





**GALAHAD** ICFS

USER DOCUMENTATION

GALAHAD Optimization Library version 5.0

#### 1 SUMMARY

Given a symmetric matrix **A**, this package **computes a symmetric, positive-definite approximation LL**<sup>T</sup> **using an incomplete Cholesky factorization**; the resulting matrix **L** is lower triangular. Subsequently, the solution **x** to the either of the linear systems  $\mathbf{L}\mathbf{x} = \mathbf{b}$  and  $\mathbf{L}^T\mathbf{x} = \mathbf{b}$  may be found for a given vector **b**.

ATTRIBUTES — Versions: GALAHAD\_ICFS\_single, GALAHAD\_ICFS\_double. Uses: GALAHAD\_SYMBOLS, GALAHAD\_SPECFILE and GALAHAD\_SPACE. Date: May 1998/December 2022. Origin: C.-J, Lin and J. J. Moré, Argonne National Laboratory, enhanced for modern fortran by N. I. M. Gould, Rutherford Appleton Laboratory. Language: Fortran 2003.

### 2 HOW TO USE THE PACKAGE

The package is available using both single and double precision reals, and either 32-bit or 64-bit integers. Access to the 32-bit integer, single precision version requires the USE statement

USE GALAHAD\_ICFS\_single

with the obvious substitution GALAHAD\_ICFS\_double, GALAHAD\_ICFS\_single\_64 and GALAHAD\_ICFS\_double\_64 for the other variants.

If it is required to use more than one of the modules at the same time, the derived types <code>ICFS\_control\_type</code>, <code>ICFS\_inform\_type</code>, <code>ICFS\_data\_type</code> and <code>NLPT\_userdata\_type</code>, (Section 2.2) and the subroutines <code>ICFS\_initialize</code>, <code>ICFS\_factorize</code>, <code>ICFS\_solve</code>, <code>ICFS\_terminate</code>, (Section 2.3) and <code>ICFS\_read\_specfile</code> (Section 2.5) must be renamed on one of the <code>USE</code> statements.

## 2.1 Real and integer kinds

We use the terms integer and real to refer to the fortran keywords REAL(rp\_) and INTEGER(ip\_), where rp\_ and ip\_ are the relevant kind values for the real and integer types employed by the particular module in use. The former are equivalent to default REAL for the single precision versions and DOUBLE PRECISION for the double precision cases, and correspond to rp\_ = real32 and rp\_ = real64, respectively, as supplied by the fortran iso\_fortran\_env module. The latter are default (32-bit) and long (64-bit) integers, and correspond to ip\_ = int32 and ip\_ = int64, respectively, again from the iso\_fortran\_env module.

## 2.2 The derived data types

Four derived data types are accessible from the package.

# 2.2.1 The derived data type for holding control parameters

The derived data type ICFS\_control\_type is used to hold controlling data. Default values may be obtained by calling ICFS\_initialize (see Section 2.3.1), while components may also be changed by calling GALAHAD\_ICFS\_read\_spec (see Section 2.5.1). The components of ICFS\_control\_type are:

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error is a scalar variable of type INTEGER(ip\_), that holds the stream number for error messages. Printing of error messages in ICFS\_analyse, ICFS\_estimate and ICFS\_terminate is suppressed if error  $\leq 0$ . The default is error = 6.

- out is a scalar variable of type INTEGER (ip\_), that holds the stream number for informational messages. Printing of informational messages in ICFS\_analyse and ICFS\_estimate is suppressed if out < 0. The default is out = 6.
- print\_level is a scalar variable of type INTEGER (ip\_), that is used to control the amount of informational output which is required. No informational output will occur if print\_level  $\leq 0$ . If print\_level > 01, details of any data errors encountered will be reported. The default is print\_level = 0.
- icfs\_vectors is a scalar variable of type INTEGER (ip\_), that holds the number of extra vectors of length *n* required by the incomplete Cholesky preconditioner. Usually, the larger the number, the better the preconditioner, but the more space and effort required to use it. Any negative value will be regarded as 0. The default is icfs\_vectors = 10.
- shift is a scalar variable of type REAL (rp\_), that holds an initial estimate of the shift  $\alpha$  used so that the incomplete factorization of  $\mathbf{A} + \mathbf{D}$  is positive definite, where  $\mathbf{D}$  is a diagonal matrix whose entries are no larger than  $\alpha$ . This value  $\alpha$  may subsequently be increased, as necessary, by the package, see inform%shift. The default is shift 0.0.
- space\_critical is a scalar variable of type default LOGICAL, that must be set .TRUE. if space is critical when allocating arrays and .FALSE. otherwise. The package may run faster if space\_critical is .FALSE. but at the possible expense of a larger storage requirement. The default is space\_critical = .FALSE..
- deallocate\_error\_fatal is a scalar variable of type default LOGICAL, that must be set .TRUE. if the user wishes to terminate execution if a deallocation fails, and .FALSE. if an attempt to continue will be made. The default is deallocate\_error\_fatal = .FALSE..
- prefix is a scalar variable of type default CHARACTER and length 30, that may be used to provide a user-selected character string to preface every line of printed output. Specifically, each line of output will be prefaced by the string prefix (2:LEN(TRIM( prefix ))-1), thus ignoreing the first and last non-null components of the supplied string. If the user does not want to preface lines by such a string, they may use the default prefix = "".

