

Note

On the Numerical Evaluation of the Ordinary Bessel Function of the Second Kind

1. INTRODUCTION

1.1. *Definitions and Relevant Properties*

The ordinary Bessel function of the first kind

$$J_\nu(z) = (z/2)^\nu \sum_{k=0}^{\infty} \frac{(-z^2/4)^k}{\Gamma(\nu + k + 1) k!} \tag{1.1}$$

and the ordinary Bessel function of the second kind

$$Y_\nu(z) = [\cos \nu\pi J_\nu(z) - J_{-\nu}(z)]/\sin \nu\pi \tag{1.2}$$

are two linearly independent solutions of the difference equation

$$f_{\nu+1} - (2\nu/z)f_\nu + f_{\nu-1} = 0. \tag{1.3}$$

This equation can be used to compute $Y_{\nu+n}$ for $n = 2, 3, \dots$ when Y_ν and $Y_{\nu+1}$ are given. In the forward direction the recurrence formula (1.3) for Y_ν is numerically stable, whereas it is unstable for J_ν (see Gautschi [1]).

The ordinary Bessel functions of the third kind are the Hankel functions

$$H_\nu^{(1)}(z) = J_\nu(z) + iY_\nu(z), H_\nu^{(2)}(z) = J_\nu(z) - iY_\nu(z). \tag{1.4}$$

Important for the representation of the Hankel functions for large $|z|$ are the functions $P(\nu, z)$ and $Q(\nu, z)$ defined by

$$H_\nu^{(1,2)}(z) = [2/(\pi z)]^{1/2} e^{\pm i\chi} [P(\nu, z) \pm iQ(\nu, z)], \tag{1.5}$$

where the $+$ sign is used for $H_\nu^{(1)}$, the $-$ sign is used for $H_\nu^{(2)}$ and

$$\chi = z - \pi(2\nu + 1)/4. \tag{1.6}$$

For large $|z|$, P and Q are slowly varying and the oscillatory behavior of $H_\nu^{(1)}$ and

$H_\nu^{(2)}$ is contained in the exponential function in (1.5). From (1.4) and (1.5) we obtain

$$\begin{aligned} Y_\nu(z) &= [2/(\pi z)]^{1/2} [P(\nu, z) \sin \chi + Q(\nu, z) \cos \chi] \\ J_\nu(z) &= [2/(\pi z)]^{1/2} [P(\nu, z) \cos \chi - Q(\nu, z) \sin \chi]. \end{aligned} \quad (1.7)$$

Again, the oscillatory behavior of J_ν and Y_ν is fully described by the circular functions in (1.7).

The connection between the ordinary Bessel functions and the modified Bessel functions follows from

$$\begin{aligned} H_\nu^{(1)}(z) &= -2i\pi^{-1}e^{-\nu\pi i/2}K_\nu(ze^{-i\pi/2}) & (-\tfrac{1}{2}\pi < \arg z \leq \pi), \\ H_\nu^{(2)}(z) &= 2i\pi^{-1}e^{\nu\pi i/2}K_\nu(ze^{i\pi/2}) & (-\pi < \arg z \leq \tfrac{1}{2}\pi). \end{aligned} \quad (1.8)$$

From the Wronskian

$$J_{\nu+1}(z) Y_\nu(z) - J_\nu(z) Y_{\nu+1}(z) = 2/(\pi z)$$

and (1.7) it easily follows that

$$P(\nu, z) P(\nu + 1, z) + Q(\nu, z) Q(\nu + 1, z) = 1. \quad (1.9)$$

1.2. Contents of the Paper

We give algorithms for the computation of Y_ν and $Y_{\nu+1}$ and we use the methods of our previous paper on the computation of K_ν and $K_{\nu+1}$ (see Temme [6]). Our results in [6] can be used for complex values of z . Here we give the explicit results for Y_ν and $Y_{\nu+1}$ and these results follow immediately from [6] by using (1.8).

