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On the Localization and Computation of Zeros of Bessel Functions

The topological degree of a continuous mapping is implemented for the calculation of the total number of the simple real zeros within any interval of the Bessel functions of first and second kind and their derivatives. A new algorithm, based on this implementation, is given for the localization and isolation of these zeros. Furthermore, a second algorithm is presented for their computation employing a modified bisection method. The only information required for this computation is the algebraic signs of function values. Moreover, lower and upper bounds of a zero can also be obtained.

MSC (1991): 33C10, 65D20, 34B30

1. Introduction

The study of the Helmholtz equation in cylindrical coordinates eventually leads to the well-known *Bessel equation*

$$x^2 u''(x) + x u'(x) + (x^2 - \nu^2) u(x) = 0, \tag{1}$$

the solutions of which are

a) the *Bessel function of order ν and argument x of the first kind* given by the series

$$J_\nu(x) = \frac{x^\nu}{2^\nu} \sum_{m=0}^{\infty} \frac{(-1)^m x^{2m}}{2^{2m} m! \Gamma(\nu + m + 1)}, \tag{2}$$

which converges for every x and

b) the *Bessel function of order ν and argument x of the second kind* expressible in terms of functions of the first kind by the formula

$$Y_\nu(x) = \begin{cases} \frac{J_\nu(x) \cos \nu\pi - J_{-\nu}(x)}{\sin \nu\pi}, & \nu \text{ non-integral,} \\ \lim_{n \rightarrow \nu} \frac{J_n(x) \cos n\pi - J_{-n}(x)}{\sin n\pi}, & \nu = 0, 1, 2, \dots \end{cases} \tag{3}$$

Regarding the zeros $j_\nu, j'_\nu, y_\nu,$ and y'_ν of the functions $J_\nu(x), J'_\nu(x), Y_\nu(x),$ and $Y'_\nu(x),$ respectively, the following holds (cf., e.g., [5, 28]): *For any real value of ν these functions have infinitely many real zeros all of which are simple, with the possible exception of $x = 0$. Especially, for $\nu \geq -1, J_\nu(x)$ has only real zeros. If $\nu \geq 0,$ all zeros of $J'_\nu(x)$ are real.*

For non-negative $\nu,$ the k -th positive zeros of the above functions are denoted correspondingly by $j_{\nu,k}, j'_{\nu,k}, y_{\nu,k},$ and $y'_{\nu,k}.$ They are interlaced according to the inequalities

$$\begin{aligned} j_{\nu,1} &< j_{\nu+1,1} < j'_{\nu,2} < j_{\nu+1,2} < j_{\nu,3} < \dots, \\ y_{\nu,1} &< y_{\nu+1,1} < y_{\nu,2} < y_{\nu+1,2} < y_{\nu,3} < \dots, \\ \nu \leq j'_{\nu,1} &< y_{\nu,1} < y'_{\nu,1} < j_{\nu,1} < j'_{\nu,2} < y_{\nu,2} < y'_{\nu,2} < j_{\nu,2} < j'_{\nu,3} < \dots \end{aligned}$$

The mathematical problem of localizing and computing zeros of Bessel functions, encountered in many fields in science and engineering, has drawn a lot of attention and has evolved to a rather specialized branch of mathematics.

In the present paper we implement the concept of the topological degree to calculate the total number of real roots of Bessel functions within a predetermined interval and to isolate each one of them. For this purpose we use PICARD's extension [16, 17, 7, 8, 21] and either Kronecker theory or KEARFOTT's degree computation method. Once a zero is isolated, it can be computed numerically, utilizing a modified bisection method, to any accuracy (subject to relative machine precision). Thus, we propose two algorithms, one for the isolation and one for the computation of a real zero of a Bessel function. Upper and lower bounds for any root can also be produced.

2. The topological degree for the localization of zeros

Definition 1: Suppose that the function $F_n = (f_1, \dots, f_n) : \mathcal{D}^n \subset \mathbb{R}^n \rightarrow \mathbb{R}^n$ is twice continuously differentiable in the domain \mathcal{D}^n the boundary of which is denoted by $b(\mathcal{D}^n).$ Suppose further that the solutions of the equation

$$F_n(x) = \mathcal{O}_n, \tag{4}$$

where $\mathcal{O}_n = (0, \dots, 0)$ denotes the origin of \mathbb{R}^n , are not located on $b(\mathcal{D}^n)$, and they are simple, i.e., the Jacobian determinant of F_n at these solutions is non-zero. Then the *topological degree of F_n at \mathcal{O}_n relative to \mathcal{D}^n* is denoted by $\text{deg}[F_n, \mathcal{D}^n, \mathcal{O}_n]$ and can be defined by the following sum:

$$\text{deg}[F_n, \mathcal{D}^n, \mathcal{O}_n] = \sum_{x \in F_n^{-1}(\mathcal{O}_n)} \text{sgn } J_{F_n}(x), \tag{5}$$

where J_{F_n} indicates the determinant of the Jacobian matrix, and sgn defines the well known sign function with values

$$\text{sgn } \psi = \begin{cases} -1, & \text{if } \psi < 0, \\ 0, & \text{if } \psi = 0, \\ 1, & \text{if } \psi > 0. \end{cases}$$

The above definition can be generalized when F_n is only continuous [6, 14].

