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### **APPENDICS**

## APPENDIX A Finite element method

The finite element is the numerical method for calculating the approximate values of the mathematical problems. By dividing the area of the problem into the elements and creating the equation of each element corresponding to the partial differential equation. The elemental equations were summarized to the system of equations for setting the boundary conditions. Gauss elimination method was used to solve the approximate values from the system of equations. These phenomena of gas streamlines assisted the operators to design the distributor in the fluidized bed, since many unique properties of fluidized beds were related to the presence of bubbles.

The problems of gas streamlines in the fluidized bed were solved by the two-dimensional finite element method. These problems of gas streamlines were derived in term of the partial differential equation.

$$\frac{\partial}{\partial x} \left( k \frac{\partial \overline{\psi}}{\partial x} \right) + \frac{\partial}{\partial y} \left( k \frac{\partial \overline{\psi}}{\partial y} \right) = 0 \tag{A1}$$

If the exact value of gas streamlines,  $\psi$ , was known and substituted into the left-hand-side of Equation A1, the result in the right-hand-side was equal to zero. Since the exact value was not solved from the partial differential equation, the approximate value,  $\psi$ , substituted into the left-hand-side of Equation A1 resulted in the residue, R, in the partial differential equation

$$\frac{\partial}{\partial x} \left( k \frac{\partial \psi}{\partial x} \right) + \frac{\partial}{\partial y} \left( k \frac{\partial \psi}{\partial y} \right) = R \tag{A2}$$

In the Galerkin method, the finite element equations were created by multiplying the residue with the weighting function, W, and integrating around the area of the element to be zero.

$$\int_{\Omega} W_i R d\Omega = 0 \tag{A3}$$

Because triangular element consisted of three nodes, therefore, it was necessary to use three equations for the solution. By substituting Equation A2 into Equation A3:

$$\int_{\Omega} W_{i} \left( \frac{\partial}{\partial x} \left( k \frac{\partial \psi}{\partial x} \right) + \frac{\partial}{\partial y} \left( k \frac{\partial \psi}{\partial y} \right) \right) d\Omega = 0$$
 (A4)

and integrating Equation A4 by Gauss's theorem:

$$\int_{\Omega} u \left( \nabla \cdot \vec{V} \right) d\Omega = \int_{\Gamma} u \left( \vec{V} \cdot \hat{n} \right) d\Gamma - \int_{\Omega} \left( \nabla u \cdot \vec{V} \right) d\Omega$$
 (A5a)

where

$$u_i = W_i \tag{A5b}$$

$$\nabla = \frac{\partial}{\partial x} \hat{i} + \frac{\partial}{\partial y} \hat{j}$$
 (A5c)

$$\vec{V}$$
 =  $k \frac{\partial \psi}{\partial x} \hat{i} + k \frac{\partial \psi}{\partial y} \hat{j}$  (A5d)

$$\hat{n} = n_x \hat{i} + n_y \hat{j}$$
 (A5e)

$$\vec{V} \cdot \hat{n} = k \frac{\partial \psi}{\partial x} n_x + k \frac{\partial \psi}{\partial y} n_y$$
 (A5f)

$$u\left(\vec{V}\cdot\hat{n}\right) = W_{i}\left(k\frac{\partial\psi}{\partial x}n_{x} + k\frac{\partial\psi}{\partial y}n_{y}\right) \tag{A5g}$$

$$\nabla u = \frac{\partial W_i}{\partial x} \hat{i} + \frac{\partial W_i}{\partial y} \hat{j}$$
 (A5h)

$$\nabla \mathbf{u} \cdot \overrightarrow{\mathbf{V}} = \frac{\partial \mathbf{W}_{i}}{\partial \mathbf{x}} \mathbf{k} \frac{\partial \mathbf{\Psi}}{\partial \mathbf{x}} + \frac{\partial \mathbf{W}_{i}}{\partial \mathbf{y}} \mathbf{k} \frac{\partial \mathbf{\Psi}}{\partial \mathbf{y}}$$
(A5i)