For the computation of J_ν the reader is referred to Gautschi [1], where an algorithm is given for the computation of $J_{\nu+n}(z)$, $n = 0, 1, 2, \dots, N$. See also Gautschi [2]. In Luke [4] rational approximations for J_ν and Y_ν are given based on Padé-representations for large $|z|$. In Luke [5] a double series of Chebyshev polynomials and values of the coefficients are given for both Y_ν, J_ν for $z \geq 5$. In Goldstein and Thaler [3] the computation of Y_ν is based on series expansions in ordinary Bessel functions of the first kind, but the treatment of small $|\nu|$ -values is not satisfactory.

2. THE COMPUTATION FOR SMALL $|z|$

In order to obtain a more symmetric representation in (1.2) we write

$$\cos \nu\pi J_\nu(z) - J_{-\nu}(z) = J_\nu(z) - J_{-\nu}(z) - 2 \sin^2(\nu\pi/2) J_\nu(z). \quad (2.1)$$

Furthermore we introduce the following notation

$$\begin{aligned} c_k &= (-z^2/4)^k/k!, \\ p_k &= (\nu/\sin \nu\pi) (z/2)^{-\nu}/\Gamma(k+1-\nu), \\ q_k &= (\nu/\sin \nu\pi) (z/2)^\nu/\Gamma(k+1+\nu), \\ f_k &= (p_k - q_k)/\nu, \\ g_k &= f_k + 2\nu^{-1} \sin^2(\nu\pi/2) q_k, \\ h_k &= -kg_k + p_k, \end{aligned}$$

where $k = 0, 1, \dots$. We have for $k = 1, 2, \dots$ the recurrence relations

$$\begin{aligned} p_k &= p_{k-1}/(k-\nu), \quad q_k = q_{k-1}/(k+\nu), \\ f_k &= (kf_{k-1} + p_{k-1} + q_{k-1})/(k^2 - \nu^2). \end{aligned}$$

Substitution of (1.1) in (1.2) and using (2.1) yields

$$Y_\nu(z) = - \sum_{k=0}^{\infty} c_k g_k. \quad (2.2)$$

Considering (2.1) with ν replaced by $\nu + 1$ and using (1.3) we have

$$\begin{aligned} &\cos(\nu + 1)\pi J_{\nu+1}(z) - J_{-\nu-1}(z) \\ &= -[J_{\nu+1}(z) - J_{-\nu-1}(z)] + (2\nu/z) J_{-\nu}(z) + 2 \sin^2(\nu\pi/2) J_{\nu+1}(z). \end{aligned}$$

We obtain by substitution of (1.1)

$$Y_{\nu+1}(z) = - (2/z) \sum_{k=0}^{\infty} c_k h_k. \quad (2.3)$$

As in [6], f_0 can be represented in such a way that it can be computed with a satisfactorily small relative error.

For small values of $|z|$ the series in (2.2) and (2.3) converge rapidly. But cancellation may occur in summing the series numerically. A strict error analysis, as for the modified Bessel function, can not easily be given, but from numerical experiments it turns out that for $|z| < 3$ the computation is stable.

3. THE COMPUTATION FOR $|z| \geq 3$

For $|z| \geq 3$ we compute $P(\nu, z)$, $P(\nu + 1, z)$, $Q(\nu, z)$ and $Q(\nu + 1, z)$, by using the functions $k_n(z)$ introduced in our previous paper [6]. For K_ν and $K_{\nu+1}$ we needed $k_0(z)$ and $k_1(z)$. From (1.8) it turns out that for the P - and Q -functions the functions $k_0(\pm iz)$ and $k_1(\pm iz)$ can be used. The application of the method in [6] is straightforward. However, the determination of the starting index N for the Miller

algorithm caused some trouble, since our error analysis in [6] was based on the case of real variables. But trying out the results of [6] for the P - and Q -functions we noticed that the determination of the starting index N can indeed be based upon the estimations given in [6].

4. ALGOL 60 PROCEDURES

The algorithms for the computation of $Y_\nu(z)$ and $Y_{\nu+1}(z)$ are given as an ALGOL 60 procedure for the case of real values of ν and z , $z > 0$. For convenience we write $\nu = a$ and $z = x$.

