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**A Second Generation Frontal
Solution Program**

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A SECOND GENERATION FRONTAL SOLUTION PROGRAM

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Synopsis

A computer program for the assembly and solution of symmetric positive definite equations as met in the Finite Element analysis based on the Frontal Solution algorithm by Irons is presented.

The program features improved direct access blocked I/O and the use of Front partitioning which makes the problem size which can be solved practically independent of the size of the computer memory. In addition the use of fast vector processors is considered which should improve CPU times considerably.

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1. INTRODUCTION

The Frontal Solution technique is based on the Gaussian Elimination method and was first published in 1970 by Irons (1).

Various authors (2,3,4,5) have since pointed out various advantages of this technique. The main advantage over band or skyline (6) solvers seems to be a simplification in the data preparation as the numbering of nodes is not restricted to minimise a band width. Also, since the node numbers are treated as "Nicknames", design changes (i.e. adding or removing elements) may be made without having to renumber the nodes. Solution time for the Frontal solver is now sensitive to the numbering of the Elements but the sequence is a natural one. The famous example of a ring structure is often mentioned in this context (1).

The main difference between the Frontal technique and a conventional band solution lies in the manner in which the structure stiffness coefficients are stored and in the order in which the equations are eliminated.

Consider, for example, a patch of 4 node/8 D.O.F. elements in Fig. 1 in which the degrees of freedom are numbered from 1 to 18.

The stiffness coefficients of Element I are stored in the order of appearance (local node numbering) and in the manner shown in Fig. 1a. Each variable has a "Destination" which determines the position of its coefficients in the Front matrix. An asterisk, *, marks the equations which are already fully summed. These variable(s) can now be eliminated before the next element is assembled by treating the other equations with:

$$c'_{i,j} = c_{i,j} - c_{n,j}^* \frac{c_{n,i}^*}{c_{n,n}^*} \quad (1)$$

Where n is the destination of the variable to be eliminated and the * denotes the coefficients which are fully summed. The reader should note that it does not matter that the coefficients $c_{i,j}$ are not in their final form since only the order in which the coefficients are added is changed.

After elimination, variables 1 and 10 cease to be "active" and the corresponding equations are transferred into buffer storage ready for output on disk. The storage locations of these equations in the Front matrix are cleared (Fig. 1b). On assembly of element II this space is re-used by Equation 3 and 6 (Fig. 1c). The Front matrix thus contains only the coefficients of "active" variables.

With the principle involved explained the reader may complete the example making the following observations:

- (1) Assembly and elimination order is governed by the order in which the variable coefficients are entered as one Element is assembled after the other. A variable is eliminated as soon as the coefficients are fully summed (i.e. on its last appearance). The position of the coefficients in the Front matrix is governed by the empty spaces available.
- (2) The storage requirement for storing the "active" coefficients is determined by the largest address used. Because of symmetry only one half of the Front matrix has to be stored and the storage requirement may be computed from

$$l = m(m + 1) / 2 \quad (2)$$

Here m is the largest "Destination" of a variable (Front width).

In the original code Irons assumes that the Front matrix resides in central storage (Comment on the program listing: "if not, buy larger computer").

It can be seen that the storage requirement increases with the square of the Front width. This puts a severe limitation on the size of problems which can be solved on a special computer.

The purpose of this paper is to present a program where this limitation has been over-come making it possible to solve large 3-D problems on a mini computer. In addition, the

transfer of data to and from disk is improved by using a blocked direct access I/O mode.

The inner-most DO-loop of the Gauss elimination is written suitable for fast vector processors which have appeared on the market in recent years.

The resulting program should not only be an improvement in solution capability but also in performance.

2. STORAGE AND BUFFERING DURING ELIMINATION

Similar to the original code by Irons, a working vector ELPA is used. ELPA is divided into 3 main areas which contain: (1) Element stiffness matrix of the Element to be assembled; (2) Front matrix or a partition of the front matrix; (3) Buffer for the equations which have ceased to be "active".

Allocation of the space for these areas is completely flexible and will depend on the type of problem solved. Whereas the space of first area is fixed by the size of the element stiffness matrix, the space allocation for areas 2 and 3 must be adjusted to give optimal solution times within the core limitations of the computer used. This will be discussed in detail later.

The number of equation coefficients in the Front matrix for a particular Front width is computed from Equation 2 and if it is greater than the space available, the partitioning algorithm has to be activated.

The number of Equations Δl_N which fit into a particular N can be computed from the inequality

$$\Delta l_N (\Delta l_N + 1) / 2 + (k_N - 1) * \Delta l_N < n_p \quad (3)$$

Here k_N is the destination of the first equation in partition N and n_p is the size of the storage space for the Front matrix. Solution of Equation 3 gives

$$\Delta l_N = \text{INT} \left[-\frac{2k_N - 1}{2} + \sqrt{\left[\frac{2k_N - 1}{2}\right]^2 + 2n_p} \right] \quad (4)$$

Where INT means the truncation of the result. This allows us to determine the limits of each partition i.e. partition N can accommodate Equation k_N to k_{N+1} .

For a coefficient k, i ($k < i$) residing in partition N, where

$$k_N < k < k_{N+1} \quad (5)$$

the address, ℓ , is computed from (see Fig. 2),

$$\ell = \ell_0 + (k - k_N) (k_N - 1) + (k - k_N) (k - k_N + 1)/2 + i \quad (6)$$

Here ℓ_0 specifies the start address of the space reserved for the Front matrix.

3. PROGRAMMING STRATEGY

As long as the Front matrix is small and fits into core the program strategy is simple and follows three basic steps for each element:

- (1) Read element stiffness matrix
- (2) Assemble its coefficients into Front matrix
- (3) Eliminate variables which are ready

The elimination essentially consists of two steps. First, the equation coefficients of the variable which is to be eliminated (n) are moved to the equation buffer. When the buffer is full its contents are written on disk and the pointer reset to the beginning of the buffer before the coefficients are moved (It should be noted at this stage that all the coefficients to the right of the minus sign in Equation 1 are now in the buffer). Then all the equations in the Front matrix are modified by Equation 1.

When the current Front matrix becomes large and no longer fits in the allocated space, partitioning is invoked automatically. The program strategy becomes more complex and follows the following basic steps:

- (1) Read Element stiffness matrix.
- (2) Swap Partition 1 into core and assemble all coefficients which are resident in this partition.
- (3) Eliminate variables which are ready in this partition (if any).
- (4) Swap Partition 2 into core and assemble further coefficients.
- (5) Modify all coefficients for the variables eliminated previously (if any) using the coefficients c_{nl}^* in the Equation buffer. This is referred to as elimination of 'old variables' in the listing.
- (6) Eliminate variables which are ready in this partition ('new variables').

Steps 4, 5 and 6 are repeated for all subsequent partitions to the last one. Note that at step 3 not all the coefficients c_{nl}^* of the equation n are available and the equation in the buffer is still incomplete. Thus, additional coefficients have to be transferred in step 5.

Our work is not completed yet since some lower partitions have not been modified due to elimination of variables in higher partitions. So we have to retrack and modify the Equations which have not yet been modified using the coefficients c_{nl}^* in the equation buffer.

4. I/O OPERATIONS DURING ELIMINATION

When the Front matrix is small and fits into the allocated space, the I/O operations are simple:

- (1) Read the assembly information for each element (The 'destination vector' is coded to indicate when each variable is ready for elimination).
- (2) Read Element stiffness matrix.
- (3) Whenever the equation buffer is full, i.e. no further equation fits, write its contents onto disk.

In the present program, the standard Fortran READ is used for operations 1 and 2. For large amounts of data the speed of the I/O operation depends greatly on the transfer mode. It has been found by the author that on a Data General Eclipse mini-computer, the data transfer is 10 times faster when the machine dependent routines RDBLK and WRBLK are used. The prerequisite for using these routines is that the number of coefficients to be transferred is divisible by the physical block size on disk (128 real numbers in this case). Since operation 3 may involve a large number of coefficients, a blocked I/O mode is used. The equation buffer is divided into a number of blocks and the space allocated for it should be a multiple of the block size. When the buffer is full it may, however, not always fill the last block completely. To avoid empty spaces on disk the last block is not written in this case, but the coefficients are rather transferred to core to the beginning of the buffer with new coefficients moved into the subsequent spaces.

