

Chapter 33

Description of program variables

33.1 Currents

```
MODULE currents
REAL, DIMENSION(1-nhalo:ncloc+nhalo,1-nhalo:nrloc+nhalo) :: udvel, uvdvel_old,&
& vdvel, vdvel_old
REAL, DIMENSION(1-nhalo:ncloc+nhalo,0:nrloc+1) :: umvel, umvel_old
REAL, DIMENSION(0:ncloc+1,1-nhalo:nrloc+nhalo) :: vmvel, vmvel_old
REAL, DIMENSION(0:ncloc+1,nrloc) :: umpred
REAL, DIMENSION(ncloc,0:nrloc+1) :: vmpred
REAL, DIMENSION(ncloc+1,nrloc) :: udfvel
REAL, DIMENSION(ncloc,nrloc+1) :: vdfvel
REAL, DIMENSION(ncloc,nrloc) :: p2dbcgradatu, p2dbcgradatv, udevint, vdevint
REAL, DIMENSION(1-nhalo:ncloc+nhalo,1-nhalo:nrloc+nhalo,nz) :: uvel, uvel_old,&
& vvel, vvel_old

REAL, DIMENSION(2-nhalo:ncloc+nhalo,nrloc,nz) :: ufvel
REAL, DIMENSION(ncloc,2-nhalo:nrloc+nhalo,nz) :: vfvel
REAL, DIMENSION(0:ncloc,0:nrloc,nz+1) :: wvel
REAL, DIMENSION(ncloc,nrloc,nz) :: p3dbcgradatu, p3dbcgradatv, wphys
```

File

currents.f90

Type

Module

Purpose

Velocity arrays

Description

p2dbcgradatu	X-component of the depth-integrated baroclinic pressure gradient $\overline{F_1^b}$	[m ² /s ²]
p2dbcgradatv	Y-component of the depth-integrated baroclinic pressure gradient $\overline{F_2^b}$	[m ² /s ²]
p3dbcgradatu	X-component of the baroclinic pressure gradient F_1^b	[m/s ²]
p3dbcgradatv	Y-component of the baroclinic pressure gradient F_2^b	[m/s ²]
udevint	Depth-integrated baroclinic terms $\overline{\delta D_{h1}} - \overline{\delta A_{h1}}$ in the U -equation	[m ² /s ²]
udfvel	X-component of the filtered depth-integrated current U_f	[m ² /s]
udvel	X-component of the depth-integrated current U	[m ² /s]
udvel_old	X-component of the depth-integrated current U at the previous 2-D time step	[m ² /s]
ufvel	X-component of the filtered 3-D current u_f	[m/s]
umpred	X-component of the depth-mean current $\overline{u^p}$ at the predictor step	[m/s]
umvel	X-component of the depth-mean current \overline{u}	[m/s]
umvel_old	X-component of the depth-mean current \overline{u} at the old barotropic (mode-splitting scheme) or baroclinic (implicit scheme) time	[m/s]
uvel	X-component of the 3-D current u	[m/s]
uvel_old	X-component of the 3-D current u at the previous corrector (baroclinic) time step	[m/s]
vdevint	Depth-integrated baroclinic terms $\overline{\delta D_{h2}} - \overline{\delta A_{h2}}$ in the V -equation	[m ² /s ²]
vdfvel	Y-component of the filtered depth-integrated current V_f	[m ² /s]
vdvel	Y-component of the depth-integrated current V	[m ² /s]
vdvel_old	Y-component of the depth-integrated current V at the previous 2-D time step	[m ² /s]
vfvel	Y-component of the filtered 3-D current v_f	[m/s]
vmpred	Y-component of the depth-mean current $\overline{v^p}$ at the predictor step	[m/s]
vmvel	Y-component of the depth-mean current \overline{v}	[m/s]

<code>vmvel_old</code>	Y-component of the depth-mean current \bar{v} at the old barotropic (mode-splitting scheme) or baroclinic (implicit scheme) time	[m/s]
<code>vvel</code>	Y-component of the 3-D current v	[m/s]
<code>vvel_old</code>	Y-component of the 3-D current v at the previous corrector (baroclinic) time step	[m/s]
<code>wphys</code>	Physical vertical velocity w	[m/s]
<code>wvel</code>	Transformed vertical velocity ω	[m/s]

33.2 Derived type definitions

MODULE `datatypes`

ExchComms

```

TYPE :: ExchComms
  LOGICAL :: sfirst
  INTEGER :: iddest, idsrce, tag
  INTEGER, DIMENSION(2) :: i1rcv, i2rcv, j1rcv, j2rcv
  INTEGER, DIMENSION(2) :: i1snd, i2snd, j1snd, j2snd
END TYPE ExchComms

```

File

`datatypes.f90`

Type

Derived type definition

Purpose

Parameters for exchange (send and receive) communications in parallel mode

Description

`iddest` Destination process in case of a send operation

`idsrce` Source process in case of a receive operation

`i1rcv` Start (local) grid index in the X-direction for the receive operation

`i1snd` Start (local) grid index in the X-direction for the send operation

`i2rcv` End (local) grid index in the X-direction for the receive operation

i2snd End (local) grid index in the X-direction for the send operation
j1rcv Start (local) grid index in the Y-direction for the receive operation
j1snd Start (local) grid index in the Y-direction for the send operation
j2rcv End (local) grid index in the Y-direction for the receive operation
j2snd End (local) grid index in the Y-direction for the send operation
sfirst Switch which determines for a local domain whether a send operation is performed before a receive operation (see Section 11.4.3.2)
tag Communication tag

FileParams

(MODULE datatypes)

TYPE :: FileParams

LOGICAL :: defined, info, opened, time_regular

CHARACTER (LEN=1) :: form, status

CHARACTER (LEN=10) :: coherens_version

CHARACTER (LEN=leniofile) :: filename, pathname

CHARACTER (LEN=lendesc) :: filedesc

CHARACTER (LEN=lentime-4) :: creation_date

INTEGER :: endfile, header_type, iostat, iunit, lenrec, maxrecs, &
 & nocoords, nodim, novars, timeid, timerec, tskips, &
 & varid, zetaid

INTEGER, DIMENSION(3) :: tlims

END TYPE FileParams

File

datatypes.f90

Type

Derived type definition

Purpose

Global file attributes

Description

coherens_version Current COHERENS version

creation_date Date and time of creation

defined .TRUE. if the file is activated (*status* is different from '0')

<code>endfile</code>	Switch to select which action needs to be taken in case an end of file condition occurs during input 0: The program aborts with an error message. 1: The program continues. No further read attempt is made. 2: The program waits until the data are available. If the data are not available after a given period of time, the program aborts.
<code>filedesc</code>	A character string with a description of the file contents
<code>filename</code>	File name
<code>form</code>	File format 'A' ASCII 'U' sequential unformatted binary 'N' netCDF 'D' direct access unformatted binary (currently not implemented)
<code>header_type</code>	Type of file header 0: No header (not allowed for netCDF files) 1: Indicates that the file contains forcing data. 2: Indicates that the file contains user-defined output data. 3: Indicates that the file contains user-defined output with output grid data only.
<code>info</code>	Header information is written to a separate information file if <code>.TRUE</code> .
<code>iostat</code>	File I/O status -1: unable to open file 0 : The file is not opened 1 : The file pointer is located before the end of the file 2 : The file pointer is located at the end of the file. An end of file condition did not yet occurred. 3 : The file pointer is located at the end of the file. An end of file condition did occur.

<code>iunit</code>	File unit number
<code>lenrec</code>	Record length (direct access files only)
<code>maxrecs</code>	Total amount of time records in case of a time series file
<code>nocoords</code>	Number of coordinate variables
<code>nodim</code>	Rank of the data variables (user output only)
<code>novars</code>	Number of data variables
<code>opened</code>	.TRUE. (.FALSE.) if the file is connected (not connected)
<code>pathname</code>	Preappended file path
<code>status</code>	File status '0' undefined, i.e. the file is not activated 'N' the file is available in a user-defined (non-COHERENS) format for input 'R' the file is available in COHERENS standard format for input 'W' the file is created for output in COHERENS standard format 'P' as 'W' but the data are appended to an already existing standard COHERENS file. The option is currently not implemented. 'T' temporary file (for internal use only). The option is currently not used.
<code>timeid</code>	NetCDF variable id of the time coordinate variable
<code>timerec</code>	Current time record number
<code>time_regular</code>	.TRUE. if the file contains time series data at regular time intervals.
<code>tlims</code>	Start/End/Step time indices used in updating the data from a forcing file
<code>tskips</code>	Currently not implemented
<code>varid</code>	ID of a variable in the file (only used internally)
<code>zetaid</code>	NetCDF variable file ID of the surface elevation coordinate variable

GridParams

```
(MODULE datatypes)
TYPE :: GridParams
  LOGICAL :: rotated
  INTEGER :: nhtype, n1dat, n2dat
  REAL :: delxdat, delydat, gridangle, longpole, x0dat, y0dat, x0rot, y0rot
END TYPE GridParams
```

File

datatypes.f90

Type

Derived type definition

Purpose

Attributes of surface grids

Description

delxdat	Spacing in the X-direction in case of a uniform rectangular grid [meters or degrees longitude]
delydat	Spacing in the Y-direction in case of a uniform rectangular grid [meters or degrees latitude]
gridangle	Rotation angle of a rotated model grid with respect to the reference grid
longpole	Longitude of the new North pole in case of a rotated spherical rectangular grid [degrees longitude]
nhtype	Type of data grid 0: single point data grid or data grid not defined 1: rectangular grid with uniform grid spacings in Cartesian or spherical coordinates (data grid is constructed from x0dat, y0dat, delxdat, delydat) 2: rectangular grid with non-uniform spacings 3: non-rectangular (curvilinear or non-structured) grid 4: data grid coincides with the model grid
n1dat	Number of grid points in the X-direction or, in case of a non-rectangular grid, the total number of grid points
n2dat	Number of grid points in the Y-direction or equal to one in case of a non-rectangular grid