The procedure *bessya* computes for $x > 0$ and $a \in \mathbb{R}$ the functions $Y_a(x)$ and $Y_{a+1}(x)$; *bessya* calls for three nonlocal procedures *sinh*, *recip gamma*, and *besspqa*. For the text of *sinh*, and *recip gamma* the reader is referred to [6]. In *besspqa* the functions $P(a, x)$, $P(a + 1, x)$, $Q(a, x)$ and $Q(a + 1, x)$ are computed. We supply *besspqa* as a separate procedure since it can also be used for the computation of the Bessel functions $J_a(x)$ and $J_{a+1}(x)$ (see (1.7)). In *bessya* the procedure *besspqa* is called for $x \geq 3$ and $|a| < .5$, but the algorithm in *besspqa* converges for all x and a ($x > 0$). It is recommended, however, to take $x > \max(|a|, 3)$. For $|a| > x$ the recurrence relations

$$\begin{aligned} P(a + 1, x) &= P(a - 1, x) - 2a/x Q(a, x) \\ Q(a + 1, x) &= Q(a - 1, x) + 2a/x P(a, x) \end{aligned}$$

can be used. These relations are valid for real a and x . They can be derived by substitution of (1.5) in (1.3). However, for $|a| + 1 > x$, computation of $J_a(x)$ and $J_{a+1}(x)$ by using (1.7) will cause a loss of correct significant digits.

The precision in the procedures *bessya* and *besspqa* can be controlled by using the variable *eps*. For *besspqa* its entry value corresponds to the desired relative accuracy in pa , $pa 1$, qa and $qa 1$. Also in *bessya* it corresponds to relative accuracy, except in the neighborhoods of zeros of $Y_a(x)$ or $Y_{a+1}(x)$. In that case ya or $ya 1$ are given with absolute accuracy *eps*.

The procedures *bessya* and *besspqa* were tested on the CD CYBER 73 of SARA, Amsterdam. For $a = 0, 0.2, 0.4$, $x = .5, 1, 2, 3, 5, 7, 10, 20, 50, 100$ and $eps = 10^{-15}$ we checked relation (1.9). The output of $|pa.pa 1 + qa.qa 1 - 1|$ is given in Table I. The procedure *bessya* was also tested in the neighborhood of $x = 3$. For $x^\pm = 3 \pm 2^{-46}$ we computed the numerical values of the expressions

$$\begin{aligned} d_0 &= \{Y_a(x^-) - Y_a(x^+)\}, \\ d_1 &= \{Y_{a+1}(x^-) - Y_{a+1}(x^+)\}. \end{aligned}$$

In Table II we give d_0 , d_1 , the maximum number of terms (n) used in (2.1), and the starting index N for the Miller algorithm.

TABLE I

| $x \backslash a$ | 0.0 | 0.2 | 0.4 |
|------------------|------------------------|------------------------|------------------------|
| 0.5 | 1.4 ₁₀ - 14 | 7.1 ₁₀ - 15 | 0.0 ₁₀ + 00 |
| 1.0 | 0.0 ₁₀ + 00 | 7.1 ₁₀ - 15 | 7.1 ₁₀ - 15 |
| 2.0 | 7.1 ₁₀ - 15 | 2.8 ₁₀ - 14 | 7.1 ₁₀ - 15 |
| 3.0 | 7.1 ₁₀ - 15 | 0.0 ₁₀ + 00 | 0.0 ₁₀ + 00 |
| 5.0 | 7.1 ₁₀ - 15 | 1.4 ₁₀ - 14 | 0.0 ₁₀ + 00 |
| 7.0 | 7.1 ₁₀ - 15 | 7.1 ₁₀ - 15 | 1.4 ₁₀ - 14 |
| 10.0 | 7.1 ₁₀ - 15 | 7.1 ₁₀ - 15 | 7.1 ₁₀ - 15 |
| 20.0 | 0.0 ₁₀ + 00 | 7.1 ₁₀ - 15 | 0.0 ₁₀ + 00 |
| 50.0 | 2.1 ₁₀ - 14 | 1.4 ₁₀ - 14 | 0.0 ₁₀ + 00 |
| 100.0 | 2.1 ₁₀ - 14 | 7.1 ₁₀ - 15 | 7.1 ₁₀ - 15 |

