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THE SOLUTION OF A CERTAIN LINEAR EQUATION SYSTEM

1. Procedure declaration. Procedure *URP* solves a group of lp systems of linear equations with lr unknowns and with various right-hand members

$$(1) \quad KX = T_r \quad (r = 1, 2, \dots, lp),$$

where the global matrix K is treated as composed of submatrices K_1, K_2, \dots, K_{lm} of dimension $n \times n$ (n — an even number) in form as shown in Fig. 1. The marked elements are summarized.

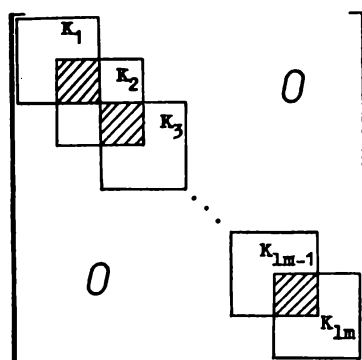


Fig. 1

The number of unknowns is $lr = .5 \times n \times (lm + 1)$. It is easy to show that the indices of the elements k_{ij} (shown as 0 in Fig. 1) in the lower part of the matrix K satisfy the relation

$$1 \leq j \leq \frac{n}{2} \left(\left[\frac{i-1}{n/2} \right] - 1 \right) \quad (i = n+1, n+2, \dots, lr),$$

while the indices of the zero elements in the upper part of the matrix K satisfy the relation

$$n + \left[\frac{i-1}{n/2} \right] \frac{n}{2} < j \leq lr \quad (i = 1, 2, \dots, lr - n),$$

where the symbol $[\cdot]$ denotes the integer part. It is a special type of strip matrix. Such systems are found in many technical problems, usually with a great number of unknowns (for example, in shell analysis). The presented algorithm is more effective (mainly due to its storage-saving properties) than those treating the matrix K as of common strip-type.

Data:

- n — degree of submatrices K_i (n must be even);
- lm — number of submatrices K_i ;
- lp — number of systems processed;
- $T[1: .5 \times n \times (lm + 1), 1: lp]$ — array of right-hand side of system (1) ($T[i, j]$ is the i -th free member of the j -th system of equations);
- CMK — procedure with the following heading: **procedure** $CMK(i, n, K)$; **integer** i, n ; **array** K ; the procedure determines the elements of a submatrix K_i for $i = 1, 2, \dots, lm$.

Results:

- $T[1: .5 \times n \times (lm + 1), 1: lp]$ — array of results for system (1) ($T[i, j]$ is the i -th unknown for the j -th system).

Other parameters:

- $exit$ — label to which the program is switched in case of singularity of the matrix K .

2. Description of the method. We assume that the element k_{rs}^i is placed at the intersection of the row r and the column s of the matrix K_i ($p = n/2$). We analyze a part of the matrix K composed of submatrices K_i and K_{i+1} ($i = 1, 2, \dots, lm - 1$):

$$\begin{array}{cccccccc}
 k_{11}^i & k_{12}^i & \dots & k_{1p}^i & k_{1,p+1}^i & \dots & k_{1n}^i & \\
 k_{21}^i & k_{22}^i & \dots & k_{2p}^i & k_{2,p+1}^i & \dots & k_{2n}^i & \\
 \dots & \dots & \dots & \dots & \dots & \dots & \dots & 0 \\
 k_{p1}^i & k_{p2}^i & \dots & k_{pp}^i & k_{p,p+1}^i & \dots & k_{pn}^i & \\
 k_{p+1,1}^i & k_{p+1,2}^i & \dots & k_{p+1,p}^i & \bar{d}_{11}^i & \dots & \bar{d}_{1p}^i & k_{1,p+1}^{i+1} & \dots & k_{1n}^{i+1} \\
 k_{p+2,1}^i & k_{p+2,2}^i & \dots & k_{p+2,p}^i & \bar{d}_{21}^i & \dots & \bar{d}_{2p}^i & k_{2,p+1}^{i+1} & \dots & k_{2n}^{i+1} \\
 \dots & \dots & \dots & \dots & \dots & \dots & \dots & \dots & \dots & \dots \\
 k_{n1}^i & k_{n2}^i & \dots & k_{np}^i & \bar{d}_{p1}^i & \dots & \bar{d}_{pp}^i & k_{p,p+1}^{i+1} & \dots & k_{pn}^{i+1} \\
 & & & & k_{p+1,1}^{i+1} & \dots & k_{p+1,p}^{i+1} & k_{p+1,p+1}^{i+1} & \dots & k_{p+1,n}^{i+1} \\
 & & & & 0 & \dots & \dots & \dots & \dots & \dots \\
 & & & & & & k_{n1}^{i+1} & \dots & k_{np}^{i+1} & k_{n,p+1}^{i+1} & \dots & k_{nn}^{i+1}
 \end{array}$$