Substituting Equation A5f-I into Equation A4 and assuming that  $W_i = N_i$ :

$$\int_{\Gamma} N_{i} \left( k \frac{\partial \psi}{\partial x} n_{x} + k \frac{\partial \psi}{\partial y} n_{y} \right) d\Gamma - \int_{\Omega} \left( \frac{\partial N_{i}}{\partial x} k \frac{\partial \Psi}{\partial x} + \frac{\partial N_{i}}{\partial y} k \frac{\partial \Psi}{\partial y} \right) d\Omega \qquad i = 1, 2, 3$$

(A6)

$$\int_{\Omega} \left( \frac{\partial N_i}{\partial x} k \frac{\partial \Psi}{\partial x} + \frac{\partial N_i}{\partial y} k \frac{\partial \Psi}{\partial y} \right) d\Omega = \int_{\Gamma} N_i \left( k \frac{\partial \Psi}{\partial x} n_x + k \frac{\partial \Psi}{\partial y} n_y \right) d\Gamma \qquad i = 1, 2, 3$$
(A7)

Since there were three equations in the element, the finite element equations were written in term of matrix.

$$\int_{\Omega} \left\{ \left\{ \frac{\partial N}{\partial x} \right\} k \frac{\partial \Psi}{\partial x} + \left\{ \frac{\partial N}{\partial y} \right\} k \frac{\partial \Psi}{\partial y} \right\} d\Omega = \int_{\Gamma} \{N\} \left( k \frac{\partial \Psi}{\partial x} n_x + k \frac{\partial \Psi}{\partial y} n_y \right) d\Gamma \quad (A8)$$

In each element, the distribution of gas streamlines in triangular element was assumed that

$$\psi = \psi(x_i, y_i) = \alpha_1 + \alpha_2 x_i + \alpha_3 y_i \quad i = 1,2,3 \quad (A9a)$$

or 
$$\psi(x, y) = [N_1(x, y) \ N_2(x, y) \ N_3(x, y)] \begin{cases} \psi_1 \\ \psi_2 \\ \psi_3 \end{cases}$$
 (A9b)

$$N_i = \frac{1}{2A}(a_i + b_i x + c_i y)$$
  $i = 1,2,3$  (A9c)

A = Area of triangular element

$$A = \frac{1}{2} \left[ x_2 (y_3 - y_1) + x_1 (y_2 - y_3) + x_3 (y_1 - y_2) \right] \quad (A9d)$$

$$\psi = \psi(x,y) = \left[ N \right] \left\{ \psi \right\}$$
 (A9e)

$$\frac{\partial \psi}{\partial x} = \left[\frac{\partial N}{\partial x}\right]_{(3\times 1)}^{\{\psi\}} \tag{A9f}$$

$$\frac{\partial \psi}{\partial y} = \left[\frac{\partial N}{\partial y}\right]_{(3\times 1)}^{\{\psi\}} \tag{A9g}$$

$$\frac{\partial N_i}{\partial x} = \frac{b_i}{2 A} \tag{A9h}$$

$$\frac{\partial N_i}{\partial y} = \frac{c_i}{2 A} \tag{A9i}$$

Substituting Equation A9f and A9g into Equation A8, the finite element equations became:

$$\int_{\Omega} \left\{ \frac{\partial N}{\partial x} \right\} k \left[ \frac{\partial N}{\partial x} \right] + \left\{ \frac{\partial N}{\partial y} \right\} k \left[ \frac{\partial N}{\partial y} \right] d\Omega \left\{ \psi \right\} = \int_{\Gamma} \left\{ N \right\} \left( k \frac{\partial \psi}{\partial x} n_x + k \frac{\partial \psi}{\partial y} n_y \right) d\Gamma$$
(A10)

From Equation A10, the finite element equations were written in the simple term of:

$$[K]_{e} \{\psi\}_{e} = \{F\}_{e}$$

$$(A11)$$

where subscript, e, was element matrix.