TABLE II

| eps | | 5.0 ₁₀ - 06 | 5.0 ₁₀ - 09 | 5.0 ₁₀ - 12 | 5.0 ₁₀ - 14 |
|----------|-----------------|------------------------|------------------------|------------------------|------------------------|
| <i>a</i> | | | | | |
| 0.0 | <i>d0</i> | 5.2 ₁₀ - 08 | 4.3 ₁₀ - 11 | 3.4 ₁₀ - 14 | 5.3 ₁₀ - 15 |
| | <i>d1</i> | 6.4 ₁₀ - 08 | 1.8 ₁₀ - 11 | 3.6 ₁₀ - 14 | 5.3 ₁₀ - 15 |
| | (<i>n, N</i>) | (9, 17) | (11, 37) | (13, 64) | (14, 87) |
| 0.2 | <i>d0</i> | 4.8 ₁₀ - 08 | 5.3 ₁₀ - 11 | 5.0 ₁₀ - 14 | 1.8 ₁₀ - 15 |
| | <i>d1</i> | 9.4 ₁₀ - 08 | 4.9 ₁₀ - 11 | 2.2 ₁₀ - 14 | 1.3 ₁₀ - 14 |
| | (<i>n, N</i>) | (9, 17) | (11, 36) | (13, 63) | (14, 86) |
| 0.4 | <i>d0</i> | 6.8 ₁₀ - 09 | 2.2 ₁₀ - 11 | 2.1 ₁₀ - 14 | 8.9 ₁₀ - 15 |
| | <i>d1</i> | 2.3 ₁₀ - 08 | 1.1 ₁₀ - 10 | 2.5 ₁₀ - 14 | 2.3 ₁₀ - 14 |
| | (<i>n, N</i>) | (10, 15) | (11, 33) | (13, 59) | (14, 81) |
| 0.6 | <i>d0</i> | 2.0 ₁₀ - 07 | 8.2 ₁₀ - 12 | 3.4 ₁₀ - 14 | 1.6 ₁₀ - 14 |
| | <i>d1</i> | 9.9 ₁₀ - 08 | 4.8 ₁₀ - 11 | 1.6 ₁₀ - 14 | 2.4 ₁₀ - 14 |
| | (<i>n, N</i>) | (8, 15) | (11, 33) | (13, 59) | (14, 81) |
| 0.8 | <i>d0</i> | 3.5 ₁₀ - 08 | 4.7 ₁₀ - 12 | 4.1 ₁₀ - 14 | 1.1 ₁₀ - 14 |
| | <i>d1</i> | 5.7 ₁₀ - 08 | 4.7 ₁₀ - 11 | 0.0 ₁₀ + 00 | 2.1 ₁₀ - 14 |
| | (<i>n, N</i>) | (9, 17) | (11, 36) | (13, 63) | (14, 86) |
| 1.0 | <i>d0</i> | 6.4 ₁₀ - 08 | 1.8 ₁₀ - 11 | 3.2 ₁₀ - 14 | 3.6 ₁₀ - 15 |
| | <i>d1</i> | 9.5 ₁₀ - 08 | 5.5 ₁₀ - 11 | 7.1 ₁₀ - 15 | 1.4 ₁₀ - 14 |
| | (<i>n, N</i>) | (9, 17) | (11, 37) | (13, 64) | (14, 87) |

