

GEOS-Chem Reference, Volume 2: Utility Modules

GEOS-CHEM SUPPORT TEAM

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Contents

1	Routine/Function Prologues	7
1.1	Fortran: Module Interface julday_mod	7
1.1.1	julday	7
1.1.2	mint	8
1.1.3	caldate	8
1.2	Fortran: Module Interface transfer_mod	9
1.2.1	transfer_a6	12
1.2.2	transfer_3d	13
1.2.3	transfer_3d_bry	14
1.2.4	transfer_g5_ple	14
1.2.5	transfer_3d_lp1	15
1.2.6	transfer_3d_trop	15
1.2.7	transfer_zonal_r4	16
1.2.8	transfer_zonal_r8	17
1.2.9	transfer_2d_int	17
1.2.10	transfer_2d_r4	18
1.2.11	transfer_2d_r8	18
1.2.12	transfer_to_1d	19
1.2.13	lump_2_r4	19
1.2.14	lump_2_r8	20
1.2.15	lump_4_r4	21
1.2.16	lump_4_r8	21
1.2.17	init_transfer	22
1.2.18	cleanup_transfer	23
1.3	Fortran: Module Interface pressure_mod	23
1.3.1	get_ap	25
1.3.2	get_bp	25
1.3.3	set_floating_pressure	26
1.3.4	get_pedge	27
1.3.5	get_pedge_fullgrid	27
1.3.6	get_pcenter	28
1.3.7	init_pressure	28
1.3.8	cleanup_pressure	29

1.3.9	accept_external_pedge	30
1.4	Fortran: Module Interface directory_mod	30
1.5	Fortran: Module Interface charpak_mod	31
1.6	Fortran: Module Interface time_mod	32
1.6.1	set_current_time	37
1.6.2	set_begin_time	38
1.6.3	set_end_time	38
1.6.4	set_ndiagtime	39
1.6.5	set_diagb	39
1.6.6	set_diage	40
1.6.7	set_timesteps	40
1.6.8	set_ct_chem	41
1.6.9	set_hg2_diag	41
1.6.10	set_ct_conv	42
1.6.11	set_ct_dyn	42
1.6.12	set_ct_emis	42
1.6.13	set_ct_diag	43
1.6.14	set_ct_a1	43
1.6.15	set_ct_a3	44
1.6.16	set_ct_a6	44
1.6.17	set_ct_i3	44
1.6.18	set_ct_i6	45
1.6.19	set_ct_extra	45
1.6.20	set_elapsed_min	46
1.6.21	get_jd	46
1.6.22	get_elapsed_min	46
1.6.23	get_elapsed_sec	47
1.6.24	get_nymdb	47
1.6.25	get_nhmsb	48
1.6.26	get_nymde	48
1.6.27	get_nhmse	48
1.6.28	get_nymd	49
1.6.29	get_nhms	49
1.6.30	get_ndiagtime	49
1.6.31	get_time_ahead	50
1.6.32	get_month	50
1.6.33	get_day	51
1.6.34	get_year	51
1.6.35	set_histyr	51
1.6.36	get_histyr	52
1.6.37	get_hour	52
1.6.38	get_minute	52
1.6.39	get_second	53
1.6.40	get_day_of_year	53
1.6.41	get_day_of_week	53
1.6.42	get_day_of_week_lt	54
1.6.43	get_gmt	54
1.6.44	get_tau	55

1.6.45	get_taub	55
1.6.46	get_tauc	56
1.6.47	get_diagb	56
1.6.48	get_diage	56
1.6.49	get_localtime	57
1.6.50	get_season	57
1.6.51	get_ts_chem	58
1.6.52	get_ts_conv	58
1.6.53	get_ts_diag	59
1.6.54	get_ts_dyn	59
1.6.55	get_ts_emis	59
1.6.56	get_ts_unit	60
1.6.57	get_ct_chem	60
1.6.58	get_ct_conv	60
1.6.59	get_ct_dyn	61
1.6.60	get_ct_emis	61
1.6.61	get_ct_a1	61
1.6.62	get_ct_a3	62
1.6.63	get_ct_a6	62
1.6.64	get_ct_i3	62
1.6.65	get_ct_i6	63
1.6.66	get_ct_extra	63
1.6.67	get_ct_diag	63
1.6.68	get_hg2_diag	64
1.6.69	get_a1_time	64
1.6.70	get_a3_time	64
1.6.71	get_a6_time	65
1.6.72	get_i3_time	65
1.6.73	get_i6_time	66
1.6.74	get_first_a1_time	66
1.6.75	get_first_a3_time	67
1.6.76	get_first_a6_time	67
1.6.77	get_first_i3_time	68
1.6.78	get_first_i6_time	68
1.6.79	its_time_for_chem	68
1.6.80	its_time_for_conv	69
1.6.81	its_time_for_dyn	69
1.6.82	its_time_for_emis	69
1.6.83	its_time_for_unit	70
1.6.84	its_time_for_diag	70
1.6.85	its_time_for_a1	71
1.6.86	its_time_for_a3	71
1.6.87	its_time_for_a6	71
1.6.88	its_time_for_a6update	72
1.6.89	its_time_for_i3	72
1.6.90	its_time_for_i6	73
1.6.91	its_time_for_unzip	73
1.6.92	its_time_for_del	74

1.6.93	its_time_for_exit	74
1.6.94	its_time_for_bpch	74
1.6.95	its_a_leapyear	75
1.6.96	its_a_new_year	75
1.6.97	its_a_new_month	76
1.6.98	its_midmonth	77
1.6.99	its_a_new_day	78
1.6.100	its_a_new_hour	79
1.6.101	its_a_new_season	79
1.6.102	print_current_time	79
1.6.103	timestamp_string	80
1.6.104	ymd_extract	80
1.6.105	expand_date	81
1.6.106	system_date_time	81
1.6.107	system_timestamp	82
1.6.108	timestamp_diag	82
1.6.109	get_nymd_diag	82
1.6.110	accept_external_date_time	83
1.7	Fortran: Module Interface error_mod	84
1.7.1	nan_float	86
1.7.2	nan_dble	87
1.7.3	finite_float	88
1.7.4	finite_dble	88
1.7.5	check_real_value	89
1.7.6	check_dble_value	90
1.7.7	error_stop	90
1.7.8	geos_chem_stop	91
1.7.9	alloc_err	91
1.7.10	debug_msg	92
1.7.11	safe_div	92
1.7.12	is_safe_div	93
1.7.13	safe_exp	94
1.7.14	is_safe_exp	94
1.7.15	safe_log	95
1.7.16	safe_log10	95
1.8	Fortran: Module Interface hdf_mod	96
1.8.1	open_hdf	97
1.8.2	close_hdf	98
1.8.3	write_hdf	98
1.8.4	init_hdf	99
1.8.5	cleanup_hdf	100
1.9	Fortran: Module Interface bpch2_mod	100
1.9.1	open_bpch2_for_read	102
1.9.2	open_bpch2_for_write	102
1.9.3	bpch2_hdr	103
1.9.4	bpch2	104
1.9.5	read_bpch2	104
1.9.6	get_modelname	106

1.9.7	get_name_ext	106
1.9.8	get_name_ext_2d	107
1.9.9	get_res_ext	108
1.9.10	get_halfpolar	109
1.9.11	get_tau0_6a	109
1.10	Fortran: Module Interface file_mod	110
1.10.1	ioerror	112
1.10.2	file_ex_c	113
1.10.3	file_ex_i	113
1.10.4	close_files	114
1.11	Fortran: Module Interface unix_cmds_mod	114
1.11.1	ifort_errmsg	115
1.12	Fortran: Module Interface regrid_a2a_mod	116
1.12.1	do_regrid_a2a	117
1.12.2	map_a2a	118
1.12.3	ymap	119
1.12.4	xmap	120
1.12.5	read_input_grid	121
1.13	Fortran: Module Interface grid_mod	122
1.13.1	compute_grid	123
1.13.2	set_xoffset	124
1.13.3	set_yoffset	124
1.13.4	get_xoffset	125
1.13.5	get_xoffset	125
1.13.6	get_xmid	125
1.13.7	get_xedge	126
1.13.8	get_ymid	126
1.13.9	get_yedge	127
1.13.10	get_ymid_r	127
1.13.11	get_ymid_r_w	128
1.13.12	get_yedge_r	128
1.13.13	get_ysin	129
1.13.14	get_area_m2	129
1.13.15	get_area_cm2	129
1.13.16	get_bounding_box	130
1.13.17	its_a_nested_grid	131
1.13.18	init_grid	131
1.13.19	cleanup_grid	132
1.14	Fortran: Module Interface inquireMod	132
1.14.1	findFreeLUN	132
1.14.2	I_Am_UnOPENed	133
1.15	Fortran: Module Interface global_grid_mod	133
1.15.1	compute_global_grid	134
1.15.2	get_iiipar	135
1.15.3	get_iiipar	135
1.15.4	get_xedge_g	135
1.15.5	get_yedge_g	136
1.15.6	init_global_grid	136

1.15.7 cleanup_global_grid	136
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1 Routine/Function Prologues

1.1 Fortran: Module Interface julday_mod

Module JULDAY_MOD contains routines used to convert from month/day/year to Astronomical Julian Date and back again.

