

Algorithms

ALGORITHM 35

SIEVE

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```

procedure Sieve (Nmax) Primes: (p) ;
    integer Nmax; integer array p ;
comment Sieve uses the Sieve of Eratosthenes to find all prime
    numbers not greater than a stated integer Nmax
    and stores them in array p. This array should be
    of dimension 1 by entier (2 × Nmax/ln (Nmax)) ;
begin integer n, i, j ;
    p[1] := 1 ; p[2] := 2 ; p[3] := j := 3 ;
    for n := 3 step 2 until Nmax do
begin
    i := 3 ;
    L1: go to if p[i] ≤ sqrt (n) then a1 else a2 ;
    a1: go to if n/p[i] = n ÷ p[i] then b1 else b2 ;
    b2: i := i + 1 ; go to L1 ;
    a2 : p[j] := n ; j := j + 1 ;
    b1: end end

```

ALGORITHM 36

TCHEBYCHEFF

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```

procedure tchebycheff (t, x, m, ℓ) ;
real array t, x ; integer ℓ, m ;
comment given a set of m+1 values of x stored in a one-
    dimensional array whose subscripts run from 0
    thru m at least, construct a table of  $t_n(x)$ ,  $n =$ 
    0, 1, ..., ℓ and store it in the two-dimensional
    array t, where you find  $t_n(x[m])$  as t[n, m] ;
begin integer i, k, n ;
    for k := 0 step 1 until m do begin t[0, k] := 1 ;
    t[1, k] := x[k] end ;
    for n := 2 step 1 until ℓ do for i = 0 step 1
    until m do
    t[n, i] := 2 × x[i] × t[n - 1, i] - t[n - 2, i]
end tcheby

```

ALGORITHM 37

TELESCOPE 1

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```

procedure Telescope 1 (N, L, eps, limit, c) ; value limit, L ;
integer N ; real L, eps, limit ; array c ;
comment: Telescope 1 takes an Nth degree polynomial approxi-
    mation  $\sum_{k=0}^N c_k x^k$  to a function which was valid to
    within eps ≥ 0 over an interval (0, L) and reduces
    it, if possible, to a polynomial of lower degree,
    valid to within limit > 0. The initial coefficients
     $c_k$  are replaced by the final coefficients, and the
    deleted coefficients are replaced by zero. The ini-
    tial eps is replaced by the final bound on the error.

```

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N is replaced by the degree of the reduced polynomial. N and eps must be variables.

This procedure computes the coefficients given in the Techniques Department of the ACM Communications, Vol. 1, No. 9, from the recursion formula

$$a_{k-1} = -a_k \cdot \frac{k \cdot L \cdot (2k - 1)}{2(N + k - 1) \cdot (N - k + 1)} ;$$

```

begin integer k ; array d[0:N] ;
if N < 1 then go to exit ; d[N] := -c[N] ;
for k := N step - 1 until 1 do
    d[k - 1] := -d[k] × L × k × (k - 0.5)/
    ((N + k - 1) × (N - k + 1)) ;
if eps + abs (d[0]) < limit then
    begin eps := eps + abs (d[0]) ;
    for k := N step - 1 until 0 do c[k] := c[k] + d[k] ;
    N := N - 1 ; go to start end ;
exit: end

```

ALGORITHM 38

TELESCOPE 2

K. A. BRONS

RCA Advanced Programming, Pennsauken, N. J.

```

procedure Telescope 2 (N, L, eps, limit, c) ; value limit, L ;
integer N ; real L, eps, limit ; array c ;
comment Telescope 2 takes an Nth degree polynomial ap-
    proximation  $\sum_{k=0}^N c_k x^k$  to a function which was
    valid to within eps ≥ 0 over an interval (-L, L)
    and reduces it, if possible, to a polynomial of
    lower degree, valid to within limit > 0. The initial
    coefficients  $c_k$  are replaced by the final coefficients,
    and deleted coefficients are replaced by zero. The
    initial eps is replaced by the final bound on the
    error, and N is replaced by the degree of the re-
    duced polynomial. N and eps must be variables.

```

This procedure computes the coefficients given in the Techniques Department of the ACM Communications, Vol. 1, No. 9, from the recursion formula

$$a_{k-2} = -a_k \frac{k \cdot L^2(k-1)}{(N+k-2) \cdot (N-k+2)} ;$$

