODY(J)
     BOTTOM = BOTTOM*TGBODY(J)*TGBODY(J)*TGBODY(J)
     IF(J.GT.20.AND.J.LT.25) SUM2 = SUM2 + 3.46D0
     IF(J.GT.24.AND.J.LT.29) SUM2 = SUM2 + 0.22D0
     IF(J.GT.48.AND.J.LT.53) SUM2 = SUM2 + GLINK
     BOTTOM = 1.0D0 + CJ*(SUM2 - BOTTOM)
```

```
TUBODY(J) = TOP/BOTTOM
    IF(J.GT.76.AND.J.LT.81) TUBODY(J) = HST
1230 CONTINUE
    SUM = 0.0D0
    AVE = 0.0D0
    DO 240 I = 1.NBODY
    BRMS = TGBODY(I) - TUBODY(I)
    BRMS = BRMS*BRMS
    SUM = SUM + BRMS
    AVE = AVE + TGBODY(I)
1240 CONTINUE
    BRMS = SUM/RNBODY
    BRMS = DSQRT(BRMS)
    AVE = AVE/RNBODY
    TEST = BRMS/AVE
    IF(TEST.LE.0.000001) GOTO 241
    DO 250 I = 1,NBODY
    TGBODY(I) = TUBODY(I)
1250 CONTINUE
    GOTO 219
241 CONTINUE
    DO 260 I = 1, NBODY
    IF(I.LE.4) QTOT(I) = 0.0D0
    T1BODY(I) = TUBODY(I)
1260 CONTINUE
    EETIME = ETIME + DT
    WRITE(6,88) EETIME
    DO 246 IS = 4, 104, 4
     IX = IS - 3
     WRITE(6,66) (T1BODY(I17), I17 = IX, IS)
1246 CONTINUE
    QBOD = 0.0D0
    DO 661 I = 1,104
     ALPHA = 0.96D0
     IF(I.LT.13) ALPHA = 0.05D0
     QBOD = QBOD
            + ALPHA*ABODY(I)*SIG*DBODY(I,106)*T1BODY(I)**4
1661 CONTINUE
     QBOD = QBOD/4.0D0
1679 CONTINUE
```

Compute the elapsed time.

ETIME = K\*DT

The 'ITER' DO-loop loops through the temperature calculations until convergence is obtained.

DO 400 ITER = 1, ITRMAX
RMS = 0.0D0
TSUM = 0.0D0
RMSMAX = 0.0D0

The 'I' DO-loop steps through the nodes.

DO 18 I = 1, M2

Compute the (Level/Position) coordinates of Node I.

I1 = I - 1 I1 = I1/NTH I2 = I - I1\*NTH I1 = I1 + 1

Compute the indices of the nodes surrounding Node I.

JJJ1 = I - 1
IF(I2.EQ.1) JJJ1 = JJJ1 + NTH
JJJ2 = I - NTH
JJJ3 = I + 1
IF(I2.EQ.NTH) JJJ3 = JJJ3 - NTH
JJJ4 = I + NTH
IF(JJJ2.LE.0) JJJ2 = I
IF(JJJ4.GT.M2) JJJ4 = I

Compute the heat flow rate radiated into each node during the calculation time interval DT.

Q1 = 0.0D0 DFSUM = 0.0D0

Loop through the J nodes which radiate to Node I.

DO 22 J = 1, M2

Compute (Level/Position) coordinates of Node J.

```
J1 = J - 1
J1 = J1/NTH
J2 = J - J1*NTH
J1 = J1 + 1
```

If Node J is in the first position on Level J1, i.e., if J2 = 1, the appropriate distribution factors have been read in; otherwise they must be found by a coordinate transformation.

```
IF(J2.EQ.1) GOTO 20
```

Coordinate transformation; rotate source Node J to first position on Level J1 and rotate receiver node an equal number of positions on Level I1.

```
JJ = J1

II = (I1 - 1)*NTH + I2 - J2 + 1

IF((I2 - J2).LT.0) II = II + NTH

GOTO 21
```

The distribution factor has been read in directly.

```
| 20 CONTINUE
| JJ = J1
| II = I
```

The distribution factor from Node J to Node I is the same as the distribution factor from Node JJ to Node II due to symmetry.