Kronecker's theorem [2, 6, 14] states that equation (4) has at least one root in \mathcal{D}^n if $\text{deg}[F_n, \mathcal{D}^n, \mathcal{O}_n] \neq 0$.

The definition of the topological degree actually indicates that its value is equal to the number of simple solutions of equation (4) for which the Jacobian determinant is positive, minus the number of simple solutions for which the Jacobian is negative. Evidently, if all of them give the same Jacobian sign, then the total number \mathcal{N}^r of simple roots of $F_n(x)$ can be obtained by the value of $\text{deg}[F_n, \mathcal{D}^n, \mathcal{O}_n]$. To this end PICARD has considered the following extensions of the function F_n and the domain \mathcal{D}^n :

$$F_{n+1} = (f_1, \dots, f_n, f_{n+1}) : \mathcal{D}^{n+1} \subset \mathbb{R}^{n+1} \rightarrow \mathbb{R}^{n+1}, \tag{6}$$

where $f_{n+1} = yJ_{F_n}$ and \mathcal{D}^{n+1} is the direct product of the domain \mathcal{D}^n with an arbitrary interval of the real y -axis containing the point $y = 0$. Then the following system of equations,

$$\begin{aligned} f_i(x_1, x_2, \dots, x_n) &= 0, & i &= 1, \dots, n, \\ yJ_{F_n}(x_1, x_2, \dots, x_n) &= 0, \end{aligned} \tag{7}$$

possesses the same simple roots with $F_n(x)$, provided $y = 0$. Also, it is easily seen that the Jacobian of (7) is equal to $J_{F_n(x)}^2$ which is always positive. Thus we conclude that the total number \mathcal{N}^r of solutions of equation (4) is

$$\mathcal{N}^r = \text{deg}[F_{n+1}, \mathcal{D}^{n+1}, \mathcal{O}_{n+1}]. \tag{8}$$

2.1 Kronecker integral approach

The topological degree can be represented by the Kronecker integral as follows:

$$\text{deg}[F_n, \mathcal{D}^n, \mathcal{O}_n] = \frac{1}{\Omega_n} \int \int_{b(\mathcal{D}^n)} \dots \int \frac{\sum_{i=1}^n A_i dx_1 \dots dx_{i-1} dx_{i+1} \dots dx_n}{(f_1^2 + f_2^2 + \dots + f_n^2)^{n/2}}, \tag{9}$$

where A_i define the following determinants:

$$A_i = (-1)^{n(i-1)} \begin{vmatrix} f_1 & \frac{\partial f_1}{\partial x_1} & \dots & \frac{\partial f_1}{\partial x_{i-1}} & \frac{\partial f_1}{\partial x_{i+1}} & \dots & \frac{\partial f_1}{\partial x_n} \\ f_2 & \frac{\partial f_2}{\partial x_1} & \dots & \frac{\partial f_2}{\partial x_{i-1}} & \frac{\partial f_2}{\partial x_{i+1}} & \dots & \frac{\partial f_2}{\partial x_n} \\ \vdots & \vdots & \ddots & \vdots & \vdots & \ddots & \vdots \\ f_n & \frac{\partial f_n}{\partial x_1} & \dots & \frac{\partial f_n}{\partial x_{i-1}} & \frac{\partial f_n}{\partial x_{i+1}} & \dots & \frac{\partial f_n}{\partial x_n} \end{vmatrix},$$

and Ω_n denotes the surface of a hypersphere in \mathbb{R}^n with radius one, i.e.: $\Omega_n = 2\pi^{n/2}/\Gamma(n/2)$.

The Kronecker integral has been numerically approximated with Gauss-Legendre quadrature [13].

In the present paper we study the zeros of a Bessel function, so we focus on the problem of calculating the total number of simple roots of a real function $f(x)$, defined in a predetermined interval $[a, b]$ and twice continuously differentiable in (a, b) , where a and b are arbitrarily chosen so that $f(a)f(b) \neq 0$.