```

start:  begin integer k ; real s ; array d[0: N] ;
        if N < 2 then go to exit ; d[N] := -c[N] ;
        for k := N step - 2 until 2 do
          d[k-2] := -d[k] × L ↑ 2 × k × (k-1)/
            ((N+k-2) × (N-k+2)) ;
        if (N/2) - entier (N/2) = 0 then s := d[0] else
          s := d[1]/N ;
        if eps + abs(s) < limit then begin
          eps := eps + abs(s) ;
          for k := N step - 2 until 0 do
            c[k] := c[k] + d[k] ;
          N := N - 1 ; go to start end ;
exit:    end

```

ALGORITHM 39 CORRELATION COEFFICIENTS WITH MATRIX MULTIPLICATION

PAPKEN SASSOUNI

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```

procedure NORM (x) number of rows: (m) number of columns:
              (n) normalized output: (y) standard deviations:
              (s) ;
value        m, n ; integer m, n ; array x, y, s ;
comment      Given an observation matrix [x] consisting of ob-
              servations xij on a population, NORM will cal-
              culate

```

$$y_{ij} = \frac{x_{ij} - \bar{x}_j}{\sqrt{\sum_{i=1}^m (x_{ij} - \bar{x}_j)^2}} \quad \text{for } i = 1, \dots, m \quad j = 1, \dots, n$$

and the standard deviations

$$s_j = \sqrt{\frac{\sum_{i=1}^m (x_{ij} - \bar{x}_j)^2}{m}}$$

where \bar{x}_j is the mean of observations on the j-th factor ;

```

begin  integer i, j ; real r, h, c, b ;
        r := sqrt (m) ; for j := 1 step 1 until n do
1: begin h := 0 ;
          for i := 1 step 1 until m do
            h := h + x[i, j] ; h := h/m ; b := 0 ;
          for i := 1 step 1 until m do
2: begin c := x[i, j] - h ; b := b + c ↑ 2 ; y[i, j] := c
          end 2 ;
          b := sqrt (b) ;
          for i := 1 step 1 until m do
            y[i, j] := y[i, j]/b ; s[j] := b/r
          end 1
        end NORM ;
comment The normalization is now completed, and we are
        ready to compute the correlation matrix ;
procedure TRANSMULT (y) number of rows: (m) number of
        columns: (n) symmetrical square matrix result:
        (z) ;
value   m, n ; integer m, n ; array y, z ;

```

comment This procedure multiplies two matrices, the first being the transpose of the second. The result is a symmetrical matrix with respect to the main diagonal, therefore only the lower part of it, including the main diagonal, is computed. The upper half is obtained by equating corresponding elements;

```

begin  integer i, j, k ; real h ;
        for j := 1 step 1 until n do
          for i := j step 1 until n do
            h := 0 ;
            for k := 1 step 1 until m do
              h := h + y[k, i] × y[k, j] ; z[i, j] := h ;
            if i ≠ j then z[j, i] := h
            end i
          end TRANSMULT. [z] is the square matrix of the
          correlation coefficients of the initial observation
          matrix [x]

```

ALGORITHM 40 CRITICAL PATH SCHEDULING

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```

procedure CRITICALPATH (n, I, J, DIJ, ES, LS, EF, LF, TF,
                        FF) ;
integer n ; integer array I, J, DIJ, ES, LS, EF, LF, TF,
                        FF ;

```

comment: Given the total number of jobs n of a project, the vector pair I_k, J_k representing the kth job, which is thought of as an arrow connecting event I_k to event J_k ($I_k < J_k$, $k = 1 \dots n$), and a duration vector $(DIJ)_k$, CRITICAL-PATH determines the earliest starting time $(ES)_k$, latest starting time $(LS)_k$, earliest completion time $(EF)_k$, latest completion time $(LF)_k$, the total float $(TF)_k$, and the free float $(FF)_k$. I_1 must be 1 and the I_k, J_k must be in ascending order. For example, if the first three jobs are labelled (1, 2), (1, 3), (3, 4), then the I, J vectors are (1, 1, 3) and (2, 3, 4) respectively. The critical path is given by each arrow whose total float is zero. The following non-local labels are used for exits: out1 — I_k not less than J_k ; out2 — I_k out of sequence ; out3 — I_k missing;

```

begin  integer k, index, max, min ; integer array ti, te [1:n] ;
        index := 1 ;
        for k := 1 step 1 until n do
          begin
            if I[k] ≥ J[k] then go to out1 ;
            if I[k] < index then go to out2 ;
            if I[k] > index ∧ I[k] ≠ index + 1 then go to out3 ;
            if I[k] = index + 1 then index := I[k] ;
          C: end ;
          for k := 1 step 1 until n do
            ti[k] := te[k] := 0 ;
            for k := 1 step 1 until n do
              begin
                max := ti[I[k]] + DIJ[k] ;
                if ti[J[k]] = 0 ∨ ti[J[k]] < max then
                  ti[J[k]] := max ;
              A: end ti ;
              te[J[n]] := ti[J[n]] ;
              for k := n step -1 until 1 do
                begin
                  min := te[J[k]] - DIJ[k] ;
                  if te[I[k]] = 0 ∨ te[I[k]] > min then
                    te[I[k]] := min ;
                B: end te ;
                for k := 1 step 1 until n do

```