```
21 CONTINUE
```

Calculate Z(I,J).

```
IF(J.NE.I) GOTO 75
II1 = (I1 - 1)*NTH + 1
ZIJ = AA(I)*EC*SIG*TO(I)*TO(I)*TO(I)*DFC(I1,II1)
Z(I,J) = -(EC + EB)*AA(I)*SIG*TO(I)*TO(I)*TO(I)
Z(I,J) = Z(I,J) + ZIJ - CC(I)/DT
Z(I,J) = Z(I,J) - G(JJJ1,I) - G(JJJ2,I)
Z(I,J) = Z(I,J) - G(JJJ3,I) - G(JJJ4,I)
GOTO 76

75 Z(I,J) = EC*SIG*TO(J)*TO(J)*TO(J)*AA(J)*DFC(JJ,II)
Z(I,J) = Z(I,J) + G(I,J)
```

There is no heat transfer by conduction from a node to to itself; thus if J = I, we skip the 'Q1' summation.

```
76 IF(J.EQ.I) GOTO 22
Q1 = Q1 + Z(I,J)*TN(J)
DFSUM = DFSUM + DFC(II,JJ)

22 CONTINUE
Q2 = 1.0D0 - DFSUM
Q2 = Q2*AA(I)*EC
Q2 = Q2*SIG*TSOURC*TSOURC*TSOURC
```

The heat input into the first ring of cavity nodes must include that from the F-O-V limiter; thus, if 'I' is less than or equal to 4, we must add the heat 'QBOD' from the F-O-V limiter.

```
IF(I.LE.4) Q2 = Q2 + QBOD
```

Calculate the quantity RHS(I).

```
RHS(I) = - AA(I)*EB*SIG*HST*HST*HST*HST
Q2 = Q2 + HEAT*AA(I)
RHS(I) = RHS(I) - Q2 - CC(I)*TO(I)/DT
```

Check to see if the matrix inversion method has been chosen; if so, proceed to next node.

```
IF(MAT.EQ.1) GOTO 888
```

Calculate TNEW for this node if the iterative solution has been chosen.

```
IF(I.LE.4) GOTO 188
    GOTO 189
188 CONTINUE
    IHLP = I + 48
    TNEW = (Q1 - RHS(I) -
           GLINK*T1BODY(IHLP))/(-Z(I,I) - GLINK)
    GOTO 190
189 CONTINUE
    TNEW = (Q1 - RHS(I))/(-Z(I,I))
1190 CONTINUE
    TRMSMX = (TNEW - TN(I))*(TNEW - TN(I))
    RMS = RMS + TRMSMX
    IF(TRMSMX.GT.RMSMAX) RMSMAX = TRMSMX
    TSUM = TSUM + TNEW
    TN(I) = TNEW
    GOTO 18
1888 CONTINUE
    IF(I.LE.4) GOTO 517
    GOTO 518
1517 CONTINUE
    IHLP = I + 48
    RHS(I) = RHS(I) - GLINK*T1BODY(IHLP)
     Z(I,I) = Z(I,I) - GLINK
1518 CONTINUE
18 CONTINUE
```

If the matrix solution method has been chosen, call the matrix inversion Subroutine GAUSS.

```
IF(MAT.EQ.0) GOTO 73
CALL GAUSS(Z,RHS,M2,NARSZ,IANS)
IF(IER.EQ.131) STOP
IF(IER.EQ.130) STOP
DO 288 NP = 1, M2
TN(NP) = RHS(NP)
288 CONTINUE
GOTO 270
```

Determine if convergence has been obtained.

```
73 RMS = DSQRT(RMS)

RMSMAX = DSQRT(RMSMAX)

ERROR = RMS*100.0/TSUM

ERRMAX = RMSMAX*100.0*FM2/TSUM

IF((ERRMAX.LE.ELIMIT).AND.

& (ITER.GT.ITRMIN)) GOTO 28

400 CONTINUE

410 CONTINUE

28 CONTINUE
```

Write the value of the elapsed time (s).

```
1270 WRITE(6,8) ETIME
```

Write out current cavity temperature distribution at the end of each integration step.