<code>rotated</code>	<code>.TRUE.</code> in case of a rotated rectangular model grid, <code>.FALSE.</code> otherwise.
<code>x0dat</code>	X-coordinate of the lower left (Southwest) corner in case of a regular grid [meters or longitude]
<code>x0rot</code>	new (rotated) X-coordinate of the lower left (Southwest) corner in case of a rotated spherical grid [degrees longitude]
<code>y0dat</code>	Y-coordinate of the lower left (Southwest) corner in case of a regular grid [meters or latitude]
<code>y0rot</code>	new (rotated) Y-coordinate of the lower left (Southwest) corner in case of a rotated spherical grid [fractional degrees latitude]

HRelativeCoords

```
(MODULE datatypes)
TYPE :: HRelativeCoords
  INTEGER :: icoord, jcoord
  REAL   :: xcoord, ycoord
END TYPE HRelativeCoords
```

File

datatypes.f90

Type

Derived type definition

Purpose

Horizontal relative coordinates

Reference

Section 10.4.1

Description

- `icoord` Integer part of the ξ_1 curvilinear coordinate with respect to the reference coordinate system
- `jcoord` Integer part of the ξ_2 curvilinear coordinate with respect to the reference coordinate system
- `xcoord` Fractional part of the ξ_1 curvilinear coordinate with respect to the reference coordinate system

ycoord Fractional part of the ξ_2 curvilinear coordinate with respect to the reference coordinate system

OutGridParams

```
(MODULE datatypes)
TYPE :: OutGridParams
  LOGICAL :: gridded, grid_file, land_mask, time_grid
  CHARACTER (LEN=1) :: status
  CHARACTER (LEN=lentime) :: enddate, refdate, startdate
  INTEGER :: ncout, nodim, nostats, nowetout, nrout, nstepout, nzout, &
    & time_format
  INTEGER, DIMENSION(3) :: tlims, xlims, ylims, zlims
  REAL :: deltout
END TYPE OutGridParams
```

File

datatypes.f90

Type

Derived type definition

Purpose

Parameters for user-defined output and attributes of output data grids

Description

deltout	Output time step. Unit is seconds in case of an absolute time. Otherwise, the unit is determined by the value of <code>time_format</code> .
enddate	Date and time of the last output
gridded	.TRUE. (.FALSE.) for gridded (non-gridded) output
grid_file	The coordinates of the output grid are written to a separate external file if .TRUE.
land_mask	The data are written with a land mask (i.e. without land points) if .TRUE.
ncout	X-dimension of the output grid in case of gridded data
nodim	Spatial dimension of the output grid (0,2,3)
nostats	Number of output stations in case of non-gridded data
nowetout	Number of wet output data in case a land mask is applied

<code>nROUT</code>	Y-dimension of the output grid in case of gridded data
<code>nSTEPOUT</code>	Number of output time steps
<code>nZOUT</code>	Vertical dimension of the output grid
<code>REFDATE</code>	Reference date/time used when the time coordinate is written in a numerical format in which case it is defined as the time elapsed since this reference date.
<code>STARTDATE</code>	Date and time of the first output
<code>STATUS</code>	Status of the corresponding output files 'W' data are written to a newly created file 'P' data are appended to an existing file. This option is currently not yet available.
<code>TIME_FORMAT</code>	Format of the time coordinate. The first and last case are absolute times, the other ones relative times with respect to the <code>REFDATE</code> 0: date/time in string format 1: seconds 2: minutes 3: hours 4: days 5: months (one month is taken as 30.4375 days) 6: years (one year is taken as 365.25 days) 7: date in years
<code>TIME_GRID</code>	Create a time-dependent grid if <code>.TRUE.</code> and <code>nodim = 3</code> .
<code>TLIMITS</code>	Start/end/step output time index. The step value determines the period in case of time-averaged or harmonic output.
<code>XLIMITS</code>	Start/end/step output space index in the X-direction in case of gridded data
<code>YLIMITS</code>	Start/end/step output space index in the Y-direction in case of gridded data
<code>ZLIMITS</code>	Start/end/step output space index in the vertical direction

StationLocs

(MODULE `datatypes`)

```
TYPE :: StationLocs
  INTEGER :: ipos, jpos
  CHARACTER (LEN=lenname) :: name
END TYPE StationLocs
```

File

datatypes.f90

Type

Derived type definition

Purpose

Attributes of the output station locations in case of non-gridded output.
The stations are assumed to be located at C-nodes.

Description

ipos C-node X-index of the stations

jpos C-node Y-index of the stations

name Name of the station

VariableAtts

(MODULE datatypes)

```
TYPE :: VariableAtts
```

```
  CHARACTER (LEN=lenname) :: f90_name
```

```
  CHARACTER (LEN=lendesc) :: long_name, vector_name
```

```
  CHARACTER (LEN=lenunit) :: units
```

```
  CHARACTER (LEN=lennode) :: node
```

```
  INTEGER :: data_type, ivarid, klev, nrank, numvar, oopt
```

```
  INTEGER, DIMENSION(4) :: shape
```

```
END TYPE VariableAtts
```

File

datatypes.f90

Type

Derived type definition

Purpose

Variable attributes

Description

f90_name	FORTRAN 90 name
long_name	Long descriptive name
vector_name	Associated vector name in case the variable denotes a vector component
units	Variable unit
node	Nodal location of the variable on the model grid. In case of user output, the nodal location on the output grid ('C' or 'W')
data_type	Data type of the variable (see Table 8.1)
ivarid	Variable key id (a zero means undefined)
klev	Vertical level used when <code>oopt</code> equals <code>oopt_klev</code>
nrank	Variable rank (0 for a scalar)
numvar	Variable number in case the last dimension of the array variable represents a variable dimension
oopt	Selects the operator to be applied in case the variable is selected for user output
oopt_null	No operator applied
oopt_min	Minimum value
oopt_max	Maximum value
oopt_mean	Spatially averaged value
oopt_klev	Value at a specified vertical level
oopt_max	Value at a specified depth below the surface
shape	Array shape (not defined in case of a scalar)
dep	Depth value used when <code>oopt</code> equals <code>oopt_dep</code>

VRelativeCoords

```
(MODULE datatypes)
TYPE :: VRelativeCoords
  INTEGER :: kcoord
  REAL :: zcoord
END TYPE VRelativeCoords
```

File

datatypes.f90

Type

Derived type definition

Purpose

Vertical relative coordinates

Description

`kcoord` Vertical level of the grid point along the reference grid just below the data point

`zcoord` Normalised vertical distance from the vertical level `kcoord` to the location of the data point (between 0 and 1)

33.3 Density

MODULE density

REAL, DIMENSION(0:ncloc+1,0:nrloc+1,nz) :: beta_sal, beta_temp, dens

REAL, DIMENSION(1-nhalo:ncloc+nhalo,1-nhalo:nrloc+nhalo,nz) :: sal, temp

File

density.f90

Type

Module

Purpose

Density arrays

Description

<code>beta_sal</code>	Salinity expansion coefficient β_S	[PSU ⁻¹]
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<code>beta_temp</code>	Temperature expansion coefficient β_T	[°C ⁻¹]
------------------------	---	---------------------

<code>dens</code>	Mass density ρ	[kg/m ³]
-------------------	---------------------	----------------------

<code>sal</code>	Salinity S	[PSU]
------------------	--------------	-------

<code>temp</code>	Temperature T	[°C]
-------------------	-----------------	------

33.4 Water depths

MODULE depths

REAL, DIMENSION(0:ncloc+1,0:nrloc+1) :: depmeanatc, deptotatc

REAL, DIMENSION(ncloc,nrloc) :: deptotatc_old, deptotatc_err

```

REAL, DIMENSION(0:ncloc+1,nrloc) :: depmeanatu, deptotatu_old, deptotatu_prev
REAL, DIMENSION(1-nhalo:ncloc+nhalo,0:nrloc+1) :: deptotatu
REAL, DIMENSION(ncloc,0:nrloc+1) :: depmeanatv, deptotatv_old, deptotatv_prev
REAL, DIMENSION(0:ncloc+1,1-nhalo:nrloc+nhalo) :: deptotatv
REAL, DIMENSION(ncloc+1,nrloc+1) :: depmeanatuv, deptotatuv
REAL, DIMENSION(0:nc+1,0:nr+1) :: depmeanglb
REAL, DIMENSION(0:ncloc+1,0:nrloc+1) :: dzeta, zeta, zeta_old

```

File

depths.f90

Type

Module

Purpose

Water depths and sea surface elevations

Description

depmeanatc	Mean water depth at the C-nodes	[m]
depmeanatu	Mean water depths at the U-nodes	[m]
depmeanatuv	Mean water depth at the UV-nodes	[m]
depmeanatv	Mean water depth at the V-nodes	[m]
depmeanglb	Mean water depth (global) at C- (or -UV-nodes)	[m]
deptotatc	Total water depth at the C-nodes	[m]
deptotatc_err	Total water depth error $\delta_e H$ at the C-nodes	[m]
deptotatc_old	Total water depth at the C-nodes and old baroclinic time step	[m]
deptotatu	Total water depth at the U-nodes	[m]
deptotatuv	Total water depth at the UV-nodes	[m]
deptotatu_old	Total water depth at the U-nodes and old baroclinic time step	[m]
deptotatu_prev	Total water depth at the U-nodes and previous (outer) iteration with the implicit scheme	[m]
deptotatv	Total water depth at the V-nodes	[m]
deptotatv_old	Total water depth at the V-nodes and old baroclinic time step	[m]
deptotatv_prev	Total water depth at the V-nodes and previous (outer) iteration with the implicit scheme	[m]

dzeta	Free surface correction defined as the difference between the surface elevation at the next and previous iteration (implicit scheme only)	[m]
zeta	Surface elevation	[m]
zeta_old	Surface elevation at the start of the first (outer) iteration with the implicit scheme	[m]

33.5 Diffusion

```

MODULE diffusion
REAL, DIMENSION(0:ncloc,0:nrloc) :: hdifcoef2datc
REAL, DIMENSION(ncloc+1,nrloc+1) :: hdifcoef2datuv
REAL, DIMENSION(0:ncloc,0:nrloc,nz) :: hdifcoef3datc
REAL, DIMENSION(ncloc+1,nrloc,nz) :: hdifcoef3datu
REAL, DIMENSION(ncloc,nrloc+1,nz) :: hdifcoef3datv
REAL, DIMENSION(ncloc+1,nrloc+1,nz) :: hdifcoef3datuv
REAL, DIMENSION(0:ncloc,0:nrloc,nz+1) :: kinvisc, vdifcoefmom
REAL, DIMENSION(ncloc,nrloc,nz+1) :: vdifcoefscal, vdifcoefvtk