Substituting Equation A9h and A9i into  $[K]_{e}$  in Equation A11.

$$K_{ij} = \int_{A} k \left( \frac{\partial N_{i}}{\partial x} \frac{\partial N_{j}}{\partial x} + \frac{\partial N_{i}}{\partial y} \frac{\partial N_{j}}{\partial y} \right) dx dy \quad i,j = 1,2,3$$
 (A12)

$$K_{ij} = k \int_{A} \left( \frac{b_i}{2A} \frac{b_j}{2A} + \frac{c_i}{2A} \frac{c_j}{2A} \right) dx dy$$
 (A13)

$$K_{ij} = \frac{k}{4 A^2} (b_i b_j + c_i c_j) \int_A dx dy$$
 (A14)

$$K_{ij} = \frac{k}{4 A} (b_i b_j + c_i c_j)$$
  $i,j = 1,2,3$  (A15)

or 
$$\begin{bmatrix} K \\ {}_{(3\times3)} \end{bmatrix} = k A \begin{bmatrix} B \end{bmatrix}^T \begin{bmatrix} B \\ {}_{(3\times2)} \end{bmatrix}$$
 (A16)

$$[B] = \frac{1}{2A} \begin{bmatrix} b_1 & b_2 & b_3 \\ c_1 & c_2 & c_3 \end{bmatrix}$$
 (A17)

After all of elemental matrices were created, they were combined to the system of equations:

$$\sum$$
 (element equations)  $\Rightarrow$   $[K]_{sys} \{\psi\}_{sys} = \{F\}_{sys}$  (A18)

where subscript, sys, is the system of equations.

#### APPENDIX B Gauss elimination method

121 127 1251

Gauss elimination method was used to solve the values of gas streamlines in the system of equations. The solution consisted of 'n' equations and 'n' variables, was displayed in Equation B1a-n.

$$a_{11} x_1 + a_{12} x_2 + a_{13} x_3 + \dots + a_{1n} x_n = b_1$$
 (B1a)

$$a_{21} x_1 + a_{22} x_2 + a_{23} x_3 + \cdots + a_{2n} x_n = b_2$$
 (B1b)

$$a_{31} x_1 + a_{32} x_2 + a_{33} x_3 + \dots + a_{3n} x_n = b_3$$
 (B1c)

$$a_{n1} x_1 + a_{n2} x_2 + a_{n3} x_3 + \dots + a_{nn} x_n = b_n$$
 (B1n)

By using the forward elimination, Equation B1a was divided by the coefficient of Equation B1a,  $a_{11}$ , and multiplied with the coefficient of Equation B1b,  $a_{21}$ .

$$x_1 + \frac{a_{12}}{a_{11}} x_2 + \frac{a_{13}}{a_{11}} x_3 + \dots + \frac{a_{1n}}{a_{11}} x_n = \frac{b_1}{a_{11}}$$
 (B2)

$$a_{21} X_1 + a_{21} \frac{a_{12}}{a_{11}} X_2 + a_{21} \frac{a_{13}}{a_{11}} X_3 + \dots + a_{21} \frac{a_{1n}}{a_{11}} X_n = a_{21} \frac{b_1}{a_{11}}$$
 (B3)

Subtracting Equation B1a by Equation B3, Equation B1a became:

$$\left(a_{22} - a_{21} \frac{a_{12}}{a_{11}}\right) x_2 + \left(a_{23} - a_{21} \frac{a_{13}}{a_{11}}\right) x_3 + \dots + \left(a_{2n} - a_{21} \frac{a_{1n}}{a_{11}}\right) x_n = b_2 - a_{21} \frac{b_1}{a_{11}}$$
(B4)

or 
$$a'_{22}x_2 + a'_{23}x_3 + \cdots + a'_{2n}x_n = b'_2$$
 (B5)

To repeat the step of the calculation from Equation B1a to Equation B1b, the system of equations was:

$$a_{11} x_1 + a_{12} x_2 + a_{13} x_3 + \cdots + a_{1n} x_n = b_1$$
 (B6a)

$$a'_{22}x_2 + a'_{23}x_3 + \cdots + a'_{2n}x_n = b'_2$$
 (B6b)

$$a'_{32} X_2 + a'_{33} X_3 + \cdots + a'_{3n} X_n = b'_3$$
 (B6c)

$$a'_{n2}x_2 + a'_{n3}x_3 + \cdots + a'_{nn}x_n = b'_n$$
 (B6n)

From the first round of elimination method, all values in the first column of Equation B6b-n except Equation B6a were equal to zero. The problem was solved by using forward elimination from the second round to the n-1 round of elimination method until the system of equations was proper to solve the results of streamlines by back substitution.