In this context it should be noted that in core transfers are typically a factor of 10^3 faster than out-of-core transfers (i.e. transfers to and from disk).

When the Front matrix becomes too large and no longer fits into core, I/O operations become more complex and frequent. The Front matrix has to be swapped in and out of core as required. Blocked I/O transfer now becomes essential and the programming critical since a program slowed down by too many I/O operations may no longer be competitive.

The aim is to reduce the I/O operations to a bare minimum even if this means that more in core operations are necessary to do this (see above statement about I/O transfer speeds).

The number of swapping operations on the partitions can be determined from the basic steps delineated in the last chapter and depends on the location of variables which are to be eliminated. In the worst case, we need $(2N - 1)$ swaps where N is the number of partitions currently used. This number is critical for the performance of the program as partitions usually involve a large number of coefficients. Ways to optimise the number of I/O operations are discussed in the next chapter.

A further complication arises which could endanger the economy of the program. For the basic step 5, (elimination of 'old variables') the coefficients $c_{n\ell}^*$ which are thought to reside in the equation buffer are needed. But in the meantime, I/O operation 3 (transfer of equation buffer on to disk when full) may have been carried out and the required coefficients may no longer be in core. Thus we must keep track of which coefficients are in core and which are on disk. If the coefficients are no longer in core they must be swapped into core and this requires additional I/O operations, the number of which depends greatly on the size of the equation buffer. If the buffer is very large then swapping of the equation buffer may occur only rarely.

5. OPTIMISATION OF I/O OPERATIONS AND CHOICE OF FRONT PARTITION AND EQUATION BUFFER LENGTH

When selecting the size of the buffer for the Front matrix, we must aim to avoid partitioning since it is expensive. If no partitioning is involved the equation buffer may be made small to accommodate a big Front matrix, the only restriction being that, the buffer has to be at least 2 blocks long and accommodate the largest equation. On the other hand, when partitioning of the Front is unavoidable because of core restrictions or size of problem, there is a case for decreasing the partition size in favour of a large equation buffer for the reasons explained in the last chapter.

To reduce the number of I/O operations further, a number of situations where swapping is not required is examined. Swapping of partitions is not required when:

- (A) There are no coefficients to be assembled into the partition, and no variables have been eliminated yet in the current element loop.
- (B) There are no coefficients to be assembled and the coefficients $c_{n\ell}^*$ ($k_N < \ell > k_{N+1}$) are zero.
- (C) Swapping of the equation buffer is not required when the coefficients $c_{n\ell}^*$ ($k_N < \ell > k_{N+1}$) in the Front partition are zero.

A check on conditions C and B is made by Subroutine SAVES in the program.

6. SHORT DESCRIPTION OF COMPUTER PROGRAM

The program consists of two main subroutines PREFR and SFRONT. The subroutine PREFR works out the coded destinations of the variables and writes them onto disk. This program is essentially the same as published by Irons and is included for completeness. Subroutine SFRONT performs the assembly and reduction of the structure or substructure stiffness matrix as detailed in the last chapters.

It uses the following subroutines:

MOVE to move equation coefficients to buffer.
 GAUSS to modify the coefficients in the Front matrix with Eq. (1).
 ASSEMB ... to assemble stiffness coefficients into the Front matrix.
 EMPDI to empty equation buffer on disk when full.
 RESBUF ... to reset buffer pointer.
 SWAPF to swap Front partitions in and out of core.
 SAVES to save swaping (see last chapter).
 UNCOD to uncode coded destinations (uses CODEST)
 PALI to work out partition limits (k_N, k_{N+1}).

In addition the following functions are used:

LADDR(M) is the local address in the current Front partition of a coefficient M, M.
 LADST(I,J) .. is the local address of a coefficient i,j in the Element stiffness matrix.

Subroutines are also used to clear integer and real arrays and write error messages.

The blocked I/O operations are performed by subroutine BLKIO which has a machine dependent coding. Files are opened and channel numbers assigned by FILO which is also machine dependent.

The fast vector operations are performed by subroutine SVECT. When a computer with vector processor is used, the appropriate coding as given in the Users Manual of the machine should be inserted here. For use on machines without this capability the Standard Fortran coding may be used as shown.

The computer program is listed in Appendix B. A list of some important arrays and variables is given in Appendix A. In addition, a program is included to test and demonstrate the substructure capability of the sub-routine SFRONT in Appendix C.

7. SUBSTRUCTURING WITH THE FRONTAL SOLUTION

For very large structures, it is often desirable to divide the mesh into several smaller meshes or substructures. These are treated as large elements and the boundary stiffness matrix obtained by elimination of the 'internal' degrees of freedom.

The substructuring has the following main advantages:

- (1) The process of solving the structure is a continuous one and errors may be detected at substructure level. Remedial actions need only to be taken in the particular substructure involved.
- (2) Sometimes a structure consists of many subareas having a similar geometry. Thus the stiffness matrix of a particular type of substructure may be computed only once and the main structure assembled with as frequent re-use of the substructure stiffness as possible.
- (3) For excavation type of problems in rock or soil mechanics the substructuring technique offers additional advantages. By defining the rock or soil mass in the full excavation as one large substructure and the material to be excavated at each stage as smaller substructures the analysis of each excavation stage just requires the assembly of substructure stiffnesses and the solution for the substructure boundary degrees of freedom.

Substructuring with the Front Solution is relatively simple. All that has to be done is to suspend the elimination of selected variables at the boundary of the substructure. The coefficients which remain in the Front matrix after the elimination of all other variables then constitute the stiffness coefficients for the super element. After suitable reordering, the stiffness matrices of all super elements can be obtained and assembled in the usual manner to solve for the complete structure.

Thus, substructuring involves the basic steps.

- (1) The PREFRONT subroutine read the substructure "Nicknames" into the vector NIX. This will modify the coding of the destinations of the substructure variables in such a way as to prevent their elimination.
- (2) Perform the usual assembly and elimination for all elements which make up the substructure.
- (3) Remove zero rows and columns from the Front matrix and reorder to obtain the substructure stiffness matrix in condensed form.

After this has been done for all substructures, perform the assembly and elimination in the usual way but this time involving all substructures which make up the structure to be analysed.

The substructuring capability is demonstrated with a test program in Appendix II where the substructure consists of a regular assembly of 4 node/8 degrees of freedom Elements.

8. RE-SOLUTION

Once the global stiffness matrix has been reduced and stored a re-resolution for as many load cases as desired can be made.

It is convenient to separate the resolution and back substitution part completely from the reduction of the left hand

side in order to have as much space available as possible. Because the size of the vector needed for each load case is only MAXPA no partitioning of the Front should be required even for large problems and the basic procedures are as follows:-

- (1) Read the Element right hand side (RHS) into the first part of ELPA.
- (2) Assemble into the space reserved for the Front-RHS.
- (3) Reduce RHS using the coefficients c_n^* on disk i.e. modify the Front RHS with

$$F_i^1 = F_i - \frac{c_{in}^*}{c_{n,m}^*} F_n^* \quad (7)$$

The procedure is exactly the same as a non-partitioned reduction except that vectors are involved instead of matrices.

The results are obtained in Element form by back substitution i.e.

$$x_n = \frac{1}{c_{nn}^*} \left(\sum_{i=n} c_{n,i} - F_n^* \right) \quad (8)$$

in the same manner as by Irons.

9. FURTHER FACILITIES OF THE COMPUTER PROGRAM AND DISCARDED FACILITIES

This section deals with features which are included in the present program and facilities which have not been considered but can be implemented easily.

9.1 Treatment of Constraints

In the present program a restrained degree of freedom is treated by setting the corresponding destination to zero and thereby preventing the assembly of the corresponding equation.

This is the simplest and most economical way. Various other types of constraints, as shown by Irons (1) can be easily implemented.

9.2 Computation of the Determinant of the Structure Stiffness Matrix

This is often required for vibration and stability analysis and is incorporated by additional coding in Subroutine GAUSS. After elimination the value of $\log_{10} K/$ is stored in the variable DET. In addition the frequency of the occurrence of a negative diagonal element is determined and stored in NEG. If NEG is odd the sign of the determinant is positive otherwise negative. The variables DET and NEG are in Common block/ EIGEN/.