```

File

diffusion.f90

Type

Module

Purpose

Horizontal and vertical diffusion coefficients

Description

hdifcoef2datc	Vertically integrated horizontal diffusion coefficient $\overline{\nu_H}$ at the C-nodes	[m ³ /s]
hdifcoef2datuv	Vertically integrated horizontal diffusion coefficient $\overline{\nu_H}$ at the UV-nodes	[m ³ /s]
hdifcoef3datc	Horizontal 3-D diffusion coefficient ν_H at the C-nodes	[m ² /s]
hdifcoef3datu	Horizontal 3-D diffusion coefficient λ_H at the U-nodes	[m ² /s]
hdifcoef3datuv	Horizontal 3-D diffusion coefficient ν_H at the UV-nodes	[m ² /s]

hdifcoef3datv	Horizontal 3-D diffusion coefficient λ_H at the V-nodes	[m ² /s]
kinvisc	Kinematic viscosity	[m ² /s]
vdifcoefmom	Vertical diffusion coefficient ν_T for momentum at the W-nodes	[m ² /s]
vdifcoefscal	Vertical diffusion coefficient λ_T for scalars at the W-nodes	[m ² /s]
vdifcoefk	Vertical diffusion coefficient ν_k for turbulent kinetic energy at the W-nodes	[m ² /s]

33.6 Fluxes

MODULE fluxes

```
REAL, DIMENSION(ncloc,nrloc) :: bdragcoefatu, bdragcoefatv, bfricatu, &
    & bfricativ, bstresatc, ces, chs, qlatent, &
    & qlwave, qnonsol, qsensible, ssalflux, &
    & sstresatc, ubstresatu, usstresatu, &
    & vbstresatv, vsstresatv, zeros2d
REAL, DIMENSION(0:ncloc,0:nrloc) :: bdragcoefatc, cds, zroughatc
REAL, DIMENSION(0:ncloc,nrloc) :: usstresatc
REAL, DIMENSION(ncloc,0:nrloc) :: vsstresatc
REAL, DIMENSION(0:nwind,-ntemp:ntemp,nrelhum) :: cdstab, cestab, chstab
```

File

fluxes.f90

Type

Module

Purpose

Bottom and surface flux arrays

Description

bdragcoefatc	Bottom drag coefficient C_{db} at the C-nodes	
bdragcoefatu	Bottom drag coefficient C_{db} at the U-nodes	
bdragcoefatv	Bottom drag coefficient C_{db} at the V-nodes	
bfricatu	Bottom friction velocity $u_{\ast b}$ at the U-nodes	[m/s]
bfricativ	Bottom friction velocity $u_{\ast b}$ at the V-nodes	[m/s]
bstresatc	Magnitude of the bottom stress τ_b at the C-nodes	[m ² /s ²]

cds	Surface drag coefficient C_{ds}	
cdstab	Surface drag coefficient given in tabular form as function of wind speed, air minus sea temperature difference and relative humidity. The table is used for interpolation in case a Monin-Obokhov formulation is selected.	
ces	Exchange coefficient for the latent heat flux C_e (Dalton number)	
cestab	Dalton number given in tabular as function of wind speed, air minus sea temperature difference and relative humidity. The table is used for interpolation in case a Monin-Obokhov formulation is selected.	
chs	Exchange coefficient for the sensible heat flux C_h (Stanton number)	
chstab	Stanton number given in tabular form as function of wind speed, air minus sea temperature difference and relative humidity. The table is used for interpolation in case a Monin-Obokhov formulation is selected.	
qlatent	Latent surface heat flux	[W/m ²]
qlwave	Long-wave surface heat flux	[W/m ²]
qnonsol	Non-solar surface heat flux	[W/m ²]
qsensible	Sensible surface heat flux	[W/m ²]
ssalflux	Surface salinity flux	[PSU m/s]
sstresatc	Magnitude of the surface stress τ_s at the C-nodes	[m ² /s ²]
ubstresatu	X-component of the bottom stress τ_{b1} at the U-nodes	[m ² /s ²]
usstresatc	X-component of the surface stress τ_{s1} at the C-nodes	[m ² /s ²]
usstresatu	X-component of the surface stress τ_{s1} at the U-nodes	[m ² /s ²]
vbstresatv	Y-component of the bottom stress τ_{b2} at the V-nodes	[m ² /s ²]
vstresatc	Y-component of the surface stress τ_{s2} at the C-nodes	[m ² /s ²]
vstresatv	Y-component of the surface stress τ_{s2} at the V-nodes	[m ² /s ²]
zeros2d	Work space array with zero values	

zroughatc Bottom roughness length z_0 at the C-nodes [m]

33.7 Model grid arrays

```

MODULE grid
LOGICAL, DIMENSION(nobu) :: westobu
LOGICAL, DIMENSION(nobv) :: soutobv
LOGICAL, DIMENSION(nobx) :: westobx
LOGICAL, DIMENSION(noby) :: soutoby
LOGICAL, DIMENSION(ncloc,nrloc) :: maskatc_int
LOGICAL, DIMENSION(0:ncloc+1,0:nrloc+1) :: seapoint

INTEGER, DIMENSION(nobu) :: indexobu, iobu, jobu
INTEGER, DIMENSION(nobv) :: indexobv, iobv, jobv
INTEGER, DIMENSION(nobx) :: indexobx, iobx, jobx
INTEGER, DIMENSION(noby) :: indexoby, ioby, joby
INTEGER, DIMENSION(nobuloc_ext) :: iobuloc, jobuloc
INTEGER, DIMENSION(nobvloc_ext) :: iobvloc, jobvloc
INTEGER, DIMENSION(nobxloc_ext) :: iobxloc, jobxloc
INTEGER, DIMENSION(nobyloc_ext) :: iobyloc, jobyloc
INTEGER, DIMENSION(1-nhalo:ncloc+nhalo,1-nhalo:nrloc+nhalo) :: &
& nodeatc, node2du, node2duv, node2dv
INTEGER, DIMENSION(1-nhalo:ncloc+nhalo,1-nhalo:nrloc+nhalo,nz) :: &
& nodeatu, nodeatuv, nodeatv
INTEGER, DIMENSION(1-nhalo:ncloc+nhalo,1-nhalo:nrloc+nhalo,nz+1) :: &
& nodeatuw, nodeatvw
REAL, DIMENSION(ncloc,nrloc) :: alphatc_fld, alphatu_fld, alphatv_fld, &
& gangleatc, &
& rlxobcatu, rlxobcatv
REAL, DIMENSION(ncloc+1,nrloc) :: gaccatu
REAL, DIMENSION(ncloc,nrloc+1) :: gaccatv
REAL, DIMENSION(ncloc+1,0:nrloc) :: coriolatu
REAL, DIMENSION(0:ncloc,nrloc+1) :: coriolatv
REAL, DIMENSION(0:nc+1) :: gdelxglb
REAL, DIMENSION(0:nr+1) :: gdelyglb
REAL, DIMENSION(0:ncloc+1,0:nrloc+1) :: gaccatc, gxcoord, gycoord
REAL, DIMENSION(0:ncloc,0:nrloc) :: gxlon, gylat
REAL, DIMENSION(0:nc+1,0:nr+1) :: gxcoordglb, gycoordglb
REAL, DIMENSION(0:nc+1,0:nr+1,nz+1) :: gscoordglb
REAL, DIMENSION(0:ncloc+1,0:nrloc+1,nz+1) :: gscoord, gscoordatuw, &

```

```

& gscoorlatvw, gscoorlatw
REAL, DIMENSION(0:ncloc+1,0:nrloc+1,2:nz) :: delzatlw
REAL, DIMENSION(0:ncloc+1,0:nrloc+1,nz) :: delzatl, gscoorlatc, gscoorlatu, &
& gscoorlatv
REAL, DIMENSION(nz+1) :: gsigcoord
REAL, DIMENSION(1-nhalo:ncloc+nhalo,1-nhalo:nrloc+nhalo) :: &
& delxatc, delxatu, delxatuv, delxatv, &
& delyatc, delyatu, delyatuv, delyatv
REAL, DIMENSION(0:ncloc+1,nrloc,nz) :: delzatu
REAL, DIMENSION(ncloc,0:nrloc+1,nz) :: delzatl
REAL, DIMENSION(ncloc+1,nrloc+1,nz) :: delzatu
REAL, DIMENSION(ncloc+1,nrloc,2:nz) :: delzatuw
REAL, DIMENSION(ncloc,nrloc+1,2:nz) :: delzatlw