$$a_{11} x_1 + a_{12} x_2 + a_{13} x_3 + \dots + a_{1n} x_n = b_1$$
 (B7a)

$$a'_{22}x_2 + a'_{23}x_3 + \cdots + a'_{2n}x_n = b'_2$$
 (B7b)

$$a_{33}''x_3 + \cdots + a_{3n}''x_n = b_3''$$
 (B7c)

. . .

$$a_{nn}^{(n-1)} x_n = b_{nn}^{(n-1)}$$
 (B7n)

From the system of Equation B7a-n, the result of  $x_n$  was calculated directly from Equation B7n:

$$x_n = \frac{b_n^{(n-1)}}{a_{nn}^{(n-1)}}$$
 (B8)

and the results of  $x_{n-1}, x_{n-2}, ..., x_2, x_1$  were computed by:

$$x_{i} = \frac{b_{i}^{(i-1)} - \sum_{j=i+1}^{n} a_{ij}^{(i-1)} x_{j}}{a_{ii}^{(i-1)}}$$
(B9)

## APPENDIX C Finite element code

C

To check a number of nodes.

```
C
         PROGRAM FINITE
         PARAMETER (MXNODE=300, MXELE=600)
         IMPLICIT REAL*8 (A-H,O-Z)
         DIMENSION COORD(MXNODE,2), SL(MXNODE), TEXT(20)
         DIMENSION SYSK(MXNODE, MXNODE), SYSF(MXNODE)
         DIMENSION PER(MXELE)
         INTEGER NNUM(MXELE,3), COND(MXNODE)
· C
         MXNODE
                            maximum number of nodes
  \mathbf{C}
                            maximum number of elements
         MXELE
                            coordinates of nodes
  \mathbf{C}
         COORD
  \mathbf{C}
         SL
                            streamline
         SYSK
  \mathbf{C}
                            [K]_{svs}
  \mathbf{C}
         SYSF
                            \{F\}_{svs}
  \mathbf{C}
         PER
                            permeability
  \mathbf{C}
         NNUM
                            nodal numbers were combined to one element
  \mathbf{C}
         COND
                            condition of calculation
  \mathbf{C}
         If the users knew the gas streamline in this point, select
                                                                        1.
  C
         If the users do not knew the gas streamline in this point, select
                                                                        0.
         OPEN(7,FILE='INPUT.DAT')
  \mathbf{C}
         To open 'input.dat' file to read the input data.
  101
         FORMAT(20A4)
         READ(7,101) TEXT
  \mathbf{C}
         To read character.
         READ(7,*) NODE, NELEM
         To read a number of nodes and elements.
  \mathbf{C}
         IF(NODE.GT.MXNODE) WRITE(6,102) NODE
```

102 FORMAT(/,' PLEASE INCREASE THE PARAMETER MXNODE TO' \* , I5) IF(NODE.GT.MXNODE) STOP IF(NELEM.GT.MXELE) WRITE(6,103) NELEM  $\mathbf{C}$ To check a number of elements. FORMAT(/,' PLEASE INCREASE THE PARAMETER MXELE TO' 103 \* , I5) IF(NELEM.GT.MXELE) STOP READ(7,101) TEXT  $\mathbf{C}$ To read character. DO 104 IP=1,NODE READ(7,\*) I, COND(I), (COORD(I,K), K=1,2), SL(I)To read nodal numbers, conditions, nodal coordinates and value of gas  $\mathbf{C}$  $\mathbf{C}$ streamlines. IF(I.NE.IP) WRITE(6,105) IP C To check the data of each node. 105 FORMAT(/, 'NODE NO. ',I5, 'IN DATA FILE IS MISSING') IF(I.NE.IP) STOP 104 **CONTINUE** READ(7,101) TEXT To read character. C

READ(7,\*) I, (NNUM(I,J), J=1,3),PER(I)

C To check element numbers, nodal numbers, permeability.

IF(I.NE.IE) WRITE(6,107) IE

C To check the data of each element.