INTERFACE:

```
MODULE JULDAY_MOD
```

USES:

```
IMPLICIT NONE
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
 - 20 Aug 2013 - R. Yantosca - Removed "define.h", this is now obsolete
-

1.1.1 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

1.1.2 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

1.1.3 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

1.2 Fortran: Module Interface transfer_mod

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 CMN_SIZE_MOD
USE ERROR_MOD, ONLY : ALLOC_ERR
USE ERROR_MOD, ONLY : GEOS_CHEM_STOP
```

```
IMPLICIT NONE
```

```
PRIVATE
```

PUBLIC MEMBER FUNCTIONS:

```
PUBLIC :: TRANSFER_A6
PUBLIC :: TRANSFER_2D
PUBLIC :: TRANSFER_3D
```

```

PUBLIC  :: TRANSFER_3D_Bry
PUBLIC  :: TRANSFER_3D_Lp1
PUBLIC  :: TRANSFER_3D_TROP
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, 2/2/12)

=====

GEOS-4, GEOS-5, GEOS-5.7, 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:

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

where

$P_{\text{surface}}(I,J)$ 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. $P_{\text{TOP}}=150$ hPa, and negative sigma edges at higher levels). For these grids, can still use the same formula as for GEOS-4, with one modification:

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

where

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

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

- (1) $B_p(\text{LLPAR}+1) = 0.0$ (L=LLPAR+1 is the atmosphere top)
- (2) $B_p(1) = 1.0$ (L=1 is the surface)
- (3) $P_{\text{TOP}} = A_p(\text{LLPAR}+1)$ (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 $\text{LLPAR} == \text{LGLOB}$ 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
- 02 Feb 2012 - R. Yantosca - Added modifications for GEOS-5.7.x met fields
- 28 Feb 2012 - R. Yantosca - Removed support for GEOS-3
- 01 Mar 2012 - R. Yantosca - Updated to use grid_mod.F90 for the GI model
- 20 Jul 2012 - R. Yantosca - Add routine TRANSFER_3D_Bry, which takes data sized (144,91,:) as inputs & outputs
- 20 Aug 2013 - R. Yantosca - Removed "define.h", this is now obsolete
- 29 Oct 2013 - R. Yantosca - Remove TRANSFER_3D_NOLUMP routine, we can just instead do a direct cast assignment

1.2.1 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,JJPARGLOB)      ! Input data
```

OUTPUT PARAMETERS:

```
REAL*8,  INTENT(OUT) :: OUT(LLPAR,IIPAR,JJPARG)   ! 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
 02 Feb 2012 - R. Yantosca - Treat GEOS-5.7.x the same way as MERRA
 28 Feb 2012 - R. Yantosca - Removed support for GEOS-3
 26 Sep 2013 - R. Yantosca - Renamed GEOS_57 Cpp switch to GEOS_FP

1.2.2 transfer_3d

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

02 Feb 2012 - R. Yantosca - Treat GEOS-5.7.x the same way as MERRA

28 Feb 2012 - R. Yantosca - Removed support for GEOS-3

26 Sep 2013 - R. Yantosca - Renamed GEOS_57 Cpp switch to GEOS_FP

Subroutine TRANSFER_3D_Bry transfers 3-dimensional data from a REAL*4 array to a REAL*8 array. Vertical layers are collapsed (from LGLOB to LLPAR) if necessary. This routine was specially created for 2 x 2.5 Bry data in strat_chem_mod.F90.

SUBROUTINE TRANSFER_3D_Bry(IN, OUT)

```
REAL*4,  INTENT(IN)  :: IN(144,91,LGLOB)    ! Input data
```

```
REAL*8,  INTENT(OUT) :: OUT(144,91,LLPAR)    ! Output data
```

20 Jul 2012 - R. Yantosca - Initial version, based on TRANSFER_3D
26 Sep 2013 - R. Yantosca - Renamed GEOS_57 Cpp switch to GEOS_FP

Subroutine TRANSFER_G5_PLE transfers GEOS-5/MERRA pressure edge data from the native 72-level grid to the reduced 47-level grid.

SUBROUTINE TRANSFER_G5_PLE(IN, OUT)

```
REAL*4,  INTENT(IN)  :: IN(IIPAR,JJPARG,LGLOB+1)      ! Input data
```

```
REAL*8,  INTENT(OUT) :: OUT(IIPAR,JJPARG,LLPAR+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
 02 Feb 2012 - R. Yantosca - Treat GEOS-5.7.x the same way as MERRA
 26 Sep 2013 - R. Yantosca - Renamed GEOS_57 Cpp switch to GEOS_FP

1.2.5 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
 02 Feb 2012 - R. Yantosca - Treat GEOS-5.7.x the same way as MERRA
 26 Sep 2013 - R. Yantosca - Renamed GEOS_57 Cpp switch to GEOS_FP

1.2.6 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+1) ! Input data
```

OUTPUT PARAMETERS:

```
REAL*8,  INTENT(OUT) :: OUT(IIPAR,JJPARG,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
```

1.2.7 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(JJPARG,LGLOB)      ! Input data
```

OUTPUT PARAMETERS:

```
REAL*4,  INTENT(OUT) :: OUT(JJPARG,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 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 ) 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
02 Feb 2012 - R. Yantosca - Treat GEOS-5.7.x the same way as MERRA
28 Feb 2012 - R. Yantosca - Removed support for GEOS-3
26 Sep 2013 - R. Yantosca - Renamed GEOS_57 Cpp switch to GEOS_FP
```

1.2.8 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 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 ) 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
02 Feb 2012 - R. Yantosca - Treat GEOS-5.7.x the same way as MERRA
28 Feb 2012 - R. Yantosca - Removed support for GEOS-3
26 Sep 2013 - R. Yantosca - Renamed GEOS_57 Cpp switch to GEOS_FP
```

1.2.9 transfer_2d_int

Subroutine TRANSFER_2D_INT transfers 2-D data from a REAL*4 array of dimension (IIPAR,JJPAR) to an INTEGER array of dimension (IIPAR,JJPAR).

INTERFACE:

```
SUBROUTINE TRANSFER_2D_INT( IN, OUT )
```

INPUT PARAMETERS:

```
REAL*4, INTENT(IN) :: IN(IIPAR,JJPAR)    ! 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 IO and JO are local variables. (bmy, 3/11/03)
13 Aug 2010 - R. Yantosca - Added ProTeX headers
```

1.2.10 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 IO and JO are local variables (bmy, 3/11/03)
13 Aug 2010 - R. Yantosca - Added ProTeX headers
```

1.2.11 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 IO and JO are local variables. (bmy, 3/11/03)
13 Aug 2010 - R. Yantosca - Added ProTeX headers
```

1.2.12 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
```

1.2.13 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

```

1.2.14 lump_2_r8

Function LUMP_2_R8 lumps 2 sigma levels into one thick level. Input arguments must be REAL*8.

INTERFACE:

```

FUNCTION LUMP_2_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: 2 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

```

1.2.15 lump_4_r4

Function LUMP_4_R4 lumps 4 sigma levels into one thick level. Input arguments must be REAL*4.

INTERFACE:

```
FUNCTION LUMP_4_R4( IN, L_IN, L ) RESULT( OUT )
```

USES:

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

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
```

1.2.16 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
```

1.2.17 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
02 Feb 2012 - R. Yantosca - Treat GEOS-5.7.x the same way as MERRA
28 Feb 2012 - R. Yantosca - Removed support for GEOS-3
26 Sep 2013 - R. Yantosca - Renamed GEOS_57 Cpp switch to GEOS_FP
```

1.2.18 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
```

1.3 Fortran: Module Interface pressure_mod

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
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
#if defined( ESMF_ )
PUBLIC  :: Accept_External_Pedge
#endif
```

REMARKS:

Hybrid Grid Coordinate Definition: (dsa, bmy, 8/27/02, 2/2/12)

=====

.

GEOS-4, GEOS-5, GEOS-5.7, 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. $PTOP=150$ 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)
 $A_p(L)$ = $PTOP$ = model top pressure
 $B_p(L)$ = $SIG(L)$ = 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) $PTOP = A_p(LLPAR+1)$ (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 Ap and Bp 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
02 Feb 2012 - R. Yantosca - Added modifications for GEOS-5.7.x met fields
28 Feb 2012 - R. Yantosca - Removed support for GEOS-3
31 Jul 2012 - R. Yantosca - Modifications for grid-independence
10 Aug 2012 - R. Yantosca - Remove DEVEL from #ifdef for EXTERNAL_PEDGE
11 Dec 2012 - R. Yantosca - Now make EXTERNAL_PEDGE private
11 Dec 2012 - R. Yantosca - Add new routine ACCEPT_PEDGE_FROM_ESMF to set
                        EXTERNAL_PEDGE from the ESMF environment
20 Aug 2013 - R. Yantosca - Removed "define.h", this is now obsolete
18 Sep 2013 - M. Long      - Now use #if defined( ESMF_ ) for HPC code
```

Function GET_AP returns the "A" term [hPa] for the hybrid ETA coordinate.

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

```
USE CMN_SIZE_MOD      ! Size parameters
```

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

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

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

Function GET_BP returns the "B" term [unitless] for the hybrid ETA coordinate.

```
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

```

1.3.3 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:

```

Prior to 11/4/13:
Eliminate array temporary by accepting assumed-shape argument into this
routine. (mpayer, 11/4/13)
      ! Array containing pressure with which to initialize PFLT [hPa]
      REAL*8, INTENT(IN) :: PS(IIPAR,JJPARG)

```

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

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
10 Jun 2013 - R. Yantosca - Avoid array temporaries in CHECK_VALUE
04 Nov 2013 - M. Sulprizio- Eliminate array temporary by accepting assumed-
                        shape dummy argument

```

1.3.4 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
02 Feb 2012 - R. Yantosca - Compute PEDGE for GEOS-5.7.2 the same as MERRA
10 Aug 2012 - R. Yantosca - Need to put #ifdef for EXTERNAL_PEDGE in the
                           section for GEOS-4, GEOS-5, MERRA, GEOS-5.7.x
10 Aug 2012 - R. Yantosca - Now only use Cpp switches EXTERNAL_GRID or
                           EXTERNAL_FORCING to use the GCM pressures.
                           This prevents problems when compiling G-C with
                           the DEVEL tag when using traditional main.F.
26 Sep 2013 - R. Yantosca - Renamed GEOS_57 Cpp switch to GEOS_FP
```

1.3.5 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=LLLPAR+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)
02 Feb 2012 - R. Yantosca - Compute PEDGE for GEOS-5.7.2 the same as MERRA
26 Sep 2013 - R. Yantosca - Renamed GEOS_57 Cpp switch to GEOS_FP

```

1.3.6 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

```

1.3.7 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( am_I_Root )
```

USES:

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

USE CMN_SIZE_MOD ! LLPAR, PTOP
```

INPUT PARAMETERS:

```
LOGICAL, INTENT(IN) :: am_I_Root ! Is this the root CPU?
```

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.\
02 Feb 2012 - R. Yantosca - Compute Ap and Bp for GEOS-5.7.x in the same way
      as for GEOS-5 and MERRA (grids are identical)
28 Feb 2012 - R. Yantosca - Removed support for GEOS-3
30 Jul 2012 - R. Yantosca - Now accept am_I_Root as an argument when
      running with the traditional driver main.F
26 Sep 2013 - R. Yantosca - Renamed GEOS_57 Cpp switch to GEOS_FP
```

1.3.8 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
```

1.3.9 accept_external_pedge

Subroutine ACCEPT_EXTERNAL_PEDGE sets the GEOS-Chem pressure edge variable with the values obtained from an external GCM (such as the NASA GEOS-5 GCM).

INTERFACE:

```
SUBROUTINE Accept_External_Pedge( am_I_Root, State_Met, RC )
```

USES:

