1 Purpose

s17dc returns a sequence of values for the Bessel functions \( Y_{\nu+n}(z) \) for complex \( z \), non-negative \( \nu \) and \( n = 0, 1, \ldots, N - 1 \), with an option for exponential scaling.

2 Syntax

\[
[c_y, n_z, \text{ifail}] = \text{s17dc}(\nu, z, n, \text{scal})
\]

3 Description

s17dc evaluates a sequence of values for the Bessel function \( Y_{\nu}(z) \), where \( z \) is complex, \(-\pi < \arg z \leq \pi\), and \( \nu \) is the real, non-negative order. The \( N \)-member sequence is generated for orders \( \nu \), \( \nu + 1 \), \ldots, \( \nu + N - 1 \). Optionally, the sequence is scaled by the factor \( e^{-i\text{Im}(z)} \).

Note: although the function may not be called with \( \nu \) less than zero, for negative orders the formula \( Y_{-\nu}(z) = Y_{\nu}(z) \cos(\pi \nu) + J_{\nu}(z) \sin(\pi \nu) \) may be used (for the Bessel function \( J_{\nu}(z) \), see s17de).

The function is derived from the function CBESY in Amos (1986). It is based on the relation

\[
Y_{\nu}(z) = \frac{H^{(1)}_{\nu}(z) - H^{(2)}_{\nu}(z)}{2i},
\]

where \( H^{(1)}_{\nu}(z) \) and \( H^{(2)}_{\nu}(z) \) are the Hankel functions of the first and second kinds respectively (see s17dl).

When \( N \) is greater than 1, extra values of \( Y_{\nu}(z) \) are computed using recurrence relations.

For very large \( |z| \) or \( (\nu + N - 1) \), argument reduction will cause total loss of accuracy, and so no computation is performed. For slightly smaller \( |z| \) or \( (\nu + N - 1) \), the computation is performed but results are accurate to less than half of machine precision. If \( |z| \) is very small, near the machine underflow threshold, or \( (\nu + N - 1) \) is too large, there is a risk of overflow and so no computation is performed. In all the above cases, a warning is given by the function.

4 References


5 Parameters

5.1 Compulsory Input Parameters

1:  \text{fnu} – double scalar

\( \nu \), the order of the first member of the sequence of functions.

\text{Constraint: } \text{fnu} \geq 0.0.

2:  \text{z} – complex scalar

\( z \), the argument of the functions.

\text{Constraint: } \text{z} \neq (0.0, 0.0).
3: \(n \rightarrow \text{int32 scalar}\)

\(N\), the number of members required in the sequence \(Y_{\nu}(z), Y_{\nu+1}(z), \ldots, Y_{\nu+N-1}(z)\).

Constraint: \(n \geq 1\).

4: \(\text{scal} \rightarrow \text{string}\)

The scaling option.

\(\text{scal} = 'U'\)

The results are returned unscaled.

\(\text{scal} = 'S'\)

The results are returned scaled by the factor \(e^{-|\text{Im}(z)|}\).

Constraint: \(\text{scal} = 'U'\) or \('S'\).

5.2 Optional Input Parameters

None.

5.3 Input Parameters Omitted from the MATLAB Interface

cwrk

5.4 Output Parameters

1: \(cy(n) \rightarrow \text{complex array}\)

The \(N\) required function values: \(\text{cy}(i)\) contains \(Y_{\nu+i-1}(z)\), for \(i = 1, 2, \ldots, N\).

2: \(nz \rightarrow \text{int32 scalar}\)

The number of components of \(cy\) that are set to zero due to underflow. The positions of such components in the array \(cy\) are arbitrary.

3: \(\text{ifail} \rightarrow \text{int32 scalar}\)

\(\text{ifail} = 0\) unless the function detects an error (see Section 6).

6 Error Indicators and Warnings

Errors or warnings detected by the function:

\(\text{ifail} = 1\)

On entry, \(\text{fnu} < 0.0\),

or \(z = (0.0, 0.0)\),

or \(n > 1\),

or \(\text{scal} \neq 'U'\) or \('S'\).

\(\text{ifail} = 2\)

No computation has been performed due to the likelihood of overflow, because \(\text{abs}(z)\) is less than a machine-dependent threshold value.

\(\text{ifail} = 3\)

No computation has been performed due to the likelihood of overflow, because \(\text{fnu} + n - 1\) is too large – how large depends on \(z\) as well as the overflow threshold of the machine.
ifail = 4

The computation has been performed, but the errors due to argument reduction in elementary functions make it likely that the results returned by s17dc are accurate to less than half of machine precision. This error exit may occur if either abs(z) or \( f_{nu} + n - 1 \) is greater than a machine-dependent threshold value.

ifail = 5

No computation has been performed because the errors due to argument reduction in elementary functions mean that all precision in results returned by s17dc would be lost. This error exit may occur if either abs(z) or \( f_{nu} + n - 1 \) is greater than a machine-dependent threshold value.

ifail = 6

No results are returned because the algorithm termination condition has not been met. This may occur because the parameters supplied to s17dc would have caused overflow or underflow.

7 Accuracy

All constants in s17dc are given to approximately 18 digits of precision. Calling the number of digits of precision in the floating-point arithmetic being used \( t \), then clearly the maximum number of correct digits in the results obtained is limited by \( p = \min(t, 18) \). Because of errors in argument reduction when computing elementary functions inside s17dc, the actual number of correct digits is limited, in general, by \( p - s \), where \( s \approx \max(1, \| \log_{10}|z| \|, \| \log_{10}|\nu| \|) \) represents the number of digits lost due to the argument reduction. Thus the larger the values of \( |z| \) and \( \nu \), the less the precision in the result. If s17dc is called with \( n > 1 \), then computation of function values via recurrence may lead to some further small loss of accuracy.

If function values which should nominally be identical are computed by calls to s17dc with different base values of \( \nu \) and different \( n \), the computed values may not agree exactly. Empirical tests with modest values of \( \nu \) and \( z \) have shown that the discrepancy is limited to the least significant 3 – 4 digits of precision.

8 Further Comments

The time taken for a call of s17dc is approximately proportional to the value of \( n \), plus a constant. In general it is much cheaper to call s17dc with \( n \) greater than 1, rather than to make \( N \) separate calls to s17dc.

Paradoxically, for some values of \( z \) and \( \nu \), it is cheaper to call s17dc with a larger value of \( n \) than is required, and then discard the extra function values returned. However, it is not possible to state the precise circumstances in which this is likely to occur. It is due to the fact that the base value used to start recurrence may be calculated in different regions for different \( n \), and the costs in each region may differ greatly.

Note that if the function required is \( Y_0(x) \) or \( Y_1(x) \), i.e., \( \nu = 0.0 \) or 1.0, where \( x \) is real and positive, and only a single unscaled function value is required, then it may be much cheaper to call s17ac or s17ad respectively.

9 Example

```plaintext
f_{nu} = 0;
z = complex(0.3, +0.4);
n = int32(2);
scal = 'U';
[cy, nz, ifail] = s17dc(f_{nu}, z, n, scal)
cy =
  -0.4983 + 0.6700i
  -1.0149 + 0.9485i
```
\begin{verbatim}
nz = 0
ifail = 0
\end{verbatim}