DO 106 IE=1,NELEM

- 107 FORMAT(/, 'ELEMENT NO. ', 15, 'IN DATA FILE IS MISSING')
  IF(I.NE.IE) STOP
- 106 CONTINUE

WRITE(6,108) 108 FORMAT(' THE FINITE ELEMENT PROGRAM') NEQ=NODE C A number of equations in the solution equaled the number of nodes. DO 109 I=1,NEQ SYSF(I)=0.To clear the values of  $\{F\}_{sys}$ .  $\mathbf{C}$ 109 **CONTINUE** DO 110 I=1,NEQ DO 110 J=1,NEQ SYSK(I,J)=0.C To clear the values of [K]<sub>sys</sub>. 110 **CONTINUE** CALL ELEMENT(NELEM, NNUM, COORD, PER, SYSK, \* MXNODE, MXELE) To call subroutine for calculating elemental matrices. C CALL BOUNDARY(NODE, COND, SL, SYSK, SYSF, MXNODE)  $\mathbf{C}$ To call subroutine for setting the boundary conditions. CALL GAUSS(NEQ, SYSK, SYSF, SL, MXNODE) To call subroutine for solving the results of gas streamlines.  $\mathbf{C}$ OPEN(8,FILE='Output.DAT')  $\mathbf{C}$ To open 'output.dat' file to record the results of gas streamline. WRITE(8,111) FORMAT('NODE',10X,", 'X',10X, 'Y',10X, 'STREAMLINE') 111 DO 112 IP=1,NODE WRITE(8,113) IP,COORD(IP,1),COORD(IP,2),SL(IP)

To print the nodal numbers, coordinates and streamlines.

FORMAT(I6, F10.5,F10.5,F10.5)

**CONTINUE** 

C

113

112

**STOP** 

**END** 

C-----

SUBROUTINE ELEMENT(NELEM,NNUM,COORD,PER,

\* SYSK, MXNODE, MXELE)

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

DIMENSION COORD(MXNODE,2), SYSK(MXNODE,MXNODE)

DIMENSION AKC(3,3), B(2,3), BT(3,2)

**INTEGER NNUM(MXELE,3)** 

**DIMENSION PER(NELEM)** 

$$C \qquad B \qquad = \qquad \begin{bmatrix} b_1 & b_2 & b_3 \\ c_1 & c_2 & c_3 \end{bmatrix}$$

$$C \qquad BT \qquad = \qquad \begin{bmatrix} b_1 & c_1 \\ b_2 & c_2 \\ b_3 & c_3 \end{bmatrix}$$

C Area of the element

C 
$$\psi(x,y) = [N_1(x,y) \quad N_2(x,y) \quad N_3(x,y)] \begin{bmatrix} \psi_1 \\ \psi_2 \\ \psi_3 \end{bmatrix}$$

$$C N_i = \frac{1}{2A} (a_i + b_i x + c_i y)$$

C 
$$a_1 = x_2y_3 - x_3y_2$$
  $b_1 = y_2 - y_1$   $c_1 = x_3 - x_2$ 

C 
$$a_2 = x_3y_1 - x_1y_3$$
  $b_2 = y_3 - y_1$   $c_2 = x_1 - x_3$ 

C 
$$a_3 = x_1y_2 - x_2y_1$$
  $b_3 = y_1 - y_2$   $c_3 = x_2 - x_1$ 

DO 201 IE=1,NELEM

II=NNUM(IE,1)

JJ=NNUM(IE,2)

KK=NNUM(IE,3)

C Combining nodal numbers into the one element.

XG1=COORD(II,1)

```
XG2=COORD(JJ,1)
       XG3=COORD(KK,1)
       YG1=COORD(II,2)
       YG2=COORD(JJ,2)
       YG3=COORD(KK,2)
\mathbf{C}
       Coordinates of node1, 2, 3 of each element
       AREA=0.5*(XG2*(YG3-YG1) + XG1*(YG2-YG3) +
    * XG3*(YG1-YG2))
       IF(AREA.LE.0.) WRITE(6,202) IE
       To check nodal connection of each element.
C
202
       FORMAT(/,'!!! ERROR!!! ELEMENT NO.', I5,
    * 'HAS NEGATIVE OR ZERO AREA', /,
    * ' --- CHECK F.E. MODEL FOR NODAL COORDINATES',
    * 'AND ELEMENT NODAL CONNECTIONS ---'
       IF(AREA.LE.0.) STOP
       B1=YG2-YG3
       B2=YG3-YG1
       B3=YG1-YG2
       C1=XG3-XG2
       C2=XG1-XG3
       C3=XG2-XG1
       DO 203 I=1,2
       DO 203 J=1,3
       B(I,J)=0.
       To clear the values of b<sub>i</sub>, c<sub>i</sub>.
C
203
       CONTINUE
       B(1,1)=B1
       B(1,2)=B2
       B(1,3)=B3
```