```
USE GIGC_ErrCode_Mod
USE GIGC_State_Met_Mod, ONLY : MetState
```

INPUT PARAMETERS:

```
LOGICAL,          INTENT(IN)  :: am_I_Root    ! Are we on root CPU?
TYPE(MetState),   INTENT(IN)  :: State_Met    ! Meteorology state object
!OUTPUT ARGUMENTS:
INTEGER,          INTENT(OUT) :: RC           ! Success or failure?
```

REMARKS:

This routine is a setter for EXTERNAL_PEDGE. It allows us to keep the EXTERNAL_PEDGE array PRIVATE to this module, which is good programming practice.

REVISION HISTORY:

06 Dec 2012 - Initial version

1.4 Fortran: Module Interface directory_mod

Module DIRECTORY_MOD contains the directory path variables used by GEOS-Chem.

INTERFACE:

```
MODULE DIRECTORY_MOD
```

USES:

```
IMPLICIT NONE
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_4_DIR      ! Subdir for GEOS-4 met data
CHARACTER(LEN=255) :: GEOS_5_DIR      ! Subdir for GEOS-5 met data
CHARACTER(LEN=255) :: GEOS_FP_DIR     ! Subdir for GEOS-FP 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=225) :: BC_DIR          ! Dir w/ mean BC files
CHARACTER(LEN=255) :: OC_DIR          ! Dir w/ mean OC 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
CHARACTER(LEN=255) :: TPBC_DIR_SE     ! Dir w/ TPCORE BC's for SEAC4RS grid
CHARACTER(LEN=255) :: POP_EMITDIR     ! Dir w/ POPS emissions files

```

REMARKS:

```

%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%
%% NOTE: THIS MODULE IS NOW OBSOLETE AND IS SLATED TO BE REMOVED! %%
%% WE NOW USE THE DIR NAMES FROM Input_Opt INSTEAD OF FROM HERE. %%
%% -- Bob Yantosca, 20 Aug 2013 %%
%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%

```

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
20 Jan 2011 - CL Friedman - Added OC_DIR and BC_DIR
01 Feb 2012 - R. Yantosca - Added GEOS_57_DIR for GEOS-5.7.x met fields
28 Feb 2012 - R. Yantosca - Removed support for GEOS-3
29 May 2012 - S. Kim - Added directory for SEAC4RS BCs
20 Aug 2013 - R. Yantosca - Removed "define.h", this is now obsolete
26 Sep 2013 - R. Yantosca - Renamed GEOS_57_DIR to GEOS_FP_DIR

```

1.5 Fortran: Module Interface charpak_mod

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

INTERFACE:

```
MODULE CHARPAK_MOD
```

USES:

```
IMPLICIT NONE
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", "txtext.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
- 20 Aug 2013 - R. Yantosca - Removed "define.h", this is now obsolete

1.6 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

PRIVATE TYPES:

PRIVATE

PRIVATE :: HG2_DIAG !hma, 25 Oct 2011

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_I3
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_DAY_OF_WEEK_LT
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
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_I3
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_I3_TIME
PUBLIC  :: GET_I6_TIME
PUBLIC  :: GET_FIRST_A1_TIME
PUBLIC  :: GET_FIRST_A3_TIME
PUBLIC  :: GET_FIRST_A6_TIME
PUBLIC  :: GET_FIRST_I3_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_I3
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
! Add for FLAMBE compatibility (skim, 6/21/13)
PUBLIC  :: ITS_A_NEW_HOUR
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
PUBLIC  :: GET_Hg2_DIAG !hma, 25 Oct 2011
PUBLIC  :: SET_Hg2_DIAG !hma, 25 Oct 2011
[eml
PUBLIC  :: SET_HISTYR
PUBLIC  :: GET_HISTYR
eml]
#if defined( APM )
PUBLIC  :: ITS_TIME_FOR_A6UPDATE
#endif

#if defined( ESMF_ )
PUBLIC  :: Accept_External_Date_Time
#endif

```

REMARKS:

References:

-
- (1) "Practical Astronomy with Your Calculator", 3rd Ed.
Peter Duffett-Smith, Cambridge UP, 1992, p9.
 - (2) Rounding algorithm from: Hultquist, P.F, "Numerical Methods for
Engineers and Computer Scientists", Benjamin/Cummings, Menlo Park CA,
1988, p. 20.
 - (3) Truncation algorithm from: <http://en.wikipedia.org/wiki/Truncation>

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
 25 Oct 2011 - H. Amos - bring Hg2 gas-particle partitioning code into
 v9-01-02
 02 Feb 2012 - R. Yantosca - Added modifications for GEOS-5.7.x met fields
 03 Feb 2012 - R. Yantosca - Added I3 fields timestep variable & routines
 28 Feb 2012 - R. Yantosca - Removed support for GEOS-3
 01 Mar 2012 - R. Yantosca - GET_LOCALTIME now takes (I,J,L,GMT) as args
 14 Jun 2013 - R. Yantosca - Now compute day of week in SET_CURRENT_TIME
 14 Jun 2013 - R. Yantosca - Added comments to module variable declarations
 20 Aug 2013 - R. Yantosca - Removed "define.h", this is now obsolete
 18 Sep 2013 - M. Long - Now use #if defined(ESMF_) for HPC code

1.6.1 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. NOTE: All time variables are returned w/r/t Greenwich Mean Time (aka Universal Time).

INTERFACE:

SUBROUTINE SET_CURRENT_TIME

USES:

USE JULDAY_MOD, ONLY : JULDAY, CALDATE

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
 14 Jun 2013 - R. Yantosca - Now move the day of week computation here
 20 Aug 2013 - R. Yantosca - Removed "define.h", this is now obsolete

1.6.2 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

1.6.3 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
```

1.6.4 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
```

1.6.5 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
```

1.6.6 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
```

1.6.7 set_timesteps

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

INTERFACE:

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

INPUT PARAMETERS:

```
LOGICAL, INTENT(IN) :: am_I_Root    ! Is this the root CPU?
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]
```

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
 30 Jul 2012 - R. Yantosca - Now accept am_I_Root as an argument when
 running with the traditional driver main.F

1.6.8 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

1.6.9 set_hg2_diag

Subroutine SET_Hg2_DIAG increments Hg2_DIAG, the counter for the number of times AAD03_Fg and AD03_Fp are recorded. (hma 20100218)

INTERFACE:

```
SUBROUTINE SET_Hg2_DIAG( INCREMENT, RESET )
```

INPUT PARAMETERS:

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

REVISION HISTORY:

18 Feb 2012 - H. Amos - Initial version
 07 Mar 2012 - M. Payer - Added ProTeX headers

1.6.10 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
```

1.6.11 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
```

1.6.12 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

```

1.6.13 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

```

1.6.14 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

```

1.6.15 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
```

1.6.16 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
```

1.6.17 set_ct_i3

Subroutine SET_CT_I3 increments CT_I3, the counter of the number of times we have read in I-3 fields.

INTERFACE:

```
SUBROUTINE SET_CT_I3( INCREMENT, RESET )
```

INPUT PARAMETERS:

```

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

```

REVISION HISTORY:

03 Feb 2012 - R. Yantosca - Initial version, for GEOS-5.7.2

1.6.18 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

1.6.19 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

1.6.20 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
```

1.6.21 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
```

1.6.22 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

1.6.23 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

1.6.24 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

1.6.25 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
```

1.6.26 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
```

1.6.27 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
```

1.6.28 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
```

1.6.29 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
```

1.6.30 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
```

1.6.31 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
```

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
20 Aug 2013 - R. Yantosca - Removed "define.h", this is now obsolete
```

1.6.32 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
```

1.6.33 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
```

1.6.34 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
```

1.6.35 set_histyr

Function SET_HISTYR returns the year stored in HISTYR w/o needing to include the CMN_O3 common block (eml, 8/20/08)

INTERFACE:

```
SUBROUTINE SET_HISTYR( YEARIN )
```

INPUT PARAMETERS:

```
INTEGER, INTENT(IN) :: YEARIN
```

REVISION HISTORY:

```
20 Aug 2008 - E. Leibensperger - Initial version
07 Mar 2012 - M. Payer          - Added ProTeX headers
```

1.6.36 get_histyr

Function GET_HISTYR returns the year stored in HISTYR w/o needing to include the CMN_O3 common block (eml, 8/20/08)

INTERFACE:

```
FUNCTION GET_HISTYR() RESULT( HISTYEAR )
```

RETURN VALUE:

```
INTEGER :: HISTYEAR
```

REVISION HISTORY:

```
20 Aug 2008 - E. Leibensperger - Initial version
07 Mar 2012 - M. Payer          - Added ProTeX headers
```

1.6.37 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
```

1.6.38 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
```

1.6.39 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
```

1.6.40 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( THISDAYOFTYEAR )
```

RETURN VALUE:

```
INTEGER :: THISDAYOFTYEAR ! Day of year
```

REVISION HISTORY:

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

1.6.41 get_day_of_week

Function GET_DAY_OF_WEEK returns the day of the week (with respect to GREENWICH MEAN TIME) 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
```

REVISION HISTORY:

```
05 Feb 2003 - R. Yantosca - Initial Version
15 Jan 2010 - R. Yantosca - Added ProTeX headers
14 Jun 2013 - R. Yantosca - Now move computation to SET_CURRENT_TIME
```

1.6.42 get_day_of_week.lt

Function GET_DAY_OF_WEEK_LT returns the day of the week (with respect to the SOLAR LOCAL TIME AT GRID BOX [I,J,L]) as a number: Sun=0, Mon=1, Tue=2, Wed=3, Thu=4, Fri=5, Sat=6.

INTERFACE:

```
FUNCTION GET_DAY_OF_WEEK_LT( I, J, L ) RESULT( DAY_NUM )
```

USES:

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

INPUT PARAMETERS:

```
INTEGER, INTENT(IN) :: I      ! Grid box lon index
INTEGER, INTENT(IN) :: J      ! Grid box lat index
INTEGER, INTENT(IN) :: L      ! Grid box level index
```

RETURN VALUE:

```
INTEGER                :: DAY_NUM    ! Day of week, w/r/t local time
```

REMARKS:

This routine is used by various emissions routines, in order to determine whether weekday or weekend emissions need to be applied.

REVISION HISTORY:

```
13 Jun 2013 - R. Yantosca - Initial version
```

1.6.43 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
```

1.6.44 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
```

1.6.45 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
```

1.6.46 get_tau_e

Function GET_TAU_e returns TAU_e (hours since 1 Jan 1985 at the end of a GEOS-Chem run) to the calling program.