```

File

grid.f90

Type

Module

Purpose

Arrays for the model grid

Description

<code>alphatc fld</code>	α -factor used in the drying/wetting algorithm at the C-nodes	
<code>alphatu fld</code>	α -factor used in the drying/wetting algorithm at the U-nodes	
<code>alphatv fld</code>	α -factor used in the drying/wetting algorithm at the V-nodes	
<code>coriolatu</code>	Coriolis frequency at the U-nodes	[rad/s]
<code>coriolatv</code>	Coriolis frequency at the V-nodes	[rad/s]
<code>delxatc</code>	Grid spacing in the X-direction at the C-nodes	[m]
<code>delxatu</code>	Grid spacing in the X-direction at the U-nodes	[m]
<code>delxatuv</code>	Grid spacing in the X-direction at the UV-nodes	[m]
<code>delxatv</code>	Grid spacing in the X-direction at the V-nodes	[m]
<code>delyatc</code>	Grid spacing in the Y-direction at the C-nodes	[m]
<code>delyatu</code>	Grid spacing in the Y-direction at the U-nodes	[m]
<code>delyatuv</code>	Grid spacing in the Y-direction at the UV-nodes	[m]

delzadc	Vertical grid spacing at the C-nodes	[m]
delzatu	Vertical grid spacing at the U-nodes	[m]
delzatuuv	Vertical grid spacing at the UV-nodes	[m]
delzatuuw	Vertical grid spacing at the UW-nodes	[m]
delzativ	Vertical grid spacing at the V-nodes	[m]
delzativv	Vertical grid spacing at the VW-nodes	[m]
delzatuw	Vertical grid spacing at the W-nodes	[m]
delyatv	Grid spacing in the Y-direction at the V-nodes	[m]
gaccadc	Acceleration of gravity at the C-nodes	[m ² /s]
gaccatu	Acceleration of gravity at the U-nodes	[m ² /s]
gaccativ	Acceleration of gravity at the V-nodes	[m ² /s]
gangleadc	Angle between the reference and local grid X-axis at C-nodes	[degrees]
gdexglb	Grid spacings in the X-direction of a non-uniform rectangular grid (global array)	[m or degrees]
gdelyglb	Grid spacings in the Y-direction of a non-uniform rectangular grid (global array)	[m or degrees]
gscoord	Sigma coordinates of the UVW-nodes	
gscoordadc	Sigma coordinates of the C-nodes	
gscoordatu	Sigma coordinates of the U-nodes	
gscoordatuuv	Sigma coordinates of the UW-nodes	
gscoordativ	Sigma coordinates of the V-nodes	
gscoordativv	Sigma coordinates of the VW-nodes	
gscoordatuw	Sigma coordinates of the W-nodes	
gscoordglb	Sigma coordinates of the W- or UVW-nodes (global array)	
gsigcoord	Sigma coordinates in case the vertical grid is horizontally uniform (iopt_grid_vtype=1)	
gxcoord	X-coordinates of the UV-nodes (local array)	[m or degrees longitude]
gxcoordglb	X-coordinates of the UV-nodes (global array)	[m or degrees longitude]
gxlon	Longitude coordinate at the C-nodes	[rad]
gycoord	Y-coordinates of the UV-nodes (local array)	[m or degrees latitude]

<code>gycoordglb</code>	Y-coordinates of the UV-nodes (global array) [m or degrees latitude]
<code>gylat</code>	Latitude coordinate at the C-nodes [rad]
<code>indexobu</code>	Maps the local open boundary indices at the U-nodes to their corresponding global ones.
<code>indexobv</code>	Maps the local open boundary indices at the V-nodes to their corresponding global ones.
<code>indexobx</code>	Maps the local open boundary indices at the X-nodes to their corresponding global ones.
<code>indexoby</code>	Maps the local open boundary indices at the Y-nodes to their corresponding global ones.
<code>iobu</code>	Global X-index of the West/East U-node open boundaries
<code>iobuloc</code>	Local X-index of the West/East U-node open boundaries
<code>iobv</code>	Global X-index of the South/North V-node open boundaries
<code>iobvloc</code>	Local X-index of the South/North V-node open boundaries
<code>iobx</code>	Global X-index of the West/East X-node open boundaries
<code>iobxloc</code>	Local X-index of the West/East X-node open boundaries
<code>ioby</code>	Global X-index of the South/North Y-node open boundaries
<code>iobyloc</code>	Local X-index of the South/north Y-node open boundaries
<code>jobu</code>	Global Y-index of the West/East U-node open boundaries
<code>jobuloc</code>	Local Y-index of the West/East U-node open boundaries
<code>jobv</code>	Global Y-index of the South/North V-node open boundaries
<code>jobvloc</code>	Local Y-index of the South/North V-node open boundaries
<code>jobx</code>	Global Y-index of the West/East X-node open boundaries
<code>jobxloc</code>	Local Y-index of the West/East X-node open boundaries
<code>joby</code>	Global Y-index of the South/north Y-node open boundaries
<code>jobyloc</code>	Local Y-index of the South/north Y-node open boundaries

<code>maskatc_int</code>	.TRUE./FALSE. at wet/dry C-node grid points
<code>nodeatc</code>	Pointers at C-nodes 0: dry cell 1: wet cell
<code>nodeatu</code>	Pointers at U-node cell faces 0: dry (land) cell face 1: coastal or solid structure boundary 2: interior wet U-node 3: open sea boundary 4: river open boundary
<code>nodeatuv</code>	Pointer at corner (UV) nodes 0: at least two surrounding U-nodes or at least two surrounding V-nodes are dry 1: interior wet node, i.e. at most one surrounding U-node and at most one surrounding V-node is dry and none of the four surrounding velocity nodes are open boundaries 2: X-node open boundary, in which case at least one of the surrounding U-nodes is an open boundary while the other one is either a closed node or open boundary, but the node is not a Y-node open boundary 3: Y-node open boundary, in which case at least one of the surrounding V-nodes is an open boundary while the other one is either a closed node or open boundary, but the node is not an X-node open boundary 4: the node is both a X- and a Y-node open boundary
<code>nodeatuw</code>	Pointer at UW-node cell faces 0: dry (land) cell face or bottom cell (1) or surface cell (nz+1) 1: coastal or solid structure boundary 2: interior wet UW-node 3: open sea boundary 4: river open boundary
<code>nodeatv</code>	Pointers at V-node cell faces

	0: dry (land) cell face
	1: coastal or solid structure boundary
	2: interior wet V-node
	3: open sea boundary
	4: river open boundary
<code>nodeatvw</code>	Pointer at VW-node cell faces
	0: dry (land) cell face or bottom cell (1) or surface cell (nz+1)
	1: coastal or solid structure boundary
	2: interior wet VW-node
	3: open sea boundary
	4: river open boundary
<code>node2du</code>	Pointers at U-nodes for 2-D calculations
	0: dry (land) cell face
	1: coastal boundary
	2: at least one U-node interface in the vertical is wet
	3: open sea boundary
	4: river open boundary
<code>node2dv</code>	Pointers at V-nodes for 2-D calculations
	0: dry (land) cell face
	1: coastal boundary
	2: at least one V-node interface in the vertical is wet
	3: open sea boundary
	4: river open boundary
<code>node2duv</code>	Pointer at corner (UV) nodes for 2-D calculations
	0: at least two surrounding U-nodes or at least two surrounding V-nodes are dry
	1: at least one corner node in the vertical is an interior wet node, i.e. at most one surrounding U-node and at most one surrounding V-node (at the same vertical position) is dry and none of the four surrounding velocity nodes are open boundaries

	2: X-node open boundary, in which case at least one of the surrounding U-nodes is an open boundary while the other one is either a closed node or open boundary, but the node is not a Y-node open boundary
	3: Y-node open boundary, in which case at least one of the surrounding V-nodes is an open boundary while the other one is either a closed node or open boundary, but the node is not an X-node open boundary
	4: the node is both a X- and a Y-node open boundary
rlxobcatu	Relaxation factor for horizontal momentum advection at U-nodes
rlxobcatv	Relaxation factor for horizontal momentum advection at V-nodes
seapoint	.FALSE. for permanent land cells, .TRUE. otherwise
seapoint	.FALSE. for permanent land cells, .FALSE. otherwise
soutobv	.TRUE./FALSE. at South/North V-open boundaries (global array)
soutoby	.TRUE./FALSE. at South/North Y-open boundaries (global array)
westobu	.TRUE./FALSE. at West/East V-open boundaries (global array)
westobx	.TRUE./FALSE. at West/East X-open boundaries (global array)

33.8 Model grid parameters

```

MODULE gridpars
!---grid dimensions
INTEGER :: nc = 0, nr = 0, nz = 0
INTEGER :: ncloc, nc1loc, nc2loc, nrloc, nr1loc, nr2loc
!---number of open boundary points
INTEGER :: nobu, nobv, nobx, noby
INTEGER :: nosbu = 0, nosbv = 0, nrvbv = 0, nrvbv = 0
INTEGER :: nobuloc, nobvloc, nobxloc, nobyloc
INTEGER :: nobuloc_ext, nobvloc_ext, nobxloc_ext, nobyloc_ext
INTEGER :: nosbuloc, nosbvloc, nrvbuloc, nrvbvloc
!---shortcuts for regular grid spacings

```



```

LOGICAL :: dXregX, dXregY, dYregX, dYregY, dXYreg, dZregZ, rotate_gvecs
!---halo dimensions
INTEGER, PARAMETER :: nhalo = 2
INTEGER :: nhdens, nhfvel, nhscal, nhturb, nh2vel, nh3vel
!---number of open boundary points
INTEGER :: noseaatc, noseaatcloc, nowetadc, nowetadcloc
!---number of sediment/biological variables used at open boundaries and for
!   nesting
INTEGER :: maxbiovars = 0, maxsedvars = 0
!---mean grid spacing
REAL :: delgrid

```

File

gridpars.f90

Type

Module

Purpose

Model grid parameters

Description

delgrid	Domain-averaged grid spacing for both X- and Y-direction [m]
dXregX	.TRUE. if the grid spacing in the X-direction is uniform in the X-direction
dXregY	.TRUE. if the grid spacing in the X-direction is uniform in the Y-direction
dXYreg	.TRUE. if all grid spacings are uniform in both X- and Y-directions
dYregX	.TRUE. if the grid spacing in the Y-direction is uniform in the X-direction
dYregY	.TRUE. if the grid spacing in the Y-direction is uniform in the Y-direction
dZregZ	.TRUE. for a uniform vertical vertical grid (<i>iopt_grid_vtype</i> =1)
maxbiovars	Maximum number of biological variables used at open boundaries and for nesting
maxsedvars	Maximum number of sediment fractions used at open boun- daries and for nesting