#### 2.2.2 The derived data type for holding timing information

The derived data type ICFS\_time\_type is used to hold elapsed CPU and system clock times for the various parts of the calculation. The components of ICFS\_time\_type are:

- total is a scalar variable of type REAL(rp\_), that gives the total CPU time spent in the package.
- factorize is a scalar variable of type REAL (rp\_), that gives the CPU time spent factorizing the required matrices.
- solve is a scalar variable of type REAL (rp\_), that gives the CPU time spent solving the resulting triangular systems.
- clock\_total is a scalar variable of type REAL(rp\_), that gives the total elapsed system clock time spent in the package.
- clock\_factorize is a scalar variable of type REAL (rp\_), that gives the elapsed system clock time spent factorizing the required matrices.
- clock\_solve is a scalar variable of type REAL(rp\_), that gives the elapsed system clock time spent solving the resulting triangular systems.

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#### 2.2.3 The derived data type for holding informational parameters

The derived data type ICFS\_inform\_type is used to hold parameters that give information about the progress and needs of the algorithm. The components of ICFS\_inform\_type are:

- status is a scalar variable of type INTEGER(ip\_), that gives the exit status of the algorithm. See Section 2.4 for details.
- alloc\_status is a scalar variable of type INTEGER( $ip_-$ ), that gives the status of the last attempted array allocation or deallocation. This will be 0 if status = 0.
- bad\_alloc is a scalar variable of type default CHARACTER and length 80, that gives the name of the last internal array for which there were allocation or deallocation errors. This will be the null string if status = 0.
- shift is a scalar variable of type REAL (rp\_), that holds the final value of the shift  $\alpha$  used so that the incomplete factorization of A + D is positive definite, where D is a diagonal matrix whose entries are no larger than  $\alpha$ .
- time is a scalar variable of type ICFS\_time\_type whose components are used to hold elapsed CPU and system clock times for the various parts of the calculation (see Section 2.2.2).

## 2.2.4 The derived data type for holding problem data

The derived data type ICFS\_data\_type is used to hold all the data for a particular problem, or sequences of problems with the same structure, between calls of ICFS procedures. This data should be preserved, untouched, from the initial call to ICFS\_initialize to the final call to ICFS\_terminate.

## 2.3 Argument lists and calling sequences

There are four procedures for user calls (see Section 2.5 for further features):

- 1. The subroutine ICFS\_initialize is used to set default values, and initialize private data.
- 2. The subroutine ICFS\_factorize is called to form the incomplete Cholesky factor L from A.
- 3. The subroutine ICFS\_triangular\_solve is called to solve either of the triangular systems  $\mathbf{L}\mathbf{x} = \mathbf{b}$  or  $\mathbf{L}^T\mathbf{x} = \mathbf{b}$  for given vectors  $\mathbf{b}$ .
- 4. The subroutine ICFS\_terminate is provided to allow the user to automatically deallocate array components of the private data, allocated by ICFS\_factorize, at the end of the solution process. It is important to do this if the data object is re-used for another matrix with a different structure since ICFS\_initialize cannot test for this situation, and any existing associated targets will subsequently become unreachable.

We note that in addition to the above calls, the user may also call the original fortran 77 subroutines DICFS and DSTRSOL directly. For details of the necessary argument lists, see Section 6.

# 2.3.1 The initialization subroutine

Default values are provided as follows:

```
CALL ICFS_initialize( data, control, inform )
```

data is a scalar INTENT (INOUT) argument of type ICFS\_data\_type (see Section 2.2.4). It is used to hold data about the matrix and its factors.