where $\bar{d}_{iq}^i = k_{iq}^{i+1} + k_{i+p,q+p}^i$ ($t, q = 1, 2, \dots, p$).

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procedure URP(n,lm,lp,T,CMK,exit);
  value n,lm,lp;
  integer n,lm,lp;
  array T;
  procedure CMK;
  label exit;
  begin
    integer lr,lm1,np,t1,k2,r1,i,j,h,i1,j1,m,pi,pj,l;
    real g,w;
    np:=n+2;
    lr:=np*(lm+1);
    l:=n+lp;
    begin
      array A[1:lr,1:np],B[1:n,1:l],K[1:n,1:n];
      integer array P[1:l];
      t1:=k2:=np;
      r1:=0;
      CMK(1,n,K);
      for i:=1 step 1 until np do
        begin
          i1:=np+i;
          for j:=1 step 1 until np do
            B[i1,np+j]:=K[i,j];
          for j:=1 step 1 until lp do
            B[i1,n+j]:=T[i,j]
          end i;
        for i:=1 step 1 until l do
          P[i]:=i;
          lm1:=lm+1;
        for h:=2 step 1 until lm1 do

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begin
  for i=1 step 1 until np do
    begin
      i1=np+1;
      for j=1 step 1 until np do
        begin
          j1=np+j;
          B[i,j]=B[i1,j1];
          B[i,j1]=K[i,j1];
          B[i1,j]=K[i1,j];
          B[i1,j1]=K[i1,j1];
        end j;
      for j=1 step 1 until lp do
        begin
          j1=n+j;
          B[i,j1]=B[i1,j1];
          B[i1,j1]=T[t1+i,j];
        end j
      end i;
      t1=t1+np;
      if h=lm+1
        then
          begin
            k2=n;
            go to E1
          end;
      CMK(h,n,K);
    for i=1 step 1 until np do
      begin
        i1=np+i;

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    for j=1 step 1 until np do
      B[i1,np+j]=B[i1,np+j]+K[i,j]
    end i;
E1: for i=1 step 1 until k2 do
  begin
    g=.0;
    for j=i step 1 until k2 do
      begin
        w=abs(B[i,P[j]]);
        if w>g
          then
            begin
              g=w;
              m=j
            end w>g
          end j;
        if g=0
          then go to exit;
        pi=P[m];
        P[m]=P[i];
        P[i]=pi;
        g=1.0/B[i,pi];
        for j=i+1 step 1 until l do
          begin
            pj=P[j];
            w=B[i,pj]=B[i,pj]*g;
            for m=1 step 1 until i-1,i+1 step 1 until n do
              B[m,pj]=B[m,pj]-B[m,pi]*w
            end j
          end i;

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for i=1 step 1 until k2 do
  begin
    i1:=P[i]+r1;
    for j=np+1 step 1 until n do
      A[i1,j-np]:=B[i,j];
    for j=1 step 1 until lp do
      T[i1,j]:=B[i,n+j]
    end i;
    r1:=r1+np
  end h;
k2:=lr-n;
r1=0;
for i=k2 step -1 until 1 do
  begin
    for h=1 step 1 until lp do
      begin
        g=0;
        for j=1 step 1 until np do
          g=g+A[i,j]*T[i+r1+j,h];
          T[i,h]=T[i,h]-g
        end h;
        r1:=r1+1;
        if r1<np
          then go to E;
        r1=0;
      E: end i
    end
  end URP