<code>nc</code>	Number of grid cells in the X-direction on the global grid
<code>ncloc</code>	Number of grid cells in the X-direction on the local grid
<code>nc1loc</code>	Global X-index of cell (1,1) on the local grid
<code>nc2loc</code>	Global X-index of cell (ncloc,1) on the local grid
<code>nhalo</code>	Maximum size of a halo in all directions. The current value of 2 cannot be changed.
<code>nhdens</code>	Minimum required size of the halo needed for density arrays (1/2)
<code>nhfvel</code>	Minimum required size of the halo for horizontal advective currents (1/2)
<code>nhscal</code>	Minimum required size of the halo for scalar (non-density) arrays (1/2)
<code>nhturb</code>	Minimum required size of the halo for the turbulence transport arrays (0/1/2)
<code>nh2vel</code>	Minimum required size of the halo for the 2-D depth-integrated currents (1/2)
<code>nh3vel</code>	Minimum required size of the halo for the 3-D currents (1/2)
<code>nobu</code>	Global number of open boundary points at the U-nodes
<code>nobuloc</code>	Local number of open boundary points at the U-nodes
<code>nobuloc_ext</code>	Local number of interior open boundary points at the U-nodes including points within the first column of the eastern halo
<code>nobv</code>	Global number of open boundary points at the V-nodes
<code>nobvloc</code>	Local number of open boundary points at the V-nodes
<code>nobvloc_ext</code>	Local number of interior open boundary points at the V-nodes including points within the first row of the northern halo
<code>nobx</code>	Global number of open boundary points at the X-nodes
<code>nobxloc</code>	Local number of open boundary points at the X-nodes
<code>nobxloc_ext</code>	Local number of interior open boundary points at the X-nodes including points within the first column of the eastern halo
<code>noby</code>	Global number of open boundary points at the Y-nodes
<code>nobyloc</code>	Local number of open boundary points at the Y-nodes

nobyloc_ext	Local number of interior open boundary points at the Y-nodes including points within the first column of the northern halo
nosbu	Global number of open sea open boundary points at the U-nodes
nosbuloc	Local number of open sea open boundary points at the U-nodes
nosbv	Global number of open sea open boundary points at the V-nodes
nosbvloc	Local number of open sea open boundary points at the V-nodes
noseaatc	Number of sea (wet or dry) C-nodes on the global domain
noseaatcloc	Number of sea (wet or dry) C-nodes on the local domain
nowetatc	Number of active wet C-nodes on the global domain
nowetatcloc	Number of active wet C-nodes on the local domain
nr	Number of grid cells in the Y-direction on the global grid
nrloc	Number of grid cells in the Y-direction on the local grid
nrvbu	Global number of river open boundary points at the U-nodes
nrvbuloc	Local number of river open boundary points at the U-nodes
nrvbv	Global number of river open boundary points at the V-nodes
nrvbvloc	Local number of river open boundary points at the V-nodes
nr1loc	Global Y-index of cell (1,1) on the local grid
nr2loc	Global Y-index of cell (1,nrloc) on the local grid
nz	Number of grid cells in the vertical direction
rotate_gvecs	.TRUE. if vectors in the model grid are rotated with respect to the reference coordinate system

33.9 General and I/O parameters

MODULE iopars

General parameters

```

LOGICAL :: cold_start = .FALSE., next_simul = .FALSE.
INTEGER :: isimul = 0, nopenf = 0, nrecunit = 4, nosimul = 0
CHARACTER (LEN=lentitle) :: intitle, outtitle, runttitle
CHARACTER (LEN=lename), DIMENSION(MaxProgLevels) :: procname
LOGICAL :: log_fill = log_undef
INTEGER :: int_fill = int_undef
REAL :: real_fill = real_undef, real_min = real_flag

```

File

iopars.f90

Type

Module

Purpose

General parameters mainly related to all kinds of input and output

Description

<code>cold_start</code>	If <code>.TRUE.</code> , the program only executes the initialisation and finalisation phases, but does not enter the time loop.
<code>intitle</code>	Title for model forcing files
<code>int_fill</code>	Data flag for invalid integer data
<code>isimul</code>	Number of the simulation (as given by the corresponding line number in the <i>defruns</i> excluding comment lines)
<code>log_fill</code>	Data flag for invalid logical data
<code>next_simul</code>	When a new data line is read from the <i>defruns</i> file, its value is set to <code>.TRUE.</code> to start a new simulation or to <code>.FALSE.</code> to exit the program.
<code>nopenf</code>	Number of current connected files
<code>nosimul</code>	Number of simulations
<code>nrecunit</code>	Unit record size in bytes for direct access I/O (not used in the present implementation)
<code>outtitle</code>	Title for user output files
<code>procname</code>	Name of the subprogram at the current program level
<code>runttitle</code>	Simulation title
<code>real_fill</code>	Data flag for invalid real data

`real_min` Real data are considered as valid or invalid if they are greater than or lower than (or equal to) this critical threshold

Parameters for model forcing

```
(MODULE iopars)
INTEGER, DIMENSION(MaxIOTypes,2) :: maxdatafiles
INTEGER, DIMENSION(MaxCIFTypes) :: ciflinenum
TYPE (FileParams), DIMENSION(MaxCIFTypes) :: ciffiles
TYPE (FileParams), DIMENSION(MaxIOTypes,MaxIOFiles,2) :: modfiles
TYPE (GridParams), DIMENSION(MaxGridTypes,MaxGridFiles) :: surfacegrids
INTEGER, PARAMETER :: io_mppmod = 1, io_inicon = 2, io_fincon = 3, &
    & io_modgrd = 4, io_metgrd = 5, io_sstgrd = 6, &
    & io_wavgrd = 7, io_biogr = 8, io_nstgrd = 9, &
    & io_sedspc = 10, io_biospc = 11, io_luvsur = 12, &
    & io_2uvobc = 13, io_3uvobc = 14, io_salobc = 15, &
    & io_tmpobc = 16, io_sedobc = 17, io_bioobc = 18, &
    & io_rlxobc = 19, io_nstspc = 20, io_2uvnst = 21, &
    & io_3uvnst = 22, io_salnst = 23, io_tmpnst = 24, &
    & io_sednst = 25, io_bionst = 26, io_metsur = 27, &
    & io_sstsur = 28, io_wavsur = 29, io_biosur = 30, &
    & io_drycel = 31, io_thndam = 32, io_weibar = 33, &
    & io_disspc = 34, io_disloc = 35, io_disvol = 36, &
    & io_discur = 37, io_dissal = 38, io_distmp = 39

INTEGER, PARAMETER :: icif_defruns= 1, icif_model = 2, icif_bio = 3, &
    & icif_sed = 4

INTEGER, PARAMETER :: ics_phys = 1, ics_bio = 2, ics_sed = 3

INTEGER, PARAMETER :: igrd_model = 1, igrd_meteo = 2, igrd_sst = 3, &
    & igrd_waves = 4, igrd_bio = 5

CHARACTER (LEN=6), DIMENSION(MaxIOTypes) :: modfiles_desc = &
    & ('mppmod','inicon','fincon','modgrd','metgrd','sstgrd','wavgrd','biogr','&
    & 'nstgrd','sedspc','biospc','luvsur','2uvobc','3uvobc','salobc','tmpobc','&
    & 'sedobc','bioobc','rlxobc','nstspc','2uvnst','3uvnst','salnst','tmpnst','&
    & 'sednst','bionst','metsur','sstsur','biosur','wavsur'/)
```

File

iopars.f90

Type

Module

Purpose

Parameters for model forcing

Description

ciflinenum	Number of the last input line read from a CIF
ciffiles	Attributes of CIFs
icif_*	Key ids of CIFs
	icif_bio CIF with setup parameters for the biological model
	icif_defruns <i>defruns</i> file
	icif_model CIF with model setup parameters
	icif_sed CIF with setup parameters for the sediment model
ics_*	Key ids for the file index in case the first index of modfiles equals io_inicon
	ics_bio initial conditions for biological arrays (currently not implemented)
	ics_phys initial conditions for physical arrays
	ics_sed initial conditions for sediment arrays
igrd_*	Key ids used for defining surface data grids
	igrd_bio surface grid for biological data (currently not implemented)
	igrd_meteo meteorological grid
	igrd_model (horizontal) model grid
	igrd_sst SST grid
	igrd_waves surface wave grid
io_*	File descriptor key ids
	io_biogrd surface grid for biological data
	io_bionst biological output for nesting
	io_bioobc definitions of open boundary conditions for biological variables (file number equals 1) or input of open boundary data (file number larger than 1).

io_biospc	(time-independent) arrays used for the setup of a biological model (currently not implemented)
io_biosur	biological surface data (currently not implemented)
io_disspc	discharge specifier arrays
io_disloc	discharge locations
io_disvol	volume discharge data
io_discur	discharge area and direction
io_dissal	salinity discharge data
io_distmp	temperature discharge data
io_drycel	dry cell locations
io_inicon	initial conditions
io_fincon	final conditions
io_metgrd	surface meteo grid
io_metsur	meteorological forcing data
io_modgrd	model grid, bathymetry and open boundary locations
io_mppmod	domain decomopistion
io_nstgrd	locations of nested sub-grid open boundaries
io_nstspc	number of open boundary locations at sub-grid nested grids
io_rlxobc	definitions of areas for the application of the relaxation open boundary scheme
io_salnst	salinity output for nesting
io_salobc	definitions of open boundary conditions for salinity (file number equals 1) or input of open boundary data (file number larger than 1)
io_sednst	sediment output for nesting
io_sedobc	definitions of open boundary conditions for sediment variables (file number equals 1) or input of open boundary data (file number larger than 1).