```

begin
  ES[k] := ti[I[k]] ;
  LS[k] := te[J[k]] - DIJ[k] ;
  EF[k] := ti[I[k]] + DIJ[k] ;
  LF[k] := te[J[k]] ;
  TF[k] := te[J[k]] - ti[I[k]] - DIJ[k] ;
  FF[k] := ti[J[k]] - ti[I[k]] - DIJ[k] ;
end
end CRITICALPATH

```

REFERENCES

- (1) JAMES E. KELLEY, JR. AND MORGAN R. WALKER, "Critical-Path Planning and Scheduling," 1959 Proceedings of the Eastern Joint Computer Conference.
- (2) M. C. FRISHBERG, "Least Cost Estimating and Scheduling — Scheduling Phase Only," IBM 650 Program Library File No. 10.3.005.

REMARKS ON ALGORITHMS 2 AND 3 (*Comm. ACM*, February 1960), ALGORITHM 15 (*Comm. ACM*, August 1960) AND ALGORITHMS 25 AND 26 (*Comm. ACM*, November 1960)

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Algorithms 2, 15, 25 and 26 were all concerned with the calculation of zeros of arbitrary functions by successive linear or quadratic interpolation. The main limiting factor on the accuracy attainable with such procedures is the condition of the method of evaluating the function in the neighbourhood of the zeros. It is this condition which should determine the tolerance which is allowed for the relative error. With a well-conditioned method of evaluation quite a strict convergence criterion will be met, even when the function has multiple roots.

For example, a real quadratic root solver (of a type similar to Algorithm 25) has been used on ACE to find the zeros of triple-diagonal matrices T having $t_{ii} = a_i$, $t_{i+1,i} = b_{i+1}$, $t_{i,i+1} = c_{i+1}$. As an extreme case I took $a_1 = a_2 = \dots = a_6 = 0$, $a_6 = a_7 = \dots = a_{10} = 1$, $a_{11} = 2$, $b_1 = 1$, $c_1 = 0$ so that the function which was being evaluated was $x^3(x-1)^5(x-2)$. In spite of the multiplicity of the roots, the answers obtained using floating-point arithmetic with a 46-bit mantissa had errors no greater than 2^{-44} . Results of similar accuracy have been obtained for the same problem using linear interpolation in place of the quadratic. This is because the method of evaluation which was used, the two-term recurrence relation for the leading principal minors, is a very well-conditioned method of evaluation. Knowing this, I was able to set a tolerance of 2^{-42} with confidence. If the same function had been evaluated from its explicit polynomial expansion, then a tolerance of about 2^{-7} would have been necessary and the multiple roots would have obtained with very low accuracy.

To find the zero roots it is necessary to have an absolute tolerance for $|x_{i+1} - x_i|$ as well as the relative tolerance condition. It is undesirable that the preliminary detection of a zero root should be necessary. The great power of rootfinders of this type is that, since we are not saddled with the problem of calculating the derivative, we have great freedom of choice in evaluating the function itself. This freedom is encroached upon if we frame the rootfinder so that it finds the zeros of $x = f(x)$ since the true function $x - f(x)$ is arbitrarily separated into two parts. The formal advantage of using this formulation is very slight. Thus, in Certification 2 (June 1960), the calculation of the zeros of $x = \tan x$ was attempted. If the function $(-x + \tan x)$ were used with a general zero finder then, provided the method of evaluation was, for example

$$x = n\pi + y$$

$$\tan x - x = -n\pi + \frac{\frac{y^3}{3} - \frac{y^5}{30} - \dots}{\cos y},$$

the multiple zeros at $x = 0$ could be found as accurately as any of the others. With a slight modification of common sine and cosine routines, this could be evaluated as

$$-n\pi + \frac{(\sin y - y) - y(\cos y - 1)}{1 + (\cos y - 1)}$$

and the evaluation is then well-conditioned in the neighbourhood of $x = 0$. As regards the number of iterations needed, the restriction to 10 (Certification 2) is rather unreasonably small. For example, the direct evaluation of $x^{60} - 1$ is well conditioned, but starting with the values $x = 2$ and $x = 1.5$ a considerable number of iterations are needed to find the root $x = 1$. Similarly a very large number are needed for Newton's method, starting with $x = 2$. If the time for evaluating the derivative is about the same as that for evaluating the function (often it is much longer), then linear interpolation is usually faster, and quadratic interpolation much faster, than Newton.

