

```

hz := hf := 0;
for i := 1 step 1 until n do
begin
  x[i] := x[i] - dx[i]; hz := hz + abs(x[i]);
  hf := hf + abs(dx[i])
end;
if hf ≥ eps2 × hz then go to ITERATION;
ENDE: FUNCTION(x, f); S := 0; itmax := z;
for i := 1 step 1 until m do S := S + f[i] × f[i]
end TAYLOR

```

REFERENCE:

[1] BRAESS, D. Über Dämpfung bei Minimalisierungsverfahren. *Computing 1* (1966), 264-272.

ALGORITHM 316
SOLUTION OF SIMULTANEOUS NON-LINEAR EQUATIONS [C5]

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procedure *nonlinearsystem* (*n*, *maxit*, *numsig*, *singular*, *x*);
value *n*, *numsig*; **integer** *n*, *maxit*, *numsig*, *singular*; **array** *x*;
comment This procedure solves a system of *n* simultaneous nonlinear equations. The method is roughly quadratically convergent and requires only $((n^2/2) + (3n/2))$ function evaluations per iterative step as compared with $(n^2 + n)$ evaluations for Newton's Method. This results in a savings of computational effort for sufficiently complicated functions. A detailed description of the general method and proof of convergence are included in [1]. Basically the technique consists in expanding the first equation in a Taylor series about the starting guess, retaining only linear terms, equating to zero and solving for one variable, say x_k , as a linear combination of the remaining $n - 1$ variables. In the second equation, x_k is eliminated by replacing it with its linear representation found above, and again the process of expanding through linear terms, equating to zero and solving for one variable in terms of the now remaining $n - 2$ variables is performed. One continues in this fashion, eliminating one variable per equation, until for the *n*th equation, we are left with one equation in one unknown. A single Newton step is now performed, followed by back-substitution in the triangularized linear system generated for the x_i 's. A pivoting effect is achieved by choosing for elimination at any step that variable having a partial derivative of largest absolute value. The pivoting is done without physical interchange of rows or columns.

The vector of initial guesses *x*, the number of significant digits desired *numsig*, the maximum number of iterations to be used, *maxit*, and the number of equations *n*, should be set up prior to the procedure call which activates *nonlinearsystem*. After execution of the procedure, the vector *x* is the solution of the system (or best approximation thereto), *maxit* is now the number of iterations used and *singular* = 0 is an indication that a Jacobian-related matrix was singular—indicative of the process "blowing-up," whereas *singular* = 1 is an indication that no such difficulty occurred. Storage space may be saved by implementing the algorithm in a way which takes advantage of the fact that the strict lower triangle of the array *pointer* and the same number of positions in the array *coe* are not used;

```

begin integer converge, m, j, k, i, jsub, itemp, kmax, kplus, tally;
real f, hold, h, fplus, dermax, test, factor, relconvg;
integer array pointer[1:n, 1:n], isub[1:n-1];
array temp, part[1:n], coe[1:n, 1:n+1];
procedure backsubstitution (k, n, x, isub, coe, pointer);
  value k, n;
  integer k, n; integer array isub, pointer; array x, coe;

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comment This procedure back-solves a triangular linear system for improved  $x[i]$  values in terms of old ones;
begin integer km, kmax, jsub;
  for km := k step -1 until 2 do
  begin kmax := isub[km-1]; x[kmax] := 0;
    for j := km step 1 until n do
    begin jsub := pointer[km, j];
      x[kmax] := x[kmax] + coe[km-1, jsub] × x[jsub]
    end;
    x[kmax] := x[kmax] + coe[km-1, n+1]
  end;
end backsubstitution;
procedure evaluatekthfunction (x, y, k);
  integer k; real y; array x;
begin comment the body of this procedure must be provided by the user. One call of the procedure should cause the value of the kth function at the current value of the vector x to be placed in y;
end evaluatekthfunction;
converge := 1; singular := 1; relconvg := 10 ↑ (-numsig);
for m := 1 step 1 until maxit do
begin
  comment An intermediate output statement may be inserted at this point in the procedure to print the successive approximation vectors x generated by each complete iterative step;
  for j := 1 step 1 until n do pointer [1, j] := j;
  for k := 1 step 1 until n do
  begin if k > 1 then backsubstitution (k, n, x, isub, coe, pointer);
    evaluatekthfunction (x, f, k); factor := .001;
  AAA: tally := 0; for i := k step 1 until n do
  begin itemp := pointer[k, i]; hold := x[itemp];
    h := factor × hold; if h = 0 then h := .001;
    x[itemp] := hold + h;
    if k > 1 then backsubstitution (k, n, x, isub, coe, pointer);
    evaluatekthfunction (x, fplus, k);
    part[itemp] := (fplus - f)/h;
    x[itemp] := hold; if (abs(part[itemp]) = 0) ∨
      (abs(f/part[itemp]) > 1.01020) then tally := tally + 1;
  end;
  if tally ≤ n - k then go to AA; factor := factor × 10.0;
  if factor > .5 then go to SING; go to AAA;
  AA: if k < n then go to A; if abs (part[itemp]) = 0
  then go to SING;
  coe[k, n+1] := 0; kmax := itemp; go to ENDK;
  A: kmax := pointer[k, k]; dermax := abs(part[kmax]);
    kplus := k + 1;
    for i := kplus step 1 until n do
    begin jsub := pointer[k, i]; test := abs(part[jsub]);
      if test < dermax then go to B; dermax := test;
      pointer [kplus, i] := kmax; kmax := jsub;
    go to ENDI;
    B: pointer [kplus, i] := jsub;
  ENDI:
  end;
  if abs(part[kmax]) = 0 then go to SING; isub[k] := kmax;
  coe[k, n+1] := 0;
  for j := kplus step 1 until n do
  begin jsub := pointer[kplus, j];
    coe[k, jsub] := -part[jsub]/part[kmax];
    coe[k, n+1] := coe[k, n+1] + part[jsub] × x[jsub]
  end;
  ENDK:
  coe[k, n+1] := (coe[k, n+1] - f)/ part[kmax] + x[kmax]
end k;
x[kmax] := coe[n, n+1];
if n > 1 then backsubstitution (n, n, x, isub, coe, pointer);
if m = 1 then go to D;
for i := 1 step 1 until n do
  if abs((temp[i] - x[i])/x[i]) > relconvg then go to C;

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converge := converge + 1;
if converge ≥ 3 then go to TERMINATE else go to D;
C: converge := 1;
D: for i := 1 step 1 until n do temp[i] := x[i]
end m;
go to THROUGH;
TERMINATE:
maxil := m; go to THROUGH;
SING:
singular := 0;
THROUGH:
end nonlinearssystem