<code>io_sedspc</code>	(time-independent) arrays used for the setup of a sediment model (particle attributes in the COHERENS sediment model)
<code>io_sstgrd</code>	surface SST grid
<code>io_sstsur</code>	SST data
<code>io_thndam</code>	thin dams locations
<code>io_tmpnst</code>	temperature output for nesting
<code>io_tmpobc</code>	definitions of open boundary conditions for temperature (file number equals 1) or input of open boundary data (file number larger than 1)
<code>io_wavgrd</code>	surface wave grid
<code>io_wavsur</code>	surface wave data
<code>io_weibar</code>	weirs and barriers locations and parameters
<code>io_1uvsur</code>	definitions of surface boundary conditions (file number equals 1) or input of surface data (file number greater than 1) in case the model is applied in water column (1-D) mode
<code>io_2uvnst</code>	2-D mode output for nesting
<code>io_2uvobc</code>	definitions of 2-D open boundary conditions (file number equals 1) or input of open boundary data (file number larger than 1)
<code>io_3uvnst</code>	3-D mode output for nesting
<code>io_3uvobc</code>	definitions of open boundary conditions for the baroclinic currents (file number equals 1) or input of open boundary data (file number larger than 1)
<code>maxdatafiles</code>	Largest file index (second dimension of <code>modfiles</code>) for a given file descriptor (first index of <code>modfiles</code>)
<code>modfiles</code>	Attributes of model forcing files. The first index is given by a file descriptor key id of the form <code>io_*</code> (see below), the second is the file index, the third equals 1 for input or 2 for output.
<code>modfiles_desc</code>	Descriptors of forcing files
<code>surfacegrids</code>	Attributes of surface data grids. The first index is given by a grid descriptor key id of the form <code>igrd_*</code> (see above), the second is the file index (currently equal to 1)

Parameters for user-defined output

```

(MODULE iopars)
!---time series
INTEGER :: nosetstsr = 0, nostatstsr = 0, novarstsr = 0
INTEGER, DIMENSION(nosetstsr,novarstsr) :: ivarstsr
INTEGER, DIMENSION(nosetstsr,nostatstsr) :: lstatstsr
TYPE (FileParams), DIMENSION(nosetstsr) :: tsrgrd, tsr0d, tsr2d, tsr3d
TYPE (OutGridParams), DIMENSION(nosetstsr) :: tsrgpars
TYPE (StationLocs), DIMENSION(nostatstsr) :: tsrstatlocs
TYPE (VariableAtts), DIMENSION(novarstsr) :: tsrvars
!---time averages
INTEGER :: nosetsavr = 0, nostatsavr = 0, novarsavr = 0
INTEGER, DIMENSION(nosetsavr,novarsavr) :: ivarsavr
INTEGER, DIMENSION(nosetsavr,nostatsavr) :: lstatsavr
TYPE (FileParams), DIMENSION(nosetsavr) :: avrgrd, avr0d, avr2d, avr3d
TYPE (OutGridParams), DIMENSION(nosetsavr) :: avrgpars
TYPE (StationLocs), DIMENSION(nostatsavr) :: avrstatlocs
TYPE (VariableAtts), DIMENSION(novarsavr) :: avrvars
!---harmonic analysis
INTEGER :: nofreqsanal = 0, nosetsanal = 0, nostatsanal = 0, novarsanal = 0
CHARACTER (LEN=lentime), DIMENSION(nosetsanal) :: cdate_time_ref
CHARACTER (LEN=lenfreq), DIMENSION(nofreqsanal) :: harm_freq_names
INTEGER, DIMENSION(nofreqsanal) :: index_anal
INTEGER, DIMENSION(nosetsanal) :: nofreqsharm
INTEGER, DIMENSION(nosetsanal,nofreqsanal) :: ifreqsharm
INTEGER, DIMENSION(nosetsanal,novarsanal) :: ivarsanal
INTEGER, DIMENSION(nosetsanal,nostatsanal) :: lstatsanal
REAL, DIMENSION(nofreqsanal) :: harm_freq
TYPE (FileParams), DIMENSION(nosetsanal) :: analgrd, res0d, res2d, res3d
TYPE (FileParams), DIMENSION(nosetsanal,nofreqsanal) :: &
& amp0d, amp2d, amp3d, &
& pha0d, pha2d, pha3d, &
& ell2d, ell3d
TYPE (OutGridParams), DIMENSION(nosetsanal) :: analgpars
TYPE (StationLocs), DIMENSION(nostatsanal) :: analstatlocs
TYPE (VariableAtts), DIMENSION(novarsanal) :: analvars
!---elliptic parameters
INTEGER, DIMENSION(nosetsanal,14) :: ivarsell
INTEGER, DIMENSION(nosetsanal,2) :: ivecell2d, ivecell3d
TYPE (VariableAtts), DIMENSION(14) :: ellvars

```

```
!---output operators
INTEGER, PARAMETER :: oopt_null = 0, oopt_dep = 1, oopt_klev = 2, &
& oopt_min = 3, oopt_max = 4, oopt_mean = 5
```

File

iopars.f90

Type

Module

Purpose

Parameters for user-defined output

Description

<code>analgpars</code>	Attributes of the output grids for harmonic output
<code>analgrd</code>	Attributes of the grid file for harmonic output containing coordinate data only
<code>analstatlocs</code>	Attributes of all output stations for harmonic output (locations, names)
<code>analvars</code>	Attributes of all variables for harmonic output
<code>amp0d</code>	Attributes of the 0-D files for the harmonic output of amplitudes
<code>amp2d</code>	Attributes of the 2-D files for the harmonic output of amplitudes
<code>amp3d</code>	Attributes of the 3-D files for the harmonic output of amplitudes
<code>avrgpars</code>	Attributes of the output grids for time averaged output
<code>avrgrd</code>	Attributes of the grid file for time averaged output containing coordinate data only
<code>avrstatlocs</code>	Attributes of all output stations for time averaged output (locations, names)
<code>avrvars</code>	Attributes of all variables for time averaged output
<code>avr0d</code>	Attributes of the 0-D files for time averaged output
<code>avr2d</code>	Attributes of the 2-D files for time averaged output
<code>avr3d</code>	Attributes of the 3-D files for time averaged output

<code>cdate_time_ref</code>	Reference date with respect to which the analysed phases are given. If not defined, the phases are given with respect to the central analysis time or the Greenwich astronomical phase.
<code>ellvars</code>	Attributes of all possible elliptic variables (set by the program)
<code>ell2d</code>	Attributes of the 2-D files for harmonic output of elliptic parameters
<code>ell3d</code>	Attributes of the 3-D files for harmonic output of elliptic parameters
<code>harm_freq</code>	Frequencies used for analysis [radian/s]
<code>harm_freq_names</code>	Names of the frequencies used for harmonic analysis
<code>ifreqsharm</code>	Index mapping array for harmonic frequencies. The element <code>ifreqsharm(iset,ifreq)</code> maps, for set <code>iset</code> , the local frequency <code>ifreq</code> into the corresponding “global” array index as defined in the array <code>index_anal</code> .
<code>index_anal</code>	Key ids of the harmonic constituents. If zero, the corresponding frequencies need to be defined by the user as well.
<code>ivarsanal</code>	Index mapping array for harmonic output. The element <code>ivarsanal(iset,ivar)</code> maps, for set <code>iset</code> , the local variable index <code>ivar</code> into the corresponding array index in <code>analvars</code> .
<code>ivarsavr</code>	Index mapping array for time averaged output. The element <code>ivarsavr(iset,ivar)</code> maps, for set <code>iset</code> , the local variable index <code>ivar</code> into the corresponding array index in <code>avrvars</code> .
<code>ivarsell</code>	Index mapping array selecting the elliptic parameters for output. The element <code>ivarsell(iset,ivar)</code> maps, for set <code>iset</code> , the local variable index <code>ivar</code> into the corresponding array index in <code>ellvars</code> .
<code>ivarstr</code>	Index mapping array for time series output. The element <code>ivarstr(iset,ivar)</code> maps, for set <code>iset</code> , the local variable index <code>ivar</code> into the corresponding array index in <code>tsrvars</code> (see below).
<code>ivecell2d</code>	Indices of the elements in <code>analvars</code> representing the X- and Y-component of the 2-D elliptic vectors

<code>ivecell3d</code>	Indices of the elements in <code>analvars</code> representing the X- and Y-component of the 3-D elliptic vectors
<code>lstatsanal</code>	Station label for harmonic output. The element <code>lstatsanal(iset,istat)</code> maps, for set <code>iset</code> , the local index <code>istat</code> into the corresponding global array index in <code>analstatlocs</code> .
<code>lstatsavr</code>	Station label for time averaged output. The element <code>lstatsavr(iset,istat)</code> maps, for set <code>iset</code> , the local index <code>istat</code> into the corresponding global array index in <code>avrstatlocs</code> .
<code>lstatstsr</code>	Station label for time series output. The element <code>lstatstsr(iset,istat)</code> maps, for set <code>iset</code> , the local index <code>istat</code> into the corresponding global array index in <code>tsrstatlocs</code> .
<code>nofreqsanal</code>	Total number of frequencies for harmonic analysis
<code>nofreqsharm</code>	Number of frequencies used for each output set
<code>nosetsanal</code>	Number of output sets for harmonic output
<code>nosetsavr</code>	Number of output sets for time averaged output
<code>nosetstsr</code>	Number of time series output sets
<code>nostatsanal</code>	Total number of stations for harmonic output
<code>nostatsavr</code>	Total number of stations for time averaged output
<code>nostatstsr</code>	Total number of stations for time series output
<code>novarsanal</code>	Total number of variables for harmonic output
<code>novarsavr</code>	Total number of variables for time averaged output
<code>novarstsr</code>	Total number of variables for time series output
<code>oopt_*</code>	User output operators
<code>oopt_dep</code>	At a specified depth from the surface
<code>oopt_klev</code>	At a specified vertical level
<code>oopt_max</code>	Maximum value
<code>oopt_mean</code>	Spatially averaged value
<code>oopt_min</code>	Minimum value
<code>oopt_null</code>	No operator
<code>pha0d</code>	Attributes of the 0-D files for the harmonic output of phases

pha2d	Attributes of the 2-D files for the harmonic output of phases
pha3d	Attributes of the 3-D files for the harmonic output of phases
res0d	Attributes of the 0-D files for the residual harmonic output
res2d	Attributes of the 2-D files for the residual harmonic output
res3d	Attributes of the 3-D files for the residual harmonic output
tsrgpars	Attributes of the output grids for time series output
tsrgrd	Attributes of the grid file for time series output containing coordinate data only
tsrstatlocs	Attributes of all output stations for time series output (locations, names)
tsrvars	Attributes of all variables for time series output
tsr0d	Attributes of the 0-D files for time series output
tsr2d	Attributes of the 2-D files for time series output
tsr3d	Attributes of the 3-D files for time series output