INTERFACE:

```
FUNCTION GET_TAU_e() RESULT( THISTAU_e )
```

RETURN VALUE:

```
REAL*8 :: THISTAU_e ! TAU_e [hrs since 1/1/1985]
```

REVISION HISTORY:

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

1.6.47 get_diag_b

Function GET_DIAG_b returns DIAG_b (hours since 1 Jan 1985 at the start of a diagnostic interval) to the calling program.

INTERFACE:

```
FUNCTION GET_DIAG_b() RESULT( THISDIAG_b )
```

RETURN VALUE:

```
INTEGER :: THISDIAG_b ! DIAG_b [hrs since 1/1/1985]
```

REVISION HISTORY:

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

1.6.48 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 since 1/1/1985]
```

REVISION HISTORY:

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

1.6.49 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, J, L, GMT ) RESULT( THISLOCALTIME )
```

USES:

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

INPUT PARAMETERS:

```
INTEGER, INTENT(IN)      :: I           ! Longitude index
INTEGER, INTENT(IN)      :: J           ! Latitude index
INTEGER, INTENT(IN)      :: L           ! Level 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
01 Mar 2012 - R. Yantosca - Now use GET_XMID(I,J,L) from grid_mod.F90, and
                           add J, L indices to the argument list
```

1.6.50 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
```

1.6.51 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
```

1.6.52 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
```

1.6.53 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
```

1.6.54 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
```

1.6.55 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
```

1.6.56 `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
```

1.6.57 `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
```

1.6.58 `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
```

1.6.59 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
```

1.6.60 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
```

1.6.61 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
```

1.6.62 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
```

1.6.63 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
```

1.6.64 get_ct_i3

Function GET_CT_I3 returns the I-3 fields timestep counter to the calling program

INTERFACE:

```
FUNCTION GET_CT_I3() RESULT( THIS_CT_I3 )
```

RETURN VALUE:

```
INTEGER :: THIS_CT_I3    ! # of I-6 timesteps
```

REVISION HISTORY:

```
03 Feb 2012 - R. Yantosca - Initial version, for GEOS-5.7.2
```

1.6.65 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
```

1.6.66 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
```

1.6.67 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
```

1.6.68 get_hg2_diag

Function GET_Hg2_DIAG returns the DIAG timestep counter to the calling program. (hma 20100218)

INTERFACE:

```
FUNCTION GET_Hg2_DIAG() RESULT( THIS_Hg2_DIAG )
```

RETURN VALUE:

```
INTEGER :: THIS_Hg2_DIAG ! # of diagnostic timesteps
```

REVISION HISTORY:

```
18 Feb 2012 - H. Amos      - Initial version
07 Mar 2012 - M. Payer    - Added ProTeX headers
```

1.6.69 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 )
```

RETURN VALUE:

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

REVISION HISTORY:

```
19 Aug 2010 - R. Yantosca - Initial version
02 Feb 2012 - R. Yantosca - Added modifications for GEOS-5.7.x met fields
20 Aug 2013 - R. Yantosca - Removed "define.h", this is now obsolete
26 Sep 2013 - R. Yantosca - Renamed GEOS_57 Cpp switch to GEOS_FP
```

1.6.70 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 )
```

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
28 Feb 2012 - R. Yantosca - Removed support for GEOS-3
20 Aug 2013 - R. Yantosca - Removed "define.h", this is now obsolete
```

1.6.71 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)
```

1.6.72 get_i3_time

Function GET_I3_TIME returns the correct YYYYMMDD and HHMMSS values that are needed to read in the next instantaneous 3-hour (I-3) fields.

INTERFACE:

```
FUNCTION GET_I3_TIME() RESULT( DATE )
```

RETURN VALUE:

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

REMARKS:

Modified for start times other than 0 GMT.

REVISION HISTORY:

6 Feb 2012 - R. Yantosca - Initial version

1.6.73 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

1.6.74 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 )
```

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
 20 Aug 2013 - R. Yantosca - Removed "define.h", this is now obsolete

1.6.75 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 )
```

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
28 Feb 2012 - R. Yantosca - Removed support for GEOS-3
20 Aug 2013 - R. Yantosca - Removed "define.h", this is now obsolete
```

1.6.76 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 )
```

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
20 Aug 2013 - R. Yantosca - Removed "define.h", this is now obsolete
```

1.6.77 get_first_i3_time

Function GET_FIRST_I3_TIME returns the correct YYYYMMDD and HHMMSS values the first time that I-3 fields are read in from disk.

INTERFACE:

```
FUNCTION GET_FIRST_I3_TIME() RESULT( DATE )
```

RETURN VALUE:

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

REVISION HISTORY:

03 Feb 2012 - R. Yantosca - Initial version, for GEOS-5.7.2

1.6.78 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

1.6.79 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)

1.6.80 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
```

1.6.81 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
```

1.6.82 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)

1.6.83 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

1.6.84 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

1.6.85 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
```

1.6.86 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
```

1.6.87 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 )
```

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)
 20 Aug 2013 - R. Yantosca - Removed "define.h", this is now obsolete

1.6.88 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 )
```

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
 20 Aug 2013 - R. Yantosca - Removed "define.h", this is now obsolete

1.6.89 its_time_for_i3

Function ITS_TIME_FOR_I3 returns TRUE if it is time to read in I2 (instantaneous 3-hr fields) and FALSE otherwise.

INTERFACE:

```
FUNCTION ITS_TIME_FOR_I3() RESULT( FLAG )
```

RETURN VALUE:

LOGICAL :: FLAG

REVISION HISTORY:

03 Feb 2012 - R. Yantosca - Initial version, for GEOS-5.7.2

1.6.90 `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

1.6.91 `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

1.6.92 `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
```

1.6.93 `its_time_for_exit`

Function `ITS_TIME_FOR_EXIT` returns TRUE if it is the end of the GEOS-Chem simulation (i.e. $\tau = \tau_e$), 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
```

1.6.94 `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
```

1.6.95 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
```

1.6.96 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
```

1.6.97 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

```

1.6.98 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

1.6.99 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

1.6.100 its_a_new_hour

Function ITS_A_NEW_HOUR returns TRUE if it's the first timestep of a new hour (it also returns TRUE on the first timestep of the run). This is useful for setting flags for reading in data. (bmy, 4/1/04)

INTERFACE:

```
FUNCTION ITS_A_NEW_HOUR( ) RESULT( IS_NEW_HOUR )
```

RETURN VALUE:

```
LOGICAL :: IS_NEW_HOUR
```

REVISION HISTORY:

```
01 Apr 2004 - R. Yantosca - Initial Version
25 Feb 2014 - M. Sulprizio- Added ProTeX headers
```

1.6.101 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
```

1.6.102 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
```

1.6.103 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 )
```

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
20 Aug 2013 - R. Yantosca - Removed "define.h", this is now obsolete
```

1.6.104 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
```

1.6.105 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
```

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
20 Aug 2013 - R. Yantosca - Removed "define.h", this is now obsolete
```

1.6.106 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
```

1.6.107 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
```

1.6.108 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
```

1.6.109 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
```

1.6.110 accept_external_date_time

Subroutine ACCEPT_EXTERNAL_DATE_TIME sets the date and time variables in time_mod.F with the values obtained from an external GCM (such as NASA's GEOS-5 GCM). The various date & time values from the GCM are passed as arguments.

INTERFACE:

```

SUBROUTINE Accept_External_Date_Time(
&   am_I_Root,    value_NYMDb,    value_NYMDe,    value_NYMD,
&   value_NHMSb,  value_NHMSe,    value_NHMS,    value_YEAR,
&   value_MONTH,  value_DAY,      value_DAYOFYR,  value_HOUR,
&   value_MINUTE, value_SECOND,    value_UTC,     value_HELAPSED,
&   value_TS_CHEM, value_TS_CONV, value_TS_DYN,   value_TS_EMIS,
&   RC
)
```

USES:

```

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

INPUT PARAMETERS:

```

LOGICAL, INTENT(IN)  :: am_I_Root      ! Are we on the root CPU?
INTEGER, OPTIONAL    :: value_NYMDb    ! YYYY/MM/DD @ start of run
INTEGER, OPTIONAL    :: value_NYMDe    ! YYYY/MM/DD @ end of run
INTEGER, OPTIONAL    :: value_NYMD     ! YYYY/MM/DD @ current time
INTEGER, OPTIONAL    :: value_NHMSb    ! hh:mm:ss @ start of run
INTEGER, OPTIONAL    :: value_NHMSe    ! hh:mm:ss @ end of run
INTEGER, OPTIONAL    :: value_NHMS     ! hh:mm:ss @ current time
INTEGER, OPTIONAL    :: value_YEAR     ! UTC year
INTEGER, OPTIONAL    :: value_MONTH    ! UTC month
INTEGER, OPTIONAL    :: value_DAY      ! UTC day
INTEGER, OPTIONAL    :: value_DAYOFYR  ! UTC day of year
INTEGER, OPTIONAL    :: value_HOUR     ! UTC hour
INTEGER, OPTIONAL    :: value_MINUTE   ! UTC minute
INTEGER, OPTIONAL    :: value_SECOND   ! UTC second
REAL*4, OPTIONAL     :: value_UTC      ! UTC time [hrs]
REAL*4, OPTIONAL     :: value_HELAPSED ! Elapsed hours
INTEGER, OPTIONAL    :: value_TS_CHEM  ! Chemistry timestep [min]
INTEGER, OPTIONAL    :: value_TS_CONV  ! Convection timestep [min]
INTEGER, OPTIONAL    :: value_TS_DYN   ! Dynamic timestep [min]
INTEGER, OPTIONAL    :: value_TS_EMIS  ! Emissions timestep [min]

! OUTPUT ARGUMENTS:
INTEGER, INTENT(OUT) :: RC              ! Success or failure?
```

REMARKS:

The date and time values are obtained via the Extract_ subroutine in module file GEOSCHEMchem_GridCompMod.F90.

```
06 Dec 2012 - Initial version
11 Dec 2012 - R. Yantosca - Renamed to ACCEPT_EXTERNAL_DATE_TIME
18 Jun 2013 - R. Yantosca - Now compute day of week w/r/t GMT, which is
                           the same modification made in SET CURRENT TIME
```

Module `ERROR_MOD` contains error checking routines.