```

APPENDIX

We include a sample procedure *evaluatekthfunction* for the 2×2 system:

$$\left(1 - \frac{1}{4\pi}\right) (e^{2x_1} - e) + \frac{e}{\pi} x_2 - 2ex_1 = 0$$

$$\frac{1}{2} \sin(x_1 x_2) - \frac{x_2}{4\pi} - \frac{x_1}{2} = 0,$$

one solution of which is $(.5, \pi)$ see [2]

```

procedure evaluatekthfunction (x, y, k);
integer k; real y; array x;
begin switch functionnumber := F1, F2;
go to functionnumber [k];
F1: y := 2.71828183 × (.920422528 × (exp(2×x[1]-1)-1)+
x[2]/3.14159265-2×x[1]);
go to RETURN;
F2: y := .5 × sin(x[1]×x[2]) - x[2]/12.5663706 - x[1]/2;
RETURN:
end evaluatekthfunction;

```

REFERENCES:

- BROWN, K. M. A quadratically convergent method for solving simultaneous non-linear equations. Doctoral Thesis, Dept. Computer Sciences, Purdue U., Lafayette, Ind., Aug., 1966.
- BROWN, K. M., AND CONTE, S. D. The solution of simultaneous nonlinear equations. Proc. ACM 22nd Nat. Conf., pp 111-114.

ALGORITHM 317*

PERMUTATION [G6]

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```

procedure permute(n, k, v); value n, k; integer array v;
integer n, k;
comment This procedure produces in the vector v the kth
permutation on n variables. When k = 0, v takes on the value
1, 2, 3, 4, ..., n. This algorithm is not as efficient as pre-
viously published algorithms [1], [2], [3] for generating a
complete set of permutations but it is significantly better
for generating a random permutation, a property useful in
certain simulation applications. Any non-negative value of
k will produce a valid permutation. To generate a random
permutation, k should be chosen from the uniform distribu-
tion over the integers from 0 to n! - 1 inclusive;
begin integer i, g, r, x, j;
for i := 1 step 1 until n do v[i] := 0;
for i := n step -1 until 1 do
begin
q := k ÷ i; r := k - q × i; x := 0; j := n;

```

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no: if v[j] = 0 then
begin
if x = r then go to it else x := x + 1
end;
j := j - 1; go to no;
it: v[j] := i; k := q;
end
end

```

REFERENCES:

- COVEYOU, R. R., AND SULLIVAN, J. G. Algorithm 71, Permutation. *Comm. ACM* 4 (Nov. 1961), 497.
- PECK, J. E. L., AND SCHRACK, G. F. Algorithm 86, Permute. *Comm. ACM* 5 (Apr. 1962), 208.
- TROTTER, H. F. Algorithm 115, Perm. *Comm. ACM* 5 (Aug. 1962), 434.

Algorithms Policy • Revised August, 1966

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<i>in</i> symbol	<i>in</i> real	<i>ou</i> tarray	<i>in</i> teger
<i>ou</i> tsymbol	<i>ou</i> treal	<i>ou</i> tboolean	<i>ou</i> tinteger
<i>le</i> ngth	<i>in</i> array	<i>ou</i> tstring	

If only one channel is used by the program for output, it should be designated by 1 and similarly a single input channel should be designated by 2. Examples:

```

outstring (1, 'x='); outreal (1,x);
for i := 1 step 1 until n do outreal (1,A[i]);
integer (2, digit [17]);

```

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