9.3 Check on Singularity and Indefiniteness

A check on singularity and indefiniteness is made during elimination. If the diagonal coefficient is less than or equal to zero an Error message is produced. Because of machine accuracy, the diagonal coefficient will not be exactly zero even for a singular matrix. More appropriate checks have been suggested (6), that is,

(1) Singularity

$$d_j < t_j \quad (9)$$

(2) Indefiniteness

$$d_j < -t_j \quad (10)$$

where d_j is the j -th diagonal element at the j -th elimination stage and

$$t_j = 8\epsilon r_j$$

where ϵ is the smallest positive floating point number for which $(1 + \epsilon) > 1$ on the computer used and r_j is the norm of the j -th row of K . This can be easily implemented in Subroutine GAUSS if the machine accuracy ϵ is known.

9.4 Check on Accuracy of Solution

In the original code by Irons a simple roundoff criterion was included. The author has found this criterion not entirely satisfactory because it is not sensitive to right hand sides and was found to register only if the difference in stiffness is too great between elements.

A better *a priori* estimate of the matrix condition is the Euclidian condition number (6). But this also involves additional unproductive computation and may be expensive.

The author favours the *a posteriori* estimate by one step of iterative refinement of the solution because it is a more productive method giving not only an estimate of the accuracy but also an improved solution. It only involves a re-solution and matrix multiplication. The iterative refinement may be made only for one load case and not repeated for the other load cases if the condition number is satisfactory.

First, the load case is solved with the re-solution facility to give x_i^0 , the unrefined result. Then the residual forces are worked out:

$$R_i^0 = F_i - K_{ij} x_j^0 \quad (11)$$

A second resolution with R_i^0 as new right hand side will give the error on x_j^0 , Δx_j^0 .

The expression

$$\|\Delta x_j^0\| / \|x_j^0\| \quad (12)$$

provides an estimate on the accuracy of the solution x_j^0 .

10. CONCLUSIONS

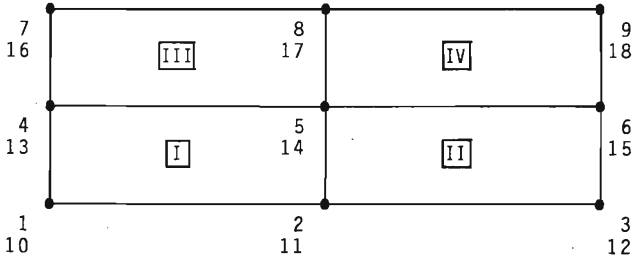
A computer program for the assembly and solution of a symmetric positive definite set of equations has been presented. The program is based on the Frontal Solution technique by Irons but uses frontal partitioning to make the problem size which can

be solved practically independent of the memory size of the computer used.

In addition, a great deal of effort has been made to optimise the I/O operations during partitioned elimination. Fast vector or pipeline processing has also been considered in the coding.

The resulting program is an improvement, not only in capability but also in performance. The program should be useful not only in mini-computer applications but also for large computers, because a reduction or optimisation of the band width is not required in the Frontal solution.

The solution time and storage requirement is influenced only by the numbering of the Elements which is a natural one.



+ non-zero entry
 P diagonal
 0 zero coeff.

EQUATION	DESTINATION	1	2	3	4	5	6	7	8
1	1*	P							
2	2	+ P							
5	3	+ + P							
4	4	+ + + P							
10	5*	+ + + + P							
11	6	+ + + + + P							
14	7	+ + + + + P							
13	8	+ + + + + + P							

(a)

EQUATION	DESTINATION	1	2	3	4	5	6	7	8
□	1	0							
2	2	0 P							
5	3	0 + P							
4	4	0 + + P							
□	5	0 0 0 0 0							
11	6	0 + + + 0 P							
14	7	0 + + + 0 + P							
13	8	0 + + + 0 + + P							

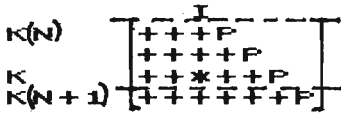
(b)

EQUATION	DESTINATION	1	2	3	4	5	6	7	8	9	10
□	1	P									
2	2	+ P									
5	3	+ + P									
4	4	0 + + P									
□	5	+ + + 0 P									
11	6	+ + + + + P									
14	7	+ + + + + P									
13	8	0 + + + 0 + + P									
11	9	+ + + 0 + + + 0 P									
12	10	+ + + 0 + + + 0 + P									

(c)

FIGURE 1 : Storage of stiffness coefficients during Frontal solution

STORAGE IN FRONT:



STORAGE IN ELPA:

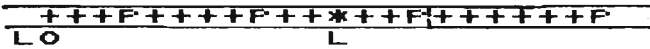


FIGURE 2 : Storage of coefficients in long vector ELPA

APPENDIX A - LIST OF IMPORTANT ARRAYS AND VARIABLES

ELPA main working space

HBWR an indicator if a partition has been written at least once

INDABL indicator on the space availability in the Front matrix also indicates in which partition space is available (coding: "+" occupied "-" ve free)

LCDEST list of coded element destinations

LUDEST list of uncoded element destinations (not entirely true since the destinations are still coded with a "-" ve sign for variables ready for elimination).

LPAL list of partition limits

LISTEQ list of start addresses of equations in the buffer or on disk. Lists address of pivot, block number and relative address in block for each element loop

KURPA current equation length

LBLK length of physical block on disk (real words)

NBLKA, NBLKE first and last block currently in the equation buffer

NPA, NPAC number of partition to be swapped into core and number of partition currently in core

NELZ, LFRBUF, LEQBUF length of buffers for element stiffness matrix, Front matrix and equation buffer (real words)
MUST BE DIVISIBLE BY THE BLOCK LENGTH.

APPENDIX B - LISTING OF THE COMPUTER PROGRAM

```

      SUBROUTINE PREFR
C-----
C   PREFRONT ROUTINE
C-----
      COMMON NIX(10000),MAXPA,NELEMZ,LDEST(62)
C-----
C   VARIABLES IN BLANK COMMON:
C     NIX   ...   WORKING SPACE
C     MAXPA ...   MAX. FRONT WIDTH
C     NELEMZ ...  NUMBER OF ELEMENTS
C     LDEST ...  ELEMENT DESTINATIONS
C-----
      COMMON /FILES/ NF6,NDIM6,NF7,NDIM7,NF8,NDIM8,NF9,NDIM9
      DIMENSION LVABL(60),MVABL(500),LCDEST(60)
      EQUIVALENCE (LPREQ,LDEST(1))
      EQUIVALENCE (KUREL,LDEST(2))
      EQUIVALENCE (LCDEST(1),LDEST(3))
      NIXEND= 2000
      CALL ICLAR(MVABL,500)
      MAXPA=1
      NIZZ= 0
C-----
C   PUT ALL ELEMENT NICKNAMES IN LONG VECTOR NIX
C-----
      DO 10 NELEM= 1,NELEMZ
      CALL GETELN(NELEM,KUREL,LVABL)
      DO 8 I=1,KUREL
      NIC= LVABL(I)
      NIZZ= NIZZ +1
      NIX(NIZZ)= -NIC
      8   CONTINUE
      NIX(NIXEND+1-NELEM)= NIZZ
      10  CONTINUE
C-----
C   PUT SUBSTRUCTURE NICKNAMES(IF ANY) AT THE END OF NIX
C-----
      CALL GETSUN(KUREL,LVABL)
      IF(KUREL .EQ. 0) GOTO 11
      NIZS= NIZZ
      DO 9 I=1,KUREL
      NIC= LVABL(I)
      NIZZ= NIZZ+1
      NIX(NIZZ)= -NIC
      9   CONTINUE
      11  CONTINUE
      KURELS= KUREL
      LCUREQ= 0
      NVABZ= 0
C-----
C   FIND DESTINATIONS
C-----
      N1= 1
      DO 26 NELEM=1,NELEMZ
      LPREQ= LCUREQ
      LCUREQ= NVABZ
      NIXE= NIXEND+1-NELEM
      NZ= NIX(NIXE)
      KUREL= NZ - N1 + 1
      DO 22 NEW= N1,NZ
      NEWA= NEW