According to Picard's extension we consider the function $F_2 = (f_1, f_2) : \mathcal{D}^2 \subset \mathbb{R}^2 \rightarrow \mathbb{R}^2$ and the corresponding system

$$f_1(x, y) = f(x) = 0, \quad f_2(x, y) = yf'(x) = 0, \tag{10}$$

where the prime denotes differentiation, and \mathcal{D}^2 is the rectangular parallelepiped $[a, b] \times [-\xi, \xi]$ in the (x, y) -plane with ξ an arbitrary positive constant. Since the roots are simple, which means $f'(x) \neq 0$ for $x \in f^{-1}(0)$, it is easily seen that the solutions of system (10) in \mathcal{D}^2 , and those of $f(x) = 0$ in (a, b) are the same. Also, since $J_{F_2} = f'^2$, the total

number of simple zeros \mathcal{N}^r of $f(x)$ in (a, b) is given by

$$\mathcal{N}^r = \text{deg} [F_2, \mathcal{P}^2, \mathcal{O}_2]. \tag{11}$$

For the computation of the topological degree of F_2 we apply Kronecker's integral (9) for $n = 2$. Using the relations

$$df_j = \frac{\partial f_j}{\partial x_1} dx_1 + \frac{\partial f_j}{\partial x_2} dx_2, \quad j = 1, 2, \text{ we obtain}$$

$$\mathcal{N}^r = \frac{1}{2\pi} \oint_{b(\mathcal{P}^2)} \frac{f_1 df_2 - f_2 df_1}{f_1^2 + f_2^2} = \frac{1}{2\pi} \oint_{b(\mathcal{P}^2)} d \arctan \left(\frac{f_2}{f_1} \right). \tag{12}$$

Replacing f_1 and f_2 by virtue of (10) and performing the integration in (12) we finally get

$$\mathcal{N}^r = -\frac{1}{\pi} \left[\xi \int_a^b \frac{f(x) f''(x) - f'^2(x)}{f^2(x) + \xi^2 f'^2(x)} dx - \arctan \left(\frac{\xi f'(b)}{f(b)} \right) + \arctan \left(\frac{\xi f'(a)}{f(a)} \right) \right]. \tag{13}$$

Remark 1: It has been explicitly shown by PICARD [16, 17] that relation (13) is independent of the value of ξ .

The above developed method is applied in the sequel for the localization of the simple real zeros of the Bessel functions of the first kind, $J_\nu(x)$. In this case relation (13) becomes

$$\mathcal{N}^r = -\frac{1}{\pi} \left[\xi \int_a^b G(x) dx - \arctan \left(\frac{\xi J'_\nu(b)}{J_\nu(b)} \right) + \arctan \left(\frac{\xi J'_\nu(a)}{J_\nu(a)} \right) \right], \tag{14}$$

where the function $G(x)$, by means of equation (1), can be expressed in the following form:

$$G(x) = \frac{\left(\frac{\nu^2}{x^2} - 1 \right) J_\nu^2(x) - \frac{1}{x} J_\nu(x) J'_\nu(x) - J_\nu'^2(x)}{J_\nu^2(x) + \xi^2 J_\nu'^2(x)}. \tag{15}$$

2.2 KEARFOTT's approach

We could use any one of the degree computation methods (see e.g. [22, 13, 10, 11, 23, 24]) to determine the total number of zeros by virtue of equation (11). Here we use KEARFOTT's method [10, 11, 12] which compares favorably to other methods, in efficiency. This method is briefly described below.

Suppose that $S^{n-1} = \langle x_1, x_2, \dots, x_n \rangle$ is an $(n-1)$ -simplex [22, 10, 11] in \mathbb{R}^n and assume $F_n = (f_1, f_2, \dots, f_n) : S^{n-1} \rightarrow \mathbb{R}^n$ is continuous. Then the *range simplex associated with S^{n-1} and F_n* , denoted by $\mathcal{R}(S^{n-1}, F_n)$, is an $n \times n$ matrix with elements ϱ_{ij} , $1 \leq i, j \leq n$, given by

$$\varrho_{ij} = \begin{cases} 1, & \text{if } f_j(x_i) \geq 0, \\ -1, & \text{if } f_j(x_i) < 0. \end{cases} \tag{16}$$

$\mathcal{R}(S^{n-1}, F_n)$ is called *usable* if one of the following conditions hold:

a) the elements ϱ_{ij} of $\mathcal{R}(S^{n-1}, F_n)$, are:

$$\varrho_{ij} = \begin{cases} 1, & \text{if } i \geq j, \\ -1, & \text{if } j = i + 1. \end{cases} \tag{17}$$

b) $\mathcal{R}(S^{n-1}, F_n)$ can be put into this form by a permutation of its rows.