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control is a scalar INTENT (OUT) argument of type ICFS\_control\_type (see Section 2.2.1). On exit, control contains default values for the components as described in Section 2.2.1. These values should only be changed after calling ICFS\_initialize.

inform is a scalar INTENT (OUT) argument of type ICFS\_inform\_type (see Section 2.2.3). A successful call to ICFS\_initialize is indicated when the component status has the value 0. For other return values of status, see Section 2.4.

#### 2.3.2 The incomplete factorization subroutine

The factorization phase, in which incomplete Cholesky factors L of A are determined, is performed as follows:

```
CALL ICFS_factorize( n, PTR, ROW, DIAG, VAL, data, control, inform )
```

- is a scalar INTENT (IN) scalar argument of type INTEGER (ip\_), that must be set to n the dimension of the matrix **A. Restrictions:** n > 0.
- PTR is a scalar INTENT(IN) rank-one array argument of type INTEGER(ip\_) and dimension n+1, whose j-th component gives the starting address for list of nonzero values and their corresponding row indices in column j of the **strict lower triangular part** of **A** (The entry  $a_{i,j}$  is in the strict lower triangular part of **A** if i > j). That is, the nonzeros in column j of the strict lower triangle of **A** must be in positions PTR(j) ..., PTR(j+1) 1 for j = 1, ... Note that PTR(n+1) points to the first position beyond that needed to store **A**.
- ROW is a scalar INTENT (IN) rank-one array argument of type INTEGER (ip\_) and dimension at least PTR (n+1) -1, that contains the row indices of the strict lower triangular part of **A** in the compressed column storage format described above.
- DIAG is a scalar INTENT (IN) rank-one array argument of type REAL (rp\_) and dimension n, whose j-th component contains the value of the j-th diagonal of A.
- VAL is a scalar INTENT (IN) rank-one array argument of type REAL (rp\_) and dimension at least PTR (n+1) -1, that contains the values of the entries strict lower triangular part of **A** that correspond to the row indices described above.
- data is a scalar INTENT (INOUT) argument of type ICFS\_data\_type (see Section 2.2.4). It is used to hold data about the problem being solved. It must not have been altered by the user since the last call to ICFS\_initialize.
- control is a scalar INTENT (IN) argument of type ICFS\_control\_type (see Section 2.2.1). Default values may be assigned by calling ICFS\_initialize prior to the first call to ICFS\_analyse.
- inform is a scalar INTENT (INOUT) argument of type ICFS\_inform\_type (see Section 2.2.3). A successful call to ICFS\_analyse is indicated when the component status has the value 0. For other return values of status, see Section 2.4.

## 2.3.3 The triangular solution subroutine

The solution phase, in which one of the triangular systems  $\mathbf{L}\mathbf{x} = \mathbf{b}$  or  $\mathbf{L}^T\mathbf{x} = \mathbf{b}$  for given vectors  $\mathbf{b}$ , is performed as follows:

```
CALL ICFS_triangular_solve( n, X, transpose, data, control, inform )
```

- n, data, control and inform are exactly as described and input to ICFS\_factorize, and must not have been changed in the interim.
- is a scalar INTENT (INOUT) rank-one array argument of type REAL (rp\_), and dimension n, that must be set on input so that X(i) contains the component  $b_i$ , i = 1, ..., n of the vector **b**. On output, this will have been overwritten by the desired vector **x**.

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transpose is a scalar INTENT (IN) argument of type default LOGICAL, that must be set .FALSE. if the user wishes to solve Lx = b and .TRUE. if instead the solution to  $L^Tx = b$  is sought.

#### 2.3.4 The termination subroutine

All previously allocated arrays are deallocated as follows:

```
CALL ICFS_terminate( data, control, inform )
```

data is a scalar INTENT (INOUT) argument of type ICFS\_data\_type exactly as for ICFS\_factorize, which must not have been altered by the user since the last call to ICFS\_initialize. On exit, array components will have been deallocated.

control is a scalar INTENT(IN) argument of type ICFS\_control\_type exactly as for ICFS\_factorize.

inform is a scalar INTENT (OUT) argument of type ICFS\_inform\_type exactly as for ICFS\_factorize. Only the component status will be set on exit, and a successful call to ICFS\_terminate is indicated when this component status has the value 0. For other return values of status, see Section 2.4.

