

Algorithms

H. J. WEGSTEIN, Editor

ALGORITHM 121

NORMDEV

DAVID SHAFER

University of Chicago, Chicago, Ill.

```

procedure NormDev(Random,A,x);
  procedure Random; real A,x;
  comment 'NormDev' uses (1) a procedure 'Random(y)' assumed to produce a random number,  $0 < y < 1$ , and (2) the constant  $A = \sqrt{2/\pi} \times \int_0^1 \exp(-x^2/2)dx$ , to produce a positive normal deviate 'x';
  begin real y;
    Random(x); if  $x > A$  then go to large;
     $x := x/A$ ;
    1: Random(y); if  $y < \exp(-x^2/2)$  then go to EndND;
      Random(x); go to 1;
    large:  $x := (x - A)/(1 - A)$ ;
    2:  $x := \sqrt{1 - 2 \times \log(x)}$ ;
      Random(y); if  $y < 1/x$  then go to EndND;
      Random(x); go to 2;
  EndND: end

```



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ALGORITHM 122

TRIDIAGONAL MATRIX

GERARD F. DIETZEL

Burroughs Corp., Pasadena, Calif.

```

procedure TRIDIAG (n,A,U);
  integer n; array A,U;
  comment This procedure reduces a real symmetric matrix A of order n to tridiagonal form (UT)AU (UT = transpose of U) by a sequence of at most  $(n-1)(n-2)/2$  binary orthogonal transformations. Also, the matrix U is calculated. [Cf. W. Givens, "Numerical computation of the characteristic values of a real symmetric matrix," Report ORNL1574 (1954), Oak Ridge Nat. Lab., Tenn., and D. E. Johansen, "A modified Givens method for the eigenvalue evaluation of large matrices," J. ACM 8, 3 (1961)];
  begin real fact,c1,c2,loc1,loc2,temp; integer i,j,j1,j2,j3,j4,n1;
  comment Set array U = identity matrix of order n;
  for i := 1 step 1 until n do
    begin
      for j := i+1 step 1 until n do U[i,j] := U[j,i] = 0;
      U[i,i] := 1.0
    end;
  comment The reduction of the matrix A begins here. Only the upper triangular elements of A are used in the computation;
  n1 := n - 2;
  for i := 1 step 1 until n1 do
    begin
      j1 := i + 1; j2 := i + 2;
      for j := j2 step 1 until n do
        begin
          if A[i,j] = 0 then go to lab;
          fact := 1 / sqrt(A[i,j1]2 + A[i,j]2);
          c1 := fact × A[i,j1]; c2 := fact × A[i,j];
          loc1 := A[j1,j1]; loc2 := A[j1,j];
          A[j1,j1] := c12 × loc1 + 2.0 × c1 × c2 × loc2 + c22 × A[j,j];
          A[j1,j] := -c1 × c2 × loc1 + (c12 - c22) × loc2 + c1 × c2 × A[j,j];
          A[j,j] := c22 × loc1 - 2.0 × c1 × c2 × loc2 + c12 × A[j,j];
          j3 := j + 1;
          for k := j3 step 1 until n do
            begin
              temp := A[j1,k];
              A[j1,k] := c1 × temp + c2 × A[j,k];
              A[j,k] := -c2 × temp + c1 × A[j,k]
            end;
          j4 := j - 1;
          for k := j2 step 1 until j4 do
            begin
              temp := A[j1,k];
              A[j1,k] := c1 × temp + c2 × A[k,j];
              A[k,j] := -c2 × temp + c1 × A[k,j]
            end;
          A[i,j1] := c1 × A[i,j1] + c2 × A[i,j];
          A[i,j] := 0;
          for k := 1 step 1 until n do

```



```

begin
  temp := U[k,j1];
  U[k,j1] := c1 × temp + c2 × U[k,j];
  U[k,j] := -c2 × temp + c1 × U[k,j];
end;
lab: end
end;
for i := 1 step 1 until n do
  for j := i+1 step 1 until n do
    A[j,i] := A[i,j]
  end
end TRIDIAG