When $\mathcal{R}(S^{n-1}, F_n)$ is usable, then the *parity* $\text{Par}(\mathcal{R}(S^{n-1}, F_n))$ is defined to be 1, if the number of the permutations of the rows required to put $\mathcal{R}(S^{n-1}, F_n)$ into the form (17) is even. If this number is odd then $\text{Par}(\mathcal{R}(S^{n-1}, F_n))$ is defined to be -1 . For all other cases, we set $\text{Par}(\mathcal{R}(S^{n-1}, F_n)) = 0$. Suppose that \mathcal{P}^n is an n -dimensional polyhedron for some $n \geq 2$, and that $\{S_i^{n-1}\}_{i=1}^m$ is a finite set of $(n-1)$ -simplexes with disjoint interiors such that $\sum_{i=1}^m S_i^{n-1} = b(\mathcal{P}^n)$; then, under some assumptions regarding S_i^{n-1} , the value of the topological degree of F_n at \mathcal{O}_n relative to \mathcal{P}^n can be obtained by the following relation:

$$\text{deg} [F_n, \mathcal{P}^n, \mathcal{O}_n] = \sum_{i=1}^m \text{Par}(\mathcal{R}(S_i^{n-1}, F_n)). \tag{18}$$

Remark 2: Kearfott's degree computation method is very efficient and has the advantage that it requires only the signs of function values to be correct.

2.3 Isolating a zero

With the above discussion in mind we give a description of our algorithm, in “pseudo-Pascal” (see e.g. [15]), independently of the degree computation method. In this algorithm $F_\nu(x)$ indicates the considered Bessel function of order ν , and (a, b) is a predetermined interval. It results in the total number \mathcal{N}^r of roots which exist within (a, b) . Besides, it isolates one of them, bisecting the initial interval, and gives a lower bound a_k and an upper one, b_k , for this root. If the isolation of the rest of them is required, the whole procedure has to be repeated successively for the remaining intervals (a, a_k) and (b_k, b) .

Algorithm *degree_isolate* (a, b, \mathcal{N}^r) ;

{**comment**: This algorithm results in the total number \mathcal{N}^r of roots of $F_\nu(x)$ in (a, b) and isolates one of them}.

procedure degree $(a_k, b_k, \mathcal{N}_k^r)$; {**comment**: computes the total number of zeros}

begin

Find \mathcal{N}_k^r , the number of zeros in (a_k, b_k) ;

end {degree}

procedure isolation $(a_k, b_k, \mathcal{N}_k^r)$; {**comment**: isolates a zero of $F_\nu(x)$ in (a_k, b_k) }

begin

while $\mathcal{N}_k^r > 1$ **do**

begin

degree $\left(a_k, \frac{a_k + b_k}{2}, \mathcal{N}_1^r\right)$;

$\mathcal{N}_2^r := \mathcal{N}_k^r - \mathcal{N}_1^r$;

if $\mathcal{N}_2^r = 0$ **then** isolation $\left(a_k, \frac{a_k + b_k}{2}, \mathcal{N}_1^r\right)$;

if $\mathcal{N}_1^r < \mathcal{N}_2^r$ **and** $\mathcal{N}_1^r \neq 0$ **then** isolation $\left(a_k, \frac{a_k + b_k}{2}, \mathcal{N}_1^r\right)$

else isolation $\left(\frac{a_k + b_k}{2}, b_k, \mathcal{N}_2^r\right)$;

end {while}

end {isolation}

begin {degree_isolate}

input a, b

degree (a, b, \mathcal{N}^r) ;

$\mathcal{N}_*^r := \mathcal{N}^r$;

if $\mathcal{N}^r > 1$ **then** isolation (a, b, \mathcal{N}^r) ;

output a, b, \mathcal{N}_*^r

end. {degree_isolate}

3. Computing roots of Bessel functions

Having isolated one root of a Bessel function within an interval, we can use a modified version of the bisection method to compute it, as described in [25, 26]. It is reported there that, in order to compute a solution of $f(x) = 0$, where $f: [a, b] \subset \mathbb{R} \rightarrow \mathbb{R}$ is continuous, the following iterative formula can be used:

$$x_{i+1} = x_i + c \operatorname{sgn} f(x_i) / 2^{i+1}, \quad i = 0, 1, \dots, \quad (19)$$

with $x_0 = a$ and $c = \operatorname{sgn} f(a) (b - a)$. The iterations (19) converge to a root $r \in (a, b)$ if for some $x_i, i = 1, 2, \dots$, there holds

$$\operatorname{sgn} f(x_0) \operatorname{sgn} f(x_i) = -1.$$

The number of iterations, η , which are required in obtaining an approximate root r^* such that $|r - r^*| \leq \varepsilon$ for some $\varepsilon \in (0, 1)$, is given by

$$\eta = \lceil \log_2 ((b - a) \varepsilon^{-1}) \rceil, \quad (20)$$

where the notation $\lceil \cdot \rceil$ refers to the smallest integer not less than the real number quoted.

