

# GEOS-Chem Reference, Vol. 2: Utility Modules

BOB YANTOSCA, MICHAEL LONG, MELISSA PAYER AND MATTHEW COOPER  
*GEOS-Chem Support Team*

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# 1 Routine/Function Prologues

## 1.1 Fortran: Module Interface bpch2\_mod

Module BPCH2\_MOD contains the routines used to read data from and write data to binary punch (BPCH) file format (v. 2.0).

### INTERFACE:

```
MODULE BPCH2_MOD
```

### USES:

```
IMPLICIT NONE
#   include "define.h"
PRIVATE
```

### PUBLIC MEMBER FUNCTIONS:

```
PUBLIC  :: OPEN_BPCH2_FOR_READ
PUBLIC  :: OPEN_BPCH2_FOR_WRITE
PUBLIC  :: BPCH2_HDR
PUBLIC  :: BPCH2
PUBLIC  :: READ_BPCH2
PUBLIC  :: GET_MODELNAME
PUBLIC  :: GET_NAME_EXT
PUBLIC  :: GET_NAME_EXT_2D
PUBLIC  :: GET_RES_EXT
PUBLIC  :: GET_HALFPOLAR
PUBLIC  :: GET_TAU0
```

```
INTERFACE GET_TAU0
  MODULE PROCEDURE GET_TAU0_6A
END INTERFACE
```

### PRIVATE MEMBER FUNCTIONS:

```
PRIVATE :: GET_TAU0_6A
```

### REVISION HISTORY:

- (1 ) Added routine GET\_TAU0 (bmy, 7/20/00)
- (2 ) Added years 1985-2001 for routine GET\_TAU0 (bmy, 8/1/00)
- (3 ) Use IOS /= 0 criterion to also check for EOF (bmy, 9/12/00)
- (4 ) Removed obsolete code in "read\_bpch2.f" (bmy, 12/18/00)
- (5 ) Correct error for 1991 TAU values in GET\_TAU0 (bnd, bmy, 1/4/01)
- (6 ) BPCH2\_MOD is now independent of any GEOS-CHEM size parameters.  
(bmy, 4/18/01)
- (7 ) Now have 2 versions of "GET\_TAU0" overloaded by an interface. The original version takes 2 arguments (MONTH, YEAR). The new version

takes 3 arguments (MONTH, DAY, YEAR). (bmy, 8/22/01)

(8 ) Updated comments (bmy, 9/4/01)

(9 ) Renamed GET\_TAU0\_3A to GET\_TAU0\_6A, and updated the GET\_TAU0 interface. Also updated comments (bmy, 9/26/01)

(10) Now use special model name for GEOS-3 w/ 30 layers (bmy, 10/9/01)

(11) Minor bug fix in GET\_TAU0\_2A. Also deleted obsolete code from 9/01. (bmy, 11/15/01)

(12) Moved routines JULDAY, MINT, CALDATE to "julian\_mod.f". Now references routine JULDAY from "julday\_mod.f". Also added code for GEOS-4/fvDAS model type. (bmy, 11/20/01)

(23) Now divide module header into MODULE PRIVATE, MODULE VARIABLES, and MODULE ROUTINES sections. Also add MODULE INTERFACES section, since we have an interface here. (bmy, 5/28/02)

(24) Added OPEN\_BPCH2\_FOR\_READ and OPEN\_BPCH2\_FOR\_WRITE. Also now reference IU\_FILE and IOERROR from "file\_mod.f". (bmy, 7/30/02)

(25) Now references "error\_mod.f". Also obsoleted routine GET\_TAU0\_2A. (bmy, 10/15/02)

(26) Made modification in READ\_BPCH2 for 1x1 nested grids (bmy, 3/11/03)

(27) Modifications for GEOS-4, 30-layer grid (bmy, 11/3/03)

(28) Added cpp switches for GEOS-4 1x125 grid (bmy, 12/1/04)

(29) Modified for GCAP and GEOS-5 met fields. Added function GET\_HALFPOLAR. (bmy, 6/28/05)

(30) Added GET\_NAME\_EXT\_2D to get filename extension for files which do not contain any vertical information (bmy, 8/16/05)

(31) Remove support for GEOS-1 and GEOS-STRAT met fields (bmy, 8/4/06)

(32) Renamed GRID30LEV to GRIDREDUCED. Also increase TEMPARRAY in READ\_BPCH2 for GEOS-5 vertical levels. (bmy, 2/16/07)

(33) Modifications for GEOS-5 nested grids (bmy, 11/6/08)

20 Nov 2009 - R. Yantosca - Added ProTeX headers

13 Aug 2010 - R. Yantosca - Added modifications for MERRA

### 1.1.1 open\_bpch2\_for\_read

Subroutine OPEN\_BPCH2\_FOR\_READ opens a binary punch file (version 2.0 format) for reading only. Also reads FTI and TITLE strings.

#### INTERFACE:

```
SUBROUTINE OPEN_BPCH2_FOR_READ( IUNIT, FILENAME, TITLE )
```

#### USES:

```
USE ERROR_MOD, ONLY : ERROR_STOP
USE FILE_MOD,  ONLY : IOERROR
```

#### INPUT PARAMETERS:

INTEGER,	INTENT(IN)	:: IUNIT	! LUN for file I/O
CHARACTER(LEN=*),	INTENT(IN)	:: FILENAME	! Name of file

**OUTPUT PARAMETERS:**

CHARACTER(LEN=80), INTENT(OUT), OPTIONAL :: TITLE ! File title string

**REVISION HISTORY:**

(1 ) Now references ERROR\_STOP from "error\_mod.f" (bmy, 10/15/02)  
 20 Nov 2009 - R. Yantosca - Added ProTeX header

---

**1.1.2 open\_bpch2\_for\_write**

Subroutine OPEN\_BPCH2\_FOR\_WRITE opens a binary punch file (version 2.0) for writing.

**INTERFACE:**

SUBROUTINE OPEN\_BPCH2\_FOR\_WRITE( IUNIT, FILENAME, TITLE )

**USES:**

USE FILE\_MOD, ONLY : IOERROR

**INPUT PARAMETERS:**

INTEGER, INTENT(IN) :: IUNIT ! LUN for file I/O  
 CHARACTER(LEN=\*), INTENT(IN) :: FILENAME ! Name of file

**OUTPUT PARAMETERS:**

CHARACTER(LEN=80), INTENT(OUT), OPTIONAL :: TITLE ! File title string

**REVISION HISTORY:**

30 Jul 2002 - R. Yantosca - Initial version  
 20 Nov 2009 - R. Yantosca - Added ProTeX header

---

**1.1.3 bpch2\_hdr**

Subroutine BPCH2\_HDR writes a header at the top of the binary punch file, version 2.0.

**INTERFACE:**

SUBROUTINE BPCH2\_HDR ( IUNIT, TITLE )

**USES:**

USE FILE\_MOD, ONLY : IOERROR

**INPUT PARAMETERS:**

INTEGER, INTENT(IN) :: IUNIT ! LUN for file I/O  
 CHARACTER(LEN=80), INTENT(IN) :: TITLE ! Top-of-file title string

**REVISION HISTORY:**

(1 ) Added this routine to "bpch\_mod.f" (bmy, 6/28/00)  
 (2 ) Use IOS /= 0 criterion to also check for EOF condition (bmy, 9/12/00)  
 (3 ) Now reference IOERROR from "file\_mod.f". (bmy, 6/26/02)  
 20 Nov 2009 - R. Yantosca - Added ProTeX header

---



### 1.1.4 bpch2

Subroutine BPCH2 writes binary punch file (version 2.0) to disk. Information about the model grid is also stored with each data block.

#### INTERFACE:

```

SUBROUTINE BPCH2( IUNIT,      MODELNAME, LONRES,  LATRES,
&                HALFPOLAR, CENTER180, CATEGORY, NTRACER,
&                UNIT,      TAU0,      TAU1,      RESERVED,
&                NI,      NJ,      NL,      IFIRST,
&                JFIRST,  LFIRST,  ARRAY )

```

#### USES:

```

USE FILE_MOD, ONLY : IOERROR

```

#### INPUT PARAMETERS:

```

! Arguments
INTEGER,          INTENT(IN) :: IUNIT           ! LUN for file I/O
CHARACTER(LEN=20), INTENT(IN) :: MODELNAME       ! Met field type
REAL*4,           INTENT(IN) :: LONRES           ! Lon resolution [deg]
REAL*4,           INTENT(IN) :: LATRES           ! Lat resolution [deg]
INTEGER,          INTENT(IN) :: HALFPOLAR        ! 1/2-size polar boxes?
INTEGER,          INTENT(IN) :: CENTER180        ! 1st box center -180?
CHARACTER(LEN=40), INTENT(IN) :: CATEGORY        ! Diag. category name
INTEGER,          INTENT(IN) :: NTRACER          ! Tracer index #
CHARACTER(LEN=40), INTENT(IN) :: UNIT            ! Unit string
REAL*8,           INTENT(IN) :: TAU0             ! TAU values @ start &
REAL*8,           INTENT(IN) :: TAU1             ! end of diag interval
CHARACTER(LEN=40), INTENT(IN) :: RESERVED        ! Extra string
INTEGER,          INTENT(IN) :: NI, NJ, NL       ! Dimensions of ARRAY
INTEGER,          INTENT(IN) :: IFIRST           ! (I,J,L) indices of
INTEGER,          INTENT(IN) :: JFIRST           ! the first grid box
INTEGER,          INTENT(IN) :: LFIRST           ! in Fortran notation
REAL*4,           INTENT(IN) :: ARRAY(NI,NJ,NL) ! Data array

```

#### REVISION HISTORY:

- (1 ) Added indices to IOERROR calls (e.g. "bpch2:1", "bpch2:2", etc.)  
(bmy, 10/4/99)
- (2 ) Added this routine to "bpch\_mod.f" (bmy, 6/28/00)
- (3 ) Use IOS /= 0 criterion to also check for EOF condition (bmy, 9/12/00)
- (4 ) Now reference IOERROR from "file\_mod.f". (bmy, 6/26/02)

### 1.1.5 read\_bpch2

Subroutine READ\_BPCH2 reads a binary punch file (v. 2.0) and extracts a data block that matches the given category, tracer, and tau value.

**INTERFACE:**

```

      SUBROUTINE READ_BPCH2( FILENAME, CATEGORY_IN, TRACER_IN,
&                           TAUO_IN,   IX,           JX,
&                           LX,         ARRAY,        QUIET )

```

**USES:**

```

      USE ERROR_MOD, ONLY : ERROR_STOP
      USE FILE_MOD,  ONLY : IU_FILE, IOERROR

```

```

      #      include "define.h"

```

**INPUT PARAMETERS:**

```

      CHARACTER(LEN=*), INTENT(IN)  :: FILENAME           ! Bpch file to read
      CHARACTER(LEN=*), INTENT(IN)  :: CATEGORY_IN        ! Diag. category name
      INTEGER,          INTENT(IN)  :: TRACER_IN          ! Tracer index #
      REAL*8,           INTENT(IN)  :: TAUO_IN            ! TAU timestamp
      INTEGER,          INTENT(IN)  :: IX, JX, LX          ! Dimensions of ARRAY
      LOGICAL, OPTIONAL, INTENT(IN) :: QUIET              ! Don't print output

```

**OUTPUT PARAMETERS:**

```

      REAL*4,          INTENT(OUT) :: ARRAY(IX,JX,LX)      ! Data array from file

```

**REVISION HISTORY:**

- (1 ) Assumes that we are reading in a global-size data block.
  - (2 ) Trap all I/O errors with subroutine IOERROR.F.
  - (3 ) Now stop with an error message if no matches are found. (bmy, 3/9/00)
  - (4 ) Added this routine to "bpch\_mod.f" (bmy, 6/28/00)
  - (5 ) Use IOS /= 0 criterion to also check for EOF condition (bmy, 9/12/00)
  - (6 ) TEMPARRAY now dimensioned to be of global size (bmy, 10/12/00)
  - (7 ) Removed obsolete code from 10/12/00 (bmy, 12/18/00)
  - (8 ) Now make TEMPARRAY independent of F77\_CMN\_SIZE parameters (bmy, 4/17/01)
  - (9 ) Removed old commented-out code (bmy, 4/20/01)
  - (10) Now reference IU\_FILE and IOERROR from "file\_mod.f". Now call  
       OPEN\_BPCH2\_FOR\_READ to open the binary punch file. Now use IU\_FILE  
       as the unit number instead of a locally-defined IUNIT. (bmy, 7/30/02)
  - (11) Now references ERROR\_STOP from "error\_mod.f" (bmy, 10/15/02)
  - (12) Now set IFIRST=1, JFIRST=1 for 1x1 nested grids. Now needs to  
       reference "define.h". Added OPTIONAL QUIET flag. (bmy, 3/14/03)
  - (13) Now separate off nested grid code in an #ifdef block using  
       NESTED\_CH or NESTED\_NA cpp switches (bmy, 12/1/04)
  - (14) Make TEMPARRAY big enough for GEOS-5 72 levels (and 73 edges)  
       (bmy, 2/15/07)
  - (15) Make TEMPARRAY large enough for 0.5 x 0.666 arrays -- but only if we  
       are doing a 0.5 x 0.666 nested simulation. (yxw, dan, bmy, 11/6/08)
  - 20 Nov 2009 - R. Yantosca - Added ProTeX header
  - 18 Dec 2009 - Aaron van D - Add NESTED\_EU flag
-

### 1.1.6 `get_modelname`

Function `GET_MODELNAME` returns the proper value of `MODELNAME` for current GEOS or GCAP met field type. `MODELNAME` is written to the binary punch file and is also used by the GAMAP package.

#### INTERFACE:

```
FUNCTION GET_MODELNAME() RESULT( MODELNAME )
```

#### USES:

```
USE CMN_SIZE_MOD
```

#### RETURN VALUE:

```
CHARACTER(LEN=20) :: MODELNAME    ! Model name for the current met field
```

#### REVISION HISTORY:

- (1 ) Now use special model name for GEOS-3 w/ 30 layers (bmy, 10/9/01)
- (2 ) Added modelname for GEOS-4/fvDAS model type (bmy, 11/20/01)
- (3 ) Added "GEOS4\_30L" for reduced GEOS-4 grid. Also now use C-preprocessor switch "GRID30LEV" instead of IF statements. (bmy, 11/3/03)
- (4 ) Updated for GCAP and GEOS-5 met fields. Rearranged coding for simplicity. (swu, bmy, 5/24/05)
- (5 ) Remove support for GEOS-1 and GEOS-STRAT met fields (bmy, 8/4/06)
- (6 ) Rename GRID30LEV to GRIDREDUCED (bmy, 2/7/07)
- 20 Nov 2009 - R. Yantosca - Added ProTeX header
- 13 Aug 2010 - R. Yantosca - Added MERRA model names

### 1.1.7 `get_name_ext`

Function `GET_NAME_EXT` returns the proper filename extension the current GEOS-Chem met field type (e.g. "geos3", "geos4", "geos5", or "gcap").

#### INTERFACE:

```
FUNCTION GET_NAME_EXT() RESULT( NAME_EXT )
```

#### USES:

```
#    include "define.h"
```

#### RETURN VALUE:

```
#if    defined( GEOS_3 )
      CHARACTER(LEN=5) :: NAME_EXT
      NAME_EXT = 'geos3'

#elif  defined( GEOS_4 )
      CHARACTER(LEN=5) :: NAME_EXT
```

```

NAME_EXT = 'geos4'

#elif defined( GEOS_5 )
  CHARACTER(LEN=5) :: NAME_EXT
  NAME_EXT = 'geos5'

#elif defined( GCAP )
  CHARACTER(LEN=4) :: NAME_EXT
  NAME_EXT = 'gcap'

#elif defined( MERRA )
  CHARACTER(LEN=5) :: NAME_EXT
  NAME_EXT = 'geos5'

#endif

```

**REVISION HISTORY:**

- (1 ) Added name string for GEOS-4/fvDAS model type (bmy, 11/20/01)
  - (2 ) Remove obsolete "geos2" model name string (bmy, 11/3/03)
  - (3 ) Modified for GCAP and GEOS-5 met fields (bmy, 5/24/05)
  - (4 ) Remove support for GEOS-1 and GEOS-STRAT met fields (bmy, 8/4/06)
  - 20 Nov 2009 - R. Yantosca - Added ProTeX header
  - 13 Aug 2010 - R. Yantosca - MERRA uses "geos5" name extension since its grid is identical to GEOS-5.
- 

**1.1.8 get\_name\_ext\_2d**

Function GET\_NAME\_EXT\_2D returns the proper filename extension for CTM model name for files which do not contain any vertical information (i.e. "geos" or "gcap").

**INTERFACE:**

```
FUNCTION GET_NAME_EXT_2D() RESULT( NAME_EXT_2D )
```

**RETURN VALUE:**

```
CHARACTER(LEN=4) :: NAME_EXT_2D
```

**REVISION HISTORY:**

- (1 ) Added name string for GEOS-4/fvDAS model type (bmy, 11/20/01)
  - (2 ) Remove obsolete "geos2" model name string (bmy, 11/3/03)
  - (3 ) Modified for GCAP and GEOS-5 met fields (bmy, 5/24/05)
  - (4 ) Remove support for GEOS-1 and GEOS-STRAT met fields (bmy, 8/4/06)
  - 20 Nov 2009 - R. Yantosca - Added ProTeX header
-

### 1.1.9 `get_res_ext`

Function `GET_RES_EXT` returns the proper filename extension for the GEOS-Chem horizontal grid resolution (e.g. "1x1", "2x25", "4x5").

#### INTERFACE:

```
FUNCTION GET_RES_EXT() RESULT( RES_EXT )
```

#### USES:

```
#    include "define.h"
```

#### RETURN VALUE:

```
#if    defined( GRID4x5 )
      CHARACTER(LEN=3) :: RES_EXT
      RES_EXT = '4x5'

#elif  defined( GRID2x25 )
      CHARACTER(LEN=4) :: RES_EXT
      RES_EXT = '2x25'

#elif  defined( GRID1x125 )
      CHARACTER(LEN=5) :: RES_EXT
      RES_EXT = '1x125'

#elif  defined( GRID1x1 )
      CHARACTER(LEN=3) :: RES_EXT
      RES_EXT = '1x1'

#elif  defined( GRID05x0666 )
      CHARACTER(LEN=7) :: RES_EXT
      RES_EXT = '05x0666'

#endif
```

#### REVISION HISTORY:

```
(1 ) Added extension for 1 x 1.25 grid (bmy, 12/1/04)
(2 ) Added extension for 0.5 x 0.666 grid (yxw, dan, bmy, 11/6/08)
20 Nov 2009 - R. Yantosca - Added ProTeX header
```

---

### 1.1.10 `get_halfpolar`

Function `GET_HALFPOLAR` returns 1 if the current grid has half-sized polar boxes (e.g. GEOS), or zero otherwise (e.g. GCAP).

#### INTERFACE:

```
FUNCTION GET_HALFPOLAR() RESULT( HALFPOLAR )
```

**USES:**

```
#      include "define.h"
```

**RETURN VALUE:**

```
INTEGER :: HALFPOLAR  ! =1 if we have half-sized polar boxes, =0 if not
```

**REVISION HISTORY:**

```
28 Jun 2005 - S. Wu & R. Yantosca - Initial version
20 Nov 2009 - R. Yantosca          - Added ProTeX header
```

**1.1.11 get\_tau0\_6a**

Function GET\_TAU0\_6A returns the corresponding TAU0 value for the first day of a given MONTH of a given YEAR. This is necessary to index monthly mean binary punch files, which are used as input to GEOS-Chem.

This function takes 3 mandatory arguments (MONTH, DAY, YEAR) and 3 optional arguments (HOUR, MIN, SEC). It is intended to replace the current 2-argument version of GET\_TAU0. The advantage being that GET\_TAU0\_6A can compute a TAU0 for any date and time in the GEOS-Chem epoch, rather than just the first day of each month. Overload this w/ an interface so that the user can also choose the version of GET\_TAU0 w/ 2 arguments (MONTH, YEAR), which is the prior version.

**INTERFACE:**

```
FUNCTION GET_TAU0_6A( MONTH, DAY, YEAR,
&                   HOUR, MIN, SEC ) RESULT( THIS_TAU0 )
```

**USES:**

```
USE ERROR_MOD, ONLY : ERROR_STOP
USE JULDAY_MOD, ONLY : JULDAY
```

**INPUT PARAMETERS:**

```
INTEGER, INTENT(IN)           :: MONTH
INTEGER, INTENT(IN)           :: DAY
INTEGER, INTENT(IN)           :: YEAR
INTEGER, INTENT(IN), OPTIONAL :: HOUR
INTEGER, INTENT(IN), OPTIONAL :: MIN
INTEGER, INTENT(IN), OPTIONAL :: SEC
```

**RETURN VALUE:**

```
REAL*8                       :: THIS_TAU0  ! TAU0 timestamp
```

**REMARKS:**

TAU0 is hours elapsed since 00:00 GMT on 01 Jan 1985.

## REVISION HISTORY:

- (1 ) 1985 is the first year of the GEOS epoch.
  - (2 ) Add TAU0 values for years 1985-2001 (bmy, 8/1/00)
  - (3 ) Correct error for 1991 TAU values. Also added 2002 and 2003.  
(bnd, bmy, 1/4/01)
  - (4 ) Updated comments (bmy, 9/26/01)
  - (5 ) Now references JULDAY from "julday\_mod.f" (bmy, 11/20/01)
  - (6 ) Now references ERROR\_STOP from "error\_mod.f" (bmy, 10/15/02)
- 20 Nov 2009 - R. Yantosca - Added ProTeX header
- 

## 1.2 Fortran: Module Interface charpak\_mod.f

Module CHARPAK\_MOD contains routines from the CHARPAK string and character manipulation package used by GEOS-Chem.

## INTERFACE:

```
MODULE CHARPAK_MOD
```

## USES:

```
IMPLICIT NONE
#   include "define.h"
PRIVATE
```

## PUBLIC MEMBER FUNCTIONS:

```
PUBLIC  :: CNTMAT
PUBLIC  :: COPYTXT
PUBLIC  :: CSTRIP
PUBLIC  :: ISDIGIT
PUBLIC  :: STRREPL
PUBLIC  :: STRSPLIT
PUBLIC  :: STRSQUEEZE
PUBLIC  :: TRANLC
PUBLIC  :: TRANUC
PUBLIC  :: TXT2INUM
PUBLIC  :: TXTEXT
```

## REMARKS:

CHARPAK routines by Robert D. Stewart, 1992. Subsequent modifications made for GEOS-CHEM by Bob Yantosca (1998, 2002, 2004).

## REVISION HISTORY:

- (1 ) Moved "cntmat.f", "copytxt.f", "cstrip.f", "fillstr.f", "txt2inum.f", "txttext.f", into this F90 module for easier bookkeeping (bmy, 10/15/01)
  - (2 ) Moved "tranuc.f" into this F90 module (bmy, 11/15/01)
  - (3 ) Now divide module header into MODULE PRIVATE, MODULE VARIABLES, and MODULE ROUTINES sections. Updated comments (bmy, 5/28/02)
  - (4 ) Wrote a new file "strrepl.f", which replaces a character pattern within a string with replacement text. Moved "tranlc.f" into this module. Replaced calls to function LENTRIM with F90 intrinsic function LEN\_TRIM. Removed function FILLSTR and replaced it w/ F90 intrinsic REPEAT. (bmy, 6/25/02)
  - (5 ) Added routine STRSPLIT as a wrapper for TXTEXT. Also added routines STRREPL and STRSQUEEZE. (bmy, 7/30/02)
  - (6 ) Added function ISDIGIT. Also replace LEN\_TRIM with LEN in routine STRREPL, to allow us to replace tabs w/ spaces. (bmy, 7/20/04)
- 20 Nov 2009 - R. Yantosca - Added ProTeX header
- 

### 1.3 Fortran: Module Interface *directory\_mod.f*

Module *DIRECTORY\_MOD* contains the directory path variables used by GEOS-Chem.