$$B(2,1)=C1$$

$$B(2,2)=C2$$

$$B(2,3)=C3$$

$$B(I,J)=B(I,J)/(2.*AREA)$$

C 
$$\frac{\partial N_i}{\partial x} = \frac{b_i}{2 A}$$
 for  $I = 1$ ,  $\frac{\partial N_i}{\partial y} = \frac{c_i}{2 A}$  for  $I=2$ 

$$BT(J,I)=B(I,J)$$

C 
$$\frac{\partial N_j}{\partial x} = \frac{b_{ji}}{2 A}$$
 for  $I = 1$ ,  $\frac{\partial N_j}{\partial y} = \frac{c_j}{2 A}$  for  $I = 2$ 

205 CONTINUE

204 CONTINUE

DO 206 I=1,3

DO 206 J=1,3

AKC(I,J)=0.

DO 207 K=1,2

$$AKC(I,J)=AKC(I,J)+BT(I,K)*B(K,J)$$

$$C \qquad AKC = \frac{1}{4A^2} (b_i b_j + c_i c_j)$$

207 CONTINUE

AKC(I,J)=PER(IE)\*AREA\*AKC(I,J)

C AKC = 
$$K_{ij} = \frac{k}{4 A} (b_i b_j + c_i c_j)$$

206 CONTINUE

C 
$$[K]_e \{\psi\}_e = \{F\}_e$$

$$(3\times 3) \quad (3\times 1)$$

CALL MATRIX(IE,NNUM,AKC,SYSK,MXNODE,MXELE)

C To call subroutine for creating the system of equations.