MODULE ERROR MOD

```

IMPLICIT NONE
PRIVATE

```

```

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_DBL
END INTERFACE

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

```

REMARKS:

```

%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%
%% NOTE: "LINUX_IFORT" and "LINUX_PGI" ARE CURRENTLY THE ONLY      %%
%% SUPPORTED COMPILER OPTIONS.  MOST SYSTEMS NOW USE A UNIX VERSION %%
%% BASED ON LINUX, SO OTHER COMPILERS (IBM/AIX, SUN/SPARC, etc.)    %%
%% ARE GENERALLY NOT USED ANYMORE.  LEAVE THE OLDER CODE HERE JUST %%
%% IN CASE WE NEED TO REVERT TO IT AGAIN IN THE FUTURE.           %%
%%      -- Bob Yantosca, 20 Aug 2013                                %%
%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%

```

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
 20 Aug 2013 - R. Yantosca - Removed "define.h", this is now obsolete

1.7.1 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:

```

#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
 20 Aug 2013 - R. Yantosca - Removed "define.h", this is now obsolete

1.7.2 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:

```
#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
 20 Aug 2013 - R. Yantosca - Removed "define.h", this is now obsolete

1.7.3 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:

```
#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
- 20 Aug 2013 - R. Yantosca - Removed "define.h", this is now obsolete

1.7.4 finite_dble

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:

```
#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
- 20 Aug 2013 - R. Yantosca - Removed "define.h", this is now obsolete

1.7.5 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=255), INTENT(IN) :: VARNAME         ! Name of variable
CHARACTER(LEN=255), 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
 10 Jun 2013 - R. Yantosca - Avoid array temporaries, use CHAR*255 args

1.7.6 check_dble_value

Subroutine CHECK_DBLE_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_DBLE_VALUE( VALUE, LOCATION, VARNAME, MESSAGE )
```

INPUT PARAMETERS:

```
REAL*8,          INTENT(IN) :: VALUE          ! Value to be checked
CHARACTER(LEN=255), INTENT(IN) :: VARNAME      ! Name of variable
CHARACTER(LEN=255), 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
 10 Jun 2013 - R. Yantosca - Avoid array temporaries, use CHAR*255 args

1.7.7 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

1.7.8 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:

```
#if defined( ESMF_ )
!-----
!      %%%% CONNECTING TO GEOS-5 GCM via ESMF INTERFACE %%%%
! Use GEOS-5 style error reporting when connecting to the GEOS-5
! GCM via the ESMF interface (bmy, 3/12/13)
!-----
USE MAPL_Mod
#   include "MAPL_Generic.h"
#endif
```

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
12 Mar 2013 - R. Yantosca - Now use GEOS-5 style traceback when using ESMF
```

1.7.9 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 )
```

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
20 Aug 2013 - R. Yantosca - Removed "define.h", this is now obsolete
```

1.7.10 debug_msg

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

INTERFACE:

```
SUBROUTINE DEBUG_MSG( MESSAGE )
```

INPUT PARAMETERS:

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

REVISION HISTORY:

```
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
20 Aug 2013 - R. Yantosca - Removed "define.h", this is now obsolete
```

1.7.11 safe_div

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.

INTERFACE:

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

INPUT PARAMETERS:

```
REAL*8,          INTENT(IN) :: N          ! Numerator
REAL*8,          INTENT(IN) :: D          ! Denominator
REAL*8,          INTENT(IN) :: ALT_NAN    ! Alternate value to be
                                           ! returned if the division
                                           ! is either NAN (0/0) or
                                           ! leads to overflow (i.e.,
                                           ! a too large number)
REAL*8, OPTIONAL, INTENT(IN) :: ALT_OVER  ! Alternate value to be
                                           ! returned if the division
                                           ! leads to overflow (default
                                           ! is ALT_NAN)
REAL*8, OPTIONAL, INTENT(IN) :: ALT_UNDER ! 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/

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

1.7.12 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

1.7.13 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

1.7.14 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

1.7.15 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

1.7.16 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

1.8 Fortran: Module Interface hdf_mod

Module HDF_MOD contains routines to write data to HDF5 files.

INTERFACE:

```
MODULE HDF_MOD
```

USES:

```
IMPLICIT NONE
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
```


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

```

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
20 Aug 2013 - R. Yantosca      - Removed "define.h", this is now obsolete

```

Subroutine OPEN_HDF creates and opens an hdf file for output.

```
#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   )
```

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

USE CMN_SIZE_MOD      ! Size 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

```
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
01 Mar 2012 - R. Yantosca      - Now use GET_XMID(I,J,L), GET_YMID(I,J,L)
                                from the new grid_mod.F90
17 Oct 2012 - M. Payer          - Bug fix: Reorder indices for GET_YMID,
                                otherwise LAT variable in HDF files will
                                be wrong (S. Philip)

```

Subroutine `CLOSE_HDF` closes an HDF5 file that is already open.

```

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

```

USE HDF5

```
INTEGER(HID_T), INTENT(INOUT) :: IU_HDF  ! HDF5 File identifier
```

```

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

```

Subroutine `WRITE_HDF` writes data to an open HDF5 file.

```
#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
17 Oct 2012 - M. Payer          - Add PBLDEPTH to fix bad HDF category error
                                (S. Philip, A. van Donkelaar)

```

1.8.4 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:

```

23 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

```

1.8.5 cleanup_hdf

Subroutine CLEANUP_HDF deallocates all module variables.

INTERFACE:

```

SUBROUTINE CLEANUP_HDF

```

REVISION HISTORY:

```

21 Dec 2009 - R. Yantosca - Initial version

```

1.9 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:

```

USE inquireMod, ONLY : findFreeLUN
USE inquireMod, ONLY : I_Am_UnOPENed

```

```

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
28 Feb 2012 - R. Yantosca - Removed support for GEOS-3
19 Jul 2012 - R. Yantosca - Bug fix in GET_NAME_EXT_2D
03 Aug 2012 - R. Yantosca - Reference file LUN routines from inquireMod.F90

```

1.9.1 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   ! File to open

```

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
06 Aug 2012 - R. Yantosca - Do not call findFreeLun() in this subroutine
                           but instead in the calling routine and pass
                           the file unit as an argument.

```

1.9.2 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
06 Aug 2012 - R. Yantosca - Do not call findFreeLun() in this subroutine
                           but instead in the calling routine and pass
                           the file unit as an argument.
```

1.9.3 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
```

1.9.4 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)

1.9.5 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 : IOERROR

```

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
- 25 May 2012 - R. Yantosca - Update TEMPARRAY for GRID025x03125

03 Aug 2012 - R. Yantosca - Move calls to findFreeLUN out of DEVEL block
 07 Aug 2012 - R. Yantosca - Now print LUN used to open file
 20 Aug 2013 - R. Yantosca - Removed "define.h", this is now obsolete
 26 Sep 2013 - R. Yantosca - Removed SEAC4RS C-preprocessor switch

1.9.6 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
 01 Feb 2012 - R. Yantosca - Added GEOS-5.7.x model names
 28 Feb 2012 - R. Yantosca - Removed support for GEOS-3
 11 Dec 2012 - R. Yantosca - Bug fix: Need to specify both EXTERNAL_GRID and EXTERNAL_FORCING Cpp switches
 26 Sep 2013 - R. Yantosca - Renamed GEOS_57 Cpp switch to GEOS_FP

1.9.7 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 )
```

RETURN VALUE:

```

#if defined( GEOS_4 )
    CHARACTER(LEN=5) :: NAME_EXT
    NAME_EXT = 'geos4'

#elif defined( GEOS_5 ) || defined( GEOS_FP )
    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 = 'merra'

#elif defined( EXTERNAL_GRID ) || ( EXTERNAL_FORCING )
    CHARACTER(LEN=5) :: NAME_EXT
    NAME_EXT = 'ext'

#endif

```

REMARKS:

NOTE: GEOS-FP and GEOS-5 met products use the same vertical grid, so we can return the name extension "geos5" for both.

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.
01 Feb 2012 - R. Yantosca - Now also define output for GEOS-5.7.x met
28 Feb 2012 - R. Yantosca - Removed support for GEOS-3
11 Dec 2012 - R. Yantosca - Bug fix: Need to specify both EXTERNAL_GRID and
                        EXTERNAL_FORCING Cpp switches
20 Aug 2013 - R. Yantosca - Removed "define.h", this is now obsolete

```

1.9.8 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    ! Return 1st 4 chars of "geos", "gcap"
```

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
19 Jul 2012 - R. Yantosca - For MERRA meterology, return "geos", which
                        indicates surface data only
```

1.9.9 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 )
```

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'

#elif defined( GRID025x03125 )
CHARACTER(LEN=9) :: RES_EXT
```

```

RES_EXT = '025x03125'

#elif defined( EXTERNAL_GRID )
  CHARACTER(LEN=8) :: RES_EXT
  RES_EXT = 'external'

#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
10 Feb 2012 - R. Yantosca - Added extension for 0.25 x 0.3125 grids
20 Aug 2013 - R. Yantosca - Removed "define.h", this is now obsolete

```

1.9.10 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 )
```

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
20 Aug 2013 - R. Yantosca - Removed "define.h", this is now obsolete

```

1.9.11 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

1.10 Fortran: Module Interface file_mod

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
      PRIVATE

```

DEFINED PARAMETERS:

```

!-----
! In the GEOS-5 GCM, the unit numbers cannot be PARAMETERS.
! Instead, use INQUIRES to find open LUNs at the point of
! request. References to most IU_* variables have now been
! made local. IU_BPCH is the only LUN that needs to be seen
! across several variables.
!-----

```

```

! Logical file unit numbers for ...
INTEGER, PUBLIC :: IU_BPCH      ! "ctm.bpch"
INTEGER, PUBLIC :: IU_FILE

```

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_CTM_CHEM 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
29 May 2010 - S. Kim        - Add IU_BC_SE for the SEAC4RS grid
06 Aug 2012 - R. Yantosca  - Remove several IU_* variables, as these have
                             now been moved to various other modules
20 Aug 2013 - R. Yantosca  - Removed "define.h", this is now obsolete

```

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)

SUBROUTINE IOERROR(ERROR_NUM, UNIT, LOCATION)

```
USE_ERROR_MOD, ONLY : GEOS_CHEM_STOP
```

```

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

```

[illegible]

- (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

20 Aug 2013 - R. Yantosca - Removed "define.h", this is now obsolete

1.10.2 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 )
```

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

20 Aug 2013 - R. Yantosca - Removed "define.h", this is now obsolete

1.10.3 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 )
```

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
20 Aug 2013 - R. Yantosca - Removed "define.h", this is now obsolete
```

1.10.4 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
29 May 2010 - S. Kim      - Now close IU_BC_SE
```

1.11 Fortran: Module Interface unix_cmds_mod

Module UNIX_CMDS_MOD contains variables which contain file suffixes and various Unix command strings.

INTERFACE:

```
MODULE UNIX_CMDS_MOD
```

USES:

```
IMPLICIT NONE
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' )
```

REMARKS:

```
%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%
%% NOTE: THIS MODULE IS NOW OBSOLETE AND IS SLATED TO BE REMOVED! %%
%% WE NOW USE THE UNIX CMDS FROM Input_Opt INSTEAD OF FROM HERE. %%
%% -- Bob Yantosca, 20 Aug 2013 %%
%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%%
```

REVISION HISTORY:

```
09 Jul 2004 - R. Yantosca - Initial version
20 Nov 2009 - R. Yantosca - Added ProTeX header
20 Aug 2013 - R. Yantosca - Removed "define.h", this is now obsolete
```

1.11.1 ifort_errmsg

Function IFORT_ERRMSG returns an error message string that corresponds to an I/O error number obtained via the IOSTAT or STAT specifiers. (This is specifically for the Intel Fortran compiler.)

INTERFACE:

```
FUNCTION IFORT_ERRMSG( ERROR_NUM ) RESULT( MSG )
```

INPUT PARAMETERS:

```
INTEGER, INTENT(IN) :: ERROR_NUM    ! Error condition from IOSTAT
```

RETURN VALUE:

```
CHARACTER(LEN=255)  :: MSG          ! Descriptive error message
```

REVISION HISTORY:

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

1.12 Fortran: Module Interface regrid_a2a_mod

Module REGRID_A2A_MOD uses an algorithm adapted from MAP_A2A code to regrid from one horizonatal grid to another.

INTERFACE:

```
MODULE REGRID_A2A_MOD
```

USES:

```
IMPLICIT NONE
PRIVATE
```

PRIVATE MEMBER FUNCTIONS:

```
PRIVATE :: XMAP
PRIVATE :: YMAP
PRIVATE :: READ_INPUT_GRID
```

PUBLIC MEMBER FUNCTIONS:

```
PUBLIC  :: DO_REGRID_A2A
PUBLIC  :: MAP_A2A
```

REVISION HISTORY:

```
13 Mar 2012 - M. Cooper    - Initial version
03 Apr 2012 - M. Payer    - Now use functions GET_AREA_CM2(I,J,L),
                           GET_YEDGE(I,J,L) and GET_YSIN(I,J,L) from the
                           new grid_mod.F90
22 May 2012 - L. Murray   - Implemented several bug fixes
23 Aug 2012 - R. Yantosca - Add capability for starting from hi-res grids
                           (generic 0.5x0.5, generic 0.25x0.25, etc.)
23 Aug 2012 - R. Yantosca - Add subroutine READ_INPUT_GRID, which reads the
                           grid parameters (lon & lat edges) w/ netCDF
27 Aug 2012 - R. Yantosca - Now parallelize key DO loops
```

1.12.1 do_regrid_a2a

Subroutine DO_REGRID_A2A regrids 2-D data in the horizontal direction.

INTERFACE:

```
SUBROUTINE DO_REGRID_A2A( FILENAME, IM, JM, INGRID, OUTGRID, IS_MASS, &
                           netCDF )
```

USES:

```
USE GRID_MOD,    ONLY : GET_XEDGE
USE GRID_MOD,    ONLY : GET_YSIN
USE GRID_MOD,    ONLY : GET_AREA_CM2
USE FILE_MOD,    ONLY : IOERROR
USE inquireMod,  ONLY : findFreeLUN
USE CMN_SIZE_MOD
USE CMN_GCTM_MOD
```

INPUT PARAMETERS:

```
! Name of file with lon and lat edge information on the INPUT GRID
CHARACTER(LEN=*), INTENT(IN)    :: FILENAME

! Number of lon centers and lat centers on the INPUT GRID
INTEGER,          INTENT(IN)    :: IM
INTEGER,          INTENT(IN)    :: JM

! Data array on the input grid
REAL*8,           INTENT(IN)    :: INGRID(IM,JM)

! IS_MASS=0 if data is units of concentration (molec/cm2/s, unitless, etc.)
! IS_MASS=1 if data is units of mass (kg/yr, etc.); we will need to convert
!           INGRID to per unit area
INTEGER,          INTENT(IN)    :: IS_MASS

! Read from netCDF file? (needed for debugging, will disappear later)
LOGICAL, OPTIONAL, INTENT(IN)   :: netCDF
```

OUTPUT PARAMETERS:

```
! Data array on the OUTPUT GRID
REAL*8,           INTENT(OUT)   :: OUTGRID(IIPAR,JJPARG)
```

REVISION HISTORY:

```
13 Mar 2012 - M. Cooper   - Initial version
22 May 2012 - L. Murray   - Bug fix: INSIN should be allocated w/ JM+1.
22 May 2012 - R. Yantosca - Updated comments, cosmetic changes
25 May 2012 - R. Yantosca - Bug fix: declare the INGRID argument as
                           INTENT(IN) to preserve the values of INGRID
```

```

                                in the calling routine
06 Aug 2012 - R. Yantosca - Now make IU_REGRID a local variable
06 Aug 2012 - R. Yantosca - Move calls to findFreeLUN out of DEVEL block
23 Aug 2012 - R. Yantosca - Now use f10.4 format for hi-res grids
23 Aug 2012 - R. Yantosca - Now can read grid info from netCDF files
27 Aug 2012 - R. Yantosca - Add parallel DO loops
03 Jan 2013 - M. Payer      - Renamed PERAREA to IS_MASS to describe parameter
                                more clearly

```

1.12.2 map_a2a

Subroutine MAP_A2A is a horizontal arbitrary grid to arbitrary grid conservative high-order mapping regridding routine by S-J Lin.

INTERFACE:

```

(1 ) INLON   (REAL*8   ) : Longitude edges of input grid
(2 ) INSIN   (REAL*8   ) : Sine of input grid latitude edges
(3 ) INGRID  (REAL*8   ) : Data array to be regridded

```

```

SUBROUTINE map_a2a( im, jm, lon1, sin1, q1, &
                   in, jn, lon2, sin2, q2, ig, iv)

```

INPUT PARAMETERS:

```

! Longitude and Latitude dimensions of INPUT grid
INTEGER, INTENT(IN)  :: im, jm

! Longitude and Latitude dimensions of OUTPUT grid
INTEGER, INTENT(IN)  :: in, jn

! IG=0: pole to pole;
! IG=1 J=1 is half-dy north of south pole
INTEGER, INTENT(IN)  :: ig

! IV=0: Regrid scalar quantity
! IV=1: Regrid vector quantity
INTEGER, INTENT(IN)  :: iv

! Longitude edges (degrees) of INPUT and OUTPUT grids
REAL*8,  INTENT(IN)  :: lon1(im+1), lon2(in+1)

! Sine of Latitude Edges (radians) of INPUT and OUTPUT grids
REAL*8,  INTENT(IN)  :: sin1(jm+1), sin2(jn+1)

! Quantity on INPUT grid
REAL*8,  INTENT(IN)  :: q1(im,jm)

```

OUTPUT PARAMETERS:

```
! Regridded quantity on OUTPUT grid
REAL*8, INTENT(OUT) :: q2(in,jn)
```

REVISION HISTORY:

- (1) Original subroutine by S-J Lin. Converted to F90 freeform format and inserted into "Geos3RegridModule" by Bob Yantosca (9/21/00)
 - (2) Added F90 type declarations to be consistent w/ TypeModule.f90. Also updated comments. (bmy, 9/21/00)
 - 21 Sep 2000 - R. Yantosca - Initial version
 - 27 Aug 2012 - R. Yantosca - Add parallel DO loops
-

1.12.3 ymap

Routine to perform area preserving mapping in N-S from an arbitrary resolution to another.

INTERFACE:

```
SUBROUTINE ymap(im, jm, sin1, q1, jn, sin2, q2, ig, iv)
```

INPUT PARAMETERS:

```
! original E-W dimension
INTEGER, INTENT(IN) :: im

! original N-S dimension
INTEGER, INTENT(IN) :: jm

! Target N-S dimension
INTEGER, INTENT(IN) :: jn

! IG=0: scalars from SP to NP (D-grid v-wind is also IG=0)
! IG=1: D-grid u-wind
INTEGER, INTENT(IN) :: ig

! IV=0: scalar;
! IV=1: vector
INTEGER, INTENT(IN) :: iv

! Original southern edge of the cell sin(lat1)
REAL*8, INTENT(IN) :: sin1(jm+1-ig)

! Original data at center of the cell
REAL*8, INTENT(IN) :: q1(im,jm)

! Target cell's southern edge sin(lat2)
REAL*8, INTENT(IN) :: sin2(jn+1-ig)
```

OUTPUT PARAMETERS:

```
! Mapped data at the target resolution
REAL*8, INTENT(OUT) :: q2(im,jn)
```

REMARKS:

```
sin1 (1) = -1 must be south pole; sin1(jm+1)=1 must be N pole.
sin1(1) < sin1(2) < sin1(3) < ... < sin1(jm) < sin1(jm+1)
sin2(1) < sin2(2) < sin2(3) < ... < sin2(jn) < sin2(jn+1)!
```

AUTHOR:

```
Developer: Prasad Kasibhatla
March 6, 2012
```

!REVISION HISTORY

```
06 Mar 2012 - P. Kasibhatla - Initial version
27 Aug 2012 - R. Yantosca   - Added parallel DO loops
27 Aug 2012 - R. Yantosca   - Change REAL*4 variables to REAL*8 to better
                               ensure numerical stability
```

1.12.4 xmap

Routine to perform area preserving mapping in E-W from an arbitrary resolution to another. Periodic domain will be assumed, i.e., the eastern wall bounding cell im is lon1(im+1) = lon1(1); Note the equal sign is true geographically.

INTERFACE:

```
SUBROUTINE xmap(im, jm, lon1, q1, in, lon2, q2)
```

INPUT PARAMETERS:

```
! Original E-W dimension
INTEGER, INTENT(IN) :: im

! Target E-W dimension
INTEGER, INTENT(IN) :: in

! Original N-S dimension
INTEGER, INTENT(IN) :: jm

! Original western edge of the cell
REAL*8, INTENT(IN) :: lon1(im+1)

! Original data at center of the cell
REAL*8, INTENT(IN) :: q1(im,jm)

! Target cell's western edge
REAL*8, INTENT(IN) :: lon2(in+1)
```

OUTPUT PARAMETERS:


```
! Mapped data at the target resolution
REAL*8,  INTENT(OUT) :: q2(in,jm)
```

REMARKS:

```
lon1(1) < lon1(2) < lon1(3) < ... < lon1(im) < lon1(im+1)
lon2(1) < lon2(2) < lon2(3) < ... < lon2(in) < lon2(in+1)
```

AUTHOR:

```
Developer: Prasad Kasibhatla
March 6, 2012
```

!REVISION HISTORY

```
06 Mar 2012 - P. Kasibhatla - Initial version
27 Aug 2012 - R. Yantosca   - Added parallel DO loops
27 Aug 2012 - R. Yantosca   - Change REAL*4 variables to REAL*8 to better
                               ensure numerical stability
```

1.12.5 read_input_grid

Routine to read variables and attributes from a netCDF file. This routine was automatically generated by the Perl script NcdfUtilities/perl/ncCodeRead.

INTERFACE:

```
SUBROUTINE READ_INPUT_GRID( IM, JM, fileName, lon_edges, lat_sines )
```

USES:

```
! Modules for netCDF read
USE m_netcdf_io_open
USE m_netcdf_io_get_dimlen
USE m_netcdf_io_read
USE m_netcdf_io_readattr
USE m_netcdf_io_close
```

```
IMPLICIT NONE
```

```
# include "netcdf.inc"
```

INPUT PARAMETERS:

```
INTEGER,          INTENT(IN)  :: IM           ! # of longitudes
INTEGER,          INTENT(IN)  :: JM           ! # of latitudes
CHARACTER(LEN=*), INTENT(IN)  :: fileName     ! File w/ grid info
```

OUTPUT PARAMETERS:

```
REAL*8,          INTENT(OUT) :: lon_edges(IM+1) ! Lon edges [degrees]
REAL*8,          INTENT(OUT) :: lat_sines(JM+1) ! SIN( latitude edges )
```

REMARKS:

Created with the ncCodeRead script of the NcdfUtilities package,
with subsequent hand-editing.

REVISION HISTORY:

23 Aug 2012 - R. Yantosca - Initial version

1.13 Fortran: Module Interface grid_mod

Module GRID_MOD contains variables and routines which are used to specify the parameters of a GEOS-Chem horizontal grid. Grid parameters are computed as 3D arrays, which are required for interfacing with a GCM.

INTERFACE:

MODULE Grid_Mod

USES:

USE CMN_GCTM_Mod ! Physical constants
USE Error_Mod ! Error-handling routines

IMPLICIT NONE
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_yEdge
PUBLIC :: Get_yEdge_r
PUBLIC :: Get_yMid
PUBLIC :: Get_yMid_r
PUBLIC :: Get_yMid_r_w
PUBLIC :: Get_ySin
PUBLIC :: Get_xOffset
PUBLIC :: Get_yOffset
PUBLIC :: Init_Grid
PUBLIC :: Its_A_Nested_Grid
PUBLIC :: Set_xOffset
PUBLIC :: Set_yOffset

```

23 Feb 2012 - R. Yantosca - Initial version, based on grid_mod.F
01 Mar 2012 - R. Yantosca - Validated for nested grids
03 Apr 2012 - M. Payer      - Added ySin for map_a2a regrid (M. Cooper)
04 Dec 2012 - R. Yantosca - Modified for GIGC running in ESMF environment
26 Feb 2013 - R. Yantosca - Fixed bug in computation of lons & lats when
                           connecting GEOS-Chem to the GEOS-5 GCM
20 Aug 2013 - R. Yantosca - Removed "define.h", this is now obsolete

```

Subroutine COMPUTE_GRID initializes the longitude, latitude and surface area arrays.

```
SUBROUTINE Compute_Grid( am_I_Root,                                &
                        I1, I2, J1,   J2,   JSP,  JNP,            &
                        L1, L2, DLON, DLAT, I_LO, J_LO, RC )
```

USE GIGC_ErrCode_Mod

```

LOGICAL, INTENT(IN)    :: am_I_Root                ! Root CPU?

! Variables with local CPU indices

INTEGER, INTENT(IN)    :: I1, I2                  ! Min lon index
INTEGER, INTENT(IN)    :: J1, J2                  ! Local lat indices
INTEGER, INTENT(IN)    :: JSP, JNP                ! Polar lat indices
INTEGER, INTENT(IN)    :: L1, L2                  ! Local lev indices
REAL*8,  INTENT(IN)    :: DLON(I2-I1+1,J2-J1+1,L2-L1+1) ! Delta lon [deg]
REAL*8,  INTENT(IN)    :: DLAT(I2-I1+1,J2-J1+1,L2-L1+1) ! Delta lat [deg]

! Variables with global CPU indices

INTEGER, INTENT(IN)    :: I_LO                    ! Min global lon
INTEGER, INTENT(IN)    :: J_LO                    ! Min global lat

```

```
INTEGER, INTENT(OUT) :: RC                                ! Success or failure?
```

- (1) Lon/lat loop indices IG, JG are global indices.
- (2) Lon/lat loop indices I, J are local to each CPU.
- (3) We do not need to have global loop indices for vertical levels, because we will always decompose the grid for MPI parallelization in longitude and/or latitude. All vertical levels must be present on each CPU for the grid-independent GEOS-Chem to function properly.

23 Feb 2012	- R. Yantosca	- Initial version, based on grid_mod.F
30 Jul 2012	- R. Yantosca	- Now accept am_I_Root as an argument when running with the traditional driver main.F
03 Dec 2012	- R. Yantosca	- Add RC to argument list
04 Dec 2012	- R. Yantosca	- Now define arrays with local CPU lon/lat bounds
07 Dec 2012	- R. Yantosca	- Bug fix: make sure the last longitude edge is computed properly. Test for IG==I2, not I==I2.
07 Dec 2012	- R. Yantosca	- Also do not apply half-polar boxes when running in ESMF environment
26 Feb 2013	- R. Yantosca	- Bug fix: now compute IND_X and IND_Y properly when connecting GEOS-Chem to the GEOS-5 GCM
21 Mar 2013	- R. Yantosca	- Add fix to prevent zero surface area at poles
21 Mar 2013	- R. Yantosca	- Rename loop indices to prevent confusion
06 Jun 2013	- M. Payer	- Add fix to compute sine of last latitude edge for MAP_A2A regridding (C. Keller)
02 Jul 2013	- R. Yantosca	- Now compute lon centers properly for GCAP, which does not have any half-sized polar boxes

Function SET_XOFFSET initializes the nested-grid longitude offset variable I0.

SUBROUTINE Set_xOffSet(X_OFFSET)

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

24 Feb 2012 - R. Yantosca - Initial version

Function SET_YOFFSET initializes the nested-grid latitude offset variable J0.

SUBROUTINE Set_yOffset(Y_OFFSET)

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

24 Feb 2012 - R. Yantosca - Initial version

1.13.4 get_xoffset

Function GET_XOFFSET returns the nested-grid longitude offset to the calling program.

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:

```
24 Feb 2012 - R. Yantosca - Initial version
```

1.13.5 get_yoffset

Function GET_YOFFSET returns the nested-grid longitude offset to the calling program.

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:

```
24 Feb 2012 - R. Yantosca - Initial version
```

1.13.6 get_xmid

Function GET_XMID returns the longitude in degrees at the center of a GEOS-Chem grid box.

INTERFACE:

```
FUNCTION Get_xMid( I, J, L ) RESULT( X )
```

INPUT PARAMETERS:

```

      INTEGER, INTENT(IN) :: I    ! Longitude index
      INTEGER, INTENT(IN) :: J    ! Latitude index
      INTEGER, INTENT(IN) :: L    ! Level index

```

RETURN VALUE:

```

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

```

REVISION HISTORY:

24 Feb 2012 - R. Yantosca - Initial version

1.13.7 get_xedge

Function GET_XEDGE returns the longitude in degrees at the western edge of a GEOS-Chem grid box.

INTERFACE:

```

      FUNCTION Get_xEdge( I, J, L ) RESULT( X )

```

INPUT PARAMETERS:

```

      INTEGER, INTENT(IN) :: I    ! Longitude index
      INTEGER, INTENT(IN) :: J    ! Latitude index
      INTEGER, INTENT(IN) :: L    ! Level index

```

RETURN VALUE:

```

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

```

REVISION HISTORY:

24 Feb 2012 - R. Yantosca - Initial version

1.13.8 get_ymid

Function GET_YMID returns the latitude in degrees at the center of a GEOS-Chem grid box.

INTERFACE:

```

      FUNCTION Get_yMid( I, J, L ) RESULT( Y )

```

INPUT PARAMETERS:

```

      INTEGER, INTENT(IN) :: I    ! Longitude index
      INTEGER, INTENT(IN) :: J    ! Latitude index
      INTEGER, INTENT(IN) :: L    ! Level index

```

RETURN VALUE:

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

REVISION HISTORY:

24 Feb 2012 - R. Yantosca - Initial version

1.13.9 get_yedge

Function GET_YEDGE returns the latitude in degrees at the southern edge of a GEOS-Chem grid box.

INTERFACE:

```
FUNCTION Get_yEdge( I, J, L ) RESULT( Y )
```

INPUT PARAMETERS:

```
INTEGER, INTENT(IN) :: I    ! Longitude index
INTEGER, INTENT(IN) :: J    ! Latitude index
INTEGER, INTENT(IN) :: L    ! Level index
```

RETURN VALUE:

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

REVISION HISTORY:

24 Feb 2012 - R. Yantosca - Initial version

1.13.10 get_ymid_r

Function GET_YMID_R returns the latitude in radians at the center of a GEOS-Chem grid box.

INTERFACE:

```
FUNCTION Get_yMid_R( I, J, L ) RESULT( Y )
```

INPUT PARAMETERS:

```
INTEGER, INTENT(IN) :: I    ! Longitude index
INTEGER, INTENT(IN) :: J    ! Latitude index
INTEGER, INTENT(IN) :: L    ! Level index
```

RETURN VALUE:

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

REVISION HISTORY:

24 Feb 2012 - R. Yantosca - Initial version

1.13.11 get_ymid_r_w

Function GET_YMID3_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( I, J, L ) RESULT( Y )
```

INPUT PARAMETERS:

```
INTEGER, INTENT(IN) :: I    ! Longitude index
INTEGER, INTENT(IN) :: J    ! Latitude index
INTEGER, INTENT(IN) :: L    ! Level index
```

RETURN VALUE:

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

REVISION HISTORY:

```
24 Feb 2012 - R. Yantosca - Initial version
01 Mar 2012 - R. Yantosca - Bracket with #ifdef for nested grids
26 Sep 2013 - R. Yantosca - Removed SEAC4RS C-preprocessor switch
```

1.13.12 get_yedge_r

Function GET_YEDGE_R returns the latitude in radians at the southern edge of a GEOS-Chem grid box.