It is evident from (19) that the only computable information required by the bisection method consists in the algebraic signs of the function f ; so it can be applied to problems with imprecise function values. Moreover, the bisection method is a globally convergent method, it always converges within the given interval, and it is optimal [19, 20] in the sense that it possesses asymptotically the best rate of convergence. Also, it can be efficiently implemented for the computation of all the zeros and extrema of a function [9].

To apply our method, as previously mentioned, we only need the algebraic signs of the function values to be correct. The following results suggest the number of terms necessary in order to determine the algebraic signs of $J_\nu(x)$.

Proposition 1: *The sign of the Bessel function J_ν for $\nu > -1$ is the same as the sign of the sum*

$$L_\nu(x) = \frac{x^\nu}{2^\nu} \sum_{m=0}^{M+K+1} (-1)^m a_m, \quad (21)$$

where a_m are given by

$$a_m = \frac{x^{2m}}{2^{2m}m! \Gamma(\nu + m + 1)}, \quad (22)$$

with

$$M = \left\lceil \frac{-\nu - 2 + \sqrt{\nu^2 + x^2}}{2} \right\rceil, \quad (23)$$

and K is defined so that the following relation holds:

$$|a_{M+K+1}| < \left| \sum_{k=0}^{M+K} (-1)^k a_k \right|. \quad (24)$$

Proof: We observe that

- a) $a_m > 0$,
- b) $\lim_{m \rightarrow \infty} a_m = 0$,
- c) $a_m \geq a_{m+1} \Leftrightarrow 4(m+1)(\nu+m+1) > x^2$.

Thus, the series

$$R_\nu(x) = \sum_{m=M}^{\infty} \frac{(-1)^m x^{2m}}{2^{2m}m! \Gamma(\nu + m + 1)} \quad (25)$$

is an alternating series which satisfies the three conditions of Leibniz's Theorem provided that

$$m > \frac{-\nu - 2 + \sqrt{\nu^2 + x^2}}{2}. \quad (26)$$

The Alternating Series Estimation Theorem ensures that the absolute value of the $(M + K + 1)$ -th term is larger than that of the remainder of the series $R_\nu(x)$. Furthermore, the signs of the $(M + K + 1)$ -th term and of the remainder of the series $R_\nu(x)$ are the same. Now, by assumption (24) (which is fulfilled for some term a_i since the sequence $\{a_i\}$ tends to zero when $i \rightarrow \infty$), the sign of the Bessel function $J_\nu(x)$ is the same as the corresponding sign of $L_\nu(x)$. Thus the proposition is proved.

We can now give a description of our algorithm that computes, within a predetermined accuracy ε , a real root of the Bessel function $J_\nu(x)$, which has been isolated in the interval (a_k, b_k) .

Algorithm *compute_zero* (a, b);

{**comment:** This algorithm computes a zero of $J_\nu(x)$ in (a, b) . It uses (19) and requires $F(x), \varepsilon$ }.

procedure sign (x, s); {**comment:** finds the sign s of $J_\nu(x)$ }

begin

$M := \lceil (-\nu - 2 + \sqrt{\nu^2 + x^2})/2 \rceil$;

$S := 0$;

for $m := 0$ **to** M **do** $S := S + (-1)^m x^{2m+\nu} / (2^{2m+\nu} m! \Gamma(\nu + m + 1))$;

$m := M + 1$;

$T := (-1)^m x^{2m+\nu} / (2^{2m+\nu} m! \Gamma(\nu + m + 1))$;

while $|T| \geq |S|$ **do**

begin

$S := S + T$;

$m := m + 1$;

$T := (-1)^m x^{2m+\nu} / (2^{2m+\nu} m! \Gamma(\nu + m + 1))$;

end {while}

Find $s := \text{sign}(S + T)$;

end {sign}

begin {compute_zero}

input a_k, b_k, ε

$x_0 := a_k$;

sign (x_0, s_0);

$c := s_0(b_k - a_k)$;

for $i := 0$ **to** $\lceil \log_2((b_k - a_k) \varepsilon^{-1}) \rceil$ **do**

begin

sign (x_i, s_i);

$x_{i+1} := x_i + s_i c / 2^{i+1}$;

end

output x_{i+1}

end. {compute_zero}

Remark 3: In order to obtain zeros of other Bessel functions, the above algorithm has to be adapted accordingly.