201 CONTINUE

**RETURN END SUBROUTINE** MATRIX(IE,NNUM,AKC,SYSK,MXNODE,MXELE) IMPLICIT REAL\*8 (A-H,O-Z) **DIMENSION AKC(3,3)** DIMENSION SYSK(MXNODE, MXNODE) INTEGER NNUM(MXELE,3) NNODE=3 DO 301 IR=1,NNODE DO 302 IC=1,NNODE IROW=NNUM(IE,IR) ICOL=NNUM(IE,IC) SYSK(IROW,ICOL)=SYSK(IROW,ICOL)+AKC(IR,IC)  $\sum$  (element equations)  $\Rightarrow$   $[K]_{sys} \{\psi\}_{sys} = \{F\}_{sys}$ C 302 **CONTINUE** 301 **CONTINUE RETURN END** SUBROUTINE BOUNDARY(NODE, COND, SL, SYSK, \* SYSF, MXNODE) IMPLICIT REAL\*8 (A-H,O-Z) DIMENSION SYSK(MXNODE, MXNODE), SYSF(MXNODE), \* SL(MXNODE) INTEGER COND(MXNODE) DO 401 IEQ=1,NODE IF(COND(IEQ).EQ.0) GOTO 401

```
C
       To check conditions of calculation
       DO 402 IR=1,NODE
       IF(IR.EQ.IEQ) GOTO 402
       SYSF(IR)=SYSF(IR) - SYSK(IR,IEQ)*SL(IEQ)
       To eliminate the values of \{F\}_{sys} after the boundary conditions were
C
C
       applied in the system of equations.
       SYSK(IR,IEQ)=0.
       The values of [K]<sub>svs</sub> in 'IR' row equaled zero except K<sub>IR,IEQ</sub>.
C
       CONTINUE
402
       DO 403 IC=1,NODE
       SYSK(IEQ,IC)=0.
       The values of [K]<sub>svs</sub> in 'IC' column equaled zero.
C
       CONTINUE
403
       SYSK(IEQ,IEQ)=1.
       K_{IO,IEO} in [K]_{sys} = 1
C
       SYSF(IEQ)=SL(IEQ)
       F_{IEO} in \{F\}_{sys} = 1
C
       CONTINUE
401
       END
       SUBROUTINE GAUSS(NEQ, SYSK, SYSF, SL, MXNODE)
       IMPLICIT REAL*8 (A-H,O-Z)
       DIMENSION SYSK(MXNODE, MXNODE), SYSF(MXNODE),
     * SL(MXNODE)
       CALL SCAL(NEQ, SYSK, SYSF, MXNODE)
       To call subroutine for dividing the equations in [K]<sub>sys</sub> by using the
C
C
       maximum coefficient of each row.
       DO 501 IP=1,NEQ-1
       CALL PIVOT(NEQ, SYSK, SYSF, MXNODE, IP)
```

```
\mathsf{C}
       To call subroutine for rearranging the equations.
       DO 502 IE=IP+1,NEQ
       RATIO=SYSK(IE,IP)/SYSK(IP,IP)
       DO 503 IC=IP+1,NEQ
       SYSK(IE,IC)=SYSK(IE,IC)-RATIO*SYSK(IP,IC)
       Forward elimination of [K]<sub>svs</sub>
C
503
       CONTINUE
       SYSF(IE)=SYSF(IE)-RATIO*SYSF(IP)
C
       Forward elimination of [F]<sub>svs</sub>
502
       CONTINUE
       DO 504 IE=IP+1,NEQ
       SYSK(IE,IP)=0.
504
       CONTINUE
501
       CONTINUE
       SL(NEQ)=SYSF(NEQ)/SYSK(NEQ,NEQ)
\mathbf{C}
       Back substitution of gas streamlines at 'n' node.
       DO 505 IE=NEQ-1,1,-1
       SUM=0.
       DO 506 IC=IE+1,NEQ
       SUM=SUM+SYSK(IE,IC)*SL(IC)
       Back substitution of gas streamlines except 'n' node.
C
506
       CONTINUE
       SL(IE)=(SYSF(IE)-SUM)/SYSK(IE,IE)
505
       CONTINUE
       RETURN
       END
       SUBROUTINE SCAL(NEQ, SYSK, SYSF, MXNODE)
       IMPLICIT REAL*8 (A-H,O-Z)
```

```
DIMENSION SYSK(MXNODE, MXNODE), SYSF(MXNODE)
       DO 601 IE=1,NEQ
       BIG=ABS(SYSK(IE,1))
       To set the first coefficient of [K]<sub>sys</sub> was maximum coefficient of each
\mathbf{C}
C
       row.
       DO 602 IC=2,NEQ
       AMAX=ABS(SYSK(IE,IC))
       IF(AMAX.GT.BIG) BIG=AMAX
       To search the maximum coefficient of each row.
\mathbf{C}
602
       CONTINUE
       DO 603 IC=1,NEQ
       SYSK(IE,IC)=SYSK(IE,IC)/BIG
       To divide the coefficient of [K]<sub>sys</sub> by the maximum coefficient of
\mathbf{C}
\mathbf{C}
       each row.
603
       CONTINUE
       SYSF(IE)=SYSF(IE)/BIG
       To divide the coefficient of \{F\}_{sys} by the maximum coefficient of
\mathbf{C}
\mathbf{C}
       each row.
601
       CONTINUE
       RETURN
       END
       SUBROUTINE PIVOT(NEQ, SYSK, SYSF, MXNODE, IP)
       IMPLICIT REAL*8 (A-H,O-Z)
       DIMENSION SYSK(MXNODE, MXNODE), SYSF(MXNODE)
       JP=IP
       BIG=ABS(SYSK(IP,IP))
       DO 701 I=IP+1,NEQ
       AMAX=ABS(SYSK(I,IP))
```

IF(AMAX.GT.BIG) THEN

BIG=AMAX

JP=I

C To search the maximum coefficient of  $[K]_{sys}$  in each column.

**ENDIF** 

701 CONTINUE

IF(JP.NE.IP) THEN

DO 702 J=IP,NEQ

DUMY=SYSK(JP,J)

SYSK(JP,J)=SYSK(IP,J)

SYSK(IP,J)=DUMY

702 CONTINUE

C To convert the coefficient in  $[K]_{sys}$ .

DUMY=SYSF(JP)

SYSF(JP)=SYSF(IP)

SYSF(IP)=DUMY

To convert the coefficient in  $\{F\}_{sys}$ .

**ENDIF** 

**RETURN** 

**END** 

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