```

ALGORITHM 123

REAL ERROR FUNCTION, $\text{ERF}(x)$

MARTIN CRAWFORD AND ROBERT TECHO

Georgia Institute of Technology, Atlanta, Ga.

```

real procedure Erf(x); real x;
comment  $\Phi(x) = \text{Erf}(x) = (2/\sqrt{\pi}) \int_0^x e^{-u^2} du$  can be computed
  by using the recursive relation for derivatives with  $\Phi'(x) =$ 
   $(2/\sqrt{\pi})e^{-x^2}$ , where  $\Phi^{(n)}(x) = -2x\Phi^{(n-1)}(x) - 2(n-2)\Phi^{(n-2)}(x)$ ,
  for  $n = 2, 3, \dots$ . The Taylor's series expansions of  $\Phi(a_k)$  are
  taken about  $k+1$  points on the interval  $0 < a_k \leq x$  and summed
  to get  $\Phi(x)$ ;
begin real A, U, V, W, Y, Z, T; integer N;
  Z := 0; 1: if  $x \neq 0$  then
begin if  $0.5 < \text{abs}(x)$  then A := - sign(x) × 0.5
  else A := - x;
  U := V := 1.12837917 × exp(-x2); Y := T := -V ×
  A; N := 1;
2: if  $\text{abs}(T) \geq 10^{-10}$  then
begin N := N + 1; W := -2 × x × V - 2 × U × (N-2);
  T := T × W × A/(V × N);
  U := V; V := W; Y := Y + T; go to 2 end;
  Z := Z + Y; x := x + A; go to 1 end;
  Erf := Z end Erf

```

ALGORITHM 124

HANKEL FUNCTION

LUIS J. SCHAEFER

Purdue University, West Lafayette, Ind.

```

procedure HANKEL(N,X,H); value N,X; integer N;
real X; array H;
comment This procedure evaluates the complex valued hankel
  function of the first kind for real argument X and integral order
  N and assigns it to H. The individual Bessel- and Neuman-func-
  tion series are not evaluated separately. Both the real and
  imaginary parts are generated from the same terms;
begin real K, P, R, A, S, T, D, L; integer Q;
  A := R := 1; H[1] := H[2] := S := 0;
  for Q := 1 step 1 until N do begin R := R × Q; S := S +
  1/Q end; D := R/N;
  R := 1/R; K := X × X/4; P := (X/2)N; T := ln(K) +
  1.1544313298631;
  for Q := 0, Q+1 while Q ≤ N ∨ L ≠ H[2] do
  begin L := H[2]; H[1] := H[1] + A × K × R;
  H[2] := H[2] + A × (R × K × (T - S) - (if Q < N then D/P
  else 0));
  A := A × K/Q; R := -R/(Q+N); S := S + 1/Q + 1/(Q+N);
  if Q < N then D := D/(N-Q)
  end; H[2] := H[2] × .31830989
end

```

Notes on Programming Languages

On the Nonexistence of a Phrase Structure Grammar for ALGOL 60

Robert W. Floyd

Computer Associates, Inc., Woburn, Massachusetts

ALGOL 60 is defined partly by formal mechanisms of phrase structure grammar, partly by informally stated restrictions. It is shown that no formal mechanisms of the type used are sufficient to define ALGOL 60.

Let a phrase structure grammar be defined as a set of definitions in the Backus notation used to define ALGOL 60 [1]. A phrase structure language, then, is a language defined by such a grammar. In such a language, because of the finite number of syntactic types, all sufficiently long programs (blocks) contain a substring which in turn contains a proper substring of the same syntactic type as itself [2, 3]. Thus, the program P takes the form $QRSTU$, where RST and S are of the same syntactic type so that either may be substituted for the other, and S is a proper substring of RST . Now $QR^{(i)}ST^{(i)}U$ is a syntactically correct program for any non-negative integer i , where $R^{(i)}$ denotes i occurrences of the string R .

Example. An ALGOL 60 program might contain the primary $(c \times d)$, which in turn contains the primary c . It would be possible, therefore, to replace $(c \times d)$ by c at that point in the program, or to replace c by $(c \times d)$ obtaining $((c \times d) \times d)$, etc., without destroying the syntactic correctness of the program.

The goal of the present paper is to exhibit a set of ALGOL 60 programs of unbounded length, and show that none of them has the property described by the first paragraph. This implies that ALGOL 60 is not definable by a phrase structure grammar alone.

Consider the ALGOL 60 program

```
begin real  $x^{(n)}$ ;  $x^{(n)} := x^{(n)}$  end
```

where $x^{(n)}$ stands for n occurrences of the letter x . If ALGOL 60 is a phrase structure language, we may choose n sufficiently large to make applicable the result of the first paragraph. That is,

```
begin real  $x^{(n)}$ ;  $x^{(n)} := x^{(n)}$  end
```

takes the form $QRSTU$, and $QR^{(i)}ST^{(i)}U$ is a syntactically correct program P_i for all $i \geq 0$. A block in ALGOL must contain at least one declarator, a semicolon, and the words **begin** and **end**; since $QSU = P_0$ is a block, R and T can contain only the characters x and $:=$. Since the declarator **real** occurs only once and there are no commas, only one identifier is declared in each of P_i , and only that identifier may be used in P_i . Two cases arise:

(1) Neither R nor T contains $:=$. Then $R = x^{(j)}$ and $T = x^{(k)}$ with j and k not both zero. One cannot, however, delete x 's from P_1 in two places R and T and still have all identifiers properly declared in P_0 ; at least three deletions would have to be made.

(2) R or T contains $:=$. Since P_0 does not contain $:=$, it must take the form

```
begin real  $x^{(j)}$ ; end
```