Parameters for monitoring

```
(MODULE iopars)
!---log files
LOGICAL :: exitlog
CHARACTER (LEN=*), PARAMETER :: logfmt1 = '(I1,':':',A)', &
& logfmt2 = '(I2,':':',A)'
CHARACTER (LEN=1), PARAMETER :: logexit = 'R'
CHARACTER (len=leniofile) :: inilog_file, runlog_file
INTEGER :: iolog = 0, loglev1, loglev2, pglev, runlog_count
INTEGER, DIMENSION(levprocs) :: levprocs_ini, levprocs_run
!---error files
LOGICAL :: errchk
CHARACTER (len=leniofile) :: errlog_file
INTEGER :: errstat, ioerr = 0, maxerrors, nerrs = 0
INTEGER, DIMENSION(npworld) :: levprocs_err
INTEGER, PARAMETER :: &
& ierrno_fopen = 1, ierrno_fclose = 2, ierrno_read = 3, ierrno_write = 4,&
```

```

    & ierrno_fend = 5, ierrno_input = 6, ierrno_inival = 7, ierrno_runval = 8, &
    & ierrno_alloc = 9, ierrno_arg = 10, ierrno_comms = 11, ierrno_MPI = 12, &
    & ierrno_CDF = 13
CHARACTER (LEN=lenerrcode), PARAMETER, DIMENSION(MaxErrCodes) :: error_code
!---warning file
LOGICAL :: warnflag, warning
CHARACTER (len=leniofile) :: warlog_file
INTEGER :: iowarn = 0
!---timer
LOGICAL :: timer = .FALSE.
CHARACTER (len=leniofile) :: timing_file
INTEGER :: levtimer = 0, maxwaitsecs = 3600, nowaitsecs = 0, npcc_max, &
    & npcc_rate, timer_format = 1
INTEGER (KIND=kndilong), DIMENSION(MaxTimers) :: nopcc
INTEGER, PARAMETER :: &
    & itm_hydro = 1, itm_1dmode = 2, itm_2dmode = 3, itm_3dmode = 4, &
    & itm_dens = 5, itm_temp = 6, itm_sal = 7, itm_init = 8, &
    & itm_trans = 9, itm_adv = 10, itm_hdif = 11, itm_vdif = 12, &
    & itm_phgrad = 13, itm_input = 14, itm_output = 15, itm_inout = 16, &
    & itm_com_coll = 17, itm_com_comb = 18, itm_com_copy = 19, &
    & itm_com_dist = 20, itm_com_exch = 21, itm_com_util = 22, itm_coms = 23, &
    & itm_MPI = 24, itm_CDF = 25, itm_arrint = 26, itm_user = 27, &
    & itm_nest = 28, itm_libs = 29, itm_astro = 30, itm_bconds = 31, &
    & itm_meteo = 32, itm_structs = 33, itm_wait = 34, itm_sed = 35, itm_bio = 3
!---timer descriptions
CHARACTER (LEN=20), DIMENSION(MaxTimers) :: desctimer

File
    iopars.f90

Type
    Module

Purpose
    Parameters for monitoring

Description
    desctimer      Strings written to the timer report for each (active)
                   timer
    errchk         Enables/disables error checking
    errlog_file    Default name of the “errlog” file (appended by the
                   process id in the parallel case)

```

<code>error_code</code>	Error message string corresponding to an error code number
<code>errstat</code>	Error status number as returned by MPI, netCDF and FORTRAN calls (e.g. ALLOCATE statement)
<code>exitlog</code>	Enables/disables the writing of the “exit” message to the “log” file when the program exits a routine
<code>ierrno_*</code>	Key ids for error messages
	<code>ierrno_alloc</code> allocation error
	<code>ierrno_arg</code> missing or invalid routine argument
	<code>ierrno_CDF</code> netCDF error
	<code>ierrno_comms</code> parallel communication error
	<code>ierrno_fclose</code> an error occurred when a file is closed
	<code>ierrno_fend</code> end of file condition
	<code>ierrno_fopen</code> an error occurred when a file is opened
	<code>ierrno_inival</code> invalid value for a setup or initial scalar or array parameter
	<code>ierrno_input</code> an error occurred when reading an invalid value for a parameter or array from a data file
	<code>ierrno_MPI</code> MPI error
	<code>ierrno_read</code> read error
	<code>ierrno_runval</code> invalid value for a scalar or array parameter during the time loop
	<code>ierrno_write</code> write error
<code>inilog_file</code>	Default name of the “inilog” file (appended by the process id in the parallel case)
<code>ioerr</code>	File unit of the “errlog” file
<code>iolog</code>	File unit of the “inilog” and “runlog” files
<code>iowarn</code>	File unit of the warning file
<code>itm_*</code>	Timer key ids
	<code>itm_adv</code> advection routines
	<code>itm_arrint</code> interpolation of model grid arrays
	<code>itm_astro</code> astronomical tide
	<code>itm_bconds</code> boundary conditions

itm_bio	biology (currently not activated)
itm_CDF	total of all netCDF calls
itm_coms	total of all communication calls
itm_com_coll	collect communication calls
itm_com_comb	combine communications calls
itm_com_copy	copy communication calls
itm_com_dist	distribute communication calls
itm_com_exch	exchange communication calls
itm_com_util	(parallel) utility communication calls
itm_dens	total of density (including temperature and salinity) calculations
itm_hdif	horizontal diffusion
itm_hydro	hydrodynamics
itm_init	initialisation procedures
itm_inout	total of input and output operations
itm_input	input operations
itm_libs	calls to library routines
itm_meteo	meteorological routines
itm_MPI	total of all MPI calls
itm_nest	nesting procedures
itm_output	output operations
itm_phgrad	baroclinic pressure gradient
itm_sal	salinity
itm_sed	sediment model
itm_structs	structures
itm_temp	temperature
itm_trans	transport routines
itm_user	calls to <code>usrdef</code> routines
itm_vdif	vertical diffusion (including turbulence modules)
itm_wait	wait calls
itm_1dmode	water column mode (1-D) calculations
itm_2dmode	2-D mode calculations

	itm_3dmode	3-D mode calculations
levprocs_err	Level of error checking	<ul style="list-style-type: none"> 0: error checking disabled 1: error checking enabled during the initialisation phase only 2: error checking enabled during the whole simulation
levprocs_ini	Maximum program level for writing a tracer information message to the “inilog” file when a routine is entered. Its value is determined for each process separately.	
levprocs_run	Maximum program level for writing a tracer information message to the “runlog” file when a routine is entered. Its value is determined for each process separately.	
levtimer	Defines the information in the timer report	<ul style="list-style-type: none"> 0: no timer report is written 1: only the execution time is written 2: Execution time and time information (in percentage of total time) is written for all “timers”. In case of a parallel simulation, the percentages are given for the process with the largest amount of time, the lowest amount of time, an average value over all processes and for the master process. 3: As the previous case, but the time percentages are now given for each individual process in addition. In the serial case, behaviour is as for case 2.
logexit	Exit message string written to the “log” file on exiting of a routine	
logfmt1	Format specification for writing the tracer information to the “log” file	
logfmt2	As logfmt1 now for program levels higher than 9	
loglev1	Maximum program level for writing a tracer information message when a routine is entered	
loglev2	Maximum program level for writing a tracer information message when a routine is exited	
maxerrors	Maximum number of messages written to the “errlog” file	

maxwaitsecs	Maximum allowed time (in seconds) for suspension of the program
nerrs	Number of detected errors
nopcc	Current number of process clock counts stored for each timer
nowaitsecs	Number of seconds to suspend the execution of the program during a wait call
npcc_max	Maximum number of clock counts which can be returned by the process clock
npcc_rate	Number of process clock counts per second
pglev	Current program level
runlog_count	Determines the number of (2-D) time steps after which the log-file will be re-written
runlog_file	Default name of the “runlog” file (appended by the process id in the parallel case)
timer	.TRUE. if levtimer>0 (see above)
timer_format	Unit of the execution time in the timer report 1: seconds 2: minutes 3: hours 4: days
timing_file	Default name of the time report file
warlog_file	Default name of the “warning” file
warnflag	Enables/disables the writing of a warning (“warlog”) file by the master process (.TRUE. on the master only if warning is .TRUE.)
warning	Enables/disables the writing of warning messages

netCDF parameters

(MODULE iopars)

```
INTEGER :: char_NF90 = 0, clobber_NF90 = 0, fill_NF90 = 0, global_NF90 = 0,&
& int_NF90 = 0, noerr_NF90 = 0, nofill_NF90 = 0, nowrite_NF90 = 0,&
& offset_64bit_NF90 = 0, real_NF90 = 0, share_NF90 = 0,&
& sizehint_default_NF90 = 0, unlimited_NF90 = 0, write_NF90 = 0
```

File

iopars.f90

Type

Module

Purpose

Aliases for parameters of the netCDF library

Description

<code>char_NF90</code>	Alias for <code>NF90_char</code>
<code>clobber_NF90</code>	Alias for <code>NF90_clobber</code>
<code>fill_NF90</code>	Alias for <code>NF90_fill</code>
<code>global_NF90</code>	Alias for <code>NF90_global</code>
<code>int_NF90</code>	Alias for <code>NF90_int</code>
<code>noerr_NF90</code>	Alias for <code>NF90_noerr</code>
<code>nofill_NF90</code>	Alias for <code>NF90_nofill</code>
<code>nowrite_NF90</code>	Alias for <code>NF90_nowrite</code>
<code>offset_64bit_NF90</code>	Alias for <code>NF90_offset_64bit</code>
<code>real_NF90</code>	Alias for <code>NF90_real</code>
<code>share_NF90</code>	Alias for <code>NF90_share</code>
<code>sizehint_default_NF90</code>	Alias for <code>NF90_sizehint_default</code>
<code>unlimited_NF90</code>	Alias for <code>NF90_unlimited</code>
<code>write_NF90</code>	Alias for <code>NF90_write</code>

33.10 Meteorological arrays

MODULE meteo

```
REAL, DIMENSION(0:ncloc,0:nrloc) :: atmpres, uwindatc, vwindatc, &
                                     & airtemp, sst
```

```
REAL, DIMENSION(ncloc,nrloc) :: cloud_cover, evapminprec, &
                                     & evaporation, precipitation, relhum
```

File

meteo.f90

Type

Module

Purpose

Meteorological data variables

Description

<code>airtemp</code>	Air temperature interpolated in time and in space at the C-nodes	[°C]
<code>atmpres</code>	Atmospheric pressure interpolated in time and in space at the C-nodes	[Pa]
<code>cloud_cover</code>	Cloud cover interpolated in time and in space at the C-nodes (between 0 and 1)	
<code>evapminprec</code>	Evaporation minus precipitation rate interpolated in time and in space at the C-nodes	[kg/m ² /s]
<code>evaporation</code>	Evaporation rate interpolated in time and in space at the C-nodes	[kg/m ² /s]
<code>precipitation</code>	Precipitation rate interpolated in time and in space at the C-nodes	[kg/m ² /s]
<code>relhum</code>	Relative humidity interpolated in time and in space at the C-nodes (between 0 and 1)	
<code>sst</code>	Sea surface temperature	[°C]
<code>uwindatc</code>	X-component of the wind at 10m height interpolated in time and in space at C-nodes	[m/s]
<code>vwindatc</code>	Y-component of the wind at 10m height interpolated in time and in space at C-nodes	[m/s]

33.11 Key ids of model variables

```

MODULE modids
!---model grid
INTEGER, PARAMETER :: &
& iarr_alphatc_fld = 1, iarr_alphatu_fld = 2, iarr_alphatv_fld = 3,&
....