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Carrying out a proper elimination of the first p unknowns of this subsystem we obtain

$$\begin{array}{cccccccc}
 1 & 0 & \dots & 0 & k_{1,p+1}^{i*} & \dots & k_{1n}^{i*} & \\
 0 & 1 & \dots & 0 & k_{2,p+1}^{i*} & \dots & k_{2n}^{i*} & \\
 \cdot & \cdot & \cdot & \cdot & \cdot & \cdot & \cdot & \cdot \\
 0 & 0 & \dots & 1 & k_{p,p+1}^{i*} & \dots & k_{pn}^{i*} & \\
 0 & 0 & \dots & 0 & d_{11}^{i*} & \dots & d_{1p}^{i*} & k_{1,p+1}^{i+1} \dots k_{1n}^{i+1} \\
 0 & 0 & \dots & 0 & d_{21}^{i*} & \dots & d_{2p}^{i*} & k_{2,p+1}^{i+1} \dots k_{2n}^{i+1} \\
 \cdot & \cdot & \cdot & \cdot & \cdot & \cdot & \cdot & \cdot \\
 0 & 0 & \dots & 0 & d_{pp}^{i*} & \dots & d_{pn}^{i*} & k_{p,p+1}^{i+1} \dots k_{pn}^{i+1} \\
 & & & & k_{p+1,1}^{i+1} & \dots & k_{p+1,p}^{i+1} & k_{p+1,p+1}^{i+1} \dots k_{p+1,n}^{i+1} \\
 0 & & & & \cdot & \cdot & \cdot & \cdot \\
 & & & & k_{n1}^{i+1} & \dots & k_{np}^{i+1} & k_{n,p+1}^{i+1} \dots k_{nn}^{i+1}
 \end{array}$$

where all elements are unchanged except those signed with *. We apply the same procedure for submatrices K_1, K_2, \dots, K_{lm} . In the reverse procedure only elements k_{tq}^{i*} ($t = 1, 2, \dots, p; q = p+1, p+2, \dots, n$) are necessary for calculation of the corresponding unknowns of the i -th step.

Conclusion. In the i -th step of the initial procedure only submatrices K_i and K_{i+1} are processed. In the i -th step of the reverse procedure only the matrix $[k_{rs}]$ of dimension $p \times p$ is processed.

In the procedure, the elimination process is carried out with selection of the main element.

3. Certification. Procedure *URP* was tested for some numerical examples. In all cases the results were the same as for the well-known Gauss procedure. The calculations were performed on the ODRA 1204 computer.

Remark. The assumption that the marked elements in Fig. 1 are summarized does not limit the application of the algorithm. If needed, the appropriate matrix elements can be set to be equal to zero.

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Received on 6. 1. 1976;
 revised version on 10. 1. 1977

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ROZWIĄZANIE PEWNEGO UKŁADU RÓWNAŃ LINIOWYCH

STRESZCZENIE

Procedura *URP* rozwiązuje grupę układów równań liniowych (1) o wspólnej macierzy układu K , zbudowanej z podmacierzy kwadratowych K_1, K_2, \dots, K_{lm} o wymiarach $n \times n$ tak, jak to pokazano na rys. 1. Elementy zakreskowane na rysunku są dodawane, co nie stanowi istotnego ograniczenia w stosowaniu procedury. Jest to specjalny typ macierzy pasmowej. Tego typu układy występują w zastosowaniach technicznych (np. przy analizie powłok obrotowych) na ogół z dużą liczbą niewiadomych. Przedstawiony algorytm jest efektywniejszy (głównie ze względu na wykorzystanie pamięci maszyny cyfrowej) niż analogiczne algorytmy, traktujące K jako macierz pasmową.

Dane:

- n — stopień podmacierzy K_i (n musi być liczbą parzystą);
- lm — liczba podmacierzy K_i ;
- lp — liczba rozwiązywanych układów;
- $T[1 : .5 \times n \times (lm + 1), 1 : lp]$ — tablica prawych stron układu (1) ($T[i, j]$ jest prawą stroną w i -tym równaniu j -tego układu);
- OMK* — procedura o nagłówku: **procedure OMK** (i, n, K);
integer i, n ; array K ; procedura oblicza elementy podmacierzy K_i dla $i = 1, 2, \dots, lm$.

Wyniki:

- $T[1 : .5 \times n \times (lm + 1), 1 : lp]$ — tablica rozwiązań układu (1) ($T[i, j]$ jest i -tą niewiadomą j -tego układu równań).

Inne parametry:

- exit* — etykieta poza treścią procedury, do której następuje skok, gdy macierz układu K jest osobliwa.