In all of the algorithms, including that for Bairstow, it is useful to have some criterion which limits the permissible change from one value of the independent variable to the next [1]. This condition is met to some extent in Algorithm 25 by the condition $S4$, that $\text{abs}(fprt) < \text{abs}(x2 \times 10)$, but here the limitation is placed on the permissible increase in the value of the function from one step to the next. Algorithms 3 and 25 have tolerances on the size of the function and on the size of the remainders $r1$ and $r0$ respectively. They are very difficult tolerances to assign since these quantities may take very small values without our wishing to accept the value of x as a root. In Algorithm 3 (*Comm. ACM* June 1960) it is useful to return to the original polynomial and to iterate with each of the computed factors. This eliminates the loss of accuracy which may occur if the factors are not found in increasing order. This presumably was the case in Certification 3 when the roots of $x^5 + 7x^4 + 5x^3 + 6x^2 + 3x + 2 = 0$ were attempted. On ACE, however, all roots of this polynomial were found very accurately and convergence was very fast using single-precision, but the roots emerged in increasing order. The reference to slow convergence is puzzling. On ACE, convergence was fast for all the initial approximations to p and q which were tried. When the initial approximations used were such that the real root $x = -6.3509936103$ and the spurious zero were found first, the remaining two quadratic factors were of lower accuracy, though this was, of course, rectified by iteration in the original polynomial. When either of the other two factors was found first, then all factors were fully accurate even without iteration in the original polynomial [1].

REFERENCE

- [1] J. H. WILKINSON. The evaluation of the zeros of ill-conditioned polynomials Parts I and II. *Num. Math.* 1 (1959), 150-180.

CERTIFICATION OF ALGORITHM 4

BISECTION ROUTINE (S. Gorn, *Comm. ACM*, March 1960)

PATTY JANE RADER,* Argonne National Laboratory, Argonne, Illinois

BISEC was coded for the Royal-Precision LGP-30 computer, using an interpretive floating point system (24.2) with 28 bits of significance.

The following minor correction was found necessary.

α : go to γ_1 should be go to γ_i

* Work supported by the U. S. Atomic Energy Commission.

After this correction was made, the program ran smoothly for $F(x) = \cos x$, using the following parameters:

y_1	y_2	ϵ	α	Results
0	1	.001	.001	FLSXT
0	2	.001	.001	1.5703
1.5	2	.001	.001	1.5703
1.55	2	.1	.1	1.5500
1.5	2	.001	.1	1.5625

These combinations test all loops of the program.

* Work supported by the U. S. Atomic Energy Commission.

REMARK ON ALGORITHM 16

CROUT WITH PIVOTING (G. E. Forsythe, *Comm. ACM*, 3 (Sept. 1960), 507-8.)

HENRY C. THACHER, JR.,* Argonne National Laboratory, Argonne, Illinois

This procedure contains the following errors:

a. In SOLVE, the expression

$c[k] := c[k] - \text{INNERPRODUCT}$
 $(B[k, p], c[p], p - 1, k - 1)$

should read:

$c[k] := c[k] - \text{INNERPRODUCT}$
 $(B[k, p], c[p], p, 1, k - 1)$

b. In CROUT, the specification part should read:

array A, b, y ; **integer** n ; **integer array** pivot ;

c. In SOLVE, the specification part should read:

array B, c, z ; **integer** n ; **integer array** pivot ;

The efficiency of the algorithm will be improved by the following changes:

a. In the elimination phase of CROUT, replace

for i := k + 1 **step** 1 **until** n **do**

begin quote := 1.0/A[k, k] ; A[i, k] := quote XA[i, k] **end** ;

by

quote := 1.0/A[k, k] ; **for** i := k + 1 **step** 1 **until** n **do**
A[i, k] := quote XA[i, k] ;

b. Omit INNERPRODUCT from the formal parameter list in both CROUT and SOLVE, and declare INNERPRODUCT either locally, or globally. This avoids any reference to INNERPRODUCT in the calling sequence produced by a compiler.

It is also to be noted that a minor modification of CROUT allows it to be used to evaluate the determinant of A.

All of these suggestions are included in a later algorithm.

* Work supported by the U. S. Atomic Energy Commission.

REMARK ON ALGORITHM 25

REAL ZEROS OF AN ARBITRARY FUNCTION

(B. Leavenworth, *Comm. ACM*, November 1960)

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On attempting to use this algorithm, I discovered the two following errors:

(1) The line following the SWITCH statement should read:
for L := 1 **step** 1 **until** n **do**

(2) The line starting with the label loop: should read:

loop: dd := 1 + d ; bi = x0 × d ↑ 2 - x1 × dd ↑ 2
+ x2 × (dd + d) ;

With these two modifications incorporated the algorithm was translated into the language of the Burroughs Algebraic Compiler and has been used successfully on the Burroughs 220 Computer.

“COMPUTERS— KEY TO TOTAL SYSTEMS CONTROL” IS THEME OF 1961 EASTERN JOINT COMPUTER CONFERENCE

*Bruce G. Oldfield, Program Chairman,
Calls For Papers To Be Presented
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