```
DO 300 I5 = NTH, M2 , NTH

I9 = I5 - NTH + 1

WRITE(6,6) (TN(I3), I3 = I9, I5)

WRITE(16,6) (TN(I3), I3 = I9, I5)

300 CONTINUE

DO 299 IJK = 1, M2

TO(IJK) = TN(IJK)

IJKL = IJK + 48

IF(IJK.LE.4) LL = IJK

IF(IJK.LE.4) QTOT(LL) = QTOT(LL)

& + GLINK*(TO(IJK) - T1BODY(IJK))

1299 CONTINUE

280 CONTINUE
```

Format statements.

```
1 FORMAT(5X, I5, 2D14.5)
 2 FORMAT(5X, 3I5, D14.5)
 3 FORMAT(5X, 215, 2D14.5)
 4 FORMAT(5X, F10.4)
 5 FORMAT(5X, 2I5, D14.5)
 6 FORMAT(5X, 2(F8.3),2(F8.3),2(F8.3),
          2(F8.3),2(F8.3))
 7 FORMAT(5X, 215, 3D14.5)
 8 FORMAT(10X, ///, 5X, 'ELAPSED TIME =',
           F6.1,' SECOND(S).', //)
 9 FORMAT(5X/5X, 'CONVERGENCE NOT OBTAINED,',
           'ERROR =',E11.4,' PERCENT'/)
66 FORMAT(45X, 2(F8.3),2(F8.3),2(F8.3),2(F8.3))
88 FORMAT(44X, 'ELAPSED TIME =',F6.1,' SECOND(S).',//)
81 FORMAT(4X, I3, 3X, F4.2, 1X, F4.2, 2X, F5.3, 5X, F5.3, 5X, F5.3)
82 FORMAT(4X, 13, 3X, F6.4)
83 FORMAT(5X, I2, 2X, I3, 2X, F5.2)
92 FORMAT(5X, 215, D14.5)
93 FORMAT(//5X, ITERATION MAXIMUM ERROR LIMIT ',
           '= '.F10.8,' PERCENT.',/,'1')
 94 FORMAT('1',//////,15X,'TRANSIENT',
           'TEMPERATURE DISTRIBUTION OF A',/,15X,
           'TRW TYPE INVERTED CONE CAVITY RADIOMETER.',
           //.14X.'DEVELOPED BY J. R. MAHAN AND ',
           'L. D. ESKIN AT',/,11X,'VIRGINIA ',
           'POLYTECHNIC INSTITUTE & STATE UNIVERSITY',
           /,17X,'FOR THE NASA LANGLEY RESEARCH CENTER.')
 95 FORMAT(10X, 'THE SOLUTION HAS SWITCHED ',
           'TO THE ITERATION METHOD.')
 96 FORMAT(5X, 'THE SYSTEM VARIABLES ARE AS ',
           'FOLLOWS:' ,//10X,'A = ',F5.2 ,' MM',
   å
           /,10X,'B = ',F5.2,' MM',/,10X,'C = ',
           F5.2, MM', /, 10X, N1 = ', I2, /, 10X, N2 '
   å
           ,'= ',I2,/,10X,'NTH =',I2,/,10X,'NSHOTN = '
   &
           ,16,/,10X,'ABS = ',F3.1,/,10X'REFR = '
   å
           '= ',16,/,10X,'ABS = ',F3.1,/,10X,'REFR = '
           .F3.1./.10X.'DT = '.F4.1./.10X.'ELIMIT = '
           ,E8.2,/10X,'TEMPERATURE SOLUTION METHOD ',
   &
           ' (O=ITERATION, 1=MATRIX) = ',I2)
 97 FORMAT(F5.2,F5.2,F5.2,I2,I2,I2,I6,
           F3.1.F3.1.F4.1,E8.2,4I1)
```