#### INTERFACE:

```
MODULE DIRECTORY_MOD
```

#### USES:

```
IMPLICIT NONE
#   include "define.h"
PUBLIC
```

#### PUBLIC DATA MEMBERS:

```
CHARACTER(LEN=255) :: DATA_DIR      ! Main DAO met field directory
CHARACTER(LEN=255) :: DATA_DIR_1x1 ! Root data dir for 1x1 emissions
CHARACTER(LEN=255) :: GCAP_DIR       ! Subdir for GCAP met data
CHARACTER(LEN=255) :: GEOS_1_DIR     ! !%%% OBSOLETE %%
CHARACTER(LEN=255) :: GEOS_S_DIR     ! !%%% OBSOLETE
CHARACTER(LEN=255) :: GEOS_3_DIR     ! Subdir for GEOS-3 met data
CHARACTER(LEN=255) :: GEOS_4_DIR     ! Subdir for GEOS-4 met data
CHARACTER(LEN=255) :: GEOS_5_DIR     ! Subdir for GEOS-5 met data
CHARACTER(LEN=255) :: MERRA_DIR      ! Subdir for MERRA met data
CHARACTER(LEN=255) :: TEMP_DIR       ! Temp dir for unzipping met data
CHARACTER(LEN=255) :: RUN_DIR        ! Run directory for GEOS-Chem
CHARACTER(LEN=255) :: OH_DIR         ! Dir w/ mean OH files
CHARACTER(LEN=255) :: O3PL_DIR       ! Dir w/ archived O3 P/L rate files
CHARACTER(LEN=255) :: TPBC_DIR       ! Dir w/ TPCORE boundary conditions
```



```

CHARACTER(LEN=255) :: TPBC_DIR_NA    ! Dir w/ TPCORE BC's for NA nest grid
CHARACTER(LEN=255) :: TPBC_DIR_EU    ! Dir w/ TPCORE BC's for EU nest grid
CHARACTER(LEN=255) :: TPBC_DIR_CH    ! Dir w/ TPCORE BC's for CH nest grid

```

## REVISION HISTORY:

```

20 Jul 2004 - R. Yantosca - Initial version
25 May 2005 - R. Yantosca - Added variables GCAP_DIR and GEOS_5_DIR
24 Oct 2005 - R. Yantosca - Added DATA_DIR_1x1
20 Nov 2009 - R. Yantosca - Added ProTeX header
18 Dec 2009 - Aaron van D - Added TPBC_DIR_NA, TPBC_DIR_EU, TPBC_DIR_CH
13 Aug 2010 - R. Yantosca - Added MERRA_DIR for MERRA met fields

```

---

## 1.4 Fortran: Module Interface error\_mod.f

Module ERROR\_MOD contains error checking routines.

### INTERFACE:

```

MODULE ERROR_MOD

```

### USES:

```

IMPLICIT NONE
#   include "define.h"
PRIVATE

```

### PUBLIC MEMBER FUNCTIONS:

```

PUBLIC  :: ALLOC_ERR
PUBLIC  :: CHECK_VALUE
PUBLIC  :: DEBUG_MSG
PUBLIC  :: ERROR_STOP
PUBLIC  :: GEOS_CHEM_STOP
PUBLIC  :: IS_SAFE_DIV
PUBLIC  :: IS_SAFE_EXP
PUBLIC  :: IT_IS_NAN
PUBLIC  :: IT_IS_FINITE
PUBLIC  :: SAFE_DIV
PUBLIC  :: SAFE_EXP
PUBLIC  :: SAFE_LOG
PUBLIC  :: SAFE_LOG10

! Interface for NaN-check routines
INTERFACE IT_IS_NAN
    MODULE PROCEDURE NAN_FLOAT
    MODULE PROCEDURE NAN_DBLE
END INTERFACE

```

```

! Interface for finite-check routines
INTERFACE IT_IS_FINITE
  MODULE PROCEDURE FINITE_FLOAT
  MODULE PROCEDURE FINITE_DBLE
END INTERFACE

! Interface for check-value routines
INTERFACE CHECK_VALUE
  MODULE PROCEDURE CHECK_REAL_VALUE
  MODULE PROCEDURE CHECK_DBLE_VALUE
END INTERFACE

```

#### PRIVATE MEMBER FUNCTIONS:

```

PRIVATE :: CHECK_DBLE_VALUE
PRIVATE :: CHECK_REAL_VALUE
PRIVATE :: FINITE_DBLE
PRIVATE :: FINITE_FLOAT
PRIVATE :: NAN_DBLE
PRIVATE :: NAN_FLOAT

```

#### REVISION HISTORY:

- 08 Mar 2001 - R. Yantosca - Initial version
- (1 ) Added subroutines CHECK\_REAL\_VALUE and CHECK\_DBLE\_VALUE, which are overloaded by interface CHECK\_VALUE. This is a convenience so that you don't have to always call IT\_IS\_NAN directly. (bmy, 6/13/01)
  - (2 ) Updated comments (bmy, 9/4/01)
  - (3 ) Now use correct values for bit masking in FINITE\_FLOAT for the ALPHA platform (bmy, 11/15/01)
  - (4 ) Now divide module header into MODULE PRIVATE, MODULE VARIABLES, and MODULE ROUTINES sections. Also add MODULE INTERFACES section, since we have an interface here. (bmy, 5/28/02)
  - (5 ) Add NaN and infinity error checking for Linux platform (bmy, 3/22/02)
  - (6 ) Added routines ERROR\_STOP, GEOS\_CHEM\_STOP, and ALLOC\_ERR to this module. Also improved CHECK\_STT. (bmy, 11/27/02)
  - (7 ) Minor bug fixes in FORMAT statements. Renamed cpp switch from DEC\_COMPAQ to COMPAQ. Also added code to trap errors on SUN platform. (bmy, 3/21/03)
  - (8 ) Added patches for IBM/AIX platform (gcc, bmy, 6/27/03)
  - (9 ) Bug fixes for LINUX platform (bmy, 9/29/03)
  - (10) Now supports INTEL\_FC compiler (bmy, 10/24/03)
  - (11) Changed the name of some cpp switches in "define.h" (bmy, 12/2/03)
  - (12) Minor fix for LINUX\_IFC and LINUX\_EFC (bmy, 1/24/04)
  - (13) Do not flush buffer for LINUX\_EFC in ERROR\_STOP (bmy, 4/6/04)
  - (14) Move CHECK\_STT routine to "tracer\_mod.f" (bmy, 7/20/04)
  - (15) Added LINUX\_IFORT switch for Intel v8 and v9 compilers (bmy, 10/18/05)

(16) Now print IFORT error messages for Intel v8/v9 compiler (bmy, 11/30/05)  
 (17) Cosmetic change in DEBUG\_MSG (bmy, 4/10/06)  
 (18) Remove support for LINUX\_IFC and LINUX\_EFC compilers (bmy, 8/4/06)  
 (19) Now use intrinsic functions for IFORT, remove C routines (bmy, 8/14/07)  
 (20) Added routine SAFE\_DIV (phs, bmy, 2/26/08)  
 (21) Added routine IS\_SAFE\_DIV (phs, bmy, 6/11/08)  
 (22) Updated routine SAFE\_DIV (phs, 4/14/09)  
 (23) Remove support for SGI, COMPAQ compilers (bmy, 7/8/09)  
 20 Nov 2009 - R. Yantosca - Added ProTeX header  
 04 Jan 2010 - R. Yantosca - Added SAFE\_EXP and IS\_SAFE\_EXP functions  
 04 Jan 2010 - R. Yantosca - Added SAVE\_LOG and SAFE\_LOG10 functions

---

### 1.4.1 nan\_float

Function NAN\_FLOAT returns TRUE if a REAL\*4 number is equal to the IEEE NaN (Not-a-Number) flag. Returns FALSE otherwise.

#### INTERFACE:

```
FUNCTION NAN_FLOAT( VALUE ) RESULT( IT_IS_A_NAN )
```

#### USES:

```
#    include "define.h"

#if    defined( IBM_AIX ) || defined( IBM_XLF )
    USE IEEE_ARITHMETIC
#endif
```

#### INPUT PARAMETERS:

```
REAL*4, INTENT(IN) :: VALUE          ! Value to be tested for NaN
```

#### RETURN VALUE:

```
LOGICAL              :: IT_IS_A_NAN  ! =T if VALUE is NaN; =F otherwise
```

#### REVISION HISTORY:

(1 ) Is overloaded by interface "IT\_IS\_NAN".  
 (2 ) Now call C routine is\_nan(x) for Linux platform (bmy, 6/13/02)  
 (3 ) Eliminate IF statement in Linux section. Also now trap NaN on the Sun/Sparc platform. Rename cpp switch from DEC\_COMPAQ to COMPAQ. (bmy, 3/23/03)  
 (4 ) Added patches for IBM/AIX platform (gcc, bmy, 6/27/03)  
 (5 ) Use LINUX error-trapping for INTEL\_FC (bmy, 10/24/03)  
 (6 ) Renamed SGI to SGI\_MIPS, LINUX to LINUX\_PGI, INTEL\_FC to INTEL\_IFC, and added LINUX\_EFC. (bmy, 12/2/03)  
 (7 ) Added LINUX\_IFORT switch for Intel v8 and v9 compilers (bmy, 10/18/05)  
 (8 ) Remove support for LINUX\_IFC & LINUX\_EFC compilers (bmy, 8/4/06)

- (9 ) Now use ISNAN for Linux/IFORT compiler (bmy, 8/14/07)
  - (10) Remove support for SGI, COMPAQ compilers. Add IBM\_XLF switch.  
(bmy, 7/8/09)
- 20 Nov 2009 - R. Yantosca - Added ProTeX header

### 1.4.2 nan\_dble

Function NAN\_DBLE returns TRUE if a REAL\*8 number is equal to the IEEE NaN (Not-a-Number) flag. Returns FALSE otherwise.

#### INTERFACE:

```
FUNCTION NAN_DBLE( VALUE ) RESULT( IT_IS_A_NAN )
```

#### USES:

```
#      include "define.h"

#if    defined( IBM_AIX ) || defined( IBM_XLF )
      USE IEEE_ARITHMETIC
#endif
```

#### INPUT PARAMETERS:

```
REAL*8, INTENT(IN) :: VALUE          ! Value to be tested for NaN
```

#### RETURN VALUE:

```
LOGICAL              :: IT_IS_A_NAN  ! =T if VALUE is NaN; =F otherwise
```

#### REVISION HISTORY:

- (1 ) Is overloaded by interface "IT\_IS\_NAN".
  - (2 ) Now call C routine is\_nan(x) for Linux platform (bmy, 6/13/02)
  - (3 ) Eliminate IF statement in Linux section. Also now trap NaN on the Sun/Sparc platform. Rename cpp switch from DEC\_COMPAQ to COMPAQ. (bmy, 3/23/03)
  - (4 ) Added patches for IBM/AIX platform (gcc, bmy, 6/27/03)
  - (5 ) Use LINUX error-trapping for INTEL\_FC (bmy, 10/24/03)
  - (6 ) Renamed SGI to SGI\_MIPS, LINUX to LINUX\_PGI, INTEL\_FC to INTEL\_IFC, and added LINUX\_EFC. (bmy, 12/2/03)
  - (7 ) Added LINUX\_IFORT switch for Intel v8 and v9 compilers (bmy, 10/18/05)
  - (8 ) Remove support for LINUX\_IFC & LINUX\_EFC compilers (bmy, 8/4/06)
  - (9 ) Now use ISNAN for Linux/IFORT compiler (bmy, 8/14/07)
  - (10) Remove support for SGI, COMPAQ compilers. Add IBM\_XLF switch.  
(bmy, 7/8/09)
- 20 Nov 2009 - R. Yantosca - Added ProTeX header

### 1.4.3 finite\_float

Function FINITE\_FLOAT returns FALSE if a REAL\*4 number is equal to the IEEE Infinity flag. Returns TRUE otherwise.

#### INTERFACE:

```
FUNCTION FINITE_FLOAT( VALUE ) RESULT( IT_IS_A_FINITE )
```

#### USES:

```
#    include "define.h"

#if    defined( IBM_AIX ) || defined( IBM_XLF )
    USE IEEE_ARITHMETIC
#endif
```

#### INPUT PARAMETERS:

```
REAL*4, INTENT(IN) :: VALUE           ! Value to be tested for infinity
```

#### RETURN VALUE:

```
LOGICAL              :: IT_IS_A_FINITE ! =T if VALUE is finite; =F else
```

#### REVISION HISTORY:

- (1 ) Is overloaded by interface "IT\_IS\_FINITE".
  - (2 ) Now use correct values for bit masking (bmy, 11/15/01)
  - (3 ) Eliminate IF statement in Linux section. Also now trap Infinity on the Sun/Sparc platform. Rename cpp switch from DEC\_COMPAQ to COMPAQ. (bmy, 3/23/03)
  - (4 ) Added patches for IBM/AIX platform (gcc, bmy, 6/27/03)
  - (5 ) Bug fix: now use external C IS\_FINITE for PGI/Linux (bmy, 9/29/03)
  - (6 ) Use LINUX error-trapping for INTEL\_FC (bmy, 10/24/03)
  - (7 ) Renamed SGI to SGI\_MIPS, LINUX to LINUX\_PGI, INTEL\_FC to INTEL\_IFC, and added LINUX\_EFC. (bmy, 12/2/03)
  - (8 ) Added LINUX\_IFORT switch for Intel v8 and v9 compilers (bmy, 10/18/05)
  - (9 ) Remove support for LINUX\_IFC & LINUX\_EFC compilers (bmy, 8/4/06)
  - (10) Now use FP\_CLASS for IFORT compiler (bmy, 8/14/07)
  - (11) Remove support for SGI, COMPAQ compilers. Add IBM\_XLF switch. (bmy, 7/8/09)
- 20 Nov 2009 - R. Yantosca - Added ProTeX header

### 1.4.4 finite\_double

Function FINITE\_FLOAT returns FALSE if a REAL\*8 number is equal to the IEEE Infinity flag. Returns TRUE otherwise.

#### INTERFACE:

```
FUNCTION FINITE_DBLE( VALUE ) RESULT( IT_IS_A_FINITE )
```

**USES:**

```
#    include "define.h"

#if    defined( IBM_AIX ) || defined( IBM_XLF )
    USE IEEE_ARITHMETIC
#endif
```

**INPUT PARAMETERS:**

```
REAL*8, INTENT(IN) :: VALUE          ! Value to be tested for infinity
```

**RETURN VALUE:**

```
LOGICAL              :: IT_IS_A_FINITE  ! =T if VALUE is finite; =F else
```

**REVISION HISTORY:**

- (1 ) Is overloaded by interface "IT\_IS\_FINITE".
  - (2 ) Now use correct values for bit masking (bmy, 11/15/01)
  - (3 ) Eliminate IF statement in Linux section. Also now trap Infinity on the Sun/Sparc platform. Rename cpp switch from DEC\_COMPAQ to COMPAQ. (bmy, 3/23/03)
  - (4 ) Added patches for IBM/AIX platform (gcc, bmy, 6/27/03)
  - (5 ) Bug fix: now use external C IS\_FINITE for PGI/Linux (bmy, 9/29/03)
  - (6 ) Use LINUX error-trapping for INTEL\_FC (bmy, 10/24/03)
  - (7 ) Renamed SGI to SGI\_MIPS, LINUX to LINUX\_PGI, INTEL\_FC to INTEL\_IFC, and added LINUX\_EFC. (bmy, 12/2/03)
  - (8 ) Added LINUX\_IFORT switch for Intel v8 and v9 compilers (bmy, 10/18/05)
  - (9 ) Remove support for LINUX\_IFC & LINUX\_EFC compilers (bmy, 8/4/06)
  - (10) Now use FP\_CLASS for IFORT compiler (bmy, 8/14/07)
  - (11) Remove support for SGI, COMPAQ compilers. Add IBM\_XLF switch. (bmy, 7/8/09)
- 20 Nov 2009 - R. Yantosca - Added ProTeX header

**1.4.5 check\_real\_value**

Subroutine CHECK\_REAL\_VALUE checks to make sure a REAL\*4 value is not NaN or Infinity. This is a wrapper for the interfaces IT\_IS\_NAN and IT\_IS\_FINITE.

**INTERFACE:**

```
SUBROUTINE CHECK_REAL_VALUE( VALUE, LOCATION, VARNAME, MESSAGE )
```

**INPUT PARAMETERS:**

```
REAL*4,              INTENT(IN) :: VALUE          ! Value to be checked
CHARACTER(LEN=*), INTENT(IN) :: VARNAME          ! Name of variable
CHARACTER(LEN=*), INTENT(IN) :: MESSAGE          ! Short descriptive msg
INTEGER,              INTENT(IN) :: LOCATION(4)  ! (/ I, J, L, N /) indices
```

**REVISION HISTORY:**

13 Jun 2001 - R. Yantosca - Initial version  
 15 Oct 2002 - R. Yantosca - Now call GEOS\_CHEM\_STOP to shutdown safely  
 15 Oct 2002 - R. Yantosca - Updated comments, cosmetic changes  
 20 Nov 2009 - R. Yantosca - Added ProTeX header

---

**1.4.6 check\_double\_value**

Subroutine CHECK\_DOUBLE\_VALUE checks to make sure a REAL\*4 value is not NaN or Infinity. This is a wrapper for the interfaces IT\_IS\_NAN and IT\_IS\_FINITE.

**INTERFACE:**

```
SUBROUTINE CHECK_DOUBLE_VALUE( VALUE, LOCATION, VARNAME, MESSAGE )
```

**INPUT PARAMETERS:**

```
REAL*8,          INTENT(IN) :: VALUE          ! Value to be checked
CHARACTER(LEN=*), INTENT(IN) :: VARNAME        ! Name of variable
CHARACTER(LEN=*), INTENT(IN) :: MESSAGE        ! Short descriptive msg
INTEGER,          INTENT(IN) :: LOCATION(4)    ! (/ I, J, L, N /) indices
```

**REVISION HISTORY:**

13 Jun 2001 - R. Yantosca - Initial version  
 15 Oct 2002 - R. Yantosca - Now call GEOS\_CHEM\_STOP to shutdown safely  
 15 Oct 2002 - R. Yantosca - Updated comments, cosmetic changes  
 20 Nov 2009 - R. Yantosca - Added ProTeX header

---

**1.4.7 error\_stop**

Subroutine ERROR\_STOP is a wrapper for GEOS\_CHEM\_STOP. It prints an error message then calls GEOS\_CHEM\_STOP to free memory and quit.

**INTERFACE:**

```
SUBROUTINE ERROR_STOP( MESSAGE, LOCATION )
```

**INPUT PARAMETERS:**

```
CHARACTER(LEN=*), INTENT(IN) :: MESSAGE        ! Error msg to print
CHARACTER(LEN=*), INTENT(IN) :: LOCATION        ! Where ERROR_STOP is called
```

**REVISION HISTORY:**

15 Oct 2002 - R. Yantosca - Initial version  
 20 Nov 2009 - R. Yantosca - Added ProTeX header

---

### 1.4.8 geos\_chem\_stop

Subroutine GEOS\_CHEM\_STOP calls CLEANUP to deallocate all module arrays and then stops the run.

#### INTERFACE:

```
SUBROUTINE GEOS_CHEM_STOP
```

#### USES:

```
#    include "define.h"
```

#### REVISION HISTORY:

```
15 Oct 2002 - R. Yantosca - Initial version
20 Nov 2009 - R. Yantosca - Now EXIT works for LINUX_IFC, LINUX_EFC,
                           so remove #if block.
20 Nov 2009 - R. Yantosca - Added ProTeX header
```

---

### 1.4.9 alloc\_err

Subroutine ALLOC\_ERR prints an error message if there is not enough memory to allocate a particular allocatable array.

#### INTERFACE:

```
SUBROUTINE ALLOC_ERR( ARRAYNAME, AS )
```

#### USES:

```
#    include "define.h"
```

#### INPUT PARAMETERS:

```
CHARACTER(LEN=*), INTENT(IN) :: ARRAYNAME ! Name of array
INTEGER, OPTIONAL, INTENT(IN) :: AS       ! Error output from "STAT"
```

#### REVISION HISTORY:

```
26 Jun 2000 - R. Yantosca - Initial version, split off from "ndxx_setup.f"
15 Oct 2002 - R. Yantosca - Added to "error_mod.f"
30 Nov 2005 - R. Yantosca - Call IFORT_ERRMSG for Intel Fortran compiler
20 Nov 2009 - R. Yantosca - Added ProTeX header
```

---



Subroutine `DEBUG_MSG` prints a message to the stdout buffer and flushes. This is useful for determining the exact location where errors occur.

## SUBROUTINE DEBUG\_MSG( MESSAGE )

```
#include "define.h"
```

```
CHARACTER(LEN=*) , INTENT(IN) :: MESSAGE    ! Message to print
```

```

07 Jan 2002 - R. Yantosca - Initial version
(1 ) Now just write the message and flush the buffer (bmy, 7/5/01)
(2 ) Renamed from "paftop.f" to "debug_msg.f" (bmy, 1/7/02)
(3 ) Bundled into "error_mod.f" (bmy, 11/22/02)
(4 ) Now do not FLUSH the buffer for EFC compiler (bmy, 4/6/04)
(5 ) Now add a little space for debug output (bmy, 4/10/06)
(6 ) Remove support for LINUX_IFC & LINUX_EFC compilers (bmy, 8/4/06)
20 Nov 2009 - R. Yantosca - Added ProTeX header

```

Function `SAFE_DIV` performs "safe division", that is to prevent overflow, underflow, NaN, or infinity errors. An alternate value is returned if the division cannot be performed.

```

FUNCTION SAFE_DIV( N,          D,
&                  ALT_NAN,  ALT_OVER,
&                  ALT_UNDER ) RESULT( Q )

```

[illegible]

```

REAL*8, OPTIONAL, INTENT(IN) :: ALT_UNDER
! is ALT_NAN)
! Alternate value to be
! returned if the division
! leads to underflow
! (default is 0, but you
! could use TINY() if you
! want a non-zero result).

```

**RETURN VALUE:**

```

REAL*8 :: Q ! Output from the division

```

**REMARKS:**

For more information, see the discussion on:

[http://groups.google.com/group/comp.lang.fortran/browse\\_thread/thread/8b367f44c419fa1d/](http://groups.google.com/group/comp.lang.fortran/browse_thread/thread/8b367f44c419fa1d/)

**REVISION HISTORY:**

26 Feb 2008 - P. Le Sager & R. Yantosca - Initial version

(1) Now can return different alternate values if NAN (that is 0/0), overflow (that is a too large number), or too small (that is greater than 0 but less than smallest possible number). Default value is zero in case of underflow (phs, 4/14/09)

(2) Some compiler options flush underflows to zero (-ftz for IFort). To think about it (phs, 4/14/09)

20 Nov 2009 - R. Yantosca - Added ProTeX header

**1.4.12 is\_safe\_div**

Function IS\_SAFE\_DIV tests for "safe division", that is check if the division will overflow/underflow or hold NaN. .FALSE. is returned if the division cannot be performed. (phs, 6/11/08)

**INTERFACE:**

```

FUNCTION IS_SAFE_DIV( N, D, R4 ) RESULT( F )

```

**INPUT PARAMETERS:**

```

REAL*8, INTENT(IN) :: N ! Numerator
REAL*8, INTENT(IN) :: D ! Denominator
LOGICAL, INTENT(IN), OPTIONAL :: R4 ! Logical flag to use the limits
! of REAL*4 to define underflow
! or overflow. Extra defensive.

```

**OUTPUT PARAMETERS:**

```

LOGICAL :: F ! =F if division isn't allowed
! =T otherwise

```

**REMARKS:**

UnderFlow, OverFlow and NaN are tested for. If you need to differentiate between the three, use the SAFE\_DIV (phs, 4/14/09)

**REVISION HISTORY:**

11 Jun 2008 - P. Le Sager - Initial version  
 20 Nov 2009 - R. Yantosca - Added ProTeX header

---

**1.4.13 safe\_exp**

Function SAFE\_EXP performs a "safe exponential", that is to prevent overflow, underflow, NaN, or infinity errors when taking the value EXP( x ). An alternate value is returned if the exponential cannot be performed.

**INTERFACE:**

```
FUNCTION SAFE_EXP( X, ALT ) RESULT( VALUE )
```

**INPUT PARAMETERS:**

```
REAL*8, INTENT(IN) :: X      ! Argument of EXP
REAL*8, INTENT(IN) :: ALT    ! Alternate value to be returned
```

**RETURN VALUE:**

```
REAL*8              :: VALUE ! Output from the exponential
```

**REVISION HISTORY:**

04 Jan 2010 - R. Yantosca - Initial version

---

**1.4.14 is\_safe\_exp**

Function IS\_SAFE\_EXP returns TRUE if it is safe to take the value EXP( x ) without encountering a floating point exception. FALSE is returned if the exponential cannot be performed.

**INTERFACE:**

```
FUNCTION IS_SAFE_EXP( X ) RESULT( F )
```

**INPUT PARAMETERS:**

```
REAL*8, INTENT(IN) :: X      ! Argument to the exponential function
```

**OUTPUT PARAMETERS:**

```
LOGICAL              :: F      ! =F if exponential isn't allowed
                          ! =T otherwise
```

**REMARKS:**

Empirical testing has revealed that  $-600 < X < 600$  will not result in a floating-point exception on Sun and IFORT compilers. This is good enough for most purposes.

**REVISION HISTORY:**

04 Jan 2010 - R. Yantosca - Initial version

---

**1.4.15 safe\_log**

Function SAFE\_LOG performs a "safe natural logarithm", that is to prevent overflow, underflow, NaN, or infinity errors when taking the value  $\text{LOG}(x)$ . An alternate value is returned if the logarithm cannot be performed.

**INTERFACE:**

```
FUNCTION SAFE_LOG( X, ALT ) RESULT( VALUE )
```

**INPUT PARAMETERS:**

```
REAL*8, INTENT(IN) :: X      ! Argument of LOG
REAL*8, INTENT(IN) :: ALT    ! Alternate value to be returned
```

**RETURN VALUE:**

```
REAL*8              :: VALUE ! Output from the natural logarithm
```

**REVISION HISTORY:**

04 Jan 2010 - R. Yantosca - Initial version

---

**1.4.16 safe\_log10**

Function SAFE\_LOG10 performs a "safe log10", that is to prevent overflow, underflow, NaN, or infinity errors when taking the value  $\text{LOG10}(x)$ . An alternate value is returned if the logarithm cannot be performed.