```

```

      NIC= NIX(NEW)
      LDES= NIC
      IF(NIC .GT. 0) GOTO 20
      LDES= 1
14    CONTINUE
      IF(MVABL(LDES) .EQ. 0) GOTO 16
      LDES= LDES + 1
      IF(LDES .LE. MAXPA) GOTO 14
      MAXPA= LDES
16    CONTINUE
      MVABL(LDES)= 1
C-----
C   RECORD FIRST, LAST AND INTERM. APPEARANCES
C-----
      KOUNT= 1000
      DO 18 LAS= NEW, NIZZ
      IF(NIX(LAS) .NE. NIC) GOTO 18
      NIX(LAS)= LDES
      KOUNT= KOUNT + 1000
      LAST= LAS
18    CONTINUE
      NIX(LAS)= LDES + 1000
      LDES= LDES + KOUNT
      NIX(NEW)= LDES
20    CONTINUE
      NZ1= NEW-N1+1
      LCDEST(NZ1)= LDES
22    CONTINUE
      N1= NEW+ 1
C-----
C   UPDATE MVABL, COUNT ELIMINATED VARIABLES
C   AND WRITE DESTINATIONS ON DISK
C-----
      DO 24 KL=1, KUREL
      CALL CODEST(KL, NSTRES, LDES, LCDEST, KUREL)
      IF(NSTRES .NE. 0 .AND. NSTRES .NE. 1) GOTO 24
      MVABL(LDES)= 0
      NVABZ= NVABZ+1
24    CONTINUE
      WRITE(NF9'NELEM) (LDEST(I), I=1, NDIM9)
26    CONTINUE
C-----
C   WRITE SUPERELEMENT DESTINATIONS ON DISK
C-----
      NIZZ= NIZS
      KUREL=KURELS
      IF(KUREL .EQ. 0) GOTO 23
      DO 25 KL=1, KUREL
      NIZZ= NIZZ + 1
      LCDEST(KL)= NIX(NIZZ) - 1000
25    CONTINUE
23    CONTINUE
      NELEM= NELEMZ+1
      WRITE(NF9'NELEM) (LDEST(I), I=1, NDIM9)
      RETURN
      END

```

SUBROUTINE SFRONT

```

C-----
C
C   S U P E R F R O N T
C
C   A SECOND GENERATION FRONTAL SOLUTION PROGRAM
C
C   G.BEER   UNIVERSITY OF QUEENSLAND   1979
C-----
      COMMON ELPA(5000),MAXPA,NELEMZ,LDEST(62)
C-----
C   VARIABLES IN BLANK COMMON :
C   ELPA   ...   WORKING SPACE
C   MAXPA  ...   MAXIMUM FRONT WIDTH (FROM PREFRONT)
C   NELEMZ ...   NUMBER OF ACTIVE ELEMENTS
C   LDEST  ...   ELEMENT DESTINATIONS
C-----
      EQUIVALENCE (KUREL,LDEST(2))
      EQUIVALENCE (LCDEST(1),LDEST(3))
      COMMON /EIGEN/ DET,NEG
      COMMON /PARTL/ NST,NEND
      COMMON /PARA/  L0,L1
      COMMON /BLOKL/ LBLK
      COMMON /FILES/ NF6,NDIM6,NF7,NDIM7,NF8,NDIM8,NF9,NDIM9
      COMMON /IOCONV/ IREAD,IWRIT
      COMMON /EQL/  KURPA
      COMMON /ENDQN/ LASTBL
      COMMON /BUFSIZE/ NELZ,LFRBUF,LEQBUF,LFRBBL,LEQBBL
      COMMON /INCORE/ NBLKA,NBLKE
C-----
C   CURRENT COMPILATION IS FOR:
C
C   ELEMENT SIZE= 60 D.O.F.
C   MAXIMUM FRONT WIDTH= 500
C   MAXIMUM NUMBER OF PARTITIONS= 50
C
C   DIMENSION FOR LISTEQ= 50 + 60*3 = 230
C-----
      DIMENSION LUDEST(60)
      DIMENSION LCDEST(60)
      DIMENSION INDABL(500)
      DIMENSION HBWR(50)
      DIMENSION LPAL(51)
      DIMENSION LISTEQ(230)
C   MAX NUMBER OF PARTITIONS THIS COMPILATION:
      MAXPAR= 50
C   SIZE OF ELPA:
      LSIZE= 5000
      IF(NELZ+LFRBUF+LEQBUF .GT. LSIZE) CALL ERROR(0.,LSIZE,3)
      LFRBBL= LFRBUF/LBLK
      LEQBBL= LEQBUF/LBLK
C-----
C   ELPA ADRESSES
C-----
C   START OF FRONT MATRIX
      L0= NELZ
C   START OF EQUATION BUFFER
      L1= NELZ + LFRBUF
C   CLEAR ARRAYS AND WORK OUT PARTITION LIMITS

```



```

CALL CLEAR(ELPA,1,LSIZE)
CALL PALI(MAXPAR,MAXPA,INDABL,LPAL,NOPAR)
PRINT 3002,MAXPA,NDPAR
3002  FORMAT(/// MAXIMUM FRONT WIDTH=',I5/
1      ' MAX. NO. OF PARTITIONS=',I5//)
CALL ICLAR(HBWR,NDPAR)
DET= 0.
NEG= 0
NBLK= 0
NBLKA= 1
NBLKE= LEQBBL
IEQ= L1
NPAC= 1
DO 1 NELEM= 1,NELEMZ
TYPE 3003,NELEM
3003  FORMAT(I5)
C READ CODED ELEMENT DESTINATIONS
READ(NF9'NELEM) (LDEST(I),I=1,NDIM9)
C UNCODE AND UPDATE SPACE INDICATOR
CALL UNCOD(LCDEST,KUREL,LUDEST,INDABL,MAXPA,NDPAR)
C READ ELEMENT STIFFNESS
READ(NF6'1) (ELPA(I),I=1,NDIM6)
NVAR= 0
III= 1
DO 2 NPA=1,NDPAR
C FIRST EQUATION IN PARTITION NPA
NST= LPAL(NPA) + 1
C LAST EQUATION IN PARTITION
NEND= LPAL(NPA + 1)
IF(NEND .GT. KURPA) NEND= KURPA
NVA= 0
LISTEQ(III)= 0
I= III
-----
C ASSEMBLY
-----
DO 4 NV=1,KUREL
IRDY= 0
LDES= LUDEST(NV)
IF(LDES) 21,4,20
21  CONTINUE
LDES= -LDES
IRDY= 1
20  CONTINUE
LPA= INDABL(LDES)
IF(LPA .NE. NPA) GOTO 4
CALL SWAPP(NPA,NPAC,HBWR)
CALL ASSEMB(LDES,LUDEST,NV,KUREL)
IF(IRDY .EQ. 0) GOTO 4
NVA= NVA + 1
NVAR= NVAR + 1
LISTEQ(III)= NVA
I= I + 1
LISTEQ(I)= LDES
I= I+2
4  CONTINUE
-----
C ELIMINATION OF OLD VARIABLES (THOSE ELIMINATED IN PREVIOUS PARTI
C -----TIONS)
IF(NPA .EQ. 1) GOTO 5
IF(NVAR .EQ. 0) GOTO 5

```

```

ICYCL= 1
NEW= 0
I= 0
NPAM= NPA-1
DO 6 NP=1,NPAM
I= I + 1
NVA= LISTEQ(I)
IF(NVA .EQ. 0) GOTO 6
CALL SWAPF(NPA,NPAC,HBWR)
DO 7 N=1,NVA
I= I+1
LDES= LISTEQ(I)
I= I+1
NNBLK= LISTEQ(I)
I= I+1
LIEQ= LISTEQ(I)
IF(LASDES .EQ. LDES) GOTO 77
CALL SAVES(LDES,IEQ,ICYCL,JES)
IF(JES .EQ. 1) GOTO 7
77 CONTINUE
CALL RESBUF(IEQ,NBLK,NNBLK,LIEQ,ICYCL)
CALL MOVE(NEW,LDES,IEQ)
CALL GAUSS(NEW,LDES,IEQ)
7 CONTINUE
6 CONTINUE
5 CONTINUE
C-----
C ELIMINATION OF NEW VARIABLES (THOSE TO BE ELIMINATED IN CURRENT
C----- PARTITION)
NEW= 1
NVA= LISTEQ(III)
I= III
IF(NVA .EQ. 0) GOTO 44
DO 45 N=1,NVA
I= I + 1
LDES= LISTEQ(I)
LASDES= LDES
LEQ= IEQ-L1
IBLK= LEQ/LBLK + 1
NNBLK= NBLK + IBLK
LIEQ= LEQ - (IBLK-1)*LBLK
I= I + 1
LISTEQ(I)= NNBLK
I= I + 1
LISTEQ(I)= LIEQ
CALL EMPDI(IEQ,NBLK)
CALL MOVE(NEW,LDES,IEQ)
CALL GAUSS(NEW,LDES,IEQ)
INDABL(LDES)= -INDABL(LDES)
45 CONTINUE
C ADDRESS OF LAST COEFFICIENT IN EQUATION BUFFER
LEQ= IEQ- L1
IBLK= (LEQ-1)/LBLK + 1
LASTBL= NBLK + IBLK
NDADD= LEQ - (IBLK-1)*LBLK
44 CONTINUE
III= I+1
2 CONTINUE
LASTP= NDPAR-1
C-----
C NOW RETRACK AND MODIFY EQUATIONS IN LOWER