```
98 FORMAT(///15X, 'NUMBER OF CAVITY NODES = '
            ,13,//,15X,12,' CAVITY NODE RINGS WITH ',12,
  å
            ' NODES PER RING.',//,1X,
  å
            T27, 'SYSTEM GEOMETRY', // 1X, T41,'
  &
            ,1X,T27,'|<---->| | |'/,1X,T27,'|',T40,
            '| | |',/1X,T27,'|',T40,'| A |'/,1X,T27,'|'
            T40, '| | | '/.1X.T27, '| '.T40, '| V | './'+', T42, '.'.', '1X.T28, '*', T39, '*', T44, 'B'/, 1X.T29, '*', T38, '*', T44, '| '/.1X.T30, '*', T37, '*', T44, '| '/.
  &
  å
            1X.T31,'*',T36,'*',T44,'\'/,1X,T32,'*',T35
  &
             '*',T44,'¦',/,1X,T33,'*',T34,'*',T44,'V',/
                                      ',//,5X,'A = ',F5.2,
            ,'+',T35,'
            ' MM',5X,'B = ',F5.2,' MM',5X,'C = ',F5.2
   å
             ('1'/,'MM',
99 FORMAT(5X,/,5X,'THE AVERAGE RMS ERROR = '
            ,F10.8,' PERCENT AFTER ',12, ' ITERATIONS .'
            //,5X,'THE MAXIMUM RMS ERROR = ',F10.8,
            ' PERCECNT.',//,5X, 'THE HEAT SINK ',
   å
             'TEMPERATURE IS ',F5.1,' K.',/)
```

End of 'TRW TEMP'.

```
STOP
END
```

## 4.3. Subroutine GAUSS

This is a standard matrix inversion subroutine which could be replaced at the discretion of the user if the argument list of the replacement routine was the same. We present this subroutine without further explanation.

# 4.3.1. Subroutine Listing

```
SUBROUTINE GAUSS(A, B, N, NARSZ, IANS)
    IMPLICIT REAL*8(A-H,O-Z), INTEGER(I-N)
    INTEGER IANS(NARSZ)
    REAL*8 A(NARSZ, NARSZ), B(NARSZ)
    COMMON /EMSGC/ IER
    DO 10 I = 1,N
    IANS(I) = I
    AM = DABS(B(I))
    DO 20 J = 1, N
    BM = DABS(A(I,J))
    IF ( BM .GT . AM ) AM = BM
20 CONTINUE
    IF(AM.EQ.0.0) GOTO 901
1 10 CONTINUE
    N1 = N + 1
    DO 100 L = 1, N
    AMX = DABS(A(L,L))
    I = L
    J = L
    DO 110 I1 = L,N
    DO 120 J1 = L,N
    F1 = DABS(A(I1,J1))
    IF(AMX.GE.F1) GOTO 120
    AMX = F1
    I = I1
    J = J1
1120 CONTINUE
110 CONTINUE
    IF(I.EQ.L) GOTO 140
    DO 130 K = 1,N
    F1 = A(I,K)
    A(I,K) = A(L,K)
    A(L,K) = F1
1130 CONTINUE
```

```
F1 = B(I)
    B(I) = B(L)
    B(L) = F1
1140 CONTINUE
    IF(J.EQ.L) GOTO 160
    DO 150 K = 1, N
    F1 = A(K,J)
    A(K,J) = A(K,L)
    A(K,L) = F1
150 CONTINUE
     III = IANS(J)
     IANS(J) = IANS(L)
     IANS(L) = III
1160 CONTINUE
     F1 = A(L,L)
     IF(F1.EQ.0.0) GOTO 902
     DO 170 LP =1, N
     A(L,LP) = A(L,LP)/F1
170 CONTINUE
     B(L) = B(L)/F1
     IF(N.EQ.L) GOTO 200
     N2 = L + 1
     DO 180 M = N2, N
     F1 = A(M,L)
     IF(F1.EQ.0.0) GOTO 182
     DO 185 LP = 1, N
     A(M,LP) = A(M,LP)/F1 - A(L,LP)
1185 CONTINUE
     B(M) = B(M)/F1 - B(L)
1182 CONTINUE
 1180 CONTINUE
1100 CONTINUE
200 CONTINUE
```