**INTERFACE:**

```
FUNCTION SAFE_LOG10( X, ALT ) RESULT( VALUE )
```

**INPUT PARAMETERS:**

```
REAL*8, INTENT(IN) :: X      ! Argument of LOG10
REAL*8, INTENT(IN) :: ALT    ! Alternate value to be returned
```

**RETURN VALUE:**

```
REAL*8              :: VALUE ! Output from the natural logarithm
```

**REVISION HISTORY:**

04 Jan 2010 - R. Yantosca - Initial version

---

## 1.5 Fortran: Module Interface file\_mod.f

Module FILE\_MOD contains file unit numbers, as well as file I/O routines for GEOS-Chem. FILE\_MOD keeps all of the I/O unit numbers in a single location for convenient access.

### INTERFACE:

```
MODULE FILE_MOD
```

### USES:

```
IMPLICIT NONE
#   include "define.h"
PRIVATE
```

### DEFINED PARAMETERS:

```
! Logical file unit numbers for ...
INTEGER, PUBLIC, PARAMETER :: IU_RST      = 1   ! Tracer restart file
INTEGER, PUBLIC, PARAMETER :: IU_CHEMDAT = 7   ! "chem.dat"
INTEGER, PUBLIC, PARAMETER :: IU_FASTJ    = 8   ! FAST-J input files
INTEGER, PUBLIC, PARAMETER :: IU_GEOS     = 10  ! "input.geos"
INTEGER, PUBLIC, PARAMETER :: IU_BPCH     = 11  ! "ctm.bpch"
INTEGER, PUBLIC, PARAMETER :: IU_ND20     = 12  ! "rate.YYYYMMDD"
INTEGER, PUBLIC, PARAMETER :: IU_ND48     = 13  ! ND48 output
INTEGER, PUBLIC, PARAMETER :: IU_ND49     = 14  ! "tsYYYYMMDD.bpch"
INTEGER, PUBLIC, PARAMETER :: IU_ND50     = 15  ! "ts24h.bpch"
INTEGER, PUBLIC, PARAMETER :: IU_ND51     = 16  ! "ts10_12am.bpch" etc.
INTEGER, PUBLIC, PARAMETER :: IU_ND51b    = 23  ! for ND51b diagnostic
INTEGER, PUBLIC, PARAMETER :: IU_ND52     = 17  ! ND52 output (NRT only)
INTEGER, PUBLIC, PARAMETER :: IU_PLANE    = 18  ! "plane.log"
INTEGER, PUBLIC, PARAMETER :: IU_BC       = 19  ! TPCORE BC files
INTEGER, PUBLIC, PARAMETER :: IU_BC_NA    = 20  ! TPCORE BC files: NA grid
INTEGER, PUBLIC, PARAMETER :: IU_BC_EU    = 21  ! TPCORE BC files: EU grid
INTEGER, PUBLIC, PARAMETER :: IU_BC_CH    = 22  ! TPCORE BC files: CH grid
INTEGER, PUBLIC, PARAMETER :: IU_FILE     = 65  ! Generic file
INTEGER, PUBLIC, PARAMETER :: IU_TP       = 69  ! "YYYYMMDD.tropp.*"
INTEGER, PUBLIC, PARAMETER :: IU_PH       = 70  ! "YYYYMMDD.phis.*"
INTEGER, PUBLIC, PARAMETER :: IU_I6       = 71  ! "YYYYMMDD.i6.*"
INTEGER, PUBLIC, PARAMETER :: IU_A6       = 72  ! "YYYYMMDD.a6.*"
INTEGER, PUBLIC, PARAMETER :: IU_A3       = 73  ! "YYYYMMDD.a3.*"
INTEGER, PUBLIC, PARAMETER :: IU_A1       = 74  ! "YYYYMMDD.a1.*"
INTEGER, PUBLIC, PARAMETER :: IU_GWET     = 75  ! "YYYYMMDD.gwet.*"
INTEGER, PUBLIC, PARAMETER :: IU_XT       = 76  ! "YYYYMMDD.xtra.*"
INTEGER, PUBLIC, PARAMETER :: IU_CN       = 77  ! "YYYYMMDD.cn.*"
INTEGER, PUBLIC, PARAMETER :: IU_SMV2LOG  = 93  ! "smv2.log"
INTEGER, PUBLIC, PARAMETER :: IU_DEBUG    = 98  ! Reserved for debugging
INTEGER, PUBLIC, PARAMETER :: IU_OAP      = 99  ! soaprod.YYYYMMDDhh
```

### PUBLIC MEMBER FUNCTIONS:

```

PUBLIC  :: CLOSE_FILES
PUBLIC  :: FILE_EXISTS
PUBLIC  :: IOERROR

INTERFACE FILE_EXISTS
  MODULE PROCEDURE FILE_EX_C
  MODULE PROCEDURE FILE_EX_I
END INTERFACE

```

#### PRIVATE MEMBER FUNCTIONS:

```

PRIVATE :: FILE_EX_C
PRIVATE :: FILE_EX_I

```

#### REVISION HISTORY:

- (1 ) Moved "ioerror.f" into this module. (bmy, 7/1/02)
- (2 ) Now references "error\_mod.f" (bmy, 10/15/02)
- (3 ) Renamed cpp switch from DEC\_COMPAQ to COMPAQ. Also added code to trap I/O errors on SUN/Sparc platform. (bmy, 3/23/03)
- (4 ) Now added IU\_BC for nested boundary conditions as unit 18 (bmy, 3/27/03)
- (5 ) Renamed IU\_CTMCHEM to IU\_SMV2LOG (bmy, 4/21/03)
- (6 ) Now print out I/O errors for IBM and INTEL\_FC compilers (bmy, 11/6/03)
- (7 ) Changed the name of some cpp switches in "define.h" (bmy, 12/2/03)
- (8 ) Renumbered the order of the files. Also removed IU\_INPTR and IU\_INPUT since they are now obsolete. (bmy, 7/20/04)
- (9 ) Added overloaded routines FILE\_EX\_C and FILE\_EX\_I (bmy, 3/23/05)
- (10) Added LINUX\_IFORT switch for Intel v8 & v9 compilers (bmy, 10/18/05)
- (11) Added IU\_XT for GEOS3 XTRA met fields files for MEGAN (tmf, 10/20/05)
- (12) Extra modification for Intel v9 compiler (bmy, 11/2/05)
- (13) Now print IFORT error messages (bmy, 11/30/05)
- (14) Remove support for LINUX\_IFC & LINUX\_EFC compilers (bmy, 8/4/06)
- (15) Remove support for SGI & COMPAQ compilers (bmy, 7/8/09)
- 20 Nov 2009 - R. Yantosca - Added ProTeX headers
- 18 Dec 2009 - Aaron van D - Added file units IU\_BC\_NA, IU\_BC\_EU, IU\_BC\_CH
- 15 Mar 2010 - D. Henze - Add IU\_OAP for SOA restart file.
- 19 Aug 2010 - R. Yantosca - Added IU\_CN and IU\_A1 parameters for MERRA
- 19 Aug 2010 - R. Yantosca - Remove IU\_KZZ

#### 1.5.1 ioerror

Subroutine IOERROR prints out I/O error messages. The error number, file unit, location, and a brief description will be printed, and program execution will be halted. (bmy, 5/28/99, 7/4/09)

#### INTERFACE:

```

SUBROUTINE IOERROR( ERROR_NUM, UNIT, LOCATION )

```

**USES:**

```
USE ERROR_MOD, ONLY : GEOS_CHEM_STOP
```

```
#      include "define.h"                                ! C-preprocessor switches
```

**INPUT PARAMETERS:**

```
INTEGER,          INTENT(IN) :: ERROR_NUM  ! I/O error from IOSTAT
INTEGER,          INTENT(IN) :: UNIT       ! Logical unit # for file
CHARACTER(LEN=*), INTENT(IN) :: LOCATION   ! Descriptive message
```

**REVISION HISTORY:**

- (1 ) Now flush the standard output buffer before stopping.  
Also updated comments. (bmy, 2/7/00)
- (2 ) Changed ROUTINE\_NAME to LOCATION. Now also use C-library routines  
gerror and strerror() to get the error string corresponding to  
ERROR\_NUM. For SGI platform, also print the command string that  
will call the SGI "explain" command, which will yield additional  
information about the error. Updated comments, cosmetic changes.  
Now also reference "define.h". (bmy, 3/21/02)
- (3 ) Moved into "file\_mod.f". Now reference GEOS\_CHEM\_STOP from module  
"error\_mod.f". Updated comments, cosmetic changes. (bmy, 10/15/02)
- (4 ) Renamed cpp switch from DEC\_COMPAQ to COMPAQ. Also added code to  
display I/O errors on SUN platform. (bmy, 3/23/03)
- (5 ) Now call GERROR for IBM and INTEL\_FC compilers (bmy, 11/6/03)
- (6 ) Renamed SGI to SGI\_MIPS, LINUX to LINUX\_PGI, INTEL\_FC to INTEL\_IFC,  
and added LINUX\_EFC. (bmy, 12/2/03)
- (7 ) Now don't flush the buffer for LINUX\_EFC (bmy, 4/23/04)
- (8 ) Modifications for Linux/IFORT Intel v9 compiler (bmy, 11/2/05)
- (9 ) Now call IFORT\_ERRMSG to get the IFORT error messages (bmy, 11/30/05)
- (10) Remove support for LINUX\_IFC & LINUX\_EFC compilers (bmy, 8/4/06)
- (10) Remove support for SGI & COMPAQ compilers. Add IBM\_XLF switch.  
(bmy, 7/8/09)
- 20 Nov 2009 - R. Yantosca - Removed commented-out code for SGI, COMPAQ
- 20 Nov 2009 - R. Yantosca - Added ProTeX header

**1.5.2 file\_ex\_c**

Function FILE\_EX\_C returns TRUE if FILENAME exists or FALSE otherwise. This is handled in a platform-independent way. The argument is of CHARACTER type.

**INTERFACE:**

```
FUNCTION FILE_EX_C( FILENAME ) RESULT( IT_EXISTS )
```

**USES:**

```
#      include "define.h"
```

**INPUT PARAMETERS:**

```
CHARACTER(LEN=*), INTENT(IN) :: FILENAME    ! Name of file or dir to test
```

**RETURN VALUE:**

```
LOGICAL                                :: IT_EXISTS    ! =T if the file/dir exists
```

**REMARKS:**

This routine is overloaded by public interface FILE\_EXISTS.

**REVISION HISTORY:**

```
23 Mar 2005 - R. Yantosca - Initial version
20 Nov 2009 - R. Yantosca - Updated for LINUX/IFORT Intel v9 compiler
20 Nov 2009 - R. Yantosca - Added ProTeX header
```

---

**1.5.3 file\_ex\_i**

Function FILE\_EX\_I returns TRUE if FILENAME exists or FALSE otherwise. This is handled in a platform-independent way. The argument is of INTEGER type.

**INTERFACE:**

```
FUNCTION FILE_EX_I( IUNIT ) RESULT( IT_EXISTS )
```

**USES:**

```
#    include "define.h"
```

**INPUT PARAMETERS:**

```
! Arguments
INTEGER, INTENT(IN) :: IUNIT    ! LUN of file to be tested
```

**RETURN VALUE:**

```
LOGICAL                                :: IT_EXISTS    ! =T if the file/dir exists
```

**REMARKS:**

This routine is overloaded by public interface FILE\_EXISTS.

**REVISION HISTORY:**

```
23 Mar 2005 - R. Yantosca - Initial version
20 Nov 2009 - R. Yantosca - Added ProTeX header
```

---



### 1.5.4 close\_files

Subroutine CLOSE\_FILES closes files used by GEOS-Chem. This should be called only from the end of the "main.f" program.

#### INTERFACE:

```
SUBROUTINE CLOSE_FILES
```

#### REVISION HISTORY:

```
04 Mar 1998 - R. Yantosca - Initial version
27 Jun 2002 - R. Yantosca - Moved into "file_mod.f"
27 Mar 2003 - R. Yantosca - Also close IU_BC
20 Jul 2004 - R. Yantosca - Removed obsolete IU_INPUT and IU_INPTR.
20 Jul 2004 - R. Yantosca - Also renamed IU_TS to IU_ND48.
20 Oct 2005 - R. Yantosca - Also close IU_XT.
20 Nov 2009 - R. Yantosca - Added ProTeX header
18 Dec 2009 - Aaron van D - Now close files IU_BC_NA, IU_BC_EU, IU_BC_CH
19 Aug 2010 - R. Yantosca - Remove IU_KZZ
19 Aug 2010 - R. Yantosca - Now close IU_A1
```

---

## 1.6 Fortran: Module Interface grid\_mod.f

Module GRID\_MOD contains variables and routines which are used to specify the parameters of a GEOS-CHEM horizontal grid.

#### INTERFACE:

```
MODULE GRID_MOD
```

#### USES:

```
IMPLICIT NONE
#   include "define.h"
PRIVATE
```

#### PUBLIC MEMBER FUNCTIONS:

```
PUBLIC  :: CLEANUP_GRID
PUBLIC  :: COMPUTE_GRID
PUBLIC  :: GET_AREA_M2
PUBLIC  :: GET_AREA_CM2
PUBLIC  :: GET_BOUNDING_BOX
PUBLIC  :: GET_XEDGE
PUBLIC  :: GET_XMID
PUBLIC  :: GET_XOFFSET
PUBLIC  :: GET_YOFFSET
PUBLIC  :: GET_YEDGE
```

```

PUBLIC  :: GET_YEDGE_R
PUBLIC  :: GET_YMID
PUBLIC  :: GET_YMID_R
PUBLIC  :: GET_YMID_R_W
PUBLIC  :: SET_XOFFSET
PUBLIC  :: SET_YOFFSET
PUBLIC  :: ITS_A_NESTED_GRID

```

#### PRIVATE MEMBER FUNCTIONS:

```

PRIVATE :: INIT_GRID

```

#### REVISION HISTORY:

```

11 Mar 2003 - R. Yantosca - Initial version
(1 ) Fixed typos in "grid_mod.f" (bmy, 4/28/03)
(2 ) Added routine GET_BOUNDING_BOX. Now define 1x125 grid. (bmy, 12/1/04)
(3 ) Modified for GCAP 4x5 horizontal grid (swu, bmy, 5/24/05)
(4 ) Added comments re: surface area derivation (bmy, 4/20/06)
(5 ) Modifications for GEOS-5 nested grids (yxw, dan, bmy, 11/6/08)
20 Nov 2009 - R. Yantosca - Added ProTeX headers

```

---

#### 1.6.1 compute\_grid

Subroutine COMPUTE\_GRID initializes the longitude, latitude and surface area arrays.

#### INTERFACE:

```

SUBROUTINE COMPUTE_GRID

```

#### USES:

```

USE CMN_SIZE_MOD  ! Size parameters
USE CMN_GCTM_MOD  ! Physical constants

```

#### REVISION HISTORY:

```

11 Mar 2003 - R. Yantosca - Initial version
(1 ) Added fancy output (bmy, 4/26/04)
(2 ) Suppress some output lines (bmy, 7/20/04)
(3 ) Now also support 1 x 1.25 grid (bmy, 12/1/04)
(4 ) Now modified for GCAP 4x5 horizontal grid (swu, bmy, 5/24/05)
(5 ) Added comments re: surface area derivation (bmy, 4/20/06)
(6 ) Compute YMID, YEDGE for 0.5x0.666 nested grids (yxw, dan, bmy, 11/6/08)
20 Nov 2009 - R. Yantosca - Added ProTeX header

```

---

### 1.6.2 set\_xoffset

Function SET\_XOFFSET initializes the nested-grid longitude offset variable I0.

#### INTERFACE:

```
SUBROUTINE SET_XOFFSET( X_OFFSET )
```

#### INPUT PARAMETERS:

```
INTEGER, INTENT(IN) :: X_OFFSET ! Value to assign to I0
```

#### REVISION HISTORY:

```
11 Mar 2003 - R. Yantosca - Initial version
20 Nov 2009 - R. Yantosca - Added ProTeX header
```

---

### 1.6.3 set\_yoffset

Function SET\_YOFFSET initializes the nested-grid latitude offset variable J0.

#### INTERFACE:

```
SUBROUTINE SET_YOFFSET( Y_OFFSET )
```

#### INPUT PARAMETERS:

```
INTEGER, INTENT(IN) :: Y_OFFSET ! Value to assign to J0
```

#### REVISION HISTORY:

```
11 Mar 2003 - R. Yantosca - Initial version
20 Nov 2009 - R. Yantosca - Added ProTeX header
```

---

### 1.6.4 get\_xoffset

Function GET\_XOFFSET returns the nested-grid longitude offset to the calling program.  
(bmy, 3/11/03)

#### INTERFACE:

```
FUNCTION GET_XOFFSET( GLOBAL ) RESULT( X_OFFSET )
```

#### INPUT PARAMETERS:

```
! If GLOBAL is passed, then return the actual window offset.
! This is necessary for certain instances (e.g. diagnostics)
LOGICAL, INTENT(IN), OPTIONAL :: GLOBAL
```

#### RETURN VALUE:

```
INTEGER :: X_OFFSET
```

#### REVISION HISTORY:

```
11 Mar 2003 - R. Yantosca - Initial version
20 Nov 2009 - R. Yantosca - Added ProTeX header
```

---

### 1.6.5 get\_xoffset

Function GET\_XOFFSET returns the nested-grid longitude offset to the calling program.  
(bmy, 3/11/03)

#### INTERFACE:

```
FUNCTION GET_YOFFSET( GLOBAL ) RESULT( Y_OFFSET )
```

#### INPUT PARAMETERS:

```
! If GLOBAL is passed, then return the actual window offset.
! This is necessary for certain instances (e.g. diagnostics)
LOGICAL, INTENT(IN), OPTIONAL :: GLOBAL
```

#### RETURN VALUE:

```
INTEGER                                :: Y_OFFSET
```

#### REVISION HISTORY:

```
11 Mar 2003 - R. Yantosca - Initial version
20 Nov 2009 - R. Yantosca - Added ProTeX header
```

---

### 1.6.6 get\_xmid

Function GET\_XMID returns the longitude in degrees at the center of a GEOS-Chem grid box. Works for nested-grids too.

#### INTERFACE:

```
FUNCTION GET_XMID( I ) RESULT( X )
```

#### INPUT PARAMETERS:

```
INTEGER, INTENT(IN) :: I ! Longitude index
```

#### RETURN VALUE:

```
REAL*8                                :: X ! Corresponding lon value @ grid box ctr
```

#### REVISION HISTORY:

```
11 Mar 2003 - R. Yantosca - Initial version
20 Nov 2009 - R. Yantosca - Added ProTeX header
```

---

### 1.6.7 get\_xedge

Function GET\_XEDGE returns the longitude in degrees at the western edge of a GEOS-Chem grid box. Works for nested-grids too.

#### INTERFACE:

```
FUNCTION GET_XEDGE( I ) RESULT( X )
```

#### INPUT PARAMETERS:

```
INTEGER, INTENT(IN) :: I ! Longitude index
```

#### RETURN VALUE:

```
REAL*8 :: X ! Corresponding lon value @ W edge of grid box
```

#### REVISION HISTORY:

```
11 Mar 2003 - R. Yantosca - Initial version
20 Nov 2009 - R. Yantosca - Added ProTeX header
```

---

### 1.6.8 get\_ymid

Function GET\_YMID returns the latitude in degrees at the center of a GEOS-Chem grid box. Works for nested-grids too.

#### INTERFACE:

```
FUNCTION GET_YMID( J ) RESULT( Y )
```

#### INPUT PARAMETERS:

```
INTEGER, INTENT(IN) :: J ! Latitude index
```

#### RETURN VALUE:

```
REAL*8 :: Y ! Latitude value at @ grid box ctr [degrees]
```

#### REVISION HISTORY:

```
11 Mar 2003 - R. Yantosca - Initial version
20 Nov 2009 - R. Yantosca - Added ProTeX header
```

---

### 1.6.9 get\_yedge

Function GET\_YEDGE returns the latitude in degrees at the southern edge of a GEOS-Chem grid box. Works for nested-grids too.

#### INTERFACE:

```
FUNCTION GET_YEDGE( J ) RESULT( Y )
```

**INPUT PARAMETERS:**

```
INTEGER, INTENT(IN) :: J ! Latitude index
```

**RETURN VALUE:**

```
REAL*8 :: Y ! Latitude value @ S edge of grid box [degrees]
```

**REVISION HISTORY:**

```
11 Mar 2003 - R. Yantosca - Initial version
20 Nov 2009 - R. Yantosca - Added ProTeX header
```

---

**1.6.10 get\_ymid**

Function GET\_YMID\_R returns the latitude in radians at the center of a GEOS-Chem grid box. Works for nested-grids too.

**INTERFACE:**

```
FUNCTION GET_YMID_R( J ) RESULT( Y )
```

**INPUT PARAMETERS:**

```
INTEGER, INTENT(IN) :: J ! Latitude index
```

**RETURN VALUE:**

```
REAL*8 :: Y ! Latitude value at @ grid box ctr [radians]
```

**REVISION HISTORY:**

```
11 Mar 2003 - R. Yantosca - Initial version
20 Nov 2009 - R. Yantosca - Added ProTeX header
```

---

**1.6.11 get\_ymid**

Function GET\_YMID\_R\_W returns the latitude in radians at the center of a GEOS-Chem grid box for the GEOS-5 nested grid.

**INTERFACE:**

```
FUNCTION GET_YMID_R_W( J ) RESULT( Y )
```

**INPUT PARAMETERS:**

```
INTEGER, INTENT(IN) :: J ! Latitude index
```

**RETURN VALUE:**

```
REAL*8 :: Y ! Latitude value at @ grid box ctr [radians]
```

**REVISION HISTORY:**

```
06 Nov 2008 - D. Chen & R. Yantosca - Initial version
20 Nov 2009 - R. Yantosca - Added ProTeX header
```

---

**1.6.12 get\_yedge\_r**

Function GET\_YEDGE\_R returns the latitude in radians at the southern edge of a GEOS-Chem grid box. Works for nested-grids too.

**INTERFACE:**

```
FUNCTION GET_YEDGE_R( J ) RESULT( Y )
```

**INPUT PARAMETERS:**

```
INTEGER, INTENT(IN) :: J ! Latitude index
```

**RETURN VALUE:**

```
REAL*8 :: Y ! Latitude value @ S edge of grid box [radians]
```

**REVISION HISTORY:**

```
11 Mar 2003 - R. Yantosca - Initial version
20 Nov 2009 - R. Yantosca - Added ProTeX header
```

---

**1.6.13 get\_area\_m2**

Function GET\_AREA\_M2 returns the surface area [m2] of a GEOS-Chem grid box. Works for nested grids too.

**INTERFACE:**

```
FUNCTION GET_AREA_M2( J ) RESULT( A )
```

**INPUT PARAMETERS:**

```
INTEGER, INTENT(IN) :: J ! Latitude index
```

**RETURN VALUE:**

```
REAL*8 :: A ! Grid box surface area [m2]
```

**REVISION HISTORY:**

```
11 Mar 2003 - R. Yantosca - Initial version
20 Nov 2009 - R. Yantosca - Added ProTeX header
```

---

**1.6.14 get\_area\_cm2**

Function GET\_AREA\_CM2 returns the surface area [cm2] of a GEOS-Chem grid box. Works for nested grids too.

**INTERFACE:**

```
FUNCTION GET_AREA_CM2( J ) RESULT( A )
```

**INPUT PARAMETERS:**

```
INTEGER, INTENT(IN) :: J ! Latitude index
```

**RETURN VALUE:**

```
REAL*8 :: A ! Grid box surface area [cm2]
```

**REVISION HISTORY:**

```
11 Mar 2003 - R. Yantosca - Initial version
20 Nov 2009 - R. Yantosca - Added ProTeX header
```

---

**1.6.15 get\_bounding\_box**

Subroutine GET\_BOUNDING\_BOX returns the indices which specify the lower left (LL) and upper right (UR) corners of a rectangular region, given the corresponding longitude and latitude values.

**INTERFACE:**

```
SUBROUTINE GET_BOUNDING_BOX( COORDS, INDICES )
```

**USES:**

```
USE ERROR_MOD, ONLY : ERROR_STOP

USE CMN_SIZE_MOD ! Size parameters
```

**INPUT PARAMETERS:**

```
REAL*8, INTENT(IN) :: COORDS(4) ! (/LON_LL, LAT_LL, LON_UR, LAT_UR/)
```

**INPUT/OUTPUT PARAMETERS:**

```
INTEGER, INTENT(OUT) :: INDICES(4) ! (/I_LL, J_LL, I_UR, J_UR/)
```

**REVISION HISTORY:**

```
01 Dec 2004 - R. Yantosca - Initial version
20 Nov 2009 - R. Yantosca - Added ProTeX header
```

---

**1.6.16 its\_a\_nested\_grid**

Function GET\_AREA\_CM2 returns the surface area [cm2] of a GEOS-Chem grid box. Works for nested grids too.

**INTERFACE:**

```
FUNCTION ITS_A_NESTED_GRID() RESULT( IT_IS_NESTED )
```



**RETURN VALUE:**

LOGICAL :: IT\_IS\_NESTED    ! =T if it's a nested grid; =F otherwise

**REVISION HISTORY:**

11 Mar 2003 - R. Yantosca - Initial version  
 20 Nov 2009 - R. Yantosca - Added ProTeX header

---

**1.6.17    init\_grid**

Subroutine INIT\_GRID initializes variables and allocates module arrays.