```

C PARTITIONS NOT YET MODIFIED

```

C-----
IF(LASTP .EQ. 0) GOTO 8
ICYCL= 2
NEW= 0
DO 9 NPA=1, LASTP
NST= LPAL(NPA) + 1
NEND= LPAL(NPA + 1)
IF(NEND .GT. KURPA) NEND=KURPA
I= LISTEQ(1)*3 + 1
DO 10 NP=2, NDPAR
I= I + 1
NVA= LISTEQ(I)
IF(NP .GT. NPA) GOTO 99
I= I + NVA*3
GOTO 10
99 CONTINUE
IF(NVA .EQ. 0) GOTO 10
DO 11 N=1, NVA
I= I + 1
LDES= LISTEQ(I)
I= I + 1
NNBLK= LISTEQ(I)
I= I + 1
LIEQ= LISTEQ(I)
CALL RESBUF(IEQ, NBLK, NNBLK, LIEQ, ICYCL)
CALL SAVES(LDES, IEQ, ICYCL, JES)
IF(JES .EQ. 1) GOTO 11
CALL SWAFF(NPA, NPAC, HBWR)
CALL GAUSS(NEW, LDES, IEQ)
11 CONTINUE
10 CONTINUE
9 CONTINUE
8 CONTINUE
1 CONTINUE

```

C-----
C WRITE CONTENTS OF EQUATION BUFFER IF NECESSARY
C-----

```

IF(NBLK .GE. LASTBL) RETURN
NBLOKS= LASTBL-NBLK
CALL BLKIO(IWRIT, NF7, NBLK+1, NBLOKS, ELPA, L1+1)

```

C-----
C CONDENSE AND REORDER SUBSTRUCTURE STIFFNESS MATRIX IF REQUIRED
C-----

```

NELEM= NELEMZ + 1
READ(NF9'NELEM) (LDEST(I), I=1, NDIM9)
IF(KUREL .EQ. 0) RETURN
DO 200 NPA= 1, NOPAR
CALL SWAFF(NPA, NPAC, HBWR)
NPAC= 0
NST= LPAL(NPA) + 1
NEND= LPAL(NPA+1)
IF(NEND .GT. KURPA) NEND= KURPA
L=0
DO 201 I=1, KUREL
IDES= LCDEST(I)
DO 202 K=1, I
L= L+1
KDES= LCDEST(K)
LDES= MAXO(IDES, KDES)
MDES= MINO(IDES, KDES)

```

```

LPA= INDABL(LDES)
IF(LPA .NE. NPA) GOTO 202
LL= LADDR(LDES-1) + MDES
ELPA(L)= ELPA(LL)
202 CONTINUE
201 CONTINUE
200 CONTINUE
C PRINT SUBSTRUCTURE STIFFNESS
PRINT 3001,(K,K=1,KUREL)
3001 FORMAT(/// CONDENSED SUBSTRUCTURE STIFFNESS MATRIX: ///4X,20I6)
LA=1
LE=1
DO 203 K=1,KUREL
PRINT 3000,K,(ELPA(L),L=LA,LE)
LA= LE+1
LE= LA+K
203 CONTINUE
3000 FORMAT(1X,I3,20F6.2)
PRINT 3004,DET,NEG
3004 FORMAT(/// LOG10 OF DETERMINANT=' ,F15.4/
1 ' NO OF NEGATIVE PIVOTS=' ,I5//)
RETURN
END
SUBROUTINE MOVE(NEW,LDES,IEQ)

```

```

C-----
C
C TO MOVE EQUATION COEFFICIENTS FROM FRONT PARTITION TO EQUATION
C BUFFER
C NEW=1 ... NEW EQUATION
C NEW=0 ... OLD EQUATION
C
C LDES ... DESTINATION OF VARIABLE TO BE ELIMINATED
C
C IEQ ... CURRENT ADDRESS OF BUFFER PONTER
C-----

```

```

COMMON ELPA(1)
COMMON /PARTL/ NST,NEND
COMMON /EQL/ KURPA
COMMON /PARA/ L0,L1
IF(NEW .EQ. 0) GOTO 1
C --- NEW EQUATION
L= LADDR(LDES-1)
M= IEQ
DO 2 J=1,LDES
L= L+1
M=M+1
ELPA(M)= ELPA(L)
ELPA(L)= 0.
2 CONTINUE
IF(LDES .EQ. NEND) GOTO 7
N1= LDES + 1
K= 0
DO 3 J=N1,NEND
K=K + 1
L= L + LDES
M= M + 1
ELPA(M)= ELPA(L)
ELPA(L)= 0.
L= L + K
3 CONTINUE

```

```

7  CONTINUE
   NDEQN= IEQ + KURPA
   IF(M .EQ. NDEQN) GOTO 5
   M1= M+1
   DO 6 I=M1,NDEQN
   ELPA(I)= 0.
6  CONTINUE
   M= NDEQN
5  CONTINUE
   ELPA(M + 1) = LDES
   ELPA(M + 2)= KURPA
   RETURN
C  --- OLD EQUATION
1  CONTINUE
   L= LDES + L0
   M= IEQ + NST - 1
   NREST= NST
   DO 4 J=NST,NEND
   M= M+1
   ELPA(M)= ELPA(L)
   ELPA(L)= 0.
   L= L + NREST
   NREST= NREST + 1
4  CONTINUE
   RETURN
   END
   SUBROUTINE GAUSS(NEW,LDES,IEQ)
C-----
C
C   MODIFIES ALL EQUATIONS OF PARTITION NPA
C   (ELIMINATION OF VARIABLE LDES)
C
C   IEQ   ...   ADDRESS OF EQUATION BUFFER POINTER
C-----
COMMON ELPA(1)
COMMON /EIGEN/ DET,NEG
COMMON /EQL/ KURPA
COMMON /PARA/ L0,L1
COMMON /PARTL/ NST,NEND
NDIAG= IEQ + LDES
PIVOT= ELPA(NDIAG)
ELPA(NDIAG)= 0.
C  CHECK FOR SINGULARITY AND WORK OUT DETERMINANT
IF(NEW .EQ. 0) GOTO 2
PIVO= ABS(PIVOT)
DET= DET + ALOG10(PIVO)
IF(PIVO .LT. 1.E-20) CALL ERROR(PIVOT,LDES,2)
IF(PIVOT .GT. 0.) GOTO 2
NEG= NEG + 1
CALL ERROR(PIVOT,LDES,1)
2  CONTINUE
   L= L0
   MI= IEQ + NST- 1
   DO 1 I=NST,NEND
   MI= MI + 1
   CONS= ELPA(MI )
   IF(CONS .EQ. 0.) GOTO 3
   CONS= CONS/PIVOT
   M= IEQ
C  CALL VECTOR PROCESSOR

```

```

      CALL SVECT(ELPA,CONS,L,M,I)
      GOTO 1
3     CONTINUE
      L= L + I
1     CONTINUE
      ELPA(NDIAG)= PIVOT
C     MOVE BUFFER POINTER TO END OF EQUATION
      IEQ= IEQ + KURPA + 2
      RETURN
      END
      SUBROUTINE UNCOD(LCDEST,KUREL,LUDEST,INDABL,MAXPA,NDPAR)

