#### D01AKF - NAG Fortran Library Routine Document

Note. Before using this routine, please read the Users' Note for your implementation to check the interpretation of bold italicised terms and other implementation-dependent details.

## 1 Purpose

D01AKF is an adaptive integrator, especially suited to oscillating, non-singular integrands, which calculates an approximation to the integral of a function f(x) over a finite interval [a, b]:

$$I = \int_a^b f(x) \, dx.$$

# 2 Specification

```
SUBROUTINE DO1AKF(F, A, B, EPSABS, EPSREL, RESULT, ABSERR, W, LW,

IW, LIW, IFAIL)

INTEGER

LW, IW(LIW), LIW, IFAIL

real

F, A, B, EPSABS, EPSREL, RESULT, ABSERR, W(LW)

EXTERNAL

F
```

# 3 Description

D01AKF is based upon the QUADPACK routine QAG (Piessens *et al.* [3]). It is an adaptive routine, using the Gauss 30-point and Kronrod 61-point rules. A 'global' acceptance criterion (as defined by Malcolm and Simpson [1]) is used. The local error estimation is described in Piessens *et al.* [3].

Because this routine is based on integration rules of high order, it is especially suitable for non-singular oscillating integrands.

D01AKF requires the user to supply a function to evaluate the integrand at a single point.

The routine D01AUF uses an identical algorithm but requires the user to supply a subroutine to evaluate the integrand at an array of points. Therefore D01AUF will be more efficient if the evaluation can be performed in vector mode on a vector-processing machine.

D01AUF also has an additional parameter KEY which allows the user to select from six different Gauss-Kronrod rules.

### 4 References

- [1] Malcolm M A and Simpson R B (1976) Local versus global strategies for adaptive quadrature ACM Trans. Math. Software 1 129–146
- [2] Piessens R (1973) An algorithm for automatic integration Angew. Inf. 15 399–401
- [3] Piessens R, de Doncker–Kapenga E, Überhuber C and Kahaner D (1983) QUADPACK, A Subroutine Package for Automatic Integration Springer-Verlag

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#### 5 Parameters

1: F - real FUNCTION, supplied by the user.

External Procedure

F must return the value of the integrand f at a given point.

Its specification is:

 $egin{array}{ll} real & {\tt FUNCTION} \ {\tt F(X)} \\ real & {\tt X} \\ \end{array}$ 

1: X-real

On entry: the point at which the integrand f must be evaluated.

F must be declared as EXTERNAL in the (sub)program from which D01AKF is called. Parameters denoted as *Input* must **not** be changed by this procedure.

2: A — real

On entry: the lower limit of integration, a.

3: B-real

On entry: the upper limit of integration, b. It is not necessary that a < b.

4: EPSABS - real Input

On entry: the absolute accuracy required. If EPSABS is negative, the absolute value is used. See Section 7.

5: EPSREL — real Input

On entry: the relative accuracy required. If EPSREL is negative, the absolute value is used. See Section 7.

6: RESULT — real

On exit: the approximation to the integral I.

7: ABSERR — real Output

On exit: an estimate of the modulus of the absolute error, which should be an upper bound |I - RESULT|.

8: W(LW) — real array Output

On exit: details of the computation, as described in Section 8.

9: LW — INTEGER Input

On entry: the dimension of W, as declared in the (sub)program from which D01AKF is called. The value of LW (together with that of LIW below) imposes a bound on the number of sub-intervals into which the interval of integration may be divided by the routine. The number of sub-intervals cannot exceed LW/4. The more difficult the integrand, the larger LW should be.

Suggested value: a value in the range 800 to 2000 is adequate for most problems.

Constraint: LW  $\geq 4$ . See IW below.

10: IW(LIW) — INTEGER array Output

On exit: IW(1) contains the actual number of sub-intervals used. The rest of the array is used as workspace.

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11: LIW — INTEGER Input

On entry: the dimension of the array IW as declared in the (sub)program from which D01AKF is called. The number of sub-intervals into which the interval of integration may be divided cannot exceed LIW.

Suggested value: LIW = LW/4.

Constraint: LIW  $\geq 1$ .

#### 12: IFAIL — INTEGER

Input/Output

On entry: IFAIL must be set to 0, -1 or 1. Users who are unfamiliar with this parameter should refer to Chapter P01 for details.

On exit: IFAIL = 0 unless the routine detects an error or gives a warning (see Section 6).

For this routine, because the values of output parameters may be useful even if IFAIL  $\neq 0$  on exit, users are recommended to set IFAIL to -1 before entry. It is then essential to test the value of IFAIL on exit.

## 6 Error Indicators and Warnings

If on entry IFAIL = 0 or -1, explanatory error messages are output on the current error message unit (as defined by X04AAF).

Errors or warnings specified by the routine:

IFAIL = 1

The maximum number of subdivisions allowed with the given workspace has been reached without the accuracy requirements being achieved. Look at the integrand in order to determine the integration difficulties. Probably another integrator which is designed for handling the type of difficulty involved must be used. Alternatively, consider relaxing the accuracy requirements specified by EPSABS and EPSREL, or increasing the amount of workspace.

IFAIL = 2

Round-off error prevents the requested tolerance from being achieved. Consider requesting less accuracy.

IFAIL = 3

Extremely bad local integrand behaviour causes a very strong subdivision around one (or more) points of the interval. The same advice applies as in the case of IFAIL = 1.