Evidently, utilizing the above results we are able to produce bounds for any real zero of a Bessel function. The following proposition proves it.

Proposition 2: Suppose that $f : [a, b] \subset \mathbb{R} \rightarrow \mathbb{R}$ is a continuous function and (a, b) contains only one simple zero x^* of f . Consider the sequence (19) and its subsequences $\{x_i^-\}$, $\{x_i^+\}$ such that $\text{sgn } f(x_i^-) = \text{sgn } f(a)$, $\text{sgn } f(x_i^+) = -\text{sgn } f(a) = \text{sgn } f(b)$. Then, a_η and b_η , determined so that

$$a_\eta = \max_{0 \leq i \leq \eta} \{x_i^-\}, \quad b_\eta = \min_{0 \leq i \leq \eta} \{x_i^+\}, \tag{27}$$

where η is given by equation (20), for any $\varepsilon \in (0, 1)$, constitute a lower and an upper bound of the zero x^* with $b_\eta - a_\eta < \varepsilon$.

Remark 4: The function f can be replaced by any of the Bessel functions considered in this paper.

4. Applications

We have tested our algorithms with several random intervals (a, b) and various Bessel functions. In each case we have calculated the total number \mathcal{N}^r of roots of the respective Bessel function existing within (a, b) . For this calculation we have used both Kronecker's integral applied to Picard's extension and Kearfott's method. The computation of the integral of equation (14) depends on the integration method used. We have tried various integration methods. Here we exhibit the results obtained by Romberg's integration method. Also, we have observed that the speed of the numerical computation of the integral in equation (14) depends on the value of ξ . Our experience is that the fastest computation (using Romberg's method) can be obtained for $0.8 \leq \xi \leq 1.2$ while for other values the computational time increases (see Fig. 1). Kearfott's method is independent of the choice of ξ , but its speed depends on a stopping parameter p and the maximum tree depth MD [11]. In our case, we have obtained accurate results even for small values of p and MD, as for instance $p = 1$, MD = 2.

Table 1 presents the number of zeros of Bessel functions of various orders ν existing within some given intervals (a, b) , as well as the respective subintervals (a_k, b_k) , where exactly one root, r_k , exists.

In Table 2 we give the first ten zeros of several Bessel functions, chosen at random, computed by Algorithm `compute_zero`.

Table 1

| | a | b | \mathcal{N}^r | a_k | b_k | r_k |
|-------------|-----|-----|-----------------|---------|----------|--------------------|
| J_0 | 1 | 100 | 32 | 1.0000 | 4.09375 | 2.404825557695773 |
| $J_{0.3}$ | 1 | 100 | 31 | 81.4375 | 87.62500 | 84.509788949453324 |
| $J_{2.5}$ | 1 | 100 | 30 | 1.0000 | 7.18750 | 5.763459196894550 |
| J_{10} | 10 | 100 | 27 | 10.0000 | 15.62500 | 14.475500686554542 |
| $J_{13.3}$ | 10 | 100 | 25 | 10.0000 | 21.25000 | 18.131465204981082 |
| J_{50} | 50 | 100 | 11 | 50.0000 | 62.50000 | 57.116899160119190 |
| J'_0 | 1 | 100 | 31 | 13.3750 | 19.56250 | 16.500922441528084 |
| $J'_{0.3}$ | 1 | 100 | 32 | 1.0000 | 4.09375 | 1.308699363719847 |
| $J'_{2.5}$ | 1 | 100 | 31 | 1.0000 | 7.18750 | 3.959527916501095 |
| J'_{10} | 10 | 100 | 27 | 10.0000 | 15.62500 | 12.128927704415439 |
| $J'_{13.3}$ | 10 | 100 | 26 | 10.0000 | 15.62500 | 15.617873137336602 |
| J'_{50} | 50 | 100 | 11 | 50.0000 | 56.25000 | 53.502858820400364 |
| Y_0 | 1 | 100 | 31 | 13.3750 | 19.56250 | 16.470630050877633 |
| $Y_{0.3}$ | 1 | 100 | 31 | 1.0000 | 7.18750 | 4.301991992307308 |
| $Y_{2.5}$ | 1 | 100 | 31 | 1.0000 | 7.18750 | 3.632797319831763 |
| Y_{10} | 10 | 100 | 27 | 10.0000 | 15.62500 | 11.770876674955581 |
| $Y_{13.3}$ | 10 | 100 | 26 | 10.0000 | 15.62500 | 15.242738650014224 |
| Y_{50} | 50 | 100 | 11 | 50.0000 | 56.25000 | 52.997640387316651 |
| Y'_0 | 1 | 100 | 32 | 1.0000 | 4.09375 | 2.197141326031016 |
| $Y'_{0.3}$ | 1 | 100 | 31 | 81.4375 | 87.62500 | 84.503872394936382 |
| $Y'_{2.5}$ | 1 | 100 | 30 | 1.0000 | 7.18750 | 5.634296563929559 |
| Y'_{10} | 10 | 100 | 27 | 10.0000 | 15.62500 | 14.353013743699869 |
| $Y'_{13.3}$ | 10 | 100 | 25 | 10.0000 | 21.25000 | 18.006479815952494 |
| Y'_{50} | 50 | 100 | 11 | 50.0000 | 62.50000 | 56.962904275167517 |