## 2.4 Warning and error messages

A negative value of inform%status on exit from ICFS\_factorize, ICFS\_triangular\_solve or ICFS\_terminate indicates that an error has occurred. No further calls should be made until the error has been corrected. Possible values are:

- -1. An allocation error occurred. A message indicating the offending array is written on unit control%error, and the returned allocation status and a string containing the name of the offending array are held in inform%alloc\_status and inform%bad\_alloc, respectively.
- -2. A deallocation error occurred. A message indicating the offending array is written on unit control%error and the returned allocation status and a string containing the name of the offending array are held in inform%alloc\_status and inform%bad\_alloc, respectively.
- -3. The restriction 0 < n has been violated.

#### 2.5 Further features

In this section, we describe an alternative means of setting control parameters, that is components of the variable control of type ICFS\_control\_type (see Section 2.2.1), by reading an appropriate data specification file using the subroutine ICFS\_read\_specfile. This facility is useful as it allows a user to change ICFS control parameters without editing and recompiling programs that call ICFS.

A specification file, or specfile, is a data file containing a number of "specification commands". Each command occurs on a separate line, and comprises a "keyword", which is a string (in a close-to-natural language) used to identify a control parameter, and an (optional) "value", which defines the value to be assigned to the given control parameter. All keywords and values are case insensitive, keywords may be preceded by one or more blanks but values must not contain blanks, and each value must be separated from its keyword by at least one blank. Values must not contain more than 30 characters, and each line of the specfile is limited to 80 characters, including the blanks separating keyword and value.

The portion of the specification file used by  $ICFS\_read\_specfile$  must start with a "BEGIN ICFS" command and end with an "END" command. The syntax of the specifile is thus defined as follows:

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```
( .. lines ignored by ICFS_read_specfile .. )
BEGIN ICFS
    keyword value
    .....
    keyword value
END
( .. lines ignored by ICFS_read_specfile .. )
```

where keyword and value are two strings separated by (at least) one blank. The "BEGIN ICFS" and "END" delimiter command lines may contain additional (trailing) strings so long as such strings are separated by one or more blanks, so that lines such as

```
BEGIN ICFS SPECIFICATION

and

END ICFS SPECIFICATION
```

are acceptable. Furthermore, between the "BEGIN ICFS" and "END" delimiters, specification commands may occur in any order. Blank lines and lines whose first non-blank character is ! or \* are ignored. The content of a line after a ! or \* character is also ignored (as is the ! or \* character itself). This provides an easy manner to "comment out" some specification commands, or to comment specific values of certain control parameters.

The value of a control parameters may be of three different types, namely integer, logical or real. Integer and real values may be expressed in any relevant Fortran integer and floating-point formats (respectively). Permitted values for logical parameters are "ON", "TRUE", ".TRUE.", "T", "YES", "Y", or "OFF", "NO", "N", "FALSE", ".FALSE." and "F". Empty values are also allowed for logical control parameters, and are interpreted as "TRUE".

The specification file must be open for input when ICFS\_read\_specifile is called, and the associated device number passed to the routine in device (see below). Note that the corresponding file is REWINDed, which makes it possible to combine the specifications for more than one program/routine. For the same reason, the file is not closed by ICFS\_read\_specifile.

## 2.5.1 To read control parameters from a specification file

Control parameters may be read from a file as follows:

```
CALL ICFS_read_specfile( control, device )
```

control is a scalar INTENT (INOUT) argument of type ICFS\_control\_type (see Section 2.2.1). Default values should have already been set, perhaps by calling ICFS\_initialize. On exit, individual components of control may have been changed according to the commands found in the specifile. Specifile commands and the component (see Section 2.2.1) of control that each affects are given in Table 2.1.

device is a scalar INTENT (IN) argument of type INTEGER (ip\_), that must be set to the unit number on which the specifile has been opened. If device is not open, control will not be altered and execution will continue, but an error message will be printed on unit control%error.

# 2.6 Information printed

If control%print\_level is positive, information about errors encountered will be printed on unit control%out.

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command	component of control	value type
error-printout-device	%error	integer
printout-device	%out	integer
print-level	%print_level	integer
number-of-icfs-vectors	icfs_vectors	integer
initial-shift	shift	real
space-critical	%space_critical	logical
deallocate-error-fatal	%deallocate_error_fatal	logical
output-line-prefix	%prefix	character

Table 2.1: Specfile commands and associated components of control.

## 3 GENERAL INFORMATION

Use of common: None.

**Workspace:** Provided automatically by the module.

Other routines called directly: None.

Other modules used directly: The module uses the GALAHAD packages GALAHAD\_SYMBOLS, GALAHAD\_SPECFILE and GALAHAD\_SPACE.

Input/output: Output is under control of the arguments control%error, control%out and control%print\_level.

**Restrictions:** n > 0.

**Portability:** ISO Fortran 2003. The package is thread-safe.

## 4 METHOD

The package computes incomplete Cholesky factors L of A so that

$$\mathbf{A} + \mathbf{D} = \mathbf{L}\mathbf{L} + \mathbf{E}.$$

The pattern of entries in L match those in A with additional "fill-ins" controlled by the extra storage provided. The shifted diagonal D whose values do not exceed a scalar shift  $\alpha$  allows for indefinite A, and also guarantees stable factors L. Increasing the extra storage provided generally decreases the size of the error matrix E, but increases the cost of the algorithm.

#### **Reference:**

The method is described in detail in

C.J. Lin and J. J. Moreé. "Incomplete Cholesky factorizations with limited memory". SIAM J. Sci. Computing. **21** (1999) 24–45.

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## 5 EXAMPLES OF USE

Suppose that

$$\mathbf{A} = \begin{pmatrix} 2 & 1 & & & \\ 1 & 5 & 1 & & 1 \\ & 1 & 1 & 1 & \\ & & 1 & 7 & \\ & 1 & & 2 \end{pmatrix} \text{ and } \mathbf{b} = \begin{pmatrix} 3 \\ 8 \\ 3 \\ 8 \\ 3 \end{pmatrix},$$

where the missing entries in **A** are structural zeros. Then we may find suitable incomplete factors **L**, and subsequently solve  $\mathbf{L}\mathbf{y} = \mathbf{b}$  and  $\mathbf{L}^T\mathbf{x} = \mathbf{y}$  using the following code; notice that we initialize **b** in **x**, overwrite this with **y** in the first triangular solve, and finally recover **x** from **y** in the second solve:

```
! GALAHAD 4.1 - 2022-12-04 AT 09:30 GMT.
  PROGRAM ICFS_EXAMPLE
  USE GALAHAD_ICFS_double
  IMPLICIT NONE
  INTEGER, PARAMETER :: wp = KIND( 1.0D+0 )
  TYPE ( ICFS_data_type ) :: data
  TYPE ( ICFS_control_type ) control
  TYPE ( ICFS_inform_type ) :: inform
  INTEGER, PARAMETER :: n = 5
  INTEGER, PARAMETER :: ne = 5
  INTEGER, ALLOCATABLE, DIMENSION(:):: PTR, ROW
  REAL ( KIND = wp ), ALLOCATABLE, DIMENSION( : ) :: DIAG, VAL
  REAL ( KIND = wp ) :: X( n )
! allocate and set lower triangle of matrix in sparse by column form
  ALLOCATE ( PTR ( n + 1 ), DIAG ( n ), VAL ( ne ), ROW ( ne ) )
  PTR = (/1, 2, 4, 5, 5, 5/)
  ROW = (/ 2, 3, 5, 4 /)
  DIAG = (/2.0_{\text{wp}}, 5.0_{\text{wp}}, 1.0_{\text{wp}}, 7.0_{\text{wp}}, 2.0_{\text{wp}})
  VAL = (/ 1.0_wp, 1.0_wp, 1.0_wp, 1.0_wp /)
! problem setup complete
  CALL ICFS_initialize( data, control, inform )
  control%icfs_vectors = 1
! form and factorize the preconditioner, P = L L^T
  CALL ICFS_factorize( n, PTR, ROW, DIAG, VAL, data, control, inform )
  IF ( inform%status < 0 ) THEN</pre>
    WRITE(6, '(A, I0)')
                                                                                  δ
           Failure of ICFS_factorize with status = ', inform%status
    STOP
  END IF
! use the factors to solve L L^T x = b, with b input in x
  X(:n) = (/3.0_{wp}, 8.0_{wp}, 3.0_{wp}, 8.0_{wp}, 3.0_{wp}, 3.0_{wp})
  CALL ICFS_triangular_solve( n, X, .FALSE., data, control, inform )
  CALL ICFS_triangular_solve( n, X, .TRUE., data, control, inform )
  IF ( inform%status == 0 ) THEN
    WRITE( 6, "( ' ICFS - Preconditioned solution is ', 5F6.2 )" ) X
  ELSE
    WRITE(6, "('ICFS - exit status = ', IO)") inform%status
  END IF
! clean up
  CALL ICFS_terminate( data, control, inform )
  DEALLOCATE ( DIAG, VAL, ROW, PTR )
  STOP
  END PROGRAM ICFS_EXAMPLE
```

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The code produces the following output:

```
ICFS - Preconditioned solution is 1.00 1.00 1.00 1.00 1.00
```

## 6 APPENDIX

The original subroutine calls DICFS and DSTRSOL, as encoded in ICFS\_factorize and ICFS\_triangular\_solve, may also be called as part of the package. To quote from the package source,

```
Subroutine dicfs
     Given a symmetric matrix A in compressed column storage, this
     subroutine computes an incomplete Cholesky factor of A + alpha*D,
     where alpha is a shift and D is the diagonal matrix with entries
     set to the 12 norms of the columns of A.
     The subroutine statement is
       subroutine dicfs(n,nnz,a,adiag,acol_ptr,arow_ind,l,ldiag,lcol_ptr,
1
                         lrow_ind,p,alpha,iwa,wa1,wa2)
     where
       n is an integer variable.
         On entry n is the order of A.
         On exit n is unchanged.
       nnz is an integer variable.
          On entry nnz is the number of nonzeros in the strict lower
            triangular part of A.
         On exit nnz is unchanged.
       a is a real array of dimension nnz.
          On entry a must contain the strict lower triangular part
            of A in compressed column storage.
         On exit a is unchanged.
       adiag is a real array of dimension n.
         On entry adiag must contain the diagonal elements of A.
          On exit adiag is unchanged.
       acol_ptr is an integer array of dimension n + 1.
          On entry acol_ptr must contain pointers to the columns of A.
            The nonzeros in column j of A must be in positions
             acol_ptr(j), ..., acol_ptr(j+1) - 1.
         On exit acol_ptr is unchanged.
       arow ind is an integer array of dimension nnz.
          On entry arow_ind must contain row indices for the strict
            lower triangular part of A in compressed column storage.
```

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```
On exit arow_ind is unchanged.
       l is a real array of dimension nnz+n*p.
         On entry 1 need not be specified.
         On exit 1 contains the strict lower triangular part
            of L in compressed column storage.
       ldiag is a real array of dimension n.
1
         On entry ldiag need not be specified.
         On exit ldiag contains the diagonal elements of L.
       lcol_ptr is an integer array of dimension n + 1.
         On entry lcol_ptr need not be specified.
         On exit lcol_ptr contains pointers to the columns of L.
            The nonzeros in column j of L are in the
             lcol_ptr(j), ..., lcol_ptr(j+1) - 1 positions of 1.
       lrow_ind is an integer array of dimension nnz+n*p.
         On entry lrow_ind need not be specified.
         On exit lrow_ind contains row indices for the strict lower
            triangular part of L in compressed column storage.
       p is an integer variable.
1
!
         On entry p specifes the amount of memory available for the
             incomplete Cholesky factorization.
         On exit p is unchanged.
       alpha is a real variable.
         On entry alpha is the initial guess of the shift.
         On exit alpha is final shift
1
      iwa is an integer work array of dimension 3*n.
       wal is a real work array of dimension n.
       wa2 is a real work array of dimension n.
!
     Subroutine dstrsol
!
     This subroutine solves the triangular systems L^*x = r or L'^*x = r.
     The subroutine statement is
       subroutine dstrsol(n,l,ldiag,jptr,indr,r,task)
!
1
     where
       n is an integer variable.
         On entry n is the order of L.
         On exit n is unchanged.
```

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```
l is a real array of dimension *.
         On entry 1 must contain the nonzeros in the strict lower
            triangular part of L in compressed column storage.
         On exit l is unchanged.
       ldiag is a real array of dimension n.
         On entry ldiag must contain the diagonal elements of L.
1
         On exit ldiag is unchanged.
       jptr is an integer array of dimension n + 1.
         On entry jptr must contain pointers to the columns of A.
            The nonzeros in column j of A must be in positions
             jptr(j), ..., jptr(j+1) - 1.
         On exit jptr is unchanged.
       indr is an integer array of dimension *.
         On entry indr must contain row indices for the strict
            lower triangular part of L in compressed column storage.
         On exit indr is unchanged.
       r is a real array of dimension n.
         On entry r must contain the vector r.
         On exit r contains the solution vector x.
      task is a character variable of length 60.
         On entry
            task(1:1) = 'N' if we need to solve L*x = r
            task(1:1) = 'T' if we need to solve L'*x = r
         On exit task is unchanged.
     MINPACK-2 Project. October 1998.
     Argonne National Laboratory.
     Chih-Jen Lin and Jorge J. More'.
```