**INTERFACE:**

SUBROUTINE INIT\_GRID

**USES:**

USE ERROR\_MOD, ONLY : ALLOC\_ERR

USE CMN\_SIZE\_MOD

**REVISION HISTORY:**

11 Mar 2003 - R. Yantosca - Initial version  
 (1 ) Fixed typos that caused AREA\_CM2\_G and AREA\_CM2 to be initialized  
      before they were allocated. (bmy, 4/28/03)  
 (2 ) Now define IIIPAR & JJJPARG for 1 x 1.25 grid (bmy, 12/1/04)  
 (3 ) Modified for GCAP 4x5 horizontal grid (swu, bmy, 5/24/05)  
 (4 ) Modifications for 0.5 x 0.666 nested grids (dan, bmy, 11/6/08)  
 20 Nov 2009 - R. Yantosca - Added ProTeX header

---

**1.6.18    cleanup\_grid**

Subroutine CLEANUP\_GRID deallocates all module arrays.

**INTERFACE:**

SUBROUTINE CLEANUP\_GRID

**REVISION HISTORY:**

11 Mar 2003 - R. Yantosca - Initial version  
 20 Nov 2009 - D. Chen        - Now also deallocate YMID\_R\_W  
 20 Nov 2009 - R. Yantosca - Added ProTeX header

---

## 1.7 Fortran: Module Interface hdf\_mod

Module HDF\_MOD contains routines to write data to HDF5 files.

### INTERFACE:

```
MODULE HDF_MOD
```

### USES:

```
IMPLICIT NONE
#   include "define.h"
PRIVATE
```

### PUBLIC MEMBER FUNCTIONS:

```
PUBLIC :: CLEANUP_HDF
PUBLIC :: CLOSE_HDF
PUBLIC :: INIT_HDF
PUBLIC :: OPEN_HDF
PUBLIC :: WRITE_HDF
```

### PUBLIC DATA MEMBERS:

```
PUBLIC :: HDFCATEGORY
PUBLIC :: HDFDESCRIPT
PUBLIC :: HDFNAME
PUBLIC :: HDIFFNAME
PUBLIC :: HDFMOLC
PUBLIC :: HDFMWT
PUBLIC :: HDFSCALE
PUBLIC :: HDFUNIT

CHARACTER(LEN=40), ALLOCATABLE :: HDFCATEGORY(:)
CHARACTER(LEN=40), ALLOCATABLE :: HDFDESCRIPT(:)
INTEGER :: MAXDIAG
INTEGER :: MAXTRACER
INTEGER :: MAXCAT
INTEGER, ALLOCATABLE :: HDFMOLC(:, :)
REAL*4, ALLOCATABLE :: HDFMWT(:, :)
REAL*4, ALLOCATABLE :: HDFSCALE(:, :)
CHARACTER(LEN=40), ALLOCATABLE :: HDFNAME(:, :)
CHARACTER(LEN=40), ALLOCATABLE :: HDIFFNAME(:, :)
CHARACTER(LEN=40), ALLOCATABLE :: HDFUNIT(:, :)
```

### REMARKS:

If you have the HDF5 library installed on your system, then you can compile GEOS-Chem with the option:

```
make HDF5=yes
```

which will activate the HDF5-specific code in this module to enable file I/O to HDF5 format. You must also specify the HDF5 include and library paths in the Makefile\_header.mk.

The default is not to activate the HDF5-specific code.

## REVISION HISTORY:

19 Nov 2009 - A. van Donkelaar - Initial Version  
 21 Dec 2009 - R. Yantosca - Modified to block out HDF5-specific code  
 so that users w/o HDF5 library can still  
 compile & run GEOS-Chem  
 21 Dec 2009 - R. Yantosca - Updated comments

### 1.7.1 open\_hdf

Subroutine OPEN\_HDF creates and opens an hdf file for output.

## INTERFACE:

```
#if defined( USE_HDF5 )
!%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%
!%%      NOTE: Subroutine is used only when USE_HDF5 is defined!      %%
!%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%
SUBROUTINE OPEN_HDF( IU_HDF, FILENAME, IMAX, IMIN,
&                   JMAX,  JMIN,   NI,   NJ   )
```

## USES:

```
USE HDF5
USE GRID_MOD, ONLY : GET_XMID
USE GRID_MOD, ONLY : GET_YMID

USE CMN_SIZE_MOD    ! Size parameters
```

## INPUT PARAMETERS:

CHARACTER(LEN=*)	INTENT(IN)	:: FILENAME	! File name to open
INTEGER,	INTENT(IN)	:: IMIN	! Min longitude index
INTEGER,	INTENT(IN)	:: IMAX	! Max longitude index
INTEGER,	INTENT(IN)	:: JMIN	! Min latitude index
INTEGER,	INTENT(IN)	:: JMAX	! Max latitude index
INTEGER,	INTENT(IN)	:: NI	! # of longitudes
INTEGER,	INTENT(IN)	:: NJ	! # of latitudes

## INPUT/OUTPUT PARAMETERS:

INTEGER(HID_T),	INTENT(INOUT)	:: IU_HDF	! HDF5 file identifier
-----------------	---------------	-----------	------------------------

## REVISION HISTORY:

Nov 20 2009 - A. van Donkelaar - Initial Version  
 21 Dec 2009 - R. Yantosca - Modified to block out HDF5-specific code  
                                   so that users w/o HDF5 library can still  
                                   compile & run GEOS-Chem  
 21 Dec 2009 - R. Yantosca - Updated comments

---

### 1.7.2 close\_hdf

Subroutine CLOSE\_HDF closes an HDF5 file that is already open.

#### INTERFACE:

```
#if defined( USE_HDF5 )
!%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%
!%%      NOTE: Subroutine is used only when USE_HDF5 is defined!    %%
!%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%
SUBROUTINE CLOSE_HDF( IU_HDF )
```

#### USES:

USE HDF5

#### INPUT/OUTPUT PARAMETERS:

INTEGER(HID\_T), INTENT(INOUT) :: IU\_HDF ! HDF5 File identifier

#### REVISION HISTORY:

20 Nov 2009 - A. van Donkelaar - Initial Version  
 21 Dec 2009 - R. Yantosca - Modified to block out HDF5-specific code  
                                   so that users w/o HDF5 library can still  
                                   compile & run GEOS-Chem  
 21 Dec 2009 - R. Yantosca - Updated comments

---

### 1.7.3 write\_hdf

Subroutine WRITE\_HDF writes data to an open HDF5 file.

#### INTERFACE:

```
#if defined( USE_HDF5 )
!%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%
!%%      NOTE: Subroutine is used only when USE_HDF5 is defined!    %%
!%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%
SUBROUTINE WRITE_HDF( IU_HDF, N, NDCATEGORY, NDTRACER,
& NDUNIT, TAU0, TAU1, RESERVED,
& NI, NJ, NL, IFIRST,
& JFIRST, LFIRST, ARRAY )
```

**USES:**

```
USE HDF5
```

**INPUT PARAMETERS:**

```

INTEGER,          INTENT(IN) :: IU_HDF           ! HDF file unit #
INTEGER,          INTENT(IN) :: N               ! Actual tracer #
CHARACTER(LEN=40), INTENT(IN) :: NDCATEGORY      ! Diagnostic category
INTEGER,          INTENT(IN) :: NDTRACER        ! Tracer # for file
CHARACTER(LEN=40), INTENT(IN) :: NDUNIT         ! Units of data
REAL*8,           INTENT(IN) :: TAU0            ! TAU at start & end
REAL*8,           INTENT(IN) :: TAU1            !   of diag interval
CHARACTER(LEN=40), INTENT(IN) :: RESERVED       ! Descriptive string
INTEGER,          INTENT(IN) :: NI              ! # of longitudes
INTEGER,          INTENT(IN) :: NJ              ! # of latitudes
INTEGER,          INTENT(IN) :: NL              ! # of levels
INTEGER,          INTENT(IN) :: IFIRST          ! Index of 1st lon
INTEGER,          INTENT(IN) :: JFIRST          ! Index of 1st lat
INTEGER,          INTENT(IN) :: LFIRST          ! Index of 1st lev
REAL*4,           INTENT(IN) :: ARRAY(NI,NJ,NL) ! Data array

```

**REVISION HISTORY:**

```

20 Nov 2009 - A. van Donkelaar - Initial Version
21 Dec 2009 - R. Yantosca      - Modified to block out HDF5-specific code
                                so that users w/o HDF5 library can still
                                compile & run GEOS-Chem
21 Dec 2009 - R. Yantosca      - Updated comments

```

---

**1.7.4 init\_hdf**

Subroutine INIT\_HDF allocates all module variables.

**INTERFACE:**

```
SUBROUTINE INIT_HDF( GMMAXCAT, GMMAXTRACER, GMMAXDIAG )
```

**USES:**

```
USE ERROR_MOD, ONLY : ALLOC_ERR
```

**INPUT PARAMETERS:**

```

INTEGER, INTENT(IN) :: GMMAXCAT
INTEGER, INTENT(IN) :: GMMAXTRACER
INTEGER, INTENT(IN) :: GMMAXDIAG

```

**REVISION HISTORY:**

### 1.7.5 cleanup\_hdf

## INTERFACE:

**REVISION HISTORY:**

### 1.7.6 ifort\_errmsg

## INTERFACE:

### INPUT PARAMETERS:

**RETURN VALUE:**

**REVISION HISTORY:**

### 1.8 Fortran: Module Interface julday\_mod.f

**INTERFACE:**

```
MODULE JULDAY_MOD
```

## USES:

```
IMPLICIT NONE
#   include "define.h"
PRIVATE
```

## PUBLIC MEMBER FUNCTIONS:

```
PUBLIC   :: JULDAY
PUBLIC  :: CALDATE
```

## PRIVATE MEMBER FUNCTIONS:

```
PRIVATE :: MINT
```

## REVISION HISTORY:

- (1 ) Moved JULDAY, MINT, CALDATE here from "bpch2\_mod.f" (bmy, 11/20/01)
  - (2 ) Bug fix: now compute NHMS correctly. Also use REAL\*4 variables to avoid roundoff errors. (bmy, 11/26/01)
  - (3 ) Updated comments (bmy, 5/28/02)
  - (4 ) Renamed arguments for clarity (bmy, 6/26/02)
  - 20 Nov 2009 - R. Yantosca - Added ProTeX Headers
- 

### 1.8.1 julday

Function JULDAY returns the astronomical Julian day.

## INTERFACE:

```
FUNCTION JULDAY( YYYY, MM, DD ) RESULT( JULIANDAY )
```

## INPUT PARAMETERS:

```
INTEGER, INTENT(IN) :: YYYY      ! Year (must be in 4-digit format!)
INTEGER, INTENT(IN) :: MM        ! Month (1-12)
REAL*8,  INTENT(IN) :: DD        ! Day of month (may be fractional!)
```

## RETURN VALUE:

```
REAL*8              :: JULIANDAY  ! Astronomical Julian Date
```

## REMARKS:

- (1) Algorithm taken from "Practical Astronomy With Your Calculator", Third Edition, by Peter Duffett-Smith, Cambridge UP, 1992.
- (2) Requires the external function MINT.F.
- (3) JulDay will compute the correct Julian day for any BC or AD date.
- (4) For BC dates, subtract 1 from the year and append a minus sign. For example, 1 BC is 0, 2 BC is -1, etc. This is necessary for the algorithm.

**REVISION HISTORY:**

26 Nov 2001 - R. Yantosca - Initial version  
 Changed YEAR to YYYY, MONTH to MM, and DAY to DD for documentation  
 purposes. (bmy, 6/26/02)  
 20 Nov 2009 - R. Yantosca - Added ProTeX headers

---

**1.8.2 mint**

Function MINT is the modified integer function.

**INTERFACE:**

```
FUNCTION MINT( X ) RESULT ( VALUE )
```

**INPUT PARAMETERS:**

```
REAL*8, INTENT(IN) :: X
```

**RETURN VALUE:**

```
REAL*8                :: VALUE
```

**REMARKS:**

The modified integer function is defined as follows:

```
      { -INT( ABS( X ) )   for X < 0
MINT = {
      {  INT( ABS( X ) )   for X >= 0
```

**REVISION HISTORY:**

20 Nov 2001 - R. Yantosca - Initial version  
 20 Nov 2009 - R. Yantosca - Added ProTeX headers

---

**1.8.3 caldate**

Subroutine CALDATE converts an astronomical Julian day to the YYYYMMDD and HH-MMSS format.

**INTERFACE:**

```
SUBROUTINE CALDATE( JULIANDAY, YYYYMMDD, HHMMSS )
```

**INPUT PARAMETERS:**

```
! Arguments
REAL*8, INTENT(IN) :: JULIANDAY ! Astronomical Julian Date
```

**OUTPUT PARAMETERS:**



```

INTEGER, INTENT(OUT) :: YYYYMMDD    ! Date in YYYY/MM/DD format
INTEGER, INTENT(OUT) :: HHMMSS      ! Time in hh:mm:ss format

```

**REMARKS:**

Algorithm taken from "Practical Astronomy With Your Calculator",  
Third Edition, by Peter Duffett-Smith, Cambridge UP, 1992.

**REVISION HISTORY:**

- (1 ) Now compute HHMMSS correctly. Also use REAL\*4 variables HH, MM, SS  
to avoid roundoff errors. (bmy, 11/21/01)
  - (2 ) Renamed NYMD to YYYYMMDD and NHMS to HHMMSS for documentation  
purposes (bmy, 6/26/02)
- 20 Nov 2009 - R. Yantosca - Added ProTeX header

**1.9 Fortran: Module Interface pressure\_mod.f**

Module PRESSURE\_MOD contains variables and routines which specify the grid box pressures for both hybrid or pure-sigma models. This is necessary for running GEOS-Chem with the GEOS-4 or GEOS-5 hybrid grids.

**INTERFACE:**

```

MODULE PRESSURE_MOD

```

**USES:**

```

      IMPLICIT NONE
#      include "define.h"
      PRIVATE

```

**PUBLIC MEMBER FUNCTIONS:**

```

      PUBLIC :: CLEANUP_PRESSURE
      PUBLIC :: GET_AP
      PUBLIC :: GET_BP
      PUBLIC :: GET_PCENTER
      PUBLIC :: GET_PEDGE
      PUBLIC :: GET_PEDGE_FULLGRID
      PUBLIC :: INIT_PRESSURE
      PUBLIC :: SET_FLOATING_PRESSURE

```

**REMARKS:**

Hybrid Grid Coordinate Definition: (dsa, bmy, 8/27/02, 8/13/10)

=====

GEOS-4, GEOS-5, and MERRA (hybrid grids):

-----

For GEOS-4/GEOS-5/MERRA met data products, the pressure at the bottom edge of grid box (I,J,L) is defined as follows:

$$P_{edge}(I,J,L) = A_p(L) + [ B_p(L) * P_{surface}(I,J) ]$$

where

$P_{surface}(I,J)$  is the "true" surface pressure at lon,lat (I,J)  
 $A_p(L)$  has the same units as surface pressure [hPa]  
 $B_p(L)$  is a unitless constant given at level edges

$A_p(L)$  and  $B_p(L)$  are given to us by GMAO.

GEOS-3 (pure-sigma) and GCAP (hybrid grid):

-----  
 GEOS-3 is a pure-sigma grid. GCAP is a hybrid grid, but its grid is defined as if it were a pure sigma grid (i.e.  $P_{TOP}=150$  hPa, and negative sigma edges at higher levels). For these grids, can still use the same formula as for GEOS-4/GEOS-5/MERRA, with one modification:

$$P_{edge}(I,J,L) = A_p(L) + [ B_p(L) * ( P_{surface}(I,J) - P_{TOP} ) ]$$

where

$P_{surface}(I,J)$  = the "true" surface pressure at lon,lat (I,J)  
 $A_p(L)$  =  $P_{TOP}$  = model top pressure  
 $B_p(L)$  =  $SIG_E(L)$  = bottom sigma edge of level L

The following are true for GCAP, GEOS-3, GEOS-4, GEOS-5, MERRA:

- (1)  $B_p(LLPAR+1) = 0.0$  (L=LLPAR+1 is the atmosphere top)  
 (2)  $B_p(1) = 1.0$  (L=1 is the surface)  
 (3)  $P_{TOP} = A_p(LLPAR+1)$  (L=LLPAR+1 is the atmosphere top)

## REVISION HISTORY:

27 Aug 2002 - D. Abbot & R. Yantosca - Initial version

- (1 ) Be sure to check PFLT for NaN or Infinities (bmy, 8/27/02)
- (2 ) Updated comments (bmy, 5/8/03)
- (3 ) Updated format string for fvDAS (bmy, 6/19/03)
- (4 ) Bug fix: use PFLT instead of PFLT-PTOP for GEOS-4 (bmy, 10/24/03)
- (5 ) Modifications for 30L and 55L GEOS-4 grids (bmy, 11/3/03)
- (6 ) Added parallel DO-loop in SET\_FLOATING\_PRESSURE (bmy, 4/14/04)
- (7 ) Modified for GCAP and GEOS-5 grids (swu, bmy, 5/24/05)
- (8 ) Removed obsolete reference to "CMN" (bmy, 4/25/06)
- (9 ) Remove support for GEOS-1 and GEOS-STRAT met fields (bmy, 8/4/06)
- (10) Added  $A_p$  and  $B_p$  for GEOS-5 met fields (bmy, 10/30/07)

20 Nov 2009 - R. Yantosca - Added ProTeX headers  
 13 Aug 2010 - R. Yantosca - Added modifications for MERRA met fields  
 30 Aug 2010 - R. Yantosca - Updated comments

---

### 1.9.1 get\_ap

Function GET\_AP returns the "A" term [hPa] for the hybrid ETA coordinate.

#### INTERFACE:

```
FUNCTION GET_AP( L ) RESULT( AP_TEMP )
```

#### USES:

```
USE CMN_SIZE_MOD           ! Size parameters
```

#### INPUT PARAMETERS:

```
INTEGER, INTENT(IN) :: L           ! GEOS-Chem level index
```

#### RETURN VALUE:

```
REAL*8                :: AP_TEMP  ! Corresponding "A" value [hPa]  
                        ! at bottom edge of level L
```

#### REVISION HISTORY:

20 Aug 2002 - D. Abbot & R. Yantosca - Initial version  
 20 Nov 2009 - R. Yantosca - Added ProTeX header

---

### 1.9.2 get\_bp

Function GET\_BP returns the "B" term [unitless] for the hybrid ETA coordinate.

#### INTERFACE:

```
FUNCTION GET_BP( L ) RESULT( BP_TEMP )
```

#### USES:

```
USE CMN_SIZE_MOD           ! Size parameters
```

#### INPUT PARAMETERS:

```
INTEGER, INTENT(IN) :: L           ! GEOS-Chem level index
```

#### RETURN VALUE:

```
REAL*8                :: BP_TEMP  ! Corresponding "B" value [unitless]  
                        ! at bottom edge of level L
```

#### REVISION HISTORY:

20 Aug 2002 - D. Abbot & R. Yantosca - Initial version  
 20 Nov 2009 - R. Yantosca - Added ProTeX header

---

### 1.9.3 set\_floating\_pressure

Subroutine SET\_FLOATING\_PRESSURE initializes the floating pressure field PFLT with a pressure from the main program. This is needed to initialize and reset PFLT after transport.

#### INTERFACE:

```
SUBROUTINE SET_FLOATING_PRESSURE( PS )
```

#### USES:

```
USE ERROR_MOD, ONLY : CHECK_VALUE
```

```
USE CMN_SIZE_MOD ! Size parameters
```

#### INPUT PARAMETERS:

```
! Array containing pressure with which to initialize PFLT [hPa]
REAL*8, INTENT(IN) :: PS(IIPAR,JJPARG)
```

#### REVISION HISTORY:

```
27 Aug 2002 - D. Abbot, B. Field, R. Yantosca - Initial version
(1 ) Now check PFLT for NaN or Infinities (bmy, 8/27/02)
(2 ) Added parallel DO-loop (bmy, 4/14/04)
20 Nov 2009 - R. Yantosca - Added ProTeX header
```

### 1.9.4 get\_pedge

Function GET\_PEDGE returns the pressure at the bottom edge of level L. L=1 is the surface, L=LLPAR+1 is the atm top.

#### INTERFACE:

```
FUNCTION GET_PEDGE( I, J, L ) RESULT( PEDGE )
```

#### USES:

```
USE CMN_SIZE_MOD ! PTOP
```

#### INPUT PARAMETERS:

```
INTEGER, INTENT(IN) :: I ! GEOS-Chem lon index
INTEGER, INTENT(IN) :: J ! GEOS-Chem lat index
INTEGER, INTENT(IN) :: L ! GEOS-Chem level index
```

#### RETURN VALUE:

```
REAL*8 :: PEDGE ! Pressure @ bottom edge of (I,J,L) [hPa]
```

#### REVISION HISTORY:

20 Aug 2002 - D. Abbot & R. Yantosca - Initial version  
 (1 ) Bug fix: use PFLT instead of PFLT-PTOP for GEOS-4 (bmy, 10/24/03)  
 (2 ) Now treat GEOS-5 the same way as GEOS-4 (bmy, 10/30/07)  
 20 Nov 2009 - R. Yantosca - Added ProTeX header  
 13 Aug 2010 - R. Yantosca - Compute PEDGE for MERRA the same as for GEOS-5

---

### 1.9.5 get\_pedge\_fullgrid

Function GET\_PEDGE\_FULLGRID returns the pressure at the bottom edge of level L of the unreduced vertical grid. L=1 is the surface, L=LLGLOB+1 is the atm top.

#### INTERFACE:

```
FUNCTION GET_PEDGE_FULLGRID( I, J, L ) RESULT( PEDGE )
```

#### USES:

```
USE CMN_SIZE_MOD    ! PTOP
```

#### INPUT PARAMETERS:

```
INTEGER, INTENT(IN) :: I      ! GEOS-Chem lon   index
INTEGER, INTENT(IN) :: J      ! GEOS-Chem lat   index
INTEGER, INTENT(IN) :: L      ! GEOS-Chem level index
```

#### RETURN VALUE:

```
REAL*8              :: PEDGE ! Pressure @ bottom edge of (I,J,L) [hPa]
```

#### REVISION HISTORY:

```
(1 ) Modified from GET_PEDGE (cdh, 1/22/09)
```

---

### 1.9.6 get\_pcenter

Function GET\_PCENTER returns the pressure at the vertical midpoint of level L.

#### INTERFACE:

```
FUNCTION GET_PCENTER( I, J, L ) RESULT( PCENTER )
```

#### USES:

```
USE CMN_SIZE_MOD    ! PTOP
```

#### INPUT PARAMETERS:

```
INTEGER, INTENT(IN) :: I      ! GEOS-Chem lon   index
INTEGER, INTENT(IN) :: J      ! GEOS-Chem lat   index
INTEGER, INTENT(IN) :: L      ! GEOS-Chem level index
```

**RETURN VALUE:**

```
REAL*8          :: PCENTER  ! Pressure @ center of (I,J,L) [hPa]
```

**REVISION HISTORY:**

```
20 Aug 2002 - D. Abbot & R. Yantosca - Initial version
(1 ) Updated format string for fvDAS (bmy, 6/19/03)
(2 ) Removed reference to "CMN", it's obsolete (bmy, 4/25/06)
20 Nov 2009 - R. Yantosca - Added ProTeX header
```

---

**1.9.7 init\_pressure**

Subroutine INIT\_PRESSURE allocates and initializes the AP and BP arrays. It must be called in "main.f", after SIGE is defined. GEOS-4 and GEOS-5 requires the hybrid pressure system specified by the listed values of AP and BP, while earlier versions of GEOS use a pure sigma pressure system. GCAP met fields (based on GISS) also use a hybrid system.

**INTERFACE:**

```
SUBROUTINE INIT_PRESSURE
```

**USES:**

```
! References to F90 modules
USE ERROR_MOD, ONLY : ALLOC_ERR

USE CMN_SIZE_MOD  ! LLPAR, PTOP
```

**REVISION HISTORY:**

```
27 Aug 2002 - D. Abbot, S. Wu, & R. Yantosca - Initial version
(1 ) Now reference ALLOC_ERR from "error_mod.f" (bmy, 10/15/02)
(2 ) Now echo Ap, Bp to std output (bmy, 3/14/03)
(3 ) Now print LLPAR+1 levels for Ap, Bp. Remove reference to SIGE, it's
      obsolete. Also now use C-preprocessor switch GRID30LEV instead of
      IF statements to define vertical coordinates. (bmy, 11/3/03)
(4 ) Now modified for both GCAP & GEOS-5 vertical grids (swu, bmy, 5/24/05)
(5 ) Renamed GRID30LEV to GRIDREDUCED (bmy, 10/30/07)
20 Nov 2009 - R. Yantosca - Added ProTeX header
13 Aug 2010 - R. Yantosca - Compute Ap and Bp for MERRA the same way as for
                          GEOS-5. The vertical grids are identical.
30 Aug 2010 - R. Yantosca - Updated comments
30 Nov 2010 - R. Yantosca - Further improved comments about how GEOS-4 and
                          GEOS-5 vertical levels are lumped together.
```

---

### 1.9.8 cleanup\_pressure

Subroutine CLEANUP\_PRESSURE deallocates all allocated arrays at the end of a GEOS-Chem model run.