INTERFACE:

```
FUNCTION Get_yEdge_R( I, J, L ) RESULT( Y )
```

INPUT PARAMETERS:

```
INTEGER, INTENT(IN) :: I    ! Longitude index
INTEGER, INTENT(IN) :: J    ! Latitude index
INTEGER, INTENT(IN) :: L    ! Level index
```

RETURN VALUE:

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

REVISION HISTORY:

```
24 Feb 2012 - R. Yantosca - Initial version
```

1.13.13 get_ysin

Function GET_YSIN returns the sine of the southern edge of a GEOS-Chem grid box.

INTERFACE:

```
FUNCTION Get_ySin( I, J, L ) RESULT( Y )
```

INPUT PARAMETERS:

```
INTEGER, INTENT(IN) :: I    ! Longitude index
INTEGER, INTENT(IN) :: J    ! Latitude index
INTEGER, INTENT(IN) :: L    ! Level index
```

RETURN VALUE:

```
REAL*8                :: Y    ! Sine of Latitude value @ S edge of grid box
```

REVISION HISTORY:

```
03 Apr 2012 - M. Payer    - Initial version (M. Cooper)
```

1.13.14 get_area_m2

Function GET_AREA_M2 returns the surface area [m2] of a GEOS-Chem grid box.

INTERFACE:

```
FUNCTION Get_Area_m2( I, J, L ) RESULT( A )
```

INPUT PARAMETERS:

```
INTEGER, INTENT(IN) :: I    ! Longitude index
INTEGER, INTENT(IN) :: J    ! Latitude index
INTEGER, INTENT(IN) :: L    ! Level index
```

RETURN VALUE:

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

REVISION HISTORY:

```
24 Feb 2012 - R. Yantosca - Initial version
```

1.13.15 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( I, J, L ) RESULT( A )
```

INPUT PARAMETERS:

```
INTEGER, INTENT(IN) :: I    ! Longitude index
INTEGER, INTENT(IN) :: J    ! Latitude index
INTEGER, INTENT(IN) :: L    ! Level index
```

RETURN VALUE:

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

REVISION HISTORY:

24 Feb 2012 - R. Yantosca - Initial version

1.13.16 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( I1, I2, J1, J2, L, COORDS, INDICES )
```

INPUT PARAMETERS:

```
INTEGER, INTENT(IN) :: I1, I2    ! Lon indices
INTEGER, INTENT(IN) :: J1, J2    ! Lat indices
INTEGER, INTENT(IN) :: L
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/)
```

REMARKS:

For now, this only works with the surface layer (which is OK since this routine is mostly just called to find a window for surface emissions)

REVISION HISTORY:

24 Feb 2012 - R. Yantosca - Initial version
 01 Mar 2012 - R. Yantosca - Modified for grids, added input parameters

1.13.17 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:

```
24 Feb 2012 - R. Yantosca - Initial version
```

1.13.18 init_grid

Subroutine INIT_GRID initializes variables and allocates module arrays.

INTERFACE:

```
SUBROUTINE Init_Grid( am_I_Root, IM, JM, LM, RC )
```

USES:

```
USE GIGC_ErrCode_Mod
```

INPUT PARAMETERS:

```
LOGICAL, INTENT(IN)  :: am_I_Root    ! Are we on the root CPU
INTEGER, INTENT(IN)  :: IM            ! # of longitudes on this CPU
INTEGER, INTENT(IN)  :: JM            ! # of latitudes  on this CPU
INTEGER, INTENT(IN)  :: LM            ! # of levels    on this CPU
```

OUTPUT PARAMETERS:

```
INTEGER, INTENT(OUT) :: RC            ! Success or failure?
```

REVISION HISTORY:

```
24 Feb 2012 - R. Yantosca - Initial version, based on grid_mod.F
01 Mar 2012 - R. Yantosca - Now define IS_NESTED based on Cpp flags
03 Dec 2012 - R. Yantosca - Add am_I_Root, RC to argument list
04 Dec 2012 - R. Yantosca - Now dimension arrays with IM, JM, LM instead
                           of I1, J1, L1, I2, J2, L2.
18 Apr 2013 - R. Yantosca - Bug fix: in nested block, use STAT=RC
26 Sep 2013 - R. Yantosca - Removed SEAC4RS C-preprocessor switch
```

1.13.19 cleanup_grid

Subroutine CLEANUP_GRID deallocates all module arrays.

INTERFACE:

```
SUBROUTINE Cleanup_Grid
```

REVISION HISTORY:

24 Feb 2012 - R. Yantosca - Initial version, based on grid_mod.F

1.14 Fortran: Module Interface inquireMod

Module inquireMod contains functions to find free and unopened logical file units (LUNs) for Fortran I/O. **INTERFACE:**

```
MODULE inquireMod
```

USES:

```
#if defined( ESMF_ )
  ! We only need to refer to these modules if we are connecting
  ! to the GEOS-5 GCM via the ESMF/MAPL framework (bmy, 8/3/12)
  USE ESMF_Mod
  USE MAPL_Mod
#endif
```

```
IMPLICIT NONE
PRIVATE
```

PUBLIC MEMBER FUNCTIONS:

```
PUBLIC  :: findFreeLUN
PUBLIC  :: I_Am_UnOPENed
```

REVISION HISTORY:

14 Jun 2012 - E. Nielsen - Initial version
03 Aug 2012 - R. Yantosca - Block off ESMF-specific code with #ifdefs
03 Aug 2012 - R. Yantosca - Cosmetic changes

1.14.1 findFreeLUN

Inquire for an existing, but unopened, logical unit number

INTERFACE:

```
FUNCTION findFreeLUN( b ) RESULT( lun )
```

USES:

```
IMPLICIT NONE
```

INPUT PARAMETERS:

```
INTEGER, INTENT(IN), OPTIONAL :: b    ! Not really used here
```

RETURN VALUE:

```
INTEGER :: lun
```

REVISION HISTORY:

```
14 Jun 2012 - E. Nielsen - Initial version
03 Aug 2012 - R. Yantosca - Block off ESMF-specific code with #ifdefs
03 Aug 2012 - R. Yantosca - Cosmetic changes
06 Aug 2012 - R. Yantosca - Now make LUN range 11..199
```

1.14.2 I_Am_UnOPENed

Inquire as to the availability of a given logical unit

INTERFACE:

```
FUNCTION I_Am_UnOPENed( n ) RESULT( TorF )
```

USES:

```
IMPLICIT NONE
```

INPUT PARAMETERS:

```
INTEGER :: n          ! Logical unit # to test
```

RETURN VALUE:

```
LOGICAL :: TorF    ! .TRUE. means the file is unopened
```

REVISION HISTORY:

```
14 Jun 2012 - E. Nielsen - Initial version
03 Aug 2012 - R. Yantosca - Block off ESMF-specific code with #ifdefs
03 Aug 2012 - R. Yantosca - Cosmetic changes
```

1.15 Fortran: Module Interface *global_grid_mod*

Module GLOBAL_GRID_MOD contains variables and routines which are used to specify the parameters of a GEOS-Chem global horizontal grid.

INTERFACE:

```
MODULE Global_Grid_Mod
```

USES:

```
USE CMN_SIZE_Mod           ! Size parameters
USE CMN_GCTM_Mod           ! Physical constants
USE Error_Mod              ! Error-handling routines
```

```
IMPLICIT NONE
PRIVATE
```

PUBLIC MEMBER FUNCTIONS:

```
PUBLIC  :: Cleanup_Global_Grid
PUBLIC  :: Compute_Global_Grid
PUBLIC  :: Get_xEdge_G
PUBLIC  :: Get_yEdge_G
PUBLIC  :: Get_IIIIPAR
PUBLIC  :: Get_JJJP
```

REVISION HISTORY:

```
01 May 2012 - M. Payer      - Initial version, based on grid_mod.F
20 Aug 2013 - R. Yantosca - Removed "define.h", this is now obsolete
```

1.15.1 compute_global_grid

Subroutine COMPUTE_GLOBAL_GRID initializes the longitude, latitude and surface area arrays.

INTERFACE:

```
SUBROUTINE Compute_Global_Grid( am_I_Root )
```

USES:

```
USE GRID_MOD,      ONLY : GET_XOFFSET, GET_YOFFSET
```

INPUT PARAMETERS:

```
LOGICAL, INTENT(IN) :: am_I_Root  ! Is this the root CPU?
```

REVISION HISTORY:

```
01 May 2012 - M. Payer      - Initial version, based on grid_mod.F
30 Jul 2012 - R. Yantosca - Now accept am_I_Root as an argument when
                           running with the traditional driver main.F
```

1.15.2 `get_iiipar`

Function GET_IIIPAR returns IIIPAR.

INTERFACE:

```
FUNCTION Get_IIIPAR() RESULT( I )
```

RETURN VALUE:

```
INTEGER                :: I ! IIIPAR
```

REVISION HISTORY:

01 Mar 2012 - P. Kasibhatla - Initial version

1.15.3 `get_jjipar`

Function GET_JJIPAR returns JJIPAR.

INTERFACE:

```
FUNCTION Get_JJIPAR() RESULT( J )
```

RETURN VALUE:

```
INTEGER                :: J ! JJIPAR
```

REVISION HISTORY:

01 Mar 2012 - P. Kasibhatla - Initial version

1.15.4 `get_xedge_g`

Function GET_XEDGE_G returns the longitude in degrees at the western edge of a GEOS-Chem grid box

INTERFACE:

```
FUNCTION Get_xEdge_G( 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:

01 Mar 2012 - P. Kasibhatla - Initial version

1.15.5 `get_yedge_g`

Function `GET_YEDGE_G` returns the latitude in degrees at the southern edge of a GEOS-Chem grid box.

INTERFACE:

```
FUNCTION Get_Yedge_G( 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:

```
01 Mar 2012 - P. Kasibhatla - Initial version
```

1.15.6 `init_global_grid`

Subroutine `INIT_GLOBAL_GRID` initializes variables and allocates module arrays.

INTERFACE:

```
SUBROUTINE Init_Global_Grid
```

REVISION HISTORY:

```
01 May 2012 - M. Payer - Initial version, based on grid_mod.F
```

1.15.7 `cleanup_global_grid`

Subroutine `CLEANUP_GLOBAL_GRID` deallocates all module arrays.

INTERFACE:

```
SUBROUTINE Cleanup_Global_Grid
```

REVISION HISTORY:

```
01 May 2012 - M. Payer - Initial version, based on grid_mod.F
```