```

```

C-----
C   UNCODES DESTINATION VECTOR LCDEST AND
C   UPDATES SPACE INDICATOR INDABL
C-----

```

```

      DIMENSION LCDEST(KUREL)
      DIMENSION LUDEST(KUREL)
      COMMON /EQL/ KURPA
      DIMENSION INDABL(MAXPA)
      DO 1 K=1,KUREL
      IRDY= 0
      CALL CODEST(K,NSTRES,LDES,LCDEST,KUREL)
      IF(LDES .EQ. 0) GOTO 3
      IF(NSTRES .NE. 0 .AND. NSTRES.NE. 1) GOTO 2
C   --- VARIABLE LDES CAN BE ELIMINATED
      IRDY= 1
2     CONTINUE
      NPA= IABS(INDABL(LDES))
      INDABL(LDES)= NPA
3     CONTINUE
      IF(IRDY .EQ. 1) LDES= -LDES
      LUDEST(K)= LDES
1     CONTINUE
CURRENT EQUATION LENGTH
      M= MAXPA
5     CONTINUE
      IF(INDABL(M) .GT. 0) GOTO 4
      M= M-1
      GOTO 5
4     CONTINUE
      KURPA= M
      NDPAR= INDABL(KURPA)
      RETURN
      END
      SUBROUTINE PALI(MAXPAR,MAXPA,INDABL,LPAL,NDPAR)

```

```

C-----
C   SETS UP PARTITION LIMIT ARRAY LPAL
C   DEPENDING ON SIZE OF FRONT BUFFER LFRBUF
C   AND MAXPA ( MAXIMUM FRONT WIDTH )
C   TOTAL NUMBER OF PARTITIONS REQUIRED: NDPAR
C-----

```

```

      DIMENSION INDABL(MAXPA),LPAL(MAXPAR)
      COMMON /BUFSIZE/ NELZ,LFRBUF,LEQBUF
      LPAL(1)= 0
      DO 1 NPA=1,MAXPAR
      FAC= 2* LPAL(NPA) + 1
      FAC1= 2*LFRBUF
      LDSPD= SQRT(.25*FAC*FAC + FAC1) - .5*FAC
      LPAL(NPA + 1)= LPAL(NPA) + LDSPD
      IF(LPAL(NPA + 1) .GE. MAXPA) GOTO 2
1     CONTINUE

```

```

      CALL ERROR(0.,NPA,4)
      2  CONTINUE
        LPAL(NPA+1)= MAXPA
        NDPAR= NPA
C-----
C   SET UP ARRAY INDABL ; CODING: '+'-VE=OCCUPIED; '-'-VE=FREE
C-----
      DO 3 M=1,MAXPA
      DO 4 NPA=1,NDPAR
      IF(M .LE. LPAL(NPA+1)) GOTO 5
      4  CONTINUE
      5  CONTINUE
        INDABL(M)= -NPA
      3  CONTINUE
        RETURN
      END
      SUBROUTINE ASSEMB(LDES,LUDEST,NV,KUREL)
C-----
C   ASSEMBLES EQUATION LDES INTO CURRENT FRONT PARTITION
C-----
      COMMON ELPA(1)
      DIMENSION LUDEST(KUREL)
      LL= LADDR(LDES-1)
      DO 1 K=1,KUREL
      II= LUDEST(K)
      IF(II) 2,1,3
      2  CONTINUE
      II= -II
      3  CONTINUE
      IF(II .GT. LDES) GOTO 1
      L= LADST(K,NV)
      LF=LL + II
      ELPA(LF)= ELPA(LF) + ELPA(L)
      1  CONTINUE
      RETURN
      END
      FUNCTION LADDR(M)
C-----
C   COMPUTES THE ADDRESS OF COEFF M,M IN CURRENT PARTITION
C-----
      COMMON /PARA/ L0,L1
      COMMON /PARTL/ NST,NEND
      NS= NST -1
      MR= M-NS
      LADDR= MR*NS + MR*(MR+1)/2+ L0
      RETURN
      END
      FUNCTION LADST(I,J)
C-----
C   COMPUTES ADDRESS OF COEFF I,J IN ELEMENT STIFFNESS MATRIX
C-----
      II= MAX0(I,J)
      JJ= MIN0(I,J)
      LADST= JJ + II*(II-1)/2.
      RETURN
      END
      SUBROUTINE EMPDI(IEQ,NBLK)
C-----
C   EMPTIES EQUATION BUFFER ONTO DISK WHEN FULL
C-----

```

```

C   IEQ   ...   BUFFER POINTER
C   NBLK  ...   NUMBER OF BLOCKS WRITTEN
C -----
      COMMON ELPA(1)
      COMMON /PARA/ L0,L1
      COMMON /EQL/ KURPA
      COMMON /INCORE/ NBLKA,NBLKE
      COMMON /BLOKL/ LBLK
      COMMON /FILES/ NF6,NDIM6,NF7,NDIM7,NF8,NDIM8,NF9,NDIM9
      COMMON /IOCONV/ IREAD,IWRIT
      COMMON /BUFSIZE/ NELZ,LFRBUF,LEQBUF,LFRBBL,LEQBBL
CHECK IF ANOTHER EQUATION FITS
      NDEQN= IEQ + KURPA + 2 - L1
      IF(NDEQN .LT. LEQBUF) RETURN
C   ---   DOES NOT FIT >> WRITE BUFFER ONTO DISK
      NBLOKS= (IEQ-L1)/LBLK
      CALL BLKIO(IWRIT,NF7,NBLK + 1,NBLOKS,ELPA,L1+1)
C   MOVE LAST BLOCK AT THE BEGINNING OF BUFFER IF NOT COMPLETELY FULL
      LEQE= IEQ
      LEQ= NBLOKS *LBLK + L1
      IEQ= L1
      IF(LEQ .EQ. LEQE) GOTO 10
      LEQ= LEQ + 1
      DO 1 I=LEQ,LEQE
      IEQ= IEQ + 1
      ELPA(IEQ)= ELPA(I)
1   CONTINUE
10  CONTINUE
      NBLK= NBLK + NBLOKS
      NBLKA= NBLK + 1
      NBLKE= NBLK + LEQBBL
      RETURN
      END
      SUBROUTINE RESBUF(IEQ,NBLK,NNBLK,LIEQ,ICYCL)
C -----
C   RESETS BUFFER POINTER IEQ TO LIEQ IN BLOCK NNBLK
C   AND SWAPS(ICYCL=1) OR READS(ICYCL=2) BLOCKS IF NECESSARY
C -----
      COMMON ELPA(1)
      COMMON /PARA/ L0,L1
      COMMON /BLOKL/ LBLK
      COMMON /EQL/ KURPA
      COMMON /INCORE/ NBLKA,NBLKE
      COMMON /FILES/ NF6,NDIM6,NF7,NDIM7,NF8,NDIM8,NF9,NDIM9
      COMMON /IOCONV/ IREAD,IWRIT
      COMMON /BUFSIZE/ NELZ,LFRBUF,LEQBUF,LFRBBL,LEQBBL
      COMMON /ENDQN / LASTBL
C -----
C   IS EQUATION STILL IN CORE ?
C -----
      IF(NNBLK .LT. NBLKA) GOTO 2
      LEQ= LIEQ + KURPA + 2
      LEQB= (LEQ-1)/LBLK
      NDBLK= NNBLK + LEQB
      IF(NDBLK .LE. NBLKE) GOTO 1
C -----
C   NO   -   SWAP BLOCKS
C -----
2   CONTINUE
      IF(ICYCL .EQ. 2) GOTO 4