#### INTERFACE:

```
SUBROUTINE CLEANUP_PRESSURE
```

#### REVISION HISTORY:

```
20 Aug 2002 - D. Abbot & R. Yantosca - Initial version
20 Nov 2009 - R. Yantosca - Added ProTeX header
```

### 1.10 Fortran: Module Interface time\_mod

Module TIME\_MOD contains GEOS-Chem date and time variables and timesteps, and routines for accessing them.

#### INTERFACE:

```
MODULE TIME_MOD
```

#### USES:

```
IMPLICIT NONE
```

```
#    include "define.h"
PRIVATE
```

#### PUBLIC MEMBER FUNCTIONS:

```
PUBLIC  :: SET_CURRENT_TIME
PUBLIC  :: SET_BEGIN_TIME
PUBLIC  :: SET_END_TIME
PUBLIC  :: SET_NDIAGTIME
PUBLIC  :: SET_DIAGb
PUBLIC  :: SET_DIAGe
PUBLIC  :: SET_TIMESTEPS
PUBLIC  :: SET_CT_CHEM
PUBLIC  :: SET_CT_CONV
PUBLIC  :: SET_CT_DYN
PUBLIC  :: SET_CT_EMIS
PUBLIC  :: SET_CT_DIAG
PUBLIC  :: SET_CT_A1
PUBLIC  :: SET_CT_A3
PUBLIC  :: SET_CT_A6
PUBLIC  :: SET_CT_I6
PUBLIC  :: SET_CT_XTRA
```

```

PUBLIC  :: SET_ELAPSED_MIN
PUBLIC  :: GET_JD
PUBLIC  :: GET_ELAPSED_MIN
PUBLIC  :: GET_ELAPSED_SEC
PUBLIC  :: GET_NYMDb
PUBLIC  :: GET_NHMSb
PUBLIC  :: GET_NYMDe
PUBLIC  :: GET_NHMSe
PUBLIC  :: GET_NYMD
PUBLIC  :: GET_NHMS
PUBLIC  :: GET_NDIAGTIME
PUBLIC  :: GET_TIME_AHEAD
PUBLIC  :: GET_MONTH
PUBLIC  :: GET_DAY
PUBLIC  :: GET_YEAR
PUBLIC  :: GET_HOUR
PUBLIC  :: GET_MINUTE
PUBLIC  :: GET_SECOND
PUBLIC  :: GET_DAY_OF_YEAR
PUBLIC  :: GET_DAY_OF_WEEK
PUBLIC  :: GET_GMT
PUBLIC  :: GET_TAU
PUBLIC  :: GET_TAUb
PUBLIC  :: GET_TAUe
PUBLIC  :: GET_DIAGb
PUBLIC  :: GET_DIAGe
PUBLIC  :: GET_LOCALTIME
PUBLIC  :: GET_SEASON
PUBLIC  :: GET_TS_CHEM
PUBLIC  :: GET_TS_CONV
PUBLIC  :: GET_TS_DIAG
PUBLIC  :: GET_TS_DYN
PUBLIC  :: GET_TS_EMIS
PUBLIC  :: GET_TS_UNIT

```

-----

Prior to 10/7/11:

```

PUBLIC  :: GET_TS_SUN_2

```

-----

```

PUBLIC  :: GET_CT_CHEM
PUBLIC  :: GET_CT_CONV
PUBLIC  :: GET_CT_DYN
PUBLIC  :: GET_CT_EMIS
PUBLIC  :: GET_CT_A1
PUBLIC  :: GET_CT_A3
PUBLIC  :: GET_CT_A6
PUBLIC  :: GET_CT_I6
PUBLIC  :: GET_CT_XTRA
PUBLIC  :: GET_CT_DIAG

```



```
PUBLIC  :: GET_A1_TIME
PUBLIC  :: GET_A3_TIME
PUBLIC  :: GET_A6_TIME
PUBLIC  :: GET_I6_TIME
PUBLIC  :: GET_FIRST_A1_TIME
PUBLIC  :: GET_FIRST_A3_TIME
PUBLIC  :: GET_FIRST_A6_TIME
PUBLIC  :: GET_FIRST_I6_TIME
PUBLIC  :: ITS_TIME_FOR_CHEM
PUBLIC  :: ITS_TIME_FOR_CONV
PUBLIC  :: ITS_TIME_FOR_DYN
PUBLIC  :: ITS_TIME_FOR_EMIS
PUBLIC  :: ITS_TIME_FOR_UNIT
PUBLIC  :: ITS_TIME_FOR_DIAG
PUBLIC  :: ITS_TIME_FOR_A1
PUBLIC  :: ITS_TIME_FOR_A3
PUBLIC  :: ITS_TIME_FOR_A6
PUBLIC  :: ITS_TIME_FOR_I6
PUBLIC  :: ITS_TIME_FOR_UNZIP
PUBLIC  :: ITS_TIME_FOR_DEL
PUBLIC  :: ITS_TIME_FOR_EXIT
PUBLIC  :: ITS_TIME_FOR_BPCH
PUBLIC  :: ITS_A_LEAPYEAR
PUBLIC  :: ITS_A_NEW_YEAR
PUBLIC  :: ITS_A_NEW_MONTH
PUBLIC  :: ITS_MIDMONTH
PUBLIC  :: ITS_A_NEW_DAY
PUBLIC  :: ITS_A_NEW_SEASON
PUBLIC  :: PRINT_CURRENT_TIME
PUBLIC  :: TIMESTAMP_STRING
PUBLIC  :: YMD_EXTRACT
PUBLIC  :: EXPAND_DATE
PUBLIC  :: SYSTEM_DATE_TIME
PUBLIC  :: SYSTEM_TIMESTAMP
PUBLIC  :: TIMESTAMP_DIAG
PUBLIC  :: GET_NYMD_DIAG
#if    defined( APM )
PUBLIC  :: ITS_TIME_FOR_A6UPDATE
#endif
```

## REVISION HISTORY:

- 21 Jun 2000 - R. Yantosca - Initial version
- (1 ) Updated comments (bmy, 9/4/01)
- (2 ) Added routine YMD\_EXTRACT. Also rewrote TIMECHECK using astronomical Julian day routines from "julday\_mod.f". (bmy, 11/21/01)
- (3 ) Eliminated obsolete code (bmy, 2/27/02)
- (4 ) Updated comments (bmy, 5/28/02)
- (5 ) Added routine "expand\_date". Also now reference "charpak\_mod.f".

- (bmy, 6/27/02)
- (6 ) Now references "error\_mod.f". Also added function GET\_SEASON, which returns the current season number. (bmy, 10/22/02)
  - (7 ) Now added module variables and various GET\_ and SET\_ routines to access them. Now minutes are the smallest timing unit. (bmy, 3/21/03)
  - (8 ) Bug fix in DATE\_STRING (bmy, 5/15/03)
  - (9 ) Added GET\_FIRST\_A3\_TIME and GET\_FIRST\_A6\_TIME. Also added changes for reading fvDAS fields. (bmy, 6/26/03)
  - (10) Now allow ITS\_A\_LEAPYEAR to take an optional argument. Bug fix for Linux: must use ENCODE to convert numbers to strings (bmy, 9/29/03)
  - (11) Bug fix in EXPAND\_DATE. Also add optional arguments to function TIMESTAMP\_STRNIG. (bmy, 10/28/03)
  - (12) Changed the name of some cpp switches in "define.h" (bmy, 12/2/03)
  - (13) Modified ITS\_TIME\_FOR\_A6 and GET\_FIRST\_A6\_TIME for both GEOS-4 "a\_llk\_03" and "a\_llk\_04" data versions. (bmy, 3/22/04)
  - (14) Added routines ITS\_A\_NEW\_MONTH, ITS\_A\_NEW\_YEAR, ITS\_A\_NEW\_DAY. (bmy, 4/1/04)
  - (15) Added routines ITS\_A\_NEW\_SEASON, GET\_NDIAGTIME, SET\_NDIAGTIME, and variable NDIAGTIME. (bmy, 7/20/04)
  - (17) Added routine GET\_DAY\_OF\_WEEK (bmy, 11/5/04)
  - (18) Removed obsolete FIRST variable in GET\_A3\_TIME (bmy, 12/10/04)
  - (19) Added routines SYSTEM\_DATE\_TIME and SYSTEM\_TIMESTAMP. Also modified for GCAP and GEOS-5 met fields. (swu, bmy, 5/3/05)
  - (20) GCAP/GISS met fields don't have leap years (swu, bmy, 8/29/05)
  - (21) Added counter variable & routines for XTRA fields (tmf, bmy, 10/20/05)
  - (22) Bug fix in ITS\_A\_NEW\_YEAR (bmy, 11/1/05)
  - (23) Added function ITS\_MIDMONTH. Also removed obsolete functions NYMD\_Y2K, NYMD6\_2\_NYMD8, NYMD\_STRING, DATE\_STRING. (sas, cdh, bmy, 12/15/05)
  - (24) GCAP bug fix: There are no leapyears, so transition from 2/28 to 3/1, skipping 2/29 for all years. (swu, bmy, 4/24/06)
  - (25) Remove support for GEOS-1 and GEOS-STRAT met fields (bmy, 8/4/06)
  - (26) Further bug fix to skip over Feb 29th in GCAP (phs, bmy, 10/3/06)
  - (27) Moved ITS\_TIME\_FOR\_BPCH here from "main.f" (bmy, 2/2/07)
  - (28) Add TS\_DIAG and CT\_DIAG variables to correctly output diagnostics (good time step).  
Add SET\_CT\_DIAG and GET\_CT\_DIAG to implement TS\_DIAG correctly. (ccc, 5/21/09)
  - (29) Add NYMD\_DIAG, GET\_NYMD\_DIAG, TIMESTAMP\_DIAG to get the good timestamp for diagnostic filenames (ccc, 8/12/09)
- 15 Jan 2010 - R. Yantosca - Added ProTeX headers
  - 27 Apr 2010 - R. Yantosca - Added OFFSET argument to GET\_LOCALTIME
  - 27 Apr 2010 - R. Yantosca - Added TS\_SUN\_2 to hold 1/2 of the interval for computing SUNCOS.
  - 27 Apr 2010 - R. Yantosca - Added public routine GET\_TS\_SUN\_2
  - 19 Aug 2010 - R. Yantosca - Added variable CT\_A1 and routine SET\_CT\_A1
  - 20 Aug 2010 - R. Yantosca - Added function ITS\_TIME\_FOR\_A1
  - 27 Sep 2010 - R. Yantosca - Added function GET\_FIRST\_I6\_TIME

17 Dec 2010 - R. Yantosca - Bug fix for HHMMSS=240000 in GET\_TIME\_AHEAD  
 27 Mar 2011 - R. Yantosca - Bug fix for GCAP leap year problem  
 29 Jul 2011 - R. Yantosca - Add LEAP\_YEAR\_DAYS as a SAVED module variable  
 17 Feb 2011 - R. Yantosca - Added ITS\_TIME\_FOR\_A6UPDATE for APM (G. Luo)  
 07 Oct 2011 - M. Payer - Modifications for central chemistry timestep  
 07 Oct 2011 - R. Yantosca - Remove obsolete TS\_SUN\_2, GET\_TS\_SUN\_2  
 07 Oct 2011 - R. Yantosca - Remove obsolete OFFSET argument to GET\_LOCALTIME  
 12 Oct 2011 - R. Yantosca - Modified ITS\_A\_NEW\_MONTH for central chem step

---

### 1.10.1 set\_current\_time

Subroutine SET\_CURRENT\_TIME takes in the elapsed time in minutes since the start of a GEOS-Chem simulation and sets the GEOS-Chem time variables accordingly.

#### INTERFACE:

```
SUBROUTINE SET_CURRENT_TIME
```

#### USES:

```
USE JULDAY_MOD, ONLY : JULDAY, CALDATE
```

```
#    include "define.h"
```

#### REMARKS:

The GEOS met fields are assimilated data, and therefore contain data on the leap-year days. However, the GCAP met fields are climatological GCM output, and do not have data on the leap-year days. SET\_CURRENT\_TIME computes the days according to the Astronomical Julian Date algorithms (in "julday\_mod.f"), which contain leap-year days. For GCAP, whenever a February 29th is encountered, we shall just skip ahead a day to March 1st and return the corresponding time & date values.

#### REVISION HISTORY:

05 Feb 2006 - R. Yantosca - Initial Version  
 (1 ) GCAP/GISS fields don't have leap years, so if JULDAY says it's  
     Feb 29th, reset MONTH, DAY, JD1 to Mar 1st. (swu, bmy, 8/29/05)  
 (2 ) Now references "define.h". Now add special handling to skip from  
     Feb 28th to Mar 1st for GCAP model. (swu, bmy, 4/24/06)  
 (3 ) Fix bug in case of GCAP fields for runs that start during leap year  
     and after February 29 (phs, 9/27/06)  
 15 Jan 2010 - R. Yantosca - Added ProTeX headers  
 29 Jul 2011 - R. Yantosca - Bug fix: For GCAP, we need to skip over the  
     # of leap-year-days that have already occurred  
     when going from Julian date to Y/M/D date

---

### 1.10.2 set\_begin\_time

Subroutine SET\_BEGIN\_TIME initializes NYMDb, NHMSb, and TAUb, which are the YYYYMMDD, HHMMSS, and hours since 1/1/1985 corresponding to the beginning date and time of a GEOS-Chem run.

#### INTERFACE:

```
SUBROUTINE SET_BEGIN_TIME( THISNYMDb, THISNHMSb )
```

#### USES:

```
USE ERROR_MOD, ONLY : ERROR_STOP
```

#### INPUT PARAMETERS:

```
INTEGER, INTENT(IN) :: THISNYMDb    ! YYYYMMDD @ start of G-C simulation
INTEGER, INTENT(IN) :: THISNHMSb    ! HHMMSS   @ start of G-C simulation
```

#### REVISION HISTORY:

```
20 Jul 2004 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
16 Dec 2010 - R. Yantosca - Updated error check for THISNYMDe, since
                           MERRA met data goes back prior to 1985
```

---

### 1.10.3 set\_end\_time

Subroutine SET\_END\_TIME initializes NYMDe, NHMSe, and TAUe, which are the YYYYMMDD, HHMMSS, and hours since 1/1/1985 corresponding to the ending date and time of a GEOS-Chem run.

#### INTERFACE:

```
SUBROUTINE SET_END_TIME( THISNYMDe, THISNHMSe )
```

#### USES:

```
USE ERROR_MOD, ONLY : ERROR_STOP
```

#### INPUT PARAMETERS:

```
INTEGER, INTENT(IN) :: THISNYMDe    ! YYYYMMDD @ end of G-C simulation
INTEGER, INTENT(IN) :: THISNHMSe    ! HHMMSS   @ end of G-C simulation
```

#### REVISION HISTORY:

```
20 Jul 2004 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
16 Dec 2010 - R. Yantosca - Updated error check for THISNYMDe, since
                           MERRA met data goes back prior to 1985
```

---

**1.10.4 set\_ndiagtime**

SET\_NDIAGTIME initializes NDIAGTIME, the time of day at which the binary punch file will be written out to disk.

**INTERFACE:**

```
SUBROUTINE SET_NDIAGTIME( THIS_NDIAGTIME )
```

**INPUT PARAMETERS:**

```
INTEGER, INTENT(IN) :: THIS_NDIAGTIME ! Initial NDIAGTIMEe [hrs]
```

**REVISION HISTORY:**

```
20 Jul 2004 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.5 set\_diagb**

Subroutine SET\_DIAGb initializes DIAGb, the TAU value at the start of the diagnostic averaging interval.

**INTERFACE:**

```
SUBROUTINE SET_DIAGb( THISDIAGb )
```

**INPUT PARAMETERS:**

```
REAL*8, INTENT(IN) :: THISDIAGb ! Initial DIAGb value [hrs from 1/1/85]
```

**REVISION HISTORY:**

```
21 Mar 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.6 set\_diage**

Subroutine SET\_DIAGe initializes DIAGe, the TAU value at the end of the diagnostic averaging interval.

**INTERFACE:**

```
SUBROUTINE SET_DIAGe( THISDIAGe )
```

**INPUT PARAMETERS:**

```
REAL*8, INTENT(IN) :: THISDIAGe ! Initial DIAGe value [hrs from 1/1/85]
```

**REVISION HISTORY:**

```
21 Mar 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

### 1.10.7 set\_timesteps

Subroutine SET\_TIMESTEPS initializes the timesteps for dynamics, convection, chemistry, emissions, and diagnostics. Counters are also zeroed.

#### INTERFACE:

```

      SUBROUTINE SET_TIMESTEPS( CHEMISTRY, CONVECTION, DYNAMICS,
&                               EMISSION,  UNIT_CONV,  DIAGNOS )

```

---

Prior to 10/5/11:

```

      Remove SUNCOS argument, it's obsolete (bmy, 10/5/11)
&                               SUNCOS )

```

---

#### INPUT PARAMETERS:

```

      INTEGER, INTENT(IN) :: CHEMISTRY    ! Chemistry timestep [min]
      INTEGER, INTENT(IN) :: CONVECTION   ! Convection timestep [min]
      INTEGER, INTENT(IN) :: DYNAMICS     ! Dynamic timestep [min]
      INTEGER, INTENT(IN) :: EMISSION     ! Emission timestep [min]
      INTEGER, INTENT(IN) :: UNIT_CONV    ! Unit conve timestep [min]
      INTEGER, INTENT(IN) :: DIAGNOS      ! Diagnostic timestep [min]

```

---

Prior to 10/5/11:

```

      Remove SUNCOS argument, it's obsolete (bmy, 10/5/11)
      INTEGER, INTENT(IN) :: SUNCOS       ! 1/2 of timestep for SUNCOS [min]

```

---

#### REVISION HISTORY:

```

      05 Feb 2003 - R. Yantosca - Initial Version
      (1 ) Suppress some output lines (bmy, 7/20/04)
      (2 ) Also zero CT_XTRA (tmf, bmy, 10/20/05)
      (3 ) Add TS_DIAG as the diagnostic timestep. (ccc, 5/13/09)
      15 Jan 2010 - R. Yantosca - Added ProTeX headers
      27 Apr 2010 - R. Yantosca - Now add SUNCOS argument to set 1/2 of the
                                interval for computing the cosine of the
                                solar zenith angle.
      07 Oct 2011 - R. Yantosca - Remove obsolete SUNCOS argument

```

---

### 1.10.8 set\_ct\_chem

Subroutine SET\_CT\_CHEM increments CT\_CHEM, the counter of chemistry timesteps executed thus far.

#### INTERFACE:

```

      SUBROUTINE SET_CT_CHEM( INCREMENT, RESET )

```

**INPUT PARAMETERS:**

```

LOGICAL, INTENT(IN), OPTIONAL :: INCREMENT ! Increment counter?
LOGICAL, INTENT(IN), OPTIONAL :: RESET     ! Reset counter?

```

**REVISION HISTORY:**

```

21 Mar 2009 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers

```

---

**1.10.9 set\_ct\_conv**

Subroutine SET\_CT\_CONV increments CT\_CONV, the counter of convection timesteps executed thus far.

**INTERFACE:**

```

SUBROUTINE SET_CT_CONV( INCREMENT, RESET )

```

**INPUT PARAMETERS:**

```

LOGICAL, INTENT(IN), OPTIONAL :: INCREMENT ! Increment counter?
LOGICAL, INTENT(IN), OPTIONAL :: RESET     ! Reset counter?

```

**REVISION HISTORY:**

```

21 Mar 2009 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers

```

---

**1.10.10 set\_ct\_dyn**

Subroutine SET\_CT\_DYN increments CT\_DYN, the counter of dynamical timesteps executed thus far.

**INTERFACE:**

```

SUBROUTINE SET_CT_DYN( INCREMENT, RESET )

```

**INPUT PARAMETERS:**

```

LOGICAL, INTENT(IN), OPTIONAL :: INCREMENT ! Increment counter?
LOGICAL, INTENT(IN), OPTIONAL :: RESET     ! Reset counter?

```

**REVISION HISTORY:**

```

21 Mar 2009 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers

```

---

### 1.10.11 set\_ct\_emis

Subroutine SET\_CT\_EMIS increments CT\_EMIS, the counter of emission timesteps executed thus far.

#### INTERFACE:

```
SUBROUTINE SET_CT_EMIS( INCREMENT, RESET )
```

#### INPUT PARAMETERS:

```
LOGICAL, INTENT(IN), OPTIONAL :: INCREMENT ! Increment counter?  
LOGICAL, INTENT(IN), OPTIONAL :: RESET      ! Reset counter?
```

#### REVISION HISTORY:

```
21 Mar 2009 - R. Yantosca - Initial Version  
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

### 1.10.12 set\_ct\_diag

Subroutine SET\_CT\_DIAG increments CT\_DIAG, the counter of largest timesteps executed thus far.

#### INTERFACE:

```
SUBROUTINE SET_CT_DIAG( INCREMENT, RESET )
```

#### INPUT PARAMETERS:

```
LOGICAL, INTENT(IN), OPTIONAL :: INCREMENT ! Increment counter?  
LOGICAL, INTENT(IN), OPTIONAL :: RESET      ! Reset counter?
```

#### REVISION HISTORY:

```
13 May 2009 - C. Carouge - Initial Version  
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

### 1.10.13 set\_ct\_a1

Subroutine SET\_CT\_A1 increments CT\_A1, the counter of the number of times we have read in A1 fields.

#### INTERFACE:

```
SUBROUTINE SET_CT_A1( INCREMENT, RESET )
```

#### INPUT PARAMETERS:



```

LOGICAL, INTENT(IN), OPTIONAL :: INCREMENT ! Increment counter?
LOGICAL, INTENT(IN), OPTIONAL :: RESET      ! Reset counter?

```

## REVISION HISTORY:

19 Aug 2010 - R. Yantosca - Initial version

---

### 1.10.14 set\_ct\_a3

Subroutine SET\_CT\_A3 increments CT\_A3, the counter of the number of times we have read in A-3 fields.

## INTERFACE:

```

SUBROUTINE SET_CT_A3( INCREMENT, RESET )

```

## INPUT PARAMETERS:

```

LOGICAL, INTENT(IN), OPTIONAL :: INCREMENT ! Increment counter?
LOGICAL, INTENT(IN), OPTIONAL :: RESET      ! Reset counter?

```

## REVISION HISTORY:

21 Mar 2003 - R. Yantosca - Initial Version  
 15 Jan 2010 - R. Yantosca - Added ProTeX headers

---

### 1.10.15 set\_ct\_a6

Subroutine SET\_CT\_A6 increments CT\_A6, the counter of the number of times we have read in A-6 fields.

## INTERFACE:

```

SUBROUTINE SET_CT_A6( INCREMENT, RESET )

```

## INPUT PARAMETERS:

```

LOGICAL, INTENT(IN), OPTIONAL :: INCREMENT ! Increment counter?
LOGICAL, INTENT(IN), OPTIONAL :: RESET      ! Reset counter?

```

## REVISION HISTORY:

21 Mar 2003 - R. Yantosca - Initial Version  
 15 Jan 2010 - R. Yantosca - Added ProTeX headers

---

**1.10.16 set\_ct\_i6**

Subroutine SET\_CT\_I6 increments CT\_I6, the counter of the number of times we have read in I-6 fields.

**INTERFACE:**

```
SUBROUTINE SET_CT_I6( INCREMENT, RESET )
```

**INPUT PARAMETERS:**

```
LOGICAL, INTENT(IN), OPTIONAL :: INCREMENT ! Increment counter?
LOGICAL, INTENT(IN), OPTIONAL :: RESET      ! Reset counter?
```

**REVISION HISTORY:**

```
21 Mar 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.17 set\_ct\_xtra**

Subroutine SET\_CT\_XTRA increments CT\_XTRA, the counter of the number of times we have read in GEOS-3 XTRA fields.

**INTERFACE:**

```
SUBROUTINE SET_CT_XTRA( INCREMENT, RESET )
```

**INPUT PARAMETERS:**

```
LOGICAL, INTENT(IN), OPTIONAL :: INCREMENT ! Increment counter?
LOGICAL, INTENT(IN), OPTIONAL :: RESET      ! Reset counter?
```

**REVISION HISTORY:**

```
20 Oct 2009 - T-M Fu, R. Yantosca - Initial Version
15 Jan 2010 -           R. Yantosca - Added ProTeX headers
```

---

**1.10.18 set\_elapsed\_min**

Subroutine SET\_ELAPSED\_MIN increments the number of elapsed minutes by the dynamic timestep TS\_DYN.

**INTERFACE:**

```
SUBROUTINE SET_ELAPSED_MIN
```

**REVISION HISTORY:**

```
05 Feb 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.19 get\_jd**

Function GET\_JD is a wrapper for the JULDAY routine. Given the current NYMD and NHMS values, GET\_JD will return the current astronomical Julian date.

**INTERFACE:**

```
FUNCTION GET_JD( THISNYMD, THISNHMS ) RESULT( THISJD )
```

**USES:**

```
USE JULDAY_MOD, ONLY : JULDAY
```

**INPUT PARAMETERS:**

```
INTEGER, INTENT(IN)  :: THISNYMD    ! YYYY/MM/DD value
INTEGER, INTENT(IN)  :: THISNHMS    ! hh:mm:ss  value
```

**RETURN VALUE:**

```
REAL*8                :: THISJD      ! Output value
```

**REVISION HISTORY:**

```
05 Feb 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.20 get\_elapsed\_min**

Function GET\_ELAPSED\_MIN returns the elapsed minutes since the start of a GEOS-chem run.