```
F1 = 0.0
    N3 = N-1
    DO 210 LP = 1.N3
    L = N - LP
    F1 = 0.0
    N4 = L + 1
    DO 220 K = N4, N
    F1 = F1 + B(K)*A(L,K)
220 CONTINUE
    B(L) = B(L) - F1
210 CONTINUE
    DO 250 LP = 1,N
    A(1,LP) = B(LP)
1250 CONTINUE
    DO 260 I = 1, N
    IA = IANS(I)
    B(IA) = A(1,I)
260 CONTINUE
    GOTO 999
|901 \text{ IER} = 130
     WRITE(6,666)
1666 FORMAT(/, GAUSS(F): ZERO ROW INPUT')
    GOTO 999
1902 IER = 131
     WRITE(6,777)
| 1777 FORMAT(/,' GAUSS(F): SINGULAR MATRIX')
1999 CONTINUE
     RETURN
     END
```

# 5. TRW DATA

The data read by 'TRW SHAPE', 'TRW NODES' and 'TRW TEMP' must be entered at the appropriate position at the end of each of the three programs. Note that not all of the programs use, or even read, all of the input data. However, the same data file serves all three.

5.4920.40 3.86	column column	38:	
	column		DT REFR
	column	٠.	ABS NSHOTN
	column column		NTH N2
	column column		N1 C
	 column column		

Variable	Type	<u>Field</u>	Description
A	F5.2	1-5	Height of barrel (mm)
В	F5.2	6-10	Height of entire cavity (mm)
C	F5.2	11–15	Radius of barrel (mm)
N 1	12	16-17	Number of barrel axial divisions
N2	12	18-19	Number of cone axial divisions
NTH	12	20-21	Number of circumferential divisions
NSHOTN	16	22-27	Number of volleys fired
ABS	F3.1	28-30	Absorptivity of cavity
REFR	F3.1	31-33	Reflectivity ratio
DT	F4.1	34-37	Cavity calculation time interval
ELIMIT	E8.2	38-45	Accuracy criterion using iteration
MAT	I1	46	= 1, matrix solution; = 0, iteration

				6. BODY	DATA	
1	2.58	0.05	2.425	1.020	0.000	
			2.425	0.710	0.000	
			2.425	0.570	0.000	
		0.0	2.425	0.580		
			2.425	1.900	0.000	
			2.425	9.080		
		0.0		1.980		
			2.425	1.330	0.000	
33			2.425	1.200	0.000	
37			2.425	0.120	0.000	
			2.425	0.260	0.000	
			2.425	0.530 0.003	0.000	
53	_		2.405	1.970	0.000	
	0.56		2.463	0.003		
	0.00		2.425	1.200		
	0.40		3.430	0.240		
	0.00		2.425	1.560	0.000	
73	0.52		3.430	0.310		
77	0.00	0.0	3.430	1.540	0.000	
81		0.0	3.430	0.310		
85	-	0.0	2.425	1.170	0.000	
89			3.430	0.540		
93	- •		2.425	1.340		
97			2.425	0.990		
101		0.0	2.463	0.016	•	
î	Î	i I	i 1	Î	1 001.000	41: QQBODY
!	i !	i !	i !	i !	ieorumn	41; QQBODI
					column	31; VVBODY
<b>!</b>	!	<b>!</b>			column	21; CCBODY
Ì	i	i	' <del></del>			,
!	1	l	<del></del>		column	16; EEBODY
1				<del></del>	column	11; AABODY
İ		<del></del>	······································		column	5; MBODY

Variable	Type	Field	Description
MBODY	13	5-7	Body node number
AABODY	F4.2	11-14	Body node surface area
EEBODY	F4.2	16-19	Body node emissivity
CCBODY	F5.3	21-25	Body node thermal capacity
VVBODY	F5.3	31-35	Body node volume
QQBODY	F5.3	41-45	Body node heat input

These variables are described in detail in Section 4.1.