IFAIL = 4

On entry, LW < 4, or LIW < 1.

# 7 Accuracy

The routine cannot guarantee, but in practice usually achieves, the following accuracy:

$$|I - RESULT| \le tol$$
,

where

$$tol = \max\{|\text{EPSABS}|, |\text{EPSREL}| \times |I|\},$$

and EPSABS and EPSREL are user-specified absolute and relative error tolerances. Moreover it returns the quantity ABSERR which, in normal circumstances satisfies

$$|I - \text{RESULT}| \leq \text{ABSERR} \leq tol.$$

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#### 8 Further Comments

The time taken by the routine depends on the integrand and the accuracy required.

If IFAIL  $\neq 0$  on exit, then the user may wish to examine the contents of the array W, which contains the end-points of the sub-intervals used by D01AKF along with the integral contributions and error estimates over these sub-intervals.

Specifically, for  $i=1,2,\ldots,n$ , let  $r_i$  denote the approximation to the value of the integral over the sub-interval  $[a_i,b_i]$  in the partition of [a,b] and  $e_i$  be the corresponding absolute error estimate. Then,  $\int_{a_i}^{b_i} f(x) \, dx \simeq r_i \text{ and RESULT} = \sum_{i=1}^n r_i. \text{ The value of } n \text{ is returned in IW}(1), \text{ and the values } a_i, b_i, e_i \text{ and } r_i \text{ are stored consecutively in the array W, that is:}$ 

$$\begin{array}{rcl} a_i & = & \mathrm{W}(i), \\ b_i & = & \mathrm{W}(n+i), \\ e_i & = & \mathrm{W}(2n+i) \text{ and } \\ r_i & = & \mathrm{W}(3n+i). \end{array}$$

## 9 Example

To compute

$$\int_0^{2\pi} x \sin(30x) \cos x \ dx.$$

#### 9.1 Program Text

**Note.** The listing of the example program presented below uses bold italicised terms to denote precision-dependent details. Please read the Users' Note for your implementation to check the interpretation of these terms. As explained in the Essential Introduction to this manual, the results produced may not be identical for all implementations.

```
DO1AKF Example Program Text
Mark 14 Revised. NAG Copyright 1989.
.. Parameters ..
INTEGER
                 LW, LIW
PARAMETER
                  (LW=800,LIW=LW/4)
INTEGER
                 NOUT
PARAMETER
                 (NOUT=6)
.. Scalars in Common ..
INTEGER
                 KOUNT
.. Local Scalars ..
                 A, ABSERR, B, EPSABS, EPSREL, PI, RESULT
real
INTEGER
.. Local Arrays ..
real
                 W(LW)
INTEGER
                 IW(LIW)
.. External Functions ..
                 FST, X01AAF
real
EXTERNAL
                 FST, X01AAF
.. External Subroutines ...
EXTERNAL
.. Common blocks ..
COMMON
                 /TELNUM/KOUNT
.. Executable Statements ..
WRITE (NOUT,*) 'DO1AKF Example Program Results'
PI = X01AAF(PI)
EPSABS = 0.0e0
EPSREL = 1.0e-03
A = 0.0e0
B = 2.0e0*PI
```

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```
KOUNT = O
     IFAIL = -1
     CALL DO1AKF(FST,A,B,EPSABS,EPSREL,RESULT,ABSERR,W,LW,IW,LIW,IFAIL)
     WRITE (NOUT,*)
     WRITE (NOUT,99998) 'EPSABS - absolute accuracy requested = ',
     WRITE (NOUT,99998) 'EPSREL - relative accuracy requested = ',
    + EPSREL
     WRITE (NOUT,*)
     IF (IFAIL.NE.O) WRITE (NOUT,99996) 'IFAIL = ', IFAIL
     IF (IFAIL.LE.3) THEN
        WRITE (NOUT,99997) 'RESULT - approximation to the integral = ',
        WRITE (NOUT,99998) 'ABSERR - estimate of the absolute error = '
         , ABSERR
        WRITE (NOUT, 99996) 'KOUNT - number of function evaluations = '
         , KOUNT
       WRITE (NOUT,99996) 'IW(1) - number of subintervals used = ',
         IW(1)
     END IF
     STOP
99999 FORMAT (1X,A,F10.4)
99998 FORMAT (1X,A,e9.2)
99997 FORMAT (1X,A,F9.5)
99996 FORMAT (1X,A,I4)
     END
     real FUNCTION FST(X)
     .. Scalar Arguments ..
     real
     .. Scalars in Common ..
     INTEGER
               KOUNT
     .. Intrinsic Functions ..
     INTRINSIC COS, SIN
     .. Common blocks ..
     COMMON
                     /TELNUM/KOUNT
     .. Executable Statements ..
     KOUNT = KOUNT + 1
     FST = X*(SIN(30.0e0*X))*COS(X)
     RETURN
     END
```

#### 9.2 Program Data

None.

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## 9.3 Program Results

D01AKF Example Program Results

```
A - lower limit of integration = 0.0000
B - upper limit of integration = 6.2832
EPSABS - absolute accuracy requested = 0.00E+00
EPSREL - relative accuracy requested = 0.10E-02

RESULT - approximation to the integral = -0.20967
ABSERR - estimate of the absolute error = 0.45E-13
KOUNT - number of function evaluations = 427
IW(1) - number of subintervals used = 4
```

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