**INTERFACE:**

```
FUNCTION GET_ELAPSED_MIN() RESULT( THIS_ELAPSED_MIN )
```

**RETURN VALUE:**

```
INTEGER :: THIS_ELAPSED_MIN
```

**REVISION HISTORY:**

```
05 Feb 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.21 get\_elapsed\_sec**

Function GET\_ELAPSED\_SEC returns the elapsed minutes since the start of a GEOS-Chem run to the calling program.

**INTERFACE:**

```
FUNCTION GET_ELAPSED_SEC() RESULT( THIS_ELAPSED_SEC )
```

**RETURN VALUE:**

```
INTEGER :: THIS_ELAPSED_SEC
```

**REVISION HISTORY:**

```
05 Feb 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.22 get\_nymdb**

Function GET\_NYMDB returns the NYMDB value (YYYYMMDD at the beginning of the run).

**INTERFACE:**

```
FUNCTION GET_NYMDB() RESULT( THISNYMDB )
```

**RETURN VALUE:**

```
INTEGER :: THISNYMDB
```

**REVISION HISTORY:**

```
05 Feb 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.23 get\_nhmsb**

Function GET\_NHMSb returns the NHMSb value (HHMMSS at the beginning of the run) to the calling program. (bmy, 3/21/03)

**INTERFACE:**

```
FUNCTION GET_NHMSb() RESULT( THISNHMSb )
```

**RETURN VALUE:**

```
INTEGER :: THISNHMSb
```

**REVISION HISTORY:**

```
05 Feb 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

### 1.10.24 get\_nymde

Function GET\_NYMDe returns the NYMDe value (YYYYMMDD at the end of the run) to the calling program. (bmy, 3/21/03)

#### INTERFACE:

```
FUNCTION GET_NYMDe() RESULT( THISNYMDe )
```

#### RETURN VALUE:

```
INTEGER :: THISNYMDe
```

#### REVISION HISTORY:

```
05 Feb 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

### 1.10.25 get\_nhmse

Function GET\_NHMSe returns the NHMSe value (HHMMSS at the end of the run).

#### INTERFACE:

```
FUNCTION GET_NHMSe() RESULT( THISNHMSe )
```

#### RETURN VALUE:

```
INTEGER :: THISNHMSe
```

#### REVISION HISTORY:

```
05 Feb 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

### 1.10.26 get\_nymd

Function GET\_NYMD returns the current NYMD value (YYYYMMDD).

#### INTERFACE:

```
FUNCTION GET_NYMD() RESULT( THISNYMD )
```

#### RETURN VALUE:

```
INTEGER :: THISNYMD
```

#### REVISION HISTORY:

```
05 Feb 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.27 get\_nhms**

Function GET\_NHMS returns the current NHMS value (HHMMSS).

**INTERFACE:**

```
FUNCTION GET_NHMS() RESULT( THISNHMS )
```

**RETURN VALUE:**

```
INTEGER :: THISNHMS
```

**REVISION HISTORY:**

```
05 Feb 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.28 get\_ndiagtime**

Subroutine GET\_NDIAGTIME returns to the calling program NDIAGTIME, the time of day at which the binary punch file will be written out to disk.

**INTERFACE:**

```
FUNCTION GET_NDIAGTIME() RESULT( THIS_NDIAGTIME )
```

**RETURN VALUE:**

```
INTEGER :: THIS_NDIAGTIME
```

**REVISION HISTORY:**

```
20 Jul 2004 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.29 get\_time\_ahead**

Function GET\_3h\_AHEAD returns to the calling program a 2-element vector containing the YYYYMMDD and HHMMSS values at the current time plus N\_MINS minutes.

**INTERFACE:**

```
FUNCTION GET_TIME_AHEAD( N_MINS ) RESULT( DATE )
```

**USES:**

```
USE JULDAY_MOD, ONLY : CALDATE
```

```
#    include "define.h"    ! C-preprocessor flags
```

**INPUT PARAMETERS:**

```
INTEGER, INTENT(IN) :: N_MINS    ! Minutes ahead to compute date & time
```

**RETURN VALUE:**

```
INTEGER                :: DATE(2) ! Date & time output
```

**REVISION HISTORY:**

```
21 Mar 2003 - R. Yantosca - Initial Version
(1 ) Bug fix for GCAP leap year case (phs, bmy, 12/8/06)
15 Jan 2010 - R. Yantosca - Added ProTeX headers
17 Dec 2010 - R. Yantosca - Added fix in case HHMMSS is returned as 240000
```

---

**1.10.30 get\_month**

Function GET\_MONTH returns the current GMT month.

**INTERFACE:**

```
FUNCTION GET_MONTH() RESULT( THISMONTH )
```

**RETURN VALUE:**

```
INTEGER :: THISMONTH
```

**REVISION HISTORY:**

```
05 Feb 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.31 get\_day**

Function GET\_DAY returns the current GMT day.

**INTERFACE:**

```
FUNCTION GET_DAY() RESULT( THISDAY )
```

**RETURN VALUE:**

```
INTEGER :: THISDAY
```

**REVISION HISTORY:**

```
05 Feb 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.32 get\_year**

Function GET\_YEAR returns the current GMT year.

**INTERFACE:**

```
FUNCTION GET_YEAR() RESULT( THISYEAR )
```

**RETURN VALUE:**

```
INTEGER :: THISYEAR
```

**REVISION HISTORY:**

```
05 Feb 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.33 get\_hour**

Function GET\_HOUR returns the current GMT hour.

**INTERFACE:**

```
FUNCTION GET_HOUR() RESULT( THISHOUR )
```

**RETURN VALUE:**

```
INTEGER :: THISHOUR
```

**REVISION HISTORY:**

```
05 Feb 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.34 get\_minute**

Function GET\_MINUTE returns the current GMT minutes.

**INTERFACE:**

```
FUNCTION GET_MINUTE() RESULT( THISMINUTE )
```

**RETURN VALUE:**

```
INTEGER :: THISMINUTE
```

**REVISION HISTORY:**

```
05 Feb 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---



**1.10.35 get\_second**

Function GET\_SECOND returns the current GMT seconds. calling program.

**INTERFACE:**

```
FUNCTION GET_SECOND() RESULT( THISSECOND )
```

**RETURN VALUE:**

```
INTEGER :: THISSECOND
```

**REVISION HISTORY:**

```
05 Feb 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.36 get\_day\_of\_year**

Function GET\_DAY\_OF\_YEAR returns the current day of the year (0-365 or 0-366 for leap years) to the calling program.

**INTERFACE:**

```
FUNCTION GET_DAY_OF_YEAR() RESULT( THISDAYOFEAR )
```

**RETURN VALUE:**

```
INTEGER :: THISDAYOFEAR ! Day of year
```

**REVISION HISTORY:**

```
05 Feb 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.37 get\_day\_of\_week**

Function GET\_DAY\_OF\_WEEK returns the day of the week as a number: Sun=0, Mon=1, Tue=2, Wed=3, Thu=4, Fri=5, Sat=6.

**INTERFACE:**

```
FUNCTION GET_DAY_OF_WEEK() RESULT( DAY_NUM )
```

**USES:**

```
USE JULDAY_MOD, ONLY : JULDAY
```

**RETURN VALUE:**

INTEGER :: DAY\_NUM ! Day number of week

## REMARKS:

Reference:

-----

"Practical Astronomy with Your Calculator", 3rd Ed. Peter Duffett-Smith,  
Cambridge UP, 1992, p9.

## REVISION HISTORY:

05 Feb 2003 - R. Yantosca - Initial Version  
15 Jan 2010 - R. Yantosca - Added ProTeX headers

### 1.10.38 get\_gmt

Function GET\_GMT returns the current Greenwich Mean Time to the calling program.

## INTERFACE:

FUNCTION GET\_GMT() RESULT( THISGMT )

## RETURN VALUE:

REAL\*8 :: THISGMT ! Greenwich mean time [hrs]

## REVISION HISTORY:

05 Feb 2003 - R. Yantosca - Initial Version  
15 Jan 2010 - R. Yantosca - Added ProTeX headers

### 1.10.39 get\_tau

Function GET\_TAU returns TAU (hours since 1 Jan 1985 at the start of a GEOS-Chem run) to the calling program.

## INTERFACE:

FUNCTION GET\_TAU() RESULT( THISTAU )

## RETURN VALUE:

REAL\*8 :: THISTAU ! TAUb [hrs since 1/1/1985]

## REVISION HISTORY:

05 Feb 2003 - R. Yantosca - Initial Version  
15 Jan 2010 - R. Yantosca - Added ProTeX headers

#### 1.10.40 `get_taub`

Function GET\_TAUb returns TAUb (hours since 1 Jan 1985 at the start of a GEOS-Chem run) to the calling program.

##### INTERFACE:

```
FUNCTION GET_TAUb() RESULT( THISTAUb )
```

##### RETURN VALUE:

```
REAL*8 :: THISTAUb  ! TAUb [hrs since 1/1/1985]
```

##### REVISION HISTORY:

```
05 Feb 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

#### 1.10.41 `get_tae`

Function GET\_TAUe returns TAUe (hours since 1 Jan 1985 at the end of a GEOS-Chem run) to the calling program.

##### INTERFACE:

```
FUNCTION GET_TAUe() RESULT( THISTAUe )
```

##### RETURN VALUE:

```
REAL*8 :: THISTAUe  ! TAUe [hrs since 1/1/1985]
```

##### REVISION HISTORY:

```
05 Feb 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

#### 1.10.42 `get_diagb`

Function GET\_DIAGb returns DIAGb (hours since 1 Jan 1985 at the start of a diagnostic interval) to the calling program.

##### INTERFACE:

```
FUNCTION GET_DIAGb() RESULT( THISDIAGb )
```

##### RETURN VALUE:

```
INTEGER :: THISDIAGb  ! DIAGb [hrs since 1/1/1985]
```

##### REVISION HISTORY:

```
05 Feb 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.43 get\_diage**

Function GET\_DIAGe returns DIAGe (hours since 1 Jan 1985 at the end of a diagnostic interval) to the calling program.

**INTERFACE:**

```
FUNCTION GET_DIAGe() RESULT( THISDIAGe )
```

**RETURN VALUE:**

```
INTEGER :: THISDIAGe    ! DIAGe [hrs sincd 1/1/1985]
```

**REVISION HISTORY:**

```
05 Feb 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.44 get\_localtime**

Function GET\_LOCALTIME returns the local time of a grid box to the calling program. (bmy, 2/5/03)

**INTERFACE:**

```
FUNCTION GET_LOCALTIME( I, GMT ) RESULT( THISLOCALTIME )
```

**USES:**

```
USE GRID_MOD, ONLY : GET_XMID
```

**INPUT PARAMETERS:**

```
INTEGER, INTENT(IN)           :: I           ! Longitude index
REAL*8,  INTENT(IN), OPTIONAL :: GMT         ! GMT time of day [hrs]
```

**RETURN VALUE:**

```
REAL*8                        :: THISLOCALTIME ! Local time [hrs]
```

**REMARKS:**

```
Local Time = GMT + ( longitude / 15 ) since each hour of time
corresponds to 15 degrees of longitude on the globe
```

**REVISION HISTORY:**

```
05 Feb 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
27 Apr 2010 - R. Yantosca - Add OFFSET to argument list, to allow the
                           local time to be computed at an arbitrary time
                           (e.g. at the halfway point of an interval)
05 Oct 2011 - R. Yantosca - Now add GMT as an optional argument
07 Oct 2011 - R. Yantosca - Removed obsolete OFFSET argument
```

---

**1.10.45 get\_season**

Function GET\_SEASON returns the climatological season number (1=DJF, 2=MAM, 3=JJA, 4=SON) to the calling program.

**INTERFACE:**

```
FUNCTION GET_SEASON() RESULT( THISSEASON )
```

**RETURN VALUE:**

```
INTEGER :: THISSEASON    ! Current season
```

**REVISION HISTORY:**

```
21 Mar 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.46 get\_ts\_chem**

Function GET\_TS\_CHEM returns the chemistry timestep in minutes.

**INTERFACE:**

```
FUNCTION GET_TS_CHEM() RESULT( THIS_TS_CHEM )
```

**RETURN VALUE:**

```
INTEGER :: THIS_TS_CHEM    ! ! Chemistry timestep [min]
```

**REVISION HISTORY:**

```
21 Mar 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.47 get\_ts\_conv**

Function GET\_TS\_CONV returns the convection timestep in minutes.

**INTERFACE:**

```
FUNCTION GET_TS_CONV() RESULT( THIS_TS_CONV )
```

**RETURN VALUE:**

```
INTEGER :: THIS_TS_CONV    ! Convective timestep [min]
```

**REVISION HISTORY:**

```
21 Mar 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.48 get\_ts\_diag**

Function GET\_TS\_DIAG returns the diagnostic timestep in minutes.

**INTERFACE:**

```
FUNCTION GET_TS_DIAG() RESULT( THIS_TS_DIAG )
```

**RETURN VALUE:**

```
INTEGER :: THIS_TS_DIAG    ! Diagnostic timestep [min]
```

**REVISION HISTORY:**

```
21 Mar 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.49 get\_ts\_dyn**

Function GET\_TS\_DIAG returns the diagnostic timestep in minutes.

**INTERFACE:**

```
FUNCTION GET_TS_DYN() RESULT( THIS_TS_DYN )
```

**RETURN VALUE:**

```
INTEGER :: THIS_TS_DYN    ! Dynamic timestep [min]
```

**REVISION HISTORY:**

```
21 Mar 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.50 get\_ts\_emis**

Function GET\_TS\_EMIS returns the emission timestep in minutes.

**INTERFACE:**

```
FUNCTION GET_TS_EMIS() RESULT( THIS_TS_EMIS )
```

**RETURN VALUE:**

```
INTEGER :: THIS_TS_EMIS    ! Emissions timestep [min]
```

**REVISION HISTORY:**

```
21 Mar 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.51 get\_ts\_unit**

Function GET\_TS\_UNIT returns the unit-conversion timestep in minutes.

**INTERFACE:**

```
FUNCTION GET_TS_UNIT() RESULT( THIS_TS_UNIT )
```

**RETURN VALUE:**

```
INTEGER :: THIS_TS_UNIT    ! Unit conversion timestep [min]
```

**REVISION HISTORY:**

```
21 Mar 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.52 get\_ct\_chem**

Function GET\_CT\_CHEM returns the chemistry timestep counter to the calling program.

**INTERFACE:**

```
FUNCTION GET_CT_CHEM() RESULT( THIS_CT_CHEM )
```

**RETURN VALUE:**

```
INTEGER :: THIS_CT_CHEM
```

**REVISION HISTORY:**

```
21 Mar 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.53 get\_ct\_conv**

Function GET\_CT\_CONV returns the convection timestep counter to the calling program.

**INTERFACE:**

```
FUNCTION GET_CT_CONV() RESULT( THIS_CT_CONV )
```

**RETURN VALUE:**

```
INTEGER :: THIS_CT_CONV    ! # of convection timesteps
```

**REVISION HISTORY:**

```
21 Mar 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

### 1.10.54 get\_ct\_dyn

Function GET\_CT\_CHEM returns the dynamic timestep counter to the calling program.

#### INTERFACE:

```
FUNCTION GET_CT_DYN() RESULT( THIS_CT_DYN )
```

#### RETURN VALUE:

```
INTEGER :: THIS_CT_DYN    ! # of dynamics timesteps
```

#### REVISION HISTORY:

```
21 Mar 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

### 1.10.55 get\_ct\_emis

Function GET\_CT\_CHEM returns the emissions timestep counter to the calling program.

#### INTERFACE:

```
FUNCTION GET_CT_EMIS() RESULT( THIS_CT_EMIS )
```

#### RETURN VALUE:

```
INTEGER :: THIS_CT_EMIS  ! # of emissions timesteps
```

#### REVISION HISTORY:

```
21 Mar 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

### 1.10.56 get\_ct\_a1

Function GET\_CT\_A1 returns the A1 fields timestep counter to the calling program.

#### INTERFACE:

```
FUNCTION GET_CT_A1() RESULT( THIS_CT_A1 )
```

#### RETURN VALUE:

```
INTEGER :: THIS_CT_A1    ! # of A-3 timesteps
```

#### REVISION HISTORY:

```
19 Aug 2010 - R. Yantosca - Initial version
```

---



**1.10.57 get\_ct\_a3**

Function GET\_CT\_A3 returns the A-3 fields timestep counter to the calling program.

**INTERFACE:**

```
FUNCTION GET_CT_A3() RESULT( THIS_CT_A3 )
```

**RETURN VALUE:**

```
INTEGER :: THIS_CT_A3    ! # of A-3 timesteps
```

**REVISION HISTORY:**

```
21 Mar 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.58 get\_ct\_a6**

Function GET\_CT\_A6 returns the A-6 fields timestep counter to the calling program.

**INTERFACE:**

```
FUNCTION GET_CT_A6() RESULT( THIS_CT_A6 )
```

**RETURN VALUE:**

```
INTEGER :: THIS_CT_A6    ! # of A-6 timesteps
```

**REVISION HISTORY:**

```
21 Mar 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.59 get\_ct\_i6**

Function GET\_CT\_I6 returns the I-6 fields timestep counter to the calling program

**INTERFACE:**

```
FUNCTION GET_CT_I6() RESULT( THIS_CT_I6 )
```

**RETURN VALUE:**

```
INTEGER :: THIS_CT_I6    ! # of I-6 timesteps
```

**REVISION HISTORY:**

```
21 Mar 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.60 get\_ct\_xtra**

Function GET\_CT\_XTRA returns the XTRA fields timestep counter to the calling program.

**INTERFACE:**

```
FUNCTION GET_CT_XTRA() RESULT( THIS_CT_XTRA )
```

**RETURN VALUE:**

```
INTEGER :: THIS_CT_XTRA      ! # of XTRA timesteps
```

**REVISION HISTORY:**

```
20 Oct 2005 - T-M Fu, R. Yantosca - Initial Version
15 Jan 2010 -           R. Yantosca - Added ProTeX headers
```

---

**1.10.61 get\_ct\_diag**

Function GET\_CT\_DIAG returns the DIAG timestep counter to the calling program.

**INTERFACE:**

```
FUNCTION GET_CT_DIAG() RESULT( THIS_CT_DIAG )
```

**RETURN VALUE:**

```
INTEGER :: THIS_CT_DIAG      ! # of diagnostic timesteps
!
```

**REVISION HISTORY:**

```
21 May 2009 - C. Carouge - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.62 get\_a1\_time**

Function GET\_A1\_TIME returns the correct YYYYMMDD and HHMMSS values that are needed to read in the next average 1-hour (A-1) fields.

**INTERFACE:**

```
FUNCTION GET_A1_TIME() RESULT( DATE )
```

**USES:**

```
#      include "define.h"
```

**RETURN VALUE:**

```
INTEGER :: DATE(2)      ! YYYYMMDD and HHMMSS values
```

**REVISION HISTORY:**

```
19 Aug 2010 - R. Yantosca - Initial version
```

---

**1.10.63 get\_a3\_time**

Function GET\_A3\_TIME returns the correct YYYYMMDD and HHMMSS values that are needed to read in the next average 3-hour (A-3) fields.

**INTERFACE:**

```
FUNCTION GET_A3_TIME() RESULT( DATE )
```

**USES:**

```
#    include "define.h"
```

**RETURN VALUE:**

```
INTEGER :: DATE(2)    ! YYYYMMDD and HHMMSS values
```

**REVISION HISTORY:**

```
21 Mar 2003 - R. Yantosca - Initial Version
(1 ) Now return proper time for GEOS-4/fvDAS fields (bmy, 6/19/03)
(2 ) Remove reference to FIRST variable (bmy, 12/10/04)
(3 ) Now modified for GCAP and GEOS-5 met fields (swu, bmy, 5/24/05)
(4 ) Remove support for GEOS-1 and GEOS-STRAT met fields (bmy, 8/4/06)

15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.64 get\_a6\_time**

Function GET\_A6\_TIME returns the correct YYYYMMDD and HHMMSS values that are needed to read in the next average 6-hour (A-6) fields.

**INTERFACE:**

```
FUNCTION GET_A6_TIME() RESULT( DATE )
```

**RETURN VALUE:**

```
INTEGER :: DATE(2)    ! YYYYMMDD and HHMMSS time
```

**REVISION HISTORY:**

```
21 Mar 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
17 Feb 2011 - R. Yantosca - Add modifications for APM microphysics (G. Luo)
```

---

**1.10.65 get\_i6\_time**

Function GET\_I6\_TIME returns the correct YYYYMMDD and HHMMSS values that are needed to read in the next instantaneous 6-hour (I-6) fields.

**INTERFACE:**

```
FUNCTION GET_I6_TIME() RESULT( DATE )
```

**RETURN VALUE:**

```
INTEGER :: DATE(2)    ! YYYYMMDD and HHMMSS values
```

**REMARKS:**

Modified for start times other than 0 GMT. However someone should check to make sure it works properly for the GCAP simulation. (bmy, 9/27/10)

**REVISION HISTORY:**

```
21 Mar 2003 - R. Yantosca - Initial Version
(1 ) Bug fix for GCAP: skip over Feb 29th (no leapyears). (bmy, 4/24/06)
15 Jan 2010 - R. Yantosca - Added ProTeX headers
27 Sep 2010 - R. Yantosca - Now works for start times other than 0 GMT
```

---

**1.10.66 get\_first\_a1\_time**

Function GET\_FIRST\_A1\_TIME returns the correct YYYYMMDD and HHMMSS values the first time that A-3 fields are read in from disk.

**INTERFACE:**

```
FUNCTION GET_FIRST_A1_TIME() RESULT( DATE )
```

**USES:**

```
#    include "define.h"
```

**RETURN VALUE:**

```
INTEGER :: DATE(2)    ! YYYYMMDD and HHMMSS values
```

**REVISION HISTORY:**

```
26 Jun 2003 - R. Yantosca - Initial Version
(1 ) Now modified for GCAP and GEOS-5 data (swu, bmy, 5/24/05)
(2 ) Remove support for GEOS-1 and GEOS-STRAT met fields (bmy, 8/4/06)
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.67 get\_first\_a3\_time**

Function GET\_FIRST\_A3\_TIME returns the correct YYYYMMDD and HHMMSS values the first time that A-3 fields are read in from disk.

**INTERFACE:**

```
FUNCTION GET_FIRST_A3_TIME() RESULT( DATE )
```

**USES:**

```
#    include "define.h"
```

**RETURN VALUE:**

```
INTEGER :: DATE(2)    ! YYYYMMDD and HHMMSS values
```

**REVISION HISTORY:**

```
26 Jun 2003 - R. Yantosca - Initial Version
(1 ) Now modified for GCAP and GEOS-5 data (swu, bmy, 5/24/05)
(2 ) Remove support for GEOS-1 and GEOS-STRAT met fields (bmy, 8/4/06)
15 Jan 2010 - R. Yantosca - Added ProTeX headers
27 Sep 2010 - R. Yantosca - Modified for start times other than 0 GMT
```

---

**1.10.68 get\_first\_a6\_time**

Function GET\_FIRST\_A6\_TIME returns the correct YYYYMMDD and HHMMSS values the first time that A-6 fields are read in from disk.

**INTERFACE:**

```
FUNCTION GET_FIRST_A6_TIME() RESULT( DATE )
```

**USES:**

```
#    include "define.h"
```

**RETURN VALUE:**

```
INTEGER :: DATE(2)    ! YYYYMMDD, HHMMSS values
```

**REVISION HISTORY:**

```
26 Jun 2003 - R. Yantosca - Initial Version
(1 ) Now modified for GEOS-4 "a_llk_03" and "a_llk_04" fields (bmy, 3/22/04)
(2 ) Modified for GCAP and GEOS-5 met fields (swu, bmy, 5/24/05)
15 Jan 2010 - R. Yantosca - Added ProTeX headers
27 Sep 2010 - R. Yantosca - Modified for start times other than 0 GMT
```

---

**1.10.69 get\_first\_i6\_time**

Function GET\_FIRST\_I6\_TIME returns the correct YYYYMMDD and HHMMSS values the first time that I-6 fields are read in from disk.

**INTERFACE:**

```
FUNCTION GET_FIRST_I6_TIME() RESULT( DATE )
```

**RETURN VALUE:**

```
INTEGER :: DATE(2)      ! YYYYMMDD, HHMMSS values
```

**REVISION HISTORY:**

27 Sep 2010 - R. Yantosca - Initial version

---

**1.10.70 its\_time\_for\_chem**

Function ITS\_TIME\_FOR\_CHEM returns TRUE if it is time to do chemistry, or FALSE otherwise.

**INTERFACE:**

```
FUNCTION ITS_TIME_FOR_CHEM() RESULT( FLAG )
```

**RETURN VALUE:**

```
LOGICAL :: FLAG
```

**REVISION HISTORY:**

21 Mar 2003 - R. Yantosca - Initial Version

15 Jan 2010 - R. Yantosca - Added ProTeX headers

27 Sep 2011 - M. Payer - Modifications for centralizing the chemistry  
time step (lzh)

---

**1.10.71 its\_time\_for\_conv**

Function ITS\_TIME\_FOR\_CONV returns TRUE if it is time to do convection, or FALSE otherwise.

**INTERFACE:**

```
FUNCTION ITS_TIME_FOR_CONV() RESULT( FLAG )
```

**RETURN VALUE:**

LOGICAL :: FLAG

## REVISION HISTORY:

21 Mar 2003 - R. Yantosca - Initial Version  
15 Jan 2010 - R. Yantosca - Added ProTeX headers

---

### 1.10.72 `its_time_for_dyn`

Function ITS\_TIME\_FOR\_DYN returns TRUE if it is time to do chemistry and false otherwise.

## INTERFACE:

FUNCTION ITS\_TIME\_FOR\_DYN() RESULT( FLAG )

## RETURN VALUE:

LOGICAL :: FLAG

## REVISION HISTORY:

21 Mar 2003 - R. Yantosca - Initial Version  
15 Jan 2010 - R. Yantosca - Added ProTeX headers

---

### 1.10.73 `its_time_for_emis`

Function ITS\_TIME\_FOR\_EMIS returns TRUE if it is time to do emissions, or FALSE otherwise.

## INTERFACE:

FUNCTION ITS\_TIME\_FOR\_EMIS() RESULT( FLAG )

## RETURN VALUE:

LOGICAL :: FLAG

## REVISION HISTORY:

21 Mar 2003 - R. Yantosca - Initial Version  
15 Jan 2010 - R. Yantosca - Added ProTeX headers  
07 Oct 2011 - R. Yantosca - Modifications for centralizing the chemistry  
time step (lzh)

---

**1.10.74    `its_time_for_unit`**

Function `ITS_TIME_FOR_UNIT` returns TRUE if it is time to do unit conversion, or FALSE otherwise.

**INTERFACE:**

```
FUNCTION ITS_TIME_FOR_UNIT() RESULT( FLAG )
```

**RETURN VALUE:**

```
LOGICAL :: FLAG
```

**REVISION HISTORY:**

```
21 Mar 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.75    `its_time_for_diag`**

Function `ITS_TIME_FOR_DIAG` returns TRUE if it is time to archive certain diagnostics, or FALSE otherwise.

