

**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  $Y_\nu$  and  $Y_{\nu+1}$  are given. In the forward direction the recurrence formula (1.3) for  $Y_\nu$  is numerically stable, whereas it is unstable for  $J_\nu$  (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  $J_\nu$  and  $Y_\nu$  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  $Y_\nu$  and  $Y_{\nu+1}$  and we use the methods of our previous paper on the computation of  $K_\nu$  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  $Y_\nu$  and  $Y_{\nu+1}$  and these results follow immediately from [6] by using (1.8).

For the computation of  $J_\nu$  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  $J_\nu$  and  $Y_\nu$  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  $Y_\nu, J_\nu$  for  $z \geq 5$ . In Goldstein and Thaler [3] the computation of  $Y_\nu$  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

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

Considering (2.1) with  $\nu$  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)

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

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  $K_\nu$  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  $\nu$  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 <sub>10</sub> - 14	7.1 <sub>10</sub> - 15	0.0 <sub>10</sub> + 00
1.0	0.0 <sub>10</sub> + 00	7.1 <sub>10</sub> - 15	7.1 <sub>10</sub> - 15
2.0	7.1 <sub>10</sub> - 15	2.8 <sub>10</sub> - 14	7.1 <sub>10</sub> - 15
3.0	7.1 <sub>10</sub> - 15	0.0 <sub>10</sub> + 00	0.0 <sub>10</sub> + 00
5.0	7.1 <sub>10</sub> - 15	1.4 <sub>10</sub> - 14	0.0 <sub>10</sub> + 00
7.0	7.1 <sub>10</sub> - 15	7.1 <sub>10</sub> - 15	1.4 <sub>10</sub> - 14
10.0	7.1 <sub>10</sub> - 15	7.1 <sub>10</sub> - 15	7.1 <sub>10</sub> - 15
20.0	0.0 <sub>10</sub> + 00	7.1 <sub>10</sub> - 15	0.0 <sub>10</sub> + 00
50.0	2.1 <sub>10</sub> - 14	1.4 <sub>10</sub> - 14	0.0 <sub>10</sub> + 00
100.0	2.1 <sub>10</sub> - 14	7.1 <sub>10</sub> - 15	7.1 <sub>10</sub> - 15

TABLE II

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

```

procedure bessya(a,x,eps,ya,yal); value a,x,eps; real a,x,eps,ya,yal;
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; yal := p;
    for n := 1, n + 1 while
    abs(g/(1 + abs(ya))) + abs(h/(1 + abs(yal))) > 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; yal := yal + h
    end;
    f := -ya; g := -yal/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; yal := g
end bessya;

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```

procedure besspqa(a,x,eps,pa,qa,pal,qal); value a,x,eps;
    real a,x,eps,pa,qa,pal,qal;
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 := pal := 1; qa := qal := 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;
        pal := (pa × q - qa × d)/x;
        qal := (qa × q + pa × d)/x
    end;
    if rec then
    begin x := 2/x; b := (a + 1) × x;
        for n := 1 step 1 until na do
        begin p0 := pa - qal × b; q0 := qa + pal × b;
            pa := pal; pal := p0; qa := qal, qal := q0; b := b + x
        end
    end;
    if rev then
    begin p0 := pal; pal := pa; pa := p0;
        q0 := qal; qal := qa; qa := q0
    end
end besspqa;

```

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