```



```

NBLKEN= NBLKE
IF(NBLKEN .LT. LASTBL) GOTO 3
NBLKEN= LASTBL
3 CONTINUE
NBLKS= NBLKEN-NBLKA + 1
CALL BLKIO(IWRIT,NF7,NBLKA,NBLOKS,ELPA,L1+1)
4 CONTINUE
NBLKS= LEQBBL
LIMBLK= LASTBL - NNBLK + 1
IF(NBLOKS .GT. LIMBLK) NBLOKS= LIMBLK
CALL BLKIO(IREAD,NF7,NNBLK,NBLOKS,ELPA,L1+1)
NBLKA= NNBLK
NBLKE= NNBLK + NBLOKS - 1
NBLK= NBLKA - 1
IEQ= LIEQ + L1
RETURN
1 CONTINUE
IEQ= (NNBLK-NBLKA)*LBLK + LIEQ + L1
RETURN
END
SUBROUTINE SWAPF(NPA,NPAC,HBWR)
-----
C SWAPS FRONT PARTITIONS IN AND OUT OF CORE AS REQUIRED
C
C NPA ... NEW PARTITION
C NPAC ... CURRENT PARTITION
C -----
COMMON ELPA(1)
DIMENSION HBWR(1)
COMMON /PARA/ L0,L1,L2
COMMON /FILES/ NF6,NDIM6,NF7,NDIM7,NF8,NDIM8,NF9,NDIM9
COMMON /IOCONV/ IREAD,IWRIT
COMMON /BLOKL/ LBLK
IF(NPA .EQ. NPAC) RETURN
NBLKS8= NDIM8/LBLK
IF(NPAC .EQ. 0) GOTO 1
NFROM= (NPAC-1)*NBLKS8 + 1
CALL BLKIO(IWRIT,NF8,NFROM,NBLKS8,ELPA,L0+1)
HBWR(NPAC)= 1
1 CONTINUE
IF(NPA .EQ. 0) RETURN
NFROM= (NPA-1)*NBLKS8 + 1
IF(HBWR(NPA) .EQ. 1) CALL BLKIO(IREAD,NF8,NFROM,NBLKS8,ELPA,
IF(HBWR(NPA) .EQ. 0) CALL CLEAR(ELPA,L0+1,L1)
NPAC= NPA
RETURN
END
SUBROUTINE CODEST(K,NSTRES,LDES,LCDEST,KUREL)
-----
C INTERPRETS CODED ELEMENT DESTINATIONS
C -----
DIMENSION LCDEST(KUREL)
LDES= LCDEST(K)
DO 2 NSTRES= 1,32000
IF(LDES .LT. 1000) GOTO 4
LDES= LDES - 1000
2 CONTINUE
4 CONTINUE
NSTRES= NSTRES - 2
RETURN
END

```

```

SUBROUTINE CLEAR(ARRAY,NST,NEN)
DIMENSION ARRAY(NEN)
DO 1 N=NST,NEN
ARRAY(N)= 0.
1 CONTINUE
RETURN
END
SUBROUTINE ICLAR(IARR,NEN)
DIMENSION IARR(NEN)
DO 1 N=1,NEN
IARR(N)= 0
1 CONTINUE
RETURN
END
SUBROUTINE ERROR(F,I,N)
GOTO (1,2,3) ,N
1 CONTINUE
PRINT 2000,F,I
2000 FORMAT(// ' *** NEGATIVE PIVOT (' ,E15.5,') AT DESTINATION',I5)
RETURN
2 CONTINUE
PRINT 2001,F,I
2001 FORMAT(/// ' SINGULARITY CHECK:/'
1 ' NEAR ZERO OR ZERO PIVOT (' ,E15.5,') AT DESTINATION',I5)
STOP
3 CONTINUE
PRINT 2002,I
2002 FORMAT(/// ' *** DIMENSION OF ELPA (' ,I5,') TOO SMALL')
STOP
4 CONTINUE
PRINT 2003,I
2003 FORMAT(/// ' *** MAXIMUM NUMBER OF PARTITIONS (' ,I5,') EXCEEDED')
STOP
RETURN
END
SUBROUTINE SAVES(LDES,IEQ,ICYCL,JES)

```

```

C-----
C TO SAVE ON COMPUTATION AND SWAPPING TIME
C FOR ZERO COEFFICIENTS
C-----

```

```

COMMON ELPA(1)
COMMON /PARTL/ NST,NEND
COMMON /PARA/ L0,L1,L2
IF(ICYCL .EQ. 2) GOTO 1
L= LDES + L0
NREST= NST
DO 2 J=NST,NEND
IF(ELPA(L) .NE. 0.) GOTO 3
L= L + NREST
NREST=NREST + 1
2 CONTINUE
JES=1
RETURN
3 CONTINUE
JES= 0
RETURN
1 CONTINUE
MI= IEQ + NST-1
DO 4 I= NST,NEND
MI= MI+1
IF(ELPA(MI) .NE. 0.) GOTO 5

```

```

4  CONTINUE
   JES= 1
   RETURN
5  CONTINUE
   JES=0
   RETURN
   END
   SUBROUTINE SVECT(VECTOR,CONS,I1,I2,N)
C-----
C  FAST VECTOR PROCESSING ROUTINE TO PERFORM :
C
C  VECTOR(I1)= VECTOR(I1) - VECTOR(I2)*CONS
C
C      VECTOR    ...    VECTOR
C      CONS      ...    SCALAR
C      I1    ...    START ADDRESS 1
C      I2    ...    START ADDRESS 2
C      N     ...    NUMBER OF OPERATIONS
C
C  MACHINE DEPENDENT CODING SHOULD BE USED IN ACTUAL IMPLEMENTATION
C  ON A GIVEN MACHINE.
C  CODING SHOWN IS STANDARD FORTRAN
C-----
      DIMENSION VECTOR(1)
      DO 2 J=1,N
      I1= I1+1
      I2= I2+1
      VECTOR(I1)= VECTOR(I1) - VECTOR(I2)*CONS
2  CONTINUE
   RETURN
   END

```

SUBROUTINE BLKIO(IRW,LUN,NBLA,NBLOKS,BUFFER,IADD)

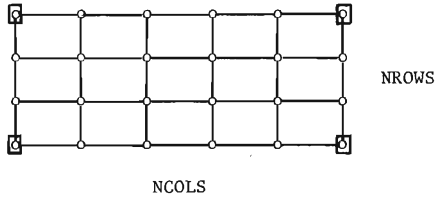
```

C-----
C  SUBROUTINE TO READ/WRITE DIRECTLY INTO BUFFER
C
C  IRW  ...  SWITCH FOR READ/WRITE
C           IRW=1  ...  READ
C           IRW=2  ...  WRITE
C  LUN  ...  LOGICAL UNIT NUMBER
C  NBLA ...  FIRST BLOCK
C  NBLOKS ...  NUMBER OF BLOCKS TO BE READ/WITTEN
C  BUFFER ...  BUFFER
C  IADD ...  START ADDRESS IN BUFFER
C
C-----
C           COMMON /IO/ IPAR(6),ISTAT(2)
C           COMMON /BLOKL/ LBLK
C           DIMENSION BUFFER(1)
C  INSERT MACHINE DEPENDENT CODING HERE
C           RETURN
C           END
C           SUBROUTINE FILO
C-----
C  SUBROUTINE TO ASSIGN CHANNEL NUMBERS
C  AND OPEN DIRECT ACCESS AND BLOCKED FILES
C
C  CHANNEL#  SIZE  CONTENTS
C  NF6       NELZ  ELEMENT STIFFNESS
C  NF9       D.O.F. + 2  ELEMENT DESTINATIONS
C  NF7       LBLK  BLOCKED EQUATIONS
C  NF8       LBLK  BLOCKED PARTITIONS
C
C-----
C           COMMON /FILES/ NF6,NDIM6,NF7,NDIM7,NF8,NDIM8,NF9,NDIM9
C           COMMON /BUFSZE/ NELZ,LFRBUF,LEQBUF
C           COMMON /IOCONV/ IREAD,IWRIT
C           COMMON /BLOKL/ LBLK
C           IREAD= 1
C           IWRIT= 2
C  INSERT MACHINE DEPENDENT CODING HERE
C           RETURN
C           END

```

APPENDIX C - TEST PROGRAM

In the following a test program is listed which can be used to test the subroutines PREFR and SFRONT. The example is a substructure condensation for a regular patch of square 4 node/8 d.o.F. Elements. The stiffness matrix of the Elements is read in and all the nodes except the 4 corner nodes of the super element are condensed out. Results can be obtained for different mesh and buffer sizes