**INTERFACE:**

```
FUNCTION ITS_TIME_FOR_DIAG() RESULT( FLAG )
```

**RETURN VALUE:**

```
LOGICAL :: FLAG
```

**REVISION HISTORY:**

```
21 Mar 2003 - R. Yantosca - Initial Version
20 Jul 2009 - C. Carouge   - Use TS_DIAG now and not 60 minutes
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.76    `its_time_for_a1`**

Function `ITS_TIME_FOR_A1` returns TRUE if it is time to read in A1 (average 1-hr fields) and FALSE otherwise.

**INTERFACE:**

```
FUNCTION ITS_TIME_FOR_A1() RESULT( FLAG )
```

**RETURN VALUE:**

```
LOGICAL :: FLAG
```

**REVISION HISTORY:**

```
20 Aug 2010 - R. Yantosca - Initial version
```

---



**1.10.77   its\_time\_for\_a3**

Function ITS\_TIME\_FOR\_A3 returns TRUE if it is time to read in A3 (average 3-hr fields) and FALSE otherwise.

**INTERFACE:**

```
FUNCTION ITS_TIME_FOR_A3() RESULT( FLAG )
```

**RETURN VALUE:**

```
LOGICAL :: FLAG
```

**REVISION HISTORY:**

```
21 Mar 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.78   its\_time\_for\_a6**

Function ITS\_TIME\_FOR\_A6 returns TRUE if it is time to read in A6 (average 6-hr fields) and FALSE otherwise.

**INTERFACE:**

```
FUNCTION ITS_TIME_FOR_A6() RESULT( FLAG )
```

**USES:**

```
#      include "define.h"
```

**RETURN VALUE:**

```
LOGICAL :: FLAG
```

**REVISION HISTORY:**

```
21 Mar 2003 - R. Yantosca - Initial Version
(1 ) Now compute when it's time to read in GEOS-4 A-6 fields. (bmy, 6/26/03)
(2 ) Now modified for GEOS-4 "a_llk_03" and "a_llk_04" fields (bmy, 3/22/04)
(3 ) Now modified for GCAP and GEOS-5 met fields (swu, bmy, 5/24/05)
15 Jan 2010 - R. Yantosca - Added ProTeX headers
17 Feb 2011 - R. Yantosca - Add modifications for APM microphysics (G. Luo)
```

---

**1.10.79    *its\_time\_for\_a6update***

Function `ITS_TIME_FOR_A6UPDATE` returns TRUE if it is time to update the A6 (average 6-hr fields) for APM microphysics, and FALSE otherwise.

**INTERFACE:**

```
FUNCTION ITS_TIME_FOR_A6UPDATE() RESULT( FLAG )
```

**USES:**

```
#      include "define.h"
```

**RETURN VALUE:**

```
LOGICAL :: FLAG
```

**REMARKS:**

This subroutine will only be compiled if you build GEOS-Chem with the APM=yes makefile option.

**REVISION HISTORY:**

```
21 Mar 2003 - R. Yantosca - Initial Version
(1 ) Now compute when it's time to read in GEOS-4 A-6 fields. (bmy, 6/26/03)
(2 ) Now modified for GEOS-4 "a_llk_03" and "a_llk_04" fields (bmy, 3/22/04)
(3 ) Now modified for GCAP and GEOS-5 met fields (swu, bmy, 5/24/05)
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.80    *its\_time\_for\_i6***

Function `ITS_TIME_FOR_I6` returns TRUE if it is time to read in I6 (instantaneous 6-hr fields) and FALSE otherwise.

**INTERFACE:**

```
FUNCTION ITS_TIME_FOR_I6() RESULT( FLAG )
```

**RETURN VALUE:**

```
LOGICAL :: FLAG
```

**REVISION HISTORY:**

```
21 Mar 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.81    `its_time_for_unzip`**

Function `ITS_TIME_FOR_UNZIP` Returns TRUE if it is time to unzip the next day's met field files, or FALSE otherwise.

**INTERFACE:**

```
FUNCTION ITS_TIME_FOR_UNZIP() RESULT( FLAG )
```

**RETURN VALUE:**

```
LOGICAL :: FLAG
```

**REVISION HISTORY:**

```
21 Mar 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.82    `its_time_for_del`**

Function `ITS_TIME_FOR_DEL` returns TRUE if it is time to delete the previous day's met field files in the temporary directory.

**INTERFACE:**

```
FUNCTION ITS_TIME_FOR_DEL() RESULT( FLAG )
```

**RETURN VALUE:**

```
LOGICAL :: FLAG
```

**REVISION HISTORY:**

```
21 Mar 2003 - R. Yantosca - Initial Version
19 Jun 2003 - R. Yantosca - Now delete files at 23 GMT each day, since the
                           last fvDAS A-3 field is 22:30 GMT and the last
                           fvDAS A-6 field is 21 GMT
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.83    `its_time_for_exit`**

Function `ITS_TIME_FOR_EXIT` returns TRUE if it is the end of the GEOS-Chem simulation (i.e.  $\tau_i = \tau_{ue}$ ), or FALSE otherwise.

**INTERFACE:**

```
FUNCTION ITS_TIME_FOR_EXIT() RESULT( FLAG )
```

**RETURN VALUE:**

LOGICAL :: FLAG

**REVISION HISTORY:**

21 Mar 2003 - R. Yantosca - Initial Version  
 15 Jan 2010 - R. Yantosca - Added ProTeX headers

---

**1.10.84 its\_time\_for\_bpch**

Function ITS\_TIME\_FOR\_BPCH returns TRUE if it's time to write output to the bpch file, or FALSE otherwise.

**INTERFACE:**

FUNCTION ITS\_TIME\_FOR\_BPCH() RESULT( DO\_BPCH )

**USES:**

USE CMN\_SIZE\_MOD ! Size parameters  
 USE CMN\_DIAG\_MOD ! NJDAY

**RETURN VALUE:**

LOGICAL :: DO\_BPCH

**REVISION HISTORY:**

02 Feb 2007 - R. Yantosca - Initial Version  
 15 Jan 2010 - R. Yantosca - Added ProTeX headers

---

**1.10.85 its\_a\_leapyear**

Function ITS\_A\_LEAPYEAR tests to see if a year is really a leapyear.

**INTERFACE:**

FUNCTION ITS\_A\_LEAPYEAR( YEAR\_IN, FORCE ) RESULT( IS\_LEAPYEAR )

**INPUT PARAMETERS:**

INTEGER, INTENT(IN), OPTIONAL :: YEAR\_IN ! Year to test if leapyear  
 LOGICAL, INTENT(IN), OPTIONAL :: FORCE ! Do not exit if using GCAP

**RETURN VALUE:**

LOGICAL :: IS\_LEAPYEAR ! =T if it's a leapyear

**REVISION HISTORY:**

17 Mar 1999 - R. Yantosca - Initial Version  
 (1 ) Now remove YEAR from ARG list; use the module variable (bmy, 3/21/03)  
 (2 ) Now add YEAR\_IN as an optional argument. If YEAR\_IN is not passed,  
       then test if the current year is a leapyear (bmy, 9/25/03)  
 (3 ) Now always return FALSE for GCAP (swu, bmy, 8/29/05)  
 (4 ) Now add FORCE argument to force ITS\_A\_LEAPYEAR to return a value  
       instead of just returning with FALSE for the GCAP met fields.  
       (swu, bmy, 4/24/06)  
 15 Jan 2010 - R. Yantosca - Added ProTeX headers

---

**1.10.86 its\_a\_new\_year**

Function ITS\_A\_NEW\_YEAR returns TRUE if it's the first timestep of the year when we have to read in annual data.

**INTERFACE:**

```
FUNCTION ITS_A_NEW_YEAR( NO_CCTS ) RESULT( IS_NEW_YEAR )
```

**INPUT PARAMETERS:**

```
LOGICAL, OPTIONAL :: NO_CCTS      ! =T reverts to previous behavior
                                   ! (i.e. w/o using central chem step)
```

**RETURN VALUE:**

```
LOGICAL           :: IS_NEW_YEAR  ! =T if it's 1st data read of year
```

**REMARKS:**

ITS\_A\_NEW\_YEAR assumes that we are using the central chemistry timestep option (i.e. do chemistry & emissions & related processes at the midpoint of each chemistry timestep). To revert to the prior behavior, set the optional flag NO\_CCTS = .TRUE.

If we are using the central chemistry timestep option (which is now the default behavior), then we must not read data at 00:00 GMT on the first day of the year, but at the center of the first chemistry timestep of the year. This is because emissions and chemistry are done at the same time. The proper time of day for reading emissions is determined by function ITS\_TIME\_FOR\_EMIS, also within time\_mod.f.

Similarly, for simulations that start at an arbitrary midmonth date and time, we must not read data at the starting date and time of the simulation, but at the midpoint of the first chemistry timestep of the simulation.

If we are not using the central chemistry timestep option (specified by

NO\_CCTS=.TRUE.), then the first data read of the month occurs at 00:00 GMT on the Jan 1st. Similarly, for those simulations that start at midmonth, the first data read will occur the starting date and time of the simulation.

## REVISION HISTORY:

01 Apr 2004 - R. Yantosca - Initial Version  
 01 Nov 2005 - R. Yantosca - Bug fix: Need month & day to be 1  
 15 Jan 2010 - R. Yantosca - Added ProTeX headers  
 14 Oct 2011 - R. Yantosca - Modified for central chemistry timestep

### 1.10.87 its\_a\_new\_month

Function ITS\_A\_NEW\_MONTH returns TRUE if it's the first timestep of the month when we have to read in monthly data.

## INTERFACE:

```
FUNCTION ITS_A_NEW_MONTH( NO_CCTS ) RESULT( IS_NEW_MONTH )
!INPUT PARAMETERS
  LOGICAL, OPTIONAL :: NO_CCTS      ! =T reverts to previous behavior
                                      ! (i.e. w/o using central chem step)
```

## RETURN VALUE:

```
  LOGICAL      :: IS_NEW_MONTH  ! =T if it's 1st data read of month
```

## REMARKS:

ITS\_A\_NEW\_MONTH assumes that we are using the central chemistry timestep option (i.e. do chemistry & emissions & related processes at the midpoint of each chemistry timestep). To revert to the prior behavior, set the optional flag NO\_CCTS = .TRUE.

If we are using the central chemistry timestep option (which is now the default behavior), then we must not read data at 00:00 GMT on the first day of the month, but at the center of the first chemistry timestep of the month. This is because emissions and chemistry are done at the same time. The proper time of day for reading emissions is determined by function ITS\_TIME\_FOR\_EMIS, also within time\_mod.f.

Similarly, for simulations that start at an arbitrary midmonth date and time, we must not read data at the starting date and time of the simulation, but at the midpoint of the first chemistry timestep of the simulation.

If we are not using the central chemistry timestep option (specified by NO\_CCTS=.TRUE.), then the first data read of the month occurs at 00:00 GMT on the first day of the month. Similarly, for those simulations that start

at midmonth, the first data read will occur the starting date and time of the simulation.

## REVISION HISTORY:

01 Apr 2004 - R. Yantosca - Initial Version  
 15 Jan 2010 - R. Yantosca - Added ProTeX headers  
 12 Oct 2011 - R. Yantosca - Modified for central chemistry timestep option

---

### 1.10.88 its\_midmonth

Function ITS\_MIDMONTH returns TRUE if it's the middle of a month.

## INTERFACE:

```
FUNCTION ITS_MIDMONTH() RESULT( IS_MIDMONTH )
```

## RETURN VALUE:

```
LOGICAL :: IS_MIDMONTH
```

## REVISION HISTORY:

10 Oct 2005 - S. Strode - Initial Version  
 15 Jan 2010 - R. Yantosca - Added ProTeX headers  
 14 Oct 2011 - R. Yantosca - Modified for central chemistry timestep

---

### 1.10.89 its\_a\_new\_day

Function ITS\_A\_NEW\_DAY returns TRUE if it's the first timestep of the day when we have to read in daily data.

## INTERFACE:

```
FUNCTION ITS_A_NEW_DAY( NO_CCTS ) RESULT( IS_NEW_DAY )
!INPUT PARAMETERS
  LOGICAL, OPTIONAL :: NO_CCTS      ! =T reverts to previous behavior
                                      ! (i.e. w/o using central chem step)
```

## RETURN VALUE:

```
LOGICAL :: IS_NEW_DAY      ! =T if it's 1st data read of day
```

## REMARKS:

ITS\_A\_NEW\_DAY assumes that we are using the central chemistry timestep option (i.e. do chemistry & emissions & related processes at the midpoint of each chemistry timestep). To revert to the prior behavior, set the optional flag NO\_CCTS = .TRUE.

If we are using the central chemistry timestep option (which is now the default behavior), then we must not read data at 00:00 GMT of each day, but at the center of the first chemistry timestep of the day. This is because emissions and chemistry are done at the same time. The proper time of day for reading emissions is determined by function `ITS_TIME_FOR_EMIS`, also within `time_mod.f`.

Similarly, for simulations that start at an arbitrary midmonth date and time, we must not read data at the starting date and time of the simulation, but at the midpoint of the first chemistry timestep of the simulation.

If we are not using the central chemistry timestep option (specified by `NO_CCTS=.TRUE.`), then the first data read of the month occurs at 00:00 GMT each day. Similarly, for those simulations that start at midmonth, the first data read will occur the starting date and time of the simulation.

## REVISION HISTORY:

01 Apr 2004 - R. Yantosca - Initial Version  
 15 Jan 2010 - R. Yantosca - Added ProTeX headers  
 14 Oct 2011 - R. Yantosca - Modified for central chemistry timestep

---

### 1.10.90 `its_a_new_season`

Function `ITS_A_NEW_SEASON` returns TRUE if it's a new season or FALSE if it's not a new season. Seasons are (1=DJF, 2=MAM, 3=JJA, 4=SON).

## INTERFACE:

```
FUNCTION ITS_A_NEW_SEASON( ) RESULT( IS_NEW_SEASON )
```

## RETURN VALUE:

```
LOGICAL :: IS_NEW_SEASON
```

## REVISION HISTORY:

20 Jul 2004 - R. Yantosca - Initial Version  
 15 Jan 2010 - R. Yantosca - Added ProTeX headers

---

### 1.10.91 `print_current_time`

Subroutine `PRINT_CURRENT_TIME` prints the date, GMT time, and elapsed hours of a GEOS-Chem simulation.

## INTERFACE:



SUBROUTINE PRINT\_CURRENT\_TIME

## REVISION HISTORY:

21 Mar 2003 - R. Yantosca - Initial Version  
 15 Jan 2010 - R. Yantosca - Added ProTeX headers

### 1.10.92 timestamp\_string

Function TIMESTAMP\_STRING returns a formatted string "YYYY/MM/DD hh:mm" for the a date and time specified by YYYYMMDD and hhmmss. If YYYYMMDD and hhmmss are omitted, then TIMESTAMP\_STRING will create a formatted string for the current date and time.

## INTERFACE:

FUNCTION TIMESTAMP\_STRING( YYYYMMDD, HHMMSS ) RESULT( TIME\_STR )

## USES:

```
#    include "define.h"
```

## INPUT PARAMETERS:

INTEGER, INTENT(IN), OPTIONAL :: YYYYMMDD    ! YYYY/MM/DD date  
 INTEGER, INTENT(IN), OPTIONAL :: HHMMSS       ! hh:mm:ss time

## RETURN VALUE:

CHARACTER(LEN=16)                                :: TIME\_STR

## REVISION HISTORY:

21 Mar 2003 - R. Yantosca - Initial Version  
 (1 ) Now use ENCODE statement for PGI/F90 on Linux (bmy, 9/29/03)  
 (2 ) Now add optional arguments YYYYMMDD and HHMMSS (bmy, 10/27/03)  
 (3 ) Renamed LINUX to LINUX\_PGI (bmy, 12/2/03)  
 15 Jan 2010 - R. Yantosca - Added ProTeX headers

### 1.10.93 ymd\_extract

Subroutine YMD\_EXTRACT extracts the year, month, and date from an integer variable in YYYYMMDD format. It can also extract the hours, minutes, and seconds from a variable in HHMMSS format.

## INTERFACE:

SUBROUTINE YMD\_EXTRACT( NYMD, Y, M, D )

**INPUT PARAMETERS:**

INTEGER, INTENT(IN) :: NYMD ! YYYY/MM/DD format date

**OUTPUT PARAMETERS:**

INTEGER, INTENT(OUT) :: Y, M, D ! Separated YYYY, MM, DD values

**REVISION HISTORY:**

21 Nov 2001 - R. Yantosca - Initial Version  
 15 Jan 2010 - R. Yantosca - Added ProTeX headers

---

**1.10.94 expand\_date**

Subroutine EXPAND\_DATE replaces "YYYYMMDD" and "hhmmss" tokens within a file-name string with the actual values.

**INTERFACE:**

SUBROUTINE EXPAND\_DATE( FILENAME, YYYYMMDD, HHMMSS )

**USES:**

USE CHARPAK\_MOD, ONLY : STRREPL

# include "define.h"

**INPUT PARAMETERS:**

INTEGER, INTENT(IN) :: YYYYMMDD ! YYYY/MM/DD date  
 INTEGER, INTENT(IN) :: HHMMSS ! hh:mm:ss time

**INPUT/OUTPUT PARAMETERS:**

CHARACTER(LEN=\*), INTENT(INOUT) :: FILENAME ! Filename to modify

**REVISION HISTORY:**

27 Jun 2002 - R. Yantosca - Initial Version  
 (1 ) Bug fix for Linux: use ENCODE statement to convert number to string  
       instead of F90 internal read. (bmy, 9/29/03)  
 (2 ) Now replace 2 and 4 digit year strings for all models (bmy, 10/23/03)  
 (3 ) Renamed LINUX to LINUX\_PGI (bmy, 12/2/03)  
 (4 ) Now do not replace "ss" with seconds, as the smallest GEOS-Chem  
       timestep is in minutes. (bmy, 7/20/04)  
 15 Jan 2010 - R. Yantosca - Added ProTeX headers

---

**1.10.95 system\_date\_time**

Subroutine SYSTEM\_DATE\_TIME returns the actual local date and time (as opposed to the model date and time).

**INTERFACE:**

```
SUBROUTINE SYSTEM_DATE_TIME( SYS_NYMD, SYS_NHMS )
```

**OUTPUT PARAMETERS:**

```
INTEGER, INTENT(OUT) :: SYS_NYMD    ! System date in YYYY/MM/DD format
INTEGER, INTENT(OUT) :: SYS_NHMS    ! System time in YYYY/MM/DD format
```

**REMARKS:**

Uses the F90 intrinsic function DATE\_AND\_TIME.

**REVISION HISTORY:**

```
02 May 2005 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.96 system\_timestamp**

Function SYSTEM\_TIMESTAMP returns a 16 character string with the system date and time in YYYY/MM/DD HH:MM format.

**INTERFACE:**

```
FUNCTION SYSTEM_TIMESTAMP() RESULT( STAMP )
```

**RETURN VALUE:**

```
CHARACTER(LEN=16) :: STAMP
```

**REVISION HISTORY:**

```
03 May 2005 - R. Yantosca - Initial version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.10.97 timestamp\_diag**

Subroutine TIMESTAMP\_DIAG save timestamps to be used in filenames for diagnostics. We do not want the time when the diagnostic is saved but the time for previous dynamic time step because midnight is considered as the beginning of next day (and not ending of previous day).

**INTERFACE:**

SUBROUTINE TIMESTAMP\_DIAG

## REVISION HISTORY:

12 Aug 2009 - C. Carouge - Initial version  
 15 Jan 2010 - R. Yantosca - Added ProTeX headers

---

### 1.10.98 get\_nymd\_diag

Function GET\_NYMD\_DIAG returns the previous NYMD value (YYYYMMDD) to the calling program. Used for diagnostic filenames.

## INTERFACE:

FUNCTION GET\_NYMD\_DIAG() RESULT( THISNYMD )

## RETURN VALUE:

INTEGER :: THISNYMD

## REVISION HISTORY:

12 Aug 2009 - C. Carouge - Initial version  
 15 Jan 2010 - R. Yantosca - Added ProTeX headers

---

### 1.11 Fortran: Module Interface transfer\_mod.f

Module TRANSFER\_MOD contains routines used to copy data from REAL\*4 to REAL\*8 arrays after being read from disk. Also, vertical levels will be collapsed in the stratosphere if necessary. This will help us to gain computational advantage.

## INTERFACE:

MODULE TRANSFER\_MOD

## USES:

USE ERROR\_MOD, ONLY : ALLOC\_ERR  
 USE ERROR\_MOD, ONLY : GEOS\_CHEM\_STOP

# include "define.h"  
 USE CMN\_SIZE\_MOD

IMPLICIT NONE

PRIVATE

**PUBLIC MEMBER FUNCTIONS:**

```

PUBLIC  :: TRANSFER_A6
PUBLIC  :: TRANSFER_2D
PUBLIC  :: TRANSFER_3D
PUBLIC  :: TRANSFER_3D_Lp1
PUBLIC  :: TRANSFER_3D_TROP
PUBLIC  :: TRANSFER_3D_NOLUMP
PUBLIC  :: TRANSFER_G5_PLE
PUBLIC  :: TRANSFER_ZONAL
PUBLIC  :: TRANSFER_TO_1D
PUBLIC  :: INIT_TRANSFER
PUBLIC  :: CLEANUP_TRANSFER

```

```

INTERFACE TRANSFER_2D
  MODULE PROCEDURE TRANSFER_2D_INT
  MODULE PROCEDURE TRANSFER_2D_R4
  MODULE PROCEDURE TRANSFER_2D_R8
END INTERFACE

```

```

INTERFACE TRANSFER_ZONAL
  MODULE PROCEDURE TRANSFER_ZONAL_R4
  MODULE PROCEDURE TRANSFER_ZONAL_R8
END INTERFACE

```

**PRIVATE MEMBER FUNCTIONS:**

```

PRIVATE :: LUMP_2
PRIVATE :: LUMP_2_R4
PRIVATE :: LUMP_2_R8
PRIVATE :: LUMP_4
PRIVATE :: LUMP_4_R4
PRIVATE :: LUMP_4_R8
PRIVATE :: TRANSFER_2D_INT
PRIVATE :: TRANSFER_2D_R4
PRIVATE :: TRANSFER_2D_R8
PRIVATE :: TRANSFER_ZONAL_R4
PRIVATE :: TRANSFER_ZONAL_R8

```

```

INTERFACE LUMP_2
  MODULE PROCEDURE LUMP_2_R4
  MODULE PROCEDURE LUMP_2_R8
END INTERFACE

```

```

INTERFACE LUMP_4
  MODULE PROCEDURE LUMP_4_R4
  MODULE PROCEDURE LUMP_4_R8
END INTERFACE

```

**REMARKS:**

Hybrid Grid Coordinate Definition: (dsa, bmy, 8/27/02, 8/13/10)

=====

GEOS-4, GEOS-5, and MERRA (hybrid grids):

-----

For GEOS-4 and GEOS-5, the pressure at the bottom edge of grid box (I,J,L) is defined as follows:

$$P_{edge}(I,J,L) = A_p(L) + [ B_p(L) * P_{surface}(I,J) ]$$

where

$P_{surface}(I,J)$  is the "true" surface pressure at lon,lat (I,J)  
 $A_p(L)$  has the same units as surface pressure [hPa]  
 $B_p(L)$  is a unitless constant given at level edges

$A_p(L)$  and  $B_p(L)$  are given to us by GMAO.