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procedure bessya(a,x,eps,ya,ya1); value a,x,eps; real a,x,eps,ya,ya1;
begin real b,c,d,e,f,g,h,p,pi,q,r,s; integer n,na; Boolean rec, rev;
  pi:= 4 × arctan(1); na:= entier(a+5); rec:= a ≥ .5;
  rev:= a < -5; if rev ∨ rec then a:= a-na;
  if a = -5 then
  begin p:= sqrt(2/pi/x); f:= p × sin(x); g:= -p × cos(x) end else
  if x < 3 then
  begin b:= x/2; d:= -ln(b); e:= a × d;
    c:= if abs(a) < 10-15 then 1/pi else a/sin(a × pi);
    s:= if abs(e) < 10-15 then 1 else sinh(e)/e;
    e:= exp(e); g:= recip gamma(a, p, q) × e; e:= (e + 1/e)/2;
    f:= 2 × c × (p × e + q × s × d); e:= a × a;
    p:= g × c; q:= 1/g/pi; c:= a × pi/2;
    r:= if abs(c) < 10-15 then 1 else sin(c)/c; r:= pi × c × r × r;
    c:= 1; d:= -b × b; ya:= f + r × q; ya1:= p;
    for n:= 1, n + 1 while
    abs(g/(1 + abs(ya))) + abs(h/(1 + abs(ya1))) > eps do
    begin f:= (f × n + p + q)/(n × n - e); c:= c × d/n;
      p:= p/(n - a); q:= q/(n + a);
      g:= c × (f + r × q); h:= c × p - n × g;
      ya:= ya + g; ya1:= ya1 + h
    end;
    f:= -ya; g:= -ya1/b
  end else
  begin b:= x - pi × (a + .5)/2; c:= cos(b); s:= sin(b);
    d:= sqrt(2/x/pi);
    besspqa(a,x,eps,p,q,b,h);
    f:= d × (p × s + q × c); g:= d × (h × s - b × c)
  end;
  if rev then
  begin x:= 2/x; na:= -na - 1;
    for n:= 0 step 1 until na do
    begin h:= x × (a - n) × f - g; g:= f; f:= h end
  end else if rec then
  begin x:= 2/x;
    for n:= 1 step 1 until na do
    begin h:= x × (a + n) × g - f; f:= g; g:= h end
  end;
  ya:= f; ya1:= g
end bessya;

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procedure besspqa(a,x,eps,pa,qa,pa1,qa1); value a,x,eps;
    real a,x,eps,pa,qa,pa1,qa1;
begin real b,c,d,e,f,g,p,p0,q,q0,r,s; integer n,na; Boolean rec,rev;
    rev := a < -.5; if rev then a := -a-1;
    rec := a ≥ .5; if rec then
    begin na := entier(a+.5); a := a - na end;
    if a = -.5 then
    begin pa := pa1 := 1; qa := qa1 := 0 end else
    begin c := .25 - a × a; b := x + x; p := 4 × arctan(1);
    e := (x × cos(a × p)/p/eps)↑2; p := 1; q := -x; r := s := 1 + x × x;
    for n := 2, n + 1 while r × n × n < e do
    begin d := (n - 1 + c/n)/s; p := (2 × n - p × d)/(n + 1);
    q := (-b + q × d)/(n + 1); s := p × p + q × q; r := r × s
    end;
    f := p := p/s; g := q := -q/s;
    for n := n, n - 1 while n > 0 do
    begin r := (n+1) × (2-p) - 2; s := b + (n+1) × q; d := (n - 1 + c/n)/
    (r × r + s × s); p := d × r; q := d × s; e := f;
    f := p × (e + 1) - g × q; g := q × (e + 1) + p × g
    end;
    f := 1 + f; d := f × f + g × g;
    pa := f/d; qa := -g/d; d := a + .5 - p; q := q + x;
    pa1 := (pa × q - qa × d)/x;
    qa1 := (qa × q + pa × d)/x
    end;
end;
if rec then
begin x := 2/x; b := (a + 1) × x;
    for n := 1 step 1 until na do
    begin p0 := pa - qa1 × b; q0 := qa + pa1 × b;
    pa := pa1; pa1 := p0; qa := qa1; qa1 := q0; b := b + x
    end
end;
if rev then
begin p0 := pa1; pa1 := pa; pa := p0;
    q0 := qa1; qa1 := qa; qa := q0
end
end
end besspqa;

```

REFERENCES

1. W. GAUTSCHI, *SIAM Rev.* **9** (1967), 24–82.
2. W. GAUTSCHI, *Comm. ACM* **7** (1964), 479–480.
3. M. GOLDSTEIN AND R. M. THALER, *Math. Tables Aids Comput.* **13** (1959), 102–108.
4. Y. L. LUKE, "The special functions and their approximations," Vols. 1 and 2, Academic Press, New York and London, 1969.
5. Y. L. LUKE, *Math. Comp.* **26** (1972), 237–240.
6. N. M. TEMME, *J. Comput. Phys.* **19** (1975).

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