□ substructure nodes

o element nodes

```

C-----
C  PROGRAM TO TEST OUT SUBROUTINE SFRONT
C-----
COMMON ELPA(5000),MAXPA,NELEMZ,LDEST(62)
COMMON /BUFSIZE/ NELZ,LFRBUF,LEQBUF,LFRBBL,LEQBBL
COMMON /MESH/ NROWS,NCOLS,NODES
COMMON /BLOKL/ LBLK
COMMON /FILES/ NF6,NDIM6,NF7,NDIM7,NF8,NDIM8,NF9,NDIM9
NCR= 7
CALL ASSIGN(NCR,'INPUT')
NELZ= 36
READ(NCR,1000) NROWS,NCOLS
1000  FORMAT(16I5)
NELEMZ= NROWS*NCOLS
NODES= (NCOLS+1)*(NROWS+1)
NDOFS= NODES*2
READ(NCR,1000) LFRBUF,LEQBUF,LBLK
CALL FILO
READ(NCR,1000) ISTIF
IF(ISTIF.NE. 1) GOTO 1
READ(NCR,1001) (ELPA(N),N=1,NELZ)
1001  FORMAT(8F10.0)
WRITE (NF6'1) (ELPA(I),I=1,NELZ)
GOTO 2
1  CONTINUE
READ(NF6'1) (ELPA(I),I=1,NELZ)
2  CONTINUE
PRINT 3000
3000  FORMAT(1H1//' *** SUBSTRUCTURE CONDENSATION EXAMPLE ***')
PRINT 3001,NROWS,NCOLS
3001  FORMAT(//' NUMBER OF ELEMENT ROWS=',I5/
1      ' NUMBER OF ELEMENT COLUMNS=',I5)
PRINT 3002,NELEMZ,NODES,NDOFS
3002  FORMAT(' NUMBER OF ELEMENTS=',I5/
1      ' NUMBER OF NODES   =',I5/
1      ' NUMBER OF D.O.F.  =',I5)
PRINT 3003,LFRBUF,LEQBUF
3003  FORMAT(' SIZE OF FRONT BUFFER   =',I5/
1      ' SIZE OF EQUATION BUFFER=',I5)
PRINT 3004,(K,K=1,8)
3004  FORMAT(//' ELEMENT STIFFNESS MATRIX: '//4X,20I6)
LA= 1
LE= 1
DO 200 K=1,8
PRINT 3005,K,(ELPA(L),L=LA,LE)
LA=LE+1
LE= LA+K
200  CONTINUE
3005  FORMAT(1X,I3,20F6.2)
CALL PREFR
CALL SFRONT
STOP
END

```

SUBROUTINE GETELN(NELEM,KUREL,LVABL)

```

C-----
C THIS IS A DUMMY SUBROUTINE FOR TESTING SFRONT
C IT CREATES CONNECTIVITY DATA FOR A REGULAR
C ASSEMBLY OF 4 NODE/8 D.O.F. ELEMENTS
C
C NROWS ... NUMBER OF ELEMENT ROWS
C NCOLS ... NUMBER OF ELEMENT COLUMNS
C NODES ... NUMBER OF NODES
C-----
      DIMENSION LVABL(8)
      COMMON /MESH/ NROWS,NCOLS,NODES
      KUREL= 8
      NROW= (NELEM-1)/NCOLS + 1
      NCOL= NELEM - NCOLS*(NROW-1)
      NCOL1= NCOLS + 1
      LVABL(1)= NCOL + (NROW-1)*NCOL1
      LVABL(2)= NCOL + 1 + (NROW-1)*NCOL1
      LVABL(3)= NCOL + 1 + NROW*NCOL1
      LVABL(4)= NCOL + NROW*NCOL1
      DO 1 N=5,8
      LVABL(N)= LVABL(N-4) + NODES
1 CONTINUE
      RETURN
      END
      SUBROUTINE GETSUN(KUREL,LVABL)

```

```

C-----
C THIS IS A DUMMY SUBROUTINE TO TEST THE SUBSTRUCTURING
C CAPABILITIES OF SFRONT,
C IT CREATES CONNECTIVITY DATA FOR A
C 4 NODE/8 D.O.F. SUPERELEMENT
C-----

```

```

      DIMENSION LVABL(8)
      COMMON /MESH/ NROWS,NCOLS,NODES
      KUREL= 8
      NCOL1= NCOLS + 1
      LVABL(1)= 1
      LVABL(2)= NCOL1
      LVABL(3)= NODES
      LVABL(4)= NCOL1*NROWS + 1
      DO 1 N=5,8
      LVABL(N)= LVABL(N-4) + NODES
1 CONTINUE
      RETURN
      END

```

*** SUBSTRUCTURE CONDENSATION EXAMPLE ***

NUMBER OF ELEMENT ROWS= 2
 NUMBER OF ELEMENT COLUMNS= 10
 NUMBER OF ELEMENTS= 20
 NUMBER OF NODES = 33
 NUMBER OF D.O.F. = 66
 SIZE OF FRONT BUFFER = 512
 SIZE OF EQUATION BUFFER= 256

ELEMENT STIFFNESS MATRIX:

	1	2	3	4	5	6	7	8
1	5.00							
2	-2.50	5.00						
3	-2.50	0.00	5.00					
4	0.00	-2.50	-2.50	5.00				
5	1.25	1.25	-1.25	-1.25	5.00			
6	-1.25	-1.25	1.25	1.25	0.00	5.00		
7	-1.25	-1.25	1.25	1.25	-2.50	-2.50	5.00	
8	1.25	1.25	-1.25	-1.25	-2.50	-2.50	0.00	5.00

MAXIMUM FRONT WIDTH= 30
 MAX. NO. OF PARTITIONS= 1

CONDENSED SUBSTRUCTURE STIFFNESS MATRIX:

	1	2	3	4	5	6	7	8
1	1.12							
2	-0.08	1.12						
3	-0.74	-0.29	1.12					
4	-0.29	-0.74	-0.08	1.12				
5	0.21	-0.01	-0.21	0.01	2.05			
6	0.01	-0.21	-0.01	0.21	0.01	2.05		
7	-0.21	0.01	0.21	-0.01	-0.05	-2.01	2.05	
8	-0.01	0.21	0.01	-0.21	-2.01	-0.05	0.01	2.05

LOG10 OF DETERMINANT= 55.9191
 NO OF NEGATIVE PIVOTS= 0

*** SUBSTRUCTURE CONDENSATION EXAMPLE ***

NUMBER OF ELEMENT ROWS= 2
 NUMBER OF ELEMENT COLUMNS= 10
 NUMBER OF ELEMENTS= 20
 NUMBER OF NODES = 33
 NUMBER OF D.O.F. = 66
 SIZE OF FRONT BUFFER = 256
 SIZE OF EQUATION BUFFER= 256

ELEMENT STIFFNESS MATRIX:

	1	2	3	4	5	6	7	8
1	5.00							
2	-2.50	5.00						
3	-2.50	0.00	5.00					
4	0.00	-2.50	-2.50	5.00				
5	1.25	1.25	-1.25	-1.25	5.00			
6	-1.25	-1.25	1.25	1.25	0.00	5.00		
7	-1.25	-1.25	1.25	1.25	-2.50	-2.50	5.00	
8	1.25	1.25	-1.25	-1.25	-2.50	-2.50	0.00	5.00

MAXIMUM FRONT WIDTH= 30
 MAX. NO. OF PARTITIONS= 2

CONDENSED SUBSTRUCTURE STIFFNESS MATRIX:

	1	2	3	4	5	6	7	8
1	1.12							
2	-0.08	1.12						
3	-0.74	-0.29	1.12					
4	-0.29	-0.74	-0.08	1.12				
5	0.21	-0.01	-0.21	0.01	2.05			
6	0.01	-0.21	-0.01	0.21	0.01	2.05		
7	-0.21	0.01	0.21	-0.01	-0.05	-2.01	2.05	
8	-0.01	0.21	0.01	-0.21	-2.01	-0.05	0.01	2.05

LOG10 OF DETERMINANT= 55.9191
 NO OF NEGATIVE PIVOTS= 0

APPENDIX D - NOMENCLATURE

<u>Symbol</u>	<u>Meaning</u>
c_{ij}	Equation coefficient
l	Address of coefficient in long vector ELPA
l_0	Start address in ELPA of space reserved for the Front matrix
m	Front width
Δl_N	Number of Equations in partition N
k_N	"destination" of first equation in partition N
n_p	Size of storage space for the Front matrix
k, i	Coefficient indexes
F_i	Coefficient on right hand side
x_n	Variable
R_i	Residual
K_{ij}	Structure stiffness coefficient
Δx_j	error on Solution for x_j

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