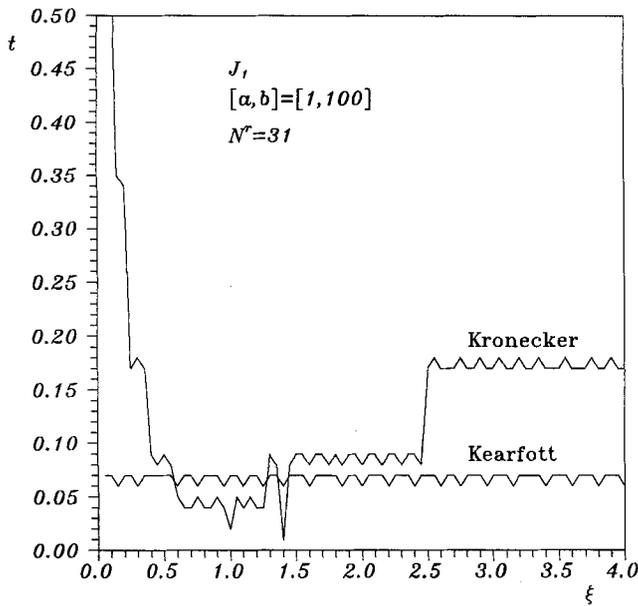


Fig. 1. Time required for the computation of the topological degree using Kronecker's integral versus ξ and comparison with Kearfott's method

Table 2. Zeros of various Bessel functions computed by Algorithm *compute_zero*

| | | | | | |
|--------|--|------------|--|-----------|--|
| J_0 | 2.404825557695773 5.520078110286310 8.653727912911011 11.791534439014282 14.930917708487784 18.071063967910918 21.211636629879254 24.352471530749303 27.493479132040244 30.634606468431967 | $J_{2.5}$ | 5.763459196894550 9.095011330476354 12.322940970566584 15.514603010886749 18.689036355362817 21.853874222709777 25.012803202289602 28.167829707993626 31.320141707447189 34.470488331285007 | J_{50} | 57.116899160119190 62.807698764835380 67.697408410764783 72.190366544011145 76.437072182667940 80.513239317465752 84.463252949061289 88.315711749199524 92.090274456391647 95.801108265953272 |
| J'_0 | 3.957678419314857 7.086051060301773 10.222345043496417 13.361097473872762 16.500922441528084 19.641309700887942 22.782028047291558 25.922957653180932 29.064030252728390 32.205204116493269 | $J'_{2.5}$ | 3.959527916501094 7.451610064214504 10.715647375791513 13.921686012308782 17.103359117208743 20.272369140216535 23.433926142067801 26.590716631086269 29.744270680556565 32.895525188224304 | J'_{50} | 53.502858820400364 60.112444427740558 65.317141149297498 69.981432989702090 74.338747166755810 78.493210918315127 82.501961830679640 86.400289204160094 90.211743342930555 93.952927913983330 |
| Y_0 | 3.831705970207513 7.015586669815619 10.173468135062721 13.323691936314222 16.470630050877633 19.615858510468247 22.760084380592769 25.903672087618386 29.046828534916836 32.189679910974398 | $Y_{2.5}$ | 3.632797319831762 7.367008971566918 10.663561390481999 13.883369775209742 17.072848832681667 20.246944819393926 23.412099866421985 26.571579024683013 29.727222608526206 32.880149822147309 | Y_{50} | 52.997640387316651 60.026319332799442 65.272723327026940 69.951692734372273 74.316349977805003 78.475187454486474 82.486831848982885 86.387212778756663 90.200199617262318 93.942572252611562 |
| Y'_0 | 2.197141326031016 5.429681040794136 8.596005868331170 11.749154830839879 14.897442128336724 18.043402276727858 21.188068934142211 24.331942571356901 27.475294980449235 30.618286491641110 | $Y'_{2.5}$ | 5.634296563929560 9.030901729624809 12.278862551656813 15.480654965172806 18.661308999166464 21.830389858741036 24.992411488536626 28.149798292577922 31.303973776577955 34.455830546321302 | Y'_{50} | 56.962904275167517 62.748881669459305 67.661781418302983 72.164827720771193 76.417109157783180 80.496796665401667 84.449230155149340 88.303453538378357 92.079360103387758 95.791251970104038 |