-----

GEOS-3 (pure-sigma) and GCAP (hybrid grid):

-----

GEOS-3 is a pure-sigma grid. GCAP is a hybrid grid, but its grid is defined as if it were a pure sigma grid (i.e.  $P_{TOP}=150$  hPa, and negative sigma edges at higher levels). For these grids, can still use the same formula as for GEOS-4, with one modification:

$$P_{edge}(I,J,L) = A_p(L) + [ B_p(L) * ( P_{surface}(I,J) - P_{TOP} ) ]$$

where

$P_{surface}(I,J)$  = the "true" surface pressure at lon,lat (I,J)  
 $A_p(L)$  =  $P_{TOP}$  = model top pressure  
 $B_p(L)$  =  $SIG_E(L)$  = bottom sigma edge of level L

-----

The following are true for GCAP, GEOS-3, GEOS-4:

- 
- (1)  $B_p(LLPAR+1) = 0.0$  (L=LLPAR+1 is the atmosphere top)
  - (2)  $B_p(1) = 1.0$  (L=1 is the surface)
  - (3)  $P_{TOP} = A_p(LLPAR+1)$  (L=LLPAR+1 is the atmosphere top)

## REVISION HISTORY:

21 Sep 2010 - M. Evans - Initial version

- (1 ) GEOS-3 Output levels were determined by Mat Evans. Groups of 2 levels and groups of 4 levels on the original grid are merged together into thick levels for the output grid. (mje, bmy, 9/26/01)
- (2 ) Assumes that  $LLPAR == LGLOB$  for GEOS-1, GEOS-STRAT (bmy, 9/26/01)
- (3 ) `EDGE_IN` needs to be provided for each model type, within an `#ifdef`

- block, in order to ensure compilation. However, EDGE\_IN is currently only used for regridding GEOS-3 data (and probably also GEOS-4 when that becomes available). (bmy, 9/26/01)
- (4 ) Add interfaces TRANSFER\_2D and TRANSFER\_ZONAL (bmy, 9/27/01)
  - (5 ) Added routine TRANSFER\_2D\_R4. Added TRANSFER\_2D\_R4 to the generic TRANSFER\_2D interface. (bmy, 1/25/02)
  - (6 ) Updated comments, cosmetic changes (bmy, 2/28/02)
  - (7 ) Bug fix: remove extraneous "," in GEOS-1 definition of EDGE\_IN array. (bmy, 3/25/02)
  - (8 ) Now divide module header into MODULE PRIVATE, MODULE VARIABLES, and MODULE ROUTINES sections. Also add MODULE INTERFACES section, since we have an interface here. (bmy, 5/28/02)
  - (9 ) Now references "pressure\_mod.f" (dsa, bdf, bmy, 8/22/02)
  - (10) Bug fix in "init\_transfer", declare variable L. Also reference GEOS\_CHEM\_STOP from "error\_mod.f" for safe stop (bmy, 10/15/02)
  - (11) Added routine TRANSFER\_3D\_TROP. Also updated comments. (bmy, 10/31/02)
  - (12) Now uses functions GET\_XOFFSET and GET\_YOFFSET from "grid\_mod.f". (bmy, 3/11/03)
  - (13) Added code to regrid GEOS-4 from 55 --> 30 levels. Renamed module variable SIGE\_IN to EDGE\_IN. (mje, bmy, 10/31/03)
  - (14) Now modified for GEOS-5 and GCAP met fields (swu, bmy, 5/24/05)
  - (15) Remove support for GEOS-1 and GEOS-STRAT met fields (bmy, 8/4/06)
  - (16) Modified for GEOS-5. Rewritten for clarity. (bmy, 10/30/07)
- 13 Aug 2010 - R. Yantosca - Added modifications for MERRA met fields  
 13 Aug 2010 - R. Yantosca - Added ProTeX headers

### 1.11.1 transfer\_a6

Subroutine TRANSFER\_A6 transfers A-6 data from a REAL\*4 array to a REAL\*8 array. Vertical layers are collapsed (from LGLOB to LLPAR) if necessary.

#### INTERFACE:

```
SUBROUTINE TRANSFER_A6( IN, OUT )
```

#### INPUT PARAMETERS:

```
REAL*4, INTENT(IN) :: IN(IIPAR,JJP,LLPAR) ! Input data
```

#### OUTPUT PARAMETERS:

```
REAL*8, INTENT(OUT) :: OUT(LLPAR,IIPAR,JJP) ! Output data
```

#### REMARKS:

#### REVISION HISTORY:

19 Sep 2001 - R. Yantosca - Initial version  
 (1 ) A-6 fields are dimensioned (LLPAR,IIPAR,JJPARG) since for Fortran efficiency, since the code loops over vertical layers L in a column located above a certain surface location (I,J). (bmy, 9/21/01)  
 (2 ) Assumes that LLPARG == LGLOB for GEOS-1, GEOS-STRAT (bmy, 9/21/01)  
 (3 ) Now use functions GET\_XOFFSET and GET\_YOFFSET from "grid\_mod.f".  
       Now IO, JO are local variables. (bmy, 3/11/03)  
 (4 ) Added code to regrid GEOS-4 from 55 --> 30 levels (mje, bmy, 10/31/03)  
 (5 ) Now modified for GEOS-5 met fields (bmy, 5/24/05)  
 (6 ) Rewritten for clarity (bmy, 2/8/07)  
 (7 ) Now get nested-grid offsets (dan, bmy, 11/6/08)  
 13 Aug 2010 - R. Yantosca - Added ProTeX headers  
 13 Aug 2010 - R. Yantosca - Treat MERRA the same way as GEOS-5, because the vertical grids are identical

---

### 1.11.2

Subroutine TRANSFER\_3D transfers 3-dimensional data from a REAL\*4 array to a REAL\*8 array. Vertical layers are collapsed (from LGLOB to LLPARG) if necessary.

#### INTERFACE:

```
SUBROUTINE TRANSFER_3D( IN, OUT )
```

#### INPUT PARAMETERS:

```
REAL*4, INTENT(IN)  :: IN(IIPAR,JJPARG,LGLOB)    ! Input data
```

#### OUTPUT PARAMETERS:

```
REAL*8, INTENT(OUT) :: OUT(IIPAR,JJPARG,LLPAR)    ! Output data
```

#### REVISION HISTORY:

19 Sep 2001 - R. Yantosca - Initial version  
 (1 ) Lump levels together in groups of 2 or 4, as dictated by Mat Evans.  
       (bmy, 9/21/01)  
 (2 ) Assumes that LLPARG == LGLOB for GEOS-1, GEOS-STRAT (bmy, 9/21/01)  
 (3 ) Now use functions GET\_XOFFSET and GET\_YOFFSET from "grid\_mod.f".  
       Now IO, JO are local variables. (bmy, 3/11/03)  
 (4 ) Added code to regrid GEOS-4 from 55 --> 30 levels (mje, bmy, 10/31/03)  
 (5 ) Now modified for GEOS-5 met fields (bmy, 5/24/05)  
 (6 ) Rewritten for clarity (bmy, 2/8/07)  
 13 Aug 2010 - R. Yantosca - Added ProTeX headers  
 13 Aug 2010 - R. Yantosca - Treat MERRA the same way as GEOS-5, because the vertical grids are identical

---



**1.11.3**

Subroutine TRANSFER\_3D transfers 3-dimensional data from a REAL\*4 array to a REAL\*8 array. Vertical layers are collapsed (from LGLOB to LLPAR) if necessary.

**INTERFACE:**

```
SUBROUTINE TRANSFER_3D_NOLUMP( IN, OUT )
```

**INPUT PARAMETERS:**

```
REAL*4,  INTENT(IN)  :: IN(IIPAR,JJPARGLOB)    ! Input data
```

**OUTPUT PARAMETERS:**

```
REAL*8,  INTENT(OUT) :: OUT(IIPAR,JJPARGLOB)    ! Output data
```

**REVISION HISTORY:**

```
04 Aug 2011 - C. Holmes - Initial version, copied from TRANSFER_3D
```

---

**1.11.4 transfer\_g5\_ple**

Subroutine TRANSFER\_G5\_PLE transfers GEOS-5/MERRA pressure edge data from the native 72-level grid to the reduced 47-level grid.

**INTERFACE:**

```
SUBROUTINE TRANSFER_G5_PLE( IN, OUT )
```

**INPUT PARAMETERS:**

```
REAL*4,  INTENT(IN)  :: IN(IIPAR,JJPARGLOB+1)    ! Input data
```

**OUTPUT PARAMETERS:**

```
REAL*8,  INTENT(OUT) :: OUT(IIPAR,JJPARGLOB+1)    ! Output data
```

**REVISION HISTORY:**

```
08 Feb 2007 - R. Yantosca - Initial version
13 Aug 2010 - R. Yantosca - Added ProTeX headers
13 Aug 2010 - R. Yantosca - Treat MERRA the same way as GEOS-5, because
                           the vertical grids are identical
```

---

### 1.11.5 transfer\_3d\_lp1

Subroutine TRANSFER\_3D\_Lp1 transfers 3-D data from a REAL\*4 array of dimension (IIPAR,JJPARGLOB+1) to a REAL\*8 array of dimension (IIPAR,JJPARGLOB+1). Regrid in the vertical if needed.

#### INTERFACE:

```
SUBROUTINE TRANSFER_3D_Lp1( IN, OUT )
```

#### INPUT PARAMETERS:

```
REAL*4, INTENT(IN) :: IN(IIPAR,JJPARGLOB+1)    ! Input data
```

#### OUTPUT PARAMETERS:

```
REAL*8, INTENT(OUT) :: OUT(IIPAR,JJPARGLOB+1)    ! Output data
```

#### REVISION HISTORY:

```
08 Feb 2007 - R. Yantosca - Initial version
13 Aug 2010 - R. Yantosca - Added ProTeX headers
13 Aug 2010 - R. Yantosca - Treat MERRA the same way as GEOS-5, because
                           the vertical grids are identical
```

---

### 1.11.6 transfer\_3d\_trop

Subroutine TRANSFER\_3D\_TROP transfers tropospheric 3-D data from a REAL\*4 array to a REAL\*8 array.

#### INTERFACE:

```
SUBROUTINE TRANSFER_3D_TROP( IN, OUT )
```

#### INPUT PARAMETERS:

```
REAL*4, INTENT(IN) :: IN(IIPAR,JJPARGLOB,LLTROP_FIX)    ! Input data
```

#### OUTPUT PARAMETERS:

```
REAL*8, INTENT(OUT) :: OUT(IIPAR,JJPARGLOB,LLTROP_FIX)    ! Output data
```

#### REVISION HISTORY:

```
19 Sep 2001 - M. Evans    - Initial version
08 Feb 2007 - R. Yantosca - Now use LLTROP_FIX instead of LLTROP, since
                           most of the offline simulations use the annual
                           mean tropopause
13 Aug 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.11.7 transfer\_zonal\_r4**

Subroutine TRANSFER\_ZONAL\_R4 transfers zonal-mean data from a REAL\*4 array to a REAL\*8 array. Vertical levels are collapsed (from LGLOB to LLPAR) if necessary. (mje, bmy, 9/21/01, 2/8/07)

**INTERFACE:**

```
SUBROUTINE TRANSFER_ZONAL_R4( IN, OUT )
```

**INPUT PARAMETERS:**

```
REAL*4, INTENT(IN) :: IN(JJPAR,LGLOB)      ! Input data
```

**OUTPUT PARAMETERS:**

```
REAL*4, INTENT(OUT) :: OUT(JJPAR,LLPAR)     ! Output data
```

**REVISION HISTORY:**

```
19 Sep 2001 - M. Evans      - Initial version
(1 ) Lump levels together in groups of 2 or 4, as dictated by Mat Evans.
      (bmy, 9/21/01)
(2 ) Assumes that LLPAR == LGLOB for GEOS-1, GEOS-STRAT (bmy, 9/21/01)
(3 ) Now use function GET_YOFFSET from "grid_mod.f". Now IO and JO are
      local variables (bmy, 3/11/03)
(4 ) Added code to regrid GEOS-4 from 55 --> 30 levels (mje, bmy, 10/31/03)
(5 ) Rewritten for clarity (bmy, 2/8/07)
13 Aug 2010 - R. Yantosca - Added ProTeX headers
13 Aug 2010 - R. Yantosca - Treat MERRA the same way as GEOS-5, because
                           the vertical grids are identical
```

---

**1.11.8 transfer\_zonal\_r8**

Subroutine TRANSFER\_ZONAL\_R8 transfers zonal mean or lat-alt data from a REAL\*4 array of dimension (JJPAR,LGLOB) to a REAL\*8 array of dimension (JJPAR,LLPAR). Regrid data in the vertical if necessary by lumping levels.

**INTERFACE:**

```
SUBROUTINE TRANSFER_ZONAL_R8( IN, OUT )
```

**INPUT PARAMETERS:**

```
REAL*4, INTENT(IN) :: IN(JJPAR,LGLOB)      ! Input data
```

**OUTPUT PARAMETERS:**

```
REAL*8, INTENT(OUT) :: OUT(JJPAR,LLPAR)     ! Output data
```

**REVISION HISTORY:**

19 Sep 2001 - R. Yantosca - Initial version  
 (1 ) Lump levels together in groups of 2 or 4, as dictated by Mat Evans.  
      (bmy, 9/21/01)  
 (2 ) Assumes that LLPAR == LGLOB for GEOS-1, GEOS-STRAT (bmy, 9/21/01)  
 (3 ) Now use function GET\_YOFFSET from "grid\_mod.f". Now I0 and J0 are  
      local variables (bmy, 3/11/03)  
 (4 ) Added code to regrid GEOS-4 from 55 --> 30 levels (mje, bmy, 10/31/03)  
 (5 ) Now modified for GEOS-5 met fields (bmy, 5/24/05)  
 13 Aug 2010 - R. Yantosca - Added ProTeX headers  
 13 Aug 2010 - R. Yantosca - Treat MERRA the same way as GEOS-5, because  
      the vertical grids are identical

---

**1.11.9 transfer\_2d\_int**

Subroutine TRANSFER\_2D\_INT transfers 2-D data from a REAL\*4 array of dimension (IIPAR,JJPARG) to an INTEGER array of dimension (IIPAR,JJPARG).

**INTERFACE:**

```
SUBROUTINE TRANSFER_2D_INT( IN, OUT )
```

**INPUT PARAMETERS:**

```
REAL*4, INTENT(IN) :: IN(IIPAR,JJPARG)    ! Input data
```

**OUTPUT PARAMETERS:**

```
INTEGER, INTENT(OUT) :: OUT(IIPAR,JJPARG)  ! Output data
```

**REVISION HISTORY:**

19 Sep 2001 - R. Yantosca - Initial version  
 (1 ) Use parallel DO loops to speed things up (bmy, 9/21/01)!  
 (2 ) Now use functions GET\_XOFFSET and GET\_YOFFSET from "grid\_mod.f".  
      Now I0 and J0 are local variables. (bmy, 3/11/03)  
 13 Aug 2010 - R. Yantosca - Added ProTeX headers

---

**1.11.10 transfer\_2d\_r4**

Subroutine TRANSFER\_2D\_R4 transfers 2-D data from a REAL\*4 array of dimension (IIPAR,JJPARG) to a REAL\*4 array of dimension (IIPAR,JJPARG).

**INTERFACE:**

```
SUBROUTINE TRANSFER_2D_R4( IN, OUT )
```

**INPUT PARAMETERS:**

```
REAL*4,  INTENT(IN)  :: IN(IIPAR,JJPARG)
```

**OUTPUT PARAMETERS:**

```
REAL*4,  INTENT(OUT) :: OUT(IIPAR,JJPARG)
```

**REVISION HISTORY:**

```
19 Sep 2001 - R. Yantosca - Initial version
(1 ) Use parallel DO loops to speed things up (bmy, 9/21/01)
(2 ) Now use functions GET_XOFFSET and GET_YOFFSET from "grid_mod.f"
      Now I0 and J0 are local variables (bmy, 3/11/03)
13 Aug 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.11.11 transfer\_2d\_r8**

Subroutine TRANSFER\_2D\_R8 transfers 2-D data from a REAL\*4 array of dimension (IIPAR,JJPARG) to a REAL\*8 array of dimension (IIPAR,JJPARG).

**INTERFACE:**

```
SUBROUTINE TRANSFER_2D_R8( IN, OUT )
```

**INPUT PARAMETERS:**

```
REAL*4,  INTENT(IN)  :: IN(IIPAR,JJPARG)    ! Input data
```

**OUTPUT PARAMETERS:**

```
REAL*8,  INTENT(OUT) :: OUT(IIPAR,JJPARG)    ! Output data
```

**REVISION HISTORY:**

```
19 Sep 2001 - R. Yantosca - Initial version
(1 ) Use parallel DO loops to speed things up (bmy, 9/21/01)
(2 ) Now use functions GET_XOFFSET and GET_YOFFSET from "grid_mod.f"
      Now I0 and J0 are local variables. (bmy, 3/11/03)
13 Aug 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.11.12 transfer\_to\_1d**

Subroutine TRANSFER\_TO\_1D transfers 2-D data from a REAL\*4 array of dimension (IIPAR,JJPARG) to 1-D a REAL\*8 array of dimension (MAXIJ), where MAXIJ = IIPAR \* JJPARG.

**INTERFACE:**

```
SUBROUTINE TRANSFER_TO_1D( IN, OUT )
```

#### INPUT PARAMETERS:

```
REAL*4,  INTENT(IN)  :: IN(IIPAR,JJPARG)  ! Input data
```

#### OUTPUT PARAMETERS:

```
REAL*8,  INTENT(OUT) :: OUT(MAXIJ)        ! Output data
```

#### REVISION HISTORY:

```
19 Sep 2001 - R. Yantosca - Initial version
(1 ) Use single-processor DO-loops for now (bmy, 9/21/01)
(2 ) Now use functions GET_XOFFSET and GET_YOFFSET from "grid_mod.f".
      Now IO and JO are local variables. (bmy, 3/11/03)
13 Aug 2010 - R. Yantosca - Added ProTeX headers
```

---

#### 1.11.13 lump\_2\_r4

Function LUMP\_2\_R4 lumps 2 sigma levels into one thick level. Input arguments must be REAL\*4.

#### INTERFACE:

```
FUNCTION LUMP_2_R4( IN, L_IN, L ) RESULT( OUT )
```

#### USES:

#### INPUT PARAMETERS:

```
REAL*4,  INTENT(IN) :: IN(L_IN)  ! Column of data on input grid
INTEGER, INTENT(IN) :: L_IN      ! Vertical dimension of the IN array
INTEGER, INTENT(IN) :: L        ! Level on input grid from which
                                ! to start regridding
```

#### RETURN VALUE:

```
REAL*4          :: OUT          ! Data on output grid: 4 lumped levels
```

#### REVISION HISTORY:

```
19 Sep 2001 - R. Yantosca - Initial version
(1 ) Now references GEOS_CHEM_STOP from "error_mod.f" (bmy, 10/15/02)
(2 ) Renamed SIGE_IN to EDGE_IN to denote that it is not always a sigma
      coordinate (as for GEOS-4). Also updated comments (bmy, 10/31/03)
13 Aug 2010 - R. Yantosca - Added ProTeX headers
```

---



**RETURN VALUE:**

```
REAL*4          :: OUT          ! Data on output grid: 4 lumped levels
```

**REVISION HISTORY:**

```
19 Sep 2001 - R. Yantosca - Initial version
(1 ) Now references GEOS_CHEM_STOP from "error_mod.f" (bmy, 10/15/02)
(2 ) Renamed SIGE_IN to EDGE_IN to denote that it is not always a sigma
      coordinate (as for GEOS-4). Also updated comments (bmy, 10/31/03)
13 Aug 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.11.16 lump\_4\_r8**

Function LUMP\_4\_R8 lumps 4 sigma levels into one thick level. Input arguments must be REAL\*8.

**INTERFACE:**

```
FUNCTION LUMP_4_R8( IN, L_IN, L ) RESULT( OUT )
```

**USES:**

```
USE ERROR_MOD, ONLY : GEOS_CHEM_STOP
```

**INPUT PARAMETERS:**

```
REAL*8,  INTENT(IN) :: IN(L_IN)  ! Column of data on input grid
INTEGER, INTENT(IN) :: L_IN      ! Vertical dimension of the IN array
INTEGER, INTENT(IN) :: L         ! Level on input grid from which
                                ! to start regridding
```

**RETURN VALUE:**

```
REAL*8          :: OUT          ! Data on output grid: 4 lumped levels
```

**REVISION HISTORY:**

```
19 Sep 2001 - R. Yantosca - Initial version
(1 ) Now references GEOS_CHEM_STOP from "error_mod.f" (bmy, 10/15/02)
(2 ) Renamed SIGE_IN to EDGE_IN to denote that it is not always a sigma
      coordinate (as for GEOS-4). Also updated comments (bmy, 10/31/03)
13 Aug 2010 - R. Yantosca - Added ProTeX headers
```

---

**1.11.17 init\_transfer**

Subroutine INIT\_TRANSFER initializes and zeroes all module variables.

**INTERFACE:**



```
SUBROUTINE INIT_TRANSFER( THIS_IO, THIS_JO )
```

## USES:

## INPUT PARAMETERS:

```
INTEGER, INTENT(IN) :: THIS_IO    ! Global X (longitude) offset
INTEGER, INTENT(IN) :: THIS_JO    ! Global Y (latitude)  offset
```

## REVISION HISTORY:

```
19 Sep 2001 - R. Yantosca - Initial version
(1 ) Removed additional "," for GEOS-1 definition of EDGE_IN (bmy, 3/25/02)
(2 ) Now use GET_BP from "pressure_mod.f" to get sigma edges for all
      grids except GEOS-3 (dsa, bdf, bmy, 8/22/02)
(3 ) Declare L as a local variable.  Also reference ALLOC_ERR from module
      "error_mod.f" (bmy, 10/15/02)
(4 ) Renamed SIGE_IN to EDGE_IN to denote that it is not always a sigma
      coordinate (as for GEOS-4).  Now assign original Ap coordinates from
      the GEOS-4 grid to the EDGE_IN array (bmy, 10/31/03)
(5 ) Now modified for GEOS-5 met fields (bmy, 5/24/05)
(6 ) Rewritten for clarity.  Remove references to "grid_mod.f" and
      "pressure_mod.f".  Now pass IO, JO from "grid_mod.f" via the arg list.
      (bmy, 2/8/07)
13 Aug 2010 - R. Yantosca - Added ProTeX headers
13 Aug 2010 - R. Yantosca - Treat MERRA the same way as GEOS-5, because
                          the vertical grids are identical
```

---

### 1.11.18 cleanup\_transfer

Subroutine CLEANUP\_TRANSFER deallocates all module variables.

## INTERFACE:

```
SUBROUTINE CLEANUP_TRANSFER
```

## REVISION HISTORY:

```
19 Sep 2001 - R. Yantosca - Initial version
31 Oct 2003 - R. Yantosca - Renamed SIGE_IN to EDGE_IN to denote that it
                          is not always a sigma coordinate (as for GEOS-4)
13 Aug 2010 - R. Yantosca - Added ProTeX headers
```

---

## 1.12 Fortran: Module Interface unix\_cmds\_mod.f

Module UNIX\_CMDS\_MOD contains variables which contain file suffixes and various Unix command strings.

### INTERFACE:

```
MODULE UNIX_CMDS_MOD
```

### USES:

```
IMPLICIT NONE
#   include "define.h"
PUBLIC
```

### PUBLIC DATA MEMBERS:

```
! Unix cmd and file suffix strings for ...
CHARACTER(LEN=255) :: BACKGROUND ! Background operator ( ' &' )
CHARACTER(LEN=255) :: REDIRECT ! Redirection operator ( ' >' )
CHARACTER(LEN=255) :: REMOVE_CMD ! File/dir remove cmd ( 'rm' )
CHARACTER(LEN=255) :: SEPARATOR ! Dir path separator ( '/' )
CHARACTER(LEN=255) :: SPACE ! Blank space ( ' ' )
CHARACTER(LEN=255) :: UNZIP_CMD ! Unzip command ( 'gzcat' )
CHARACTER(LEN=255) :: WILD_CARD ! Wild card operator ( '*' )
CHARACTER(LEN=255) :: A3_SUFFIX ! !%% OBSOLETE %%
CHARACTER(LEN=255) :: A6_SUFFIX ! !%% OBSOLETE %%
CHARACTER(LEN=255) :: I6_SUFFIX ! !%% OBSOLETE %%
CHARACTER(LEN=255) :: PH_SUFFIX ! !%% OBSOLETE %%
CHARACTER(LEN=255) :: KZZ_SUFFIX ! !%% OBSOLETE %%
CHARACTER(LEN=255) :: GRID_SUFFIX ! !%% OBSOLETE %%
CHARACTER(LEN=255) :: ZIP_SUFFIX ! Zipped file suffix ( '.gz' )
```

### REVISION HISTORY:

```
09 Jul 2004 - R. Yantosca - Initial version
20 Nov 2009 - R. Yantosca - Added ProTeX header
```