```
1
5
9
13
      0.0592
0.0412
      0.0320
      0.0340
17
      0.1384
21
      0.6240
25
      0.1240
29
      0.1760
33
      0.3440
37
      0.0920
41
      0.2000
45
      0.4000
49
      0.0192
53
      0.8000
      0.0208
57
61
      0.4400
65
      0.6800
69
      0.5600
73
      0.8800
      1.0800
77
81
      0.8800
85
      3.9200
89
      1.5600
93
      6.7600
97
      4.6400
101
      1.0800
                                           column 11; COND
                                           column 5; MBODY
```

<u>Variable</u>	Type	<u>Field</u>	Description
MBODY	I3	5-7	Body node number
COND	F6.4	11-16	Conductance between body nodes on MBODY

1234567890123456345634563456345633 123456789012345634563456333333333333333333333333333	5678901123456789012222222233322223333333334423	8.95 8.95
35 36	40	0.66 0.66
33 34		
35	43	0.02
30 33	44 45	0.02 0.04
34	46 47	0.04 0.04
36	47	0.04
33	53	0.05

34 35 37 38 39 40 41 43 44 45 47	54 55 56 41 42 44 45 47 48 54 55 55	0.05 0.05 0.05 0.00 0.00 0.00 7.11 7.11 7.11 7.11 4.03 4.03
48 49	56 57 58	4.03 0.03
50	58	0.03
51 52	59 60	0.03 0.03
53 54	61 62	7.77 7.77
55 55 56 57 58 59 60	63 64	7.77
57	65	0.06
58 59	66 67	0.06 0.06
60 61	68 65	0.06 1.99
62	66	1.99
63 64	67 68	1.99 1.99
61 62	69 70	8.44 8.44
63	71	8.44
64 65 66	72 73	8.44 3.17
66 67	74 75	3.17 3.17
68 69	76	3.17 2.59 2.59 2.59
70	76 73 74 75	2.59
71 72	75 76	2.59 2.59
69 70	77 78	8.92 8.92
71	79	8.92
72 73	80 81	8.92 2.82
73 74 75	82 83	2.82 2.82
76	84	2.82

```
77
     81
           4.51
78
     82
           4.51
79
     83
           4.51
80
     84
           4.51
77
     85
          13.46
78
     86
          13.46
79
     87
          13.46
80
     88
          13.46
81
           2.06
     89
           2.06
82
     90
83
     91
           2.06
84
     92
           2.06
     89
85
           9.17
86
     90
           9.17
87
     91
           9.17
88
     92
           9.17
85
     93
           0.72
     94
           0.72
86
87
     95
           0.72
           0.72
88
     96
           0.03
89
    101
90
    102
           0.03
91
    103
           0.03
92
    104
           0.03
93
     97
           2.47
94
     98
           2.47
95
     99
           2.47
    100
96
           2.47
                                          column 16; GBODY(MIBODY, MJBODY)
                                          _column 10; MJBODY
                                          column 6; MIBODY
```

<u>Variable</u>	Type	<u>Field</u>	Description			
MIBODY MJBODY	I2 I3	6 <b>-</b> 7 10 <b>-</b> 12	Body node 'level' index Body node 'level' index			
GBODY	F5.2	16-20	Conductance between MIBODY and MJBODY			

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1. Report No. NASA CR-172240	2. Government Accessio	n No.	3. Recip	ent's Catalog No.		
4. Title and Subtitle			5. Repor	t Date		
TRANSIENT THERMAL MODEL	ING OF THE NONSCANN	ING	Nove	ember 1983		
ERBE DETECTOR			6. Perfor	ming Organization Code		
7. Author(s)			8. Perfor	ming Organization Report No.		
J. R. Mahan						
J. R. Manan			<del></del>			
			10. Work	Unit No.		
9. Performing Organization Name and Addr				•		
Mechanical Engineering I	=		11 Contr	act or Grant No.		
Virginia Polytechnic Ins		niversity	√ I	1-16508		
Blacksburg, Virginia 24	1061			1 10000		
			13. Type	of Report and Period Covered		
12. Sponsoring Agency Name and Address			Con	Contractor Report		
National Association and	1 Coors Administrat	d a n	<b></b>			
National Aeronautics and	i Space Administrat	1011	14. Spons	oring Agency Code		
Washington, DC 20546	•					
15. Supplementary Notes		<del> </del>				
Langley technical monito	r: Dwayne E. Hinto	n	•	1		
Final Report	Ji. Dwayne B. Mineo	••		·		
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16. Abstract				·		
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17. Key Words (Suggested by Author(s))		18. Distributi	on Statement			
Thermal modeling						
Monte Carlo techniques		فسعو	Distribution			
radiometry		F	DISCLIDENTION			
Earth Radiation Budget	Experiment (ERBE)					
Harri Maaracron 200800			Subjec	t Category 43		
		. <u></u>				
19. Security Classif. (of this report)	20. Security Classif. (of this	page)	21. No. of Pages	22. Price		
Unclassified	Unclassified		107			

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