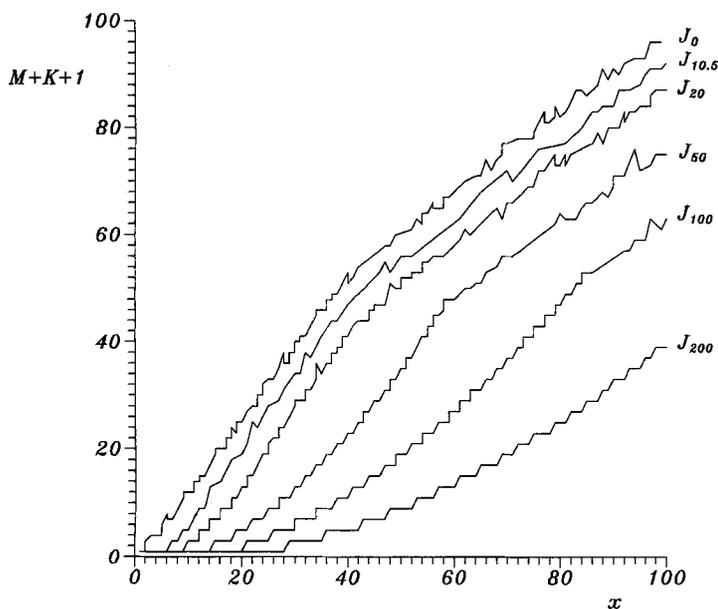


Fig. 2. The number of terms to be added for the determination of the sign of $J_\nu(x)$ versus x for $\nu = 0, 10.5, 20, 50, 100, 200$

By virtue of Proposition 1, we have obtained the number of terms of J_ν required to obtain the corresponding algebraic sign for various values of the order ν and the argument x . As it is shown in Fig. 2, the number of terms decreases as the order increases.

5. Concluding remarks

An efficient method for locating, isolating, and computing real zeros of Bessel functions is described in this paper. It has been implemented and tested, and our experience is that it behaves predictably and accurately.

The first phase of this method (algorithm *degree_isolate*) exploits topological degree theory and especially Picard's extension to calculate the total number of real roots of Bessel functions within a predetermined interval and to isolate one of them. This procedure can be repeated for the isolation of each one of the zeros in this interval.

Once a zero is isolated, the second phase (algorithm *compute_zero*) is applied for its computation to any accuracy (subject to relative machine precision). This algorithm utilizes a modified bisection method. The only computable information required consists in the algebraic signs of the function and, consequently, it is not affected by imprecise function values. Moreover, it always converges rapidly to a zero within the initially specified region independently of the starting guess. It is also a globally convergent method, it can be applied to nondifferentiable continuous functions and does not involve derivatives or approximations of such derivatives. Furthermore, the number of iterations needed to compute a zero to a predetermined accuracy is a priori known.

The rootfinding portion of our method requires the smallest amount of function value information which is its algebraic sign. Using this and the stopping criterion (20), it computes a zero of a Bessel function within a given accuracy.

We have been able to calculate the total number of zeros by computing the value of the topological degree of (11) utilizing efficient methods such as STENGER's [22], STYNES' [23, 24], or KEARFOTT's method [10, 11] (see also [1]), which are based only on sign calculations.

For large argument x , one can use the usual asymptotic expressions for Bessel functions [28]. Moreover, if one would like to compute zeros of a combination of Bessel functions which may not be an alternating series, the van Wijngaarden transformation can be used to convert this new series into an alternating one [18]. Furthermore, very accurate function values for the Bessel functions can be obtained by employing STEED's and TEMME's methods [18], or, alternatively, by means of Coulomb functions, employing BARNETT's procedure [3, 4].

When just one root is required, the isolation portion of our method can be avoided if the function values at the endpoints of the given interval are opposite (see [27] for extensions). This is so because, in this case, the modified bisection method always converges to a zero.

Since the first derivatives of Bessel functions are available, it is at the user's disposal to apply any other rootfinding method, as for instance Newton's method, to accelerate the convergence within the predetermined interval. In such a case, though, unless the starting point is close enough to a zero and certainly away from an extremum of the function, convergence is not ensured. Thus, a few initial steps of our method should be performed so that these conditions are satisfied.

At last, for any given interval (a_k, b_k) containing a single zero of a Bessel function, Proposition 2 can provide a lower and an upper bound for this zero, theoretically as close to it as one desires, but practically as close to it as the accuracy of the calculations permits.

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