```

File

modids.f90

Type

Module

Purpose

Definitions of key ids for model variables. The key id name has the form `iarr_*` where `*` is the FORTRAN name of the variable.

33.12 Nested sub-grids

```

MODULE nestgrids
INTEGER :: nonestsets = 0
INTEGER, DIMENSION(nonestsets) :: nestcoords, nohnstglbc, &
                                & nohnstglbu, nohnstglbv
INTEGER, DIMENSION(nonestsets) :: nohnstatc, nohnstatu, nohnstatv, novnst
INTEGER, DIMENSION(nonestsets) :: inst2dtype, lbhnstatc, lbhnstatu, lbhnstatv
INTEGER, DIMENSION(nonestsets) :: nobionst, nosednst
INTEGER, DIMENSION(nprocs,nonestsets) :: nohnstcprocs, nohnstuvprocs
INTEGER, DIMENSION(maxbiovars,nonestsets) :: instbio
INTEGER, DIMENSION(maxsedvars,nonestsets) :: instsed
INTEGER, DIMENSION(nprocs,MAXVAL(nohnstcprocs),nonestsets) :: indexnstc
INTEGER, DIMENSION(nprocs,MAXVAL(nohnstuvprocs),nonestsets) :: indexnstuv
TYPE (HRelativeCoords), DIMENSION(SUM(nohnstatc)) :: hnstctoc
TYPE (HRelativeCoords), DIMENSION(SUM(nohnstatu)) :: hnstctou, hnstutou
TYPE (HRelativeCoords), DIMENSION(SUM(nohnstatv)) :: hnstctov, hnstvtov
TYPE (VRelativeCoords), DIMENSION(2,2,SUM(nohnstatc),&
                                & MAXVAL(novnst)) :: vnstctoc
TYPE (VRelativeCoords), DIMENSION(2,2,SUM(nohnstatu),&
                                & MAXVAL(novnst)) :: vnstutou
TYPE (VRelativeCoords), DIMENSION(2,2,SUM(nohnstatv),&
                                & MAXVAL(novnst)) :: vnstvtov

```

File

nestgrids.f90

Type

Module

Purpose

Arrays for the setup of sub-grid nesting

Description

`hnstctoc` Relative horizontal coordinates of the C-node sub-grid points with respect to the local C-node main grid

<code>hnstctou</code>	Relative horizontal coordinates of the U-node sub-grid points with respect to the local C-node main grid
<code>hnstctov</code>	Relative horizontal coordinates of the V-node sub-grid points with respect to the local C-node main grid
<code>hnstutou</code>	Relative horizontal coordinates of the U-node sub-grid points with respect to the local U-node main grid
<code>hnstvtov</code>	Relative horizontal coordinates of the V-node sub-grid points with respect to the local V-node main grid
<code>indexnstc</code>	Index mapping array for C-node points. Element <code>indexnstc(iproc,lsub,iset)</code> equals <code>lglb-l1+1</code> where <code>iproc</code> is the process number, <code>lglb</code> the corresponding index in the global indexing system and <code>l1</code> the index in the global indexing system of the first global point in set <code>iset</code>
<code>indexnstuv</code>	Index mapping array for U- and V-points. Element <code>indexnstuv(iproc,lsub,iset)</code> equals <code>lglb-l1+1</code> where <code>iproc</code> is the process number, <code>lglb</code> the index in the global indexing system and <code>l1</code> the index in the global indexing system of the first global point in set <code>iset</code> . Note that U-points are counted before V-node points.
<code>instbio</code>	Variable indices of the biological state variables used for nesting per set
<code>instsed</code>	Fraction indices of the sediment concentrations used for nesting per set
<code>inst2dtype</code>	Type of data used for 2-D nesting 1: transports and elevations 2: elevations 3: transports
<code>lbhnstatc</code>	The C-node sub-grid points are indexed globally over all sub-grids. The element <code>lbhnstatc(iset)</code> represents the global index of the first point in the global indexing system which is located on the local sub-domain.
<code>lbhnstatu</code>	The U-node sub-grid points are indexed globally over all sub-grids. The element <code>lbhnstatu(iset)</code> represents the global index of the first point in the global indexing system which is located on the local sub-domain.
<code>lbhnstatv</code>	The V-node sub-grid points are indexed globally over all sub-grids. The element <code>lbhnstatv(iset)</code> represents the

	global index of the first point in the global indexing system which is located on the local sub-domain.
nestcoords	Type of coordinates used in the setup 1: absolute coordinates 2: relative coordinates
nobionst	Number of nested biological state variables per nested sub-grid
nohnstatc	Local number of C-node (horizontal) sub-grid open boundary points
nohnstatu	Local number of U-node (horizontal) sub-grid open boundary points
nohnstatv	Local number of V-node (horizontal) sub-grid open boundary points
nohnstcprocs	Number of C-node sub-grid points for each sub-domain and each sub-grid
nohnstglbc	Global number of C-node (horizontal) sub-grid open boundary points
nohnstglbu	Global number of U-node (horizontal) sub-grid open boundary points
nohnstglbv	Global number of V-node (horizontal) sub-grid open boundary points
nohnstuvprocs	Number of U- and V-node sub-grid points for each sub-domain and each sub-grid
nonestsets	Number of sub-grid nests
nosednst	Number of nested sediment fractions per nested sub-grid
novnst	Number of vertical levels on the sub-grid
vnstctoc	Relative vertical coordinates of the C-node sub-grid points with respect to the local main C-node grid taken at the four points surrounding each sub-grid point.
vnstctou	Relative vertical coordinates of the C-node sub-grid points with respect to the local main U-node grid taken at the four points surrounding each sub-grid point.
vnstctov	Relative vertical coordinates of the C-node sub-grid points with respect to the local main V-node grid taken at the four points surrounding each sub-grid point.

33.13 Open boundary conditions

```

MODULE obconds
!---2-D open boundary conditions
INTEGER, DIMENSION(nobu) :: iloczobu, ityp2dobu
INTEGER, DIMENSION(nobv) :: iloczobv, ityp2dobv
!nofiles = maxdatafiles(io_2uvobc,1)
INTEGER, DIMENSION(2:nofiles) :: iobc2dtype, no2dobc
INTEGER, DIMENSION(nobu+nobv,2:nofiles) :: index2dobc
REAL, DIMENSION(nobu,2) :: ud2obu, zetobu
REAL, DIMENSION(nobv,2) :: vd2obv, zetobv
REAL, DIMENSION(nobu,nconobc) :: ud2obu_amp, ud2obu_pha, &
                                & zetobu_amp, zetobu_pha
REAL, DIMENSION(nobv,nconobc) :: vd2obv_amp, vd2obv_pha, &
                                & zetobv_amp, zetobv_pha
REAL, DIMENSION(nobu,2) :: obc2uvatu, obc2uvatu_old
REAL, DIMENSION(nobv,2) :: obc2uvatv, obc2uvatv_old
!---3-D open boundary forcing
REAL, DIMENSION(nobu,nz,2) :: obc3uvatu
REAL, DIMENSION(nobv,nz,2) :: obc3uvatv
REAL, DIMENSION(nobu,nz,0:2) :: obcsalatu, obctmpatu
REAL, DIMENSION(nobv,nz,0:2) :: obcsalatv, obctmpatv
!---surface forcing (1-D application)
INTEGER :: isur1dtype
REAL :: gxslope = 0.0, gyslope = 0.0
REAL, DIMENSION(nconobc) :: gxslope_amp, gxslope_pha, gyslope_amp, &
                                & gyslope_pha, zeta_amp, zeta_pha

```

File

obconds.f90

Type

Module

Purpose

Parameters and arrays used to define open boundary conditions for the 2-D and 3-D mode or surface boundary conditions (surface slope and elevations) for 1-D applications

Description

gxslope	X-component of the (barotropic) pressure gradient in case of a 1-D application	[m/s ²]
----------------	--	---------------------

<code>gxslope_amp</code>	Amplitudes of the X-component of the (barotropic) pressure gradient in case of a 1-D application [m/s ²]
<code>gxslope pha</code>	Phases of the X-component of the (barotropic) pressure gradient in case of a 1-D application [rad]
<code>gyslope</code>	Y-component of the (barotropic) pressure gradient in case of a 1-D application [m/s ²]
<code>gyslope_amp</code>	Amplitudes of the Y-component of the (barotropic) pressure gradient in case of a 1-D application [m/s ²]
<code>gyslope pha</code>	Phases of the Y-component of the (barotropic) pressure gradient in case of a 1-D application [rad]
<code>iloczobu</code>	If the elevation has to be specified at a U-open boundary, the array selects the position of the specified elevation with respect to the open boundary. 0: not required 1: at the open boundary U-node 2: at the “nearest” C-node outside the domain
<code>iloczobv</code>	The same as <code>iloczobu</code> now at V-open boundaries
<code>index2dobc</code>	Each 2-D open boundary data file contains a sub-set of open boundary data points. The element <code>index2dobc(idat,ifil)</code> maps, for file <code>ifil</code> , the local data point <code>idat</code> into a corresponding global open boundary index (between <code>1:nobu</code> for U- and <code>nobu+1:nobu+nobv</code> for V-open boundaries). The physical size of the first dimension for file <code>ifil</code> equals <code>no2dobc(ifil)</code> .
<code>iobc2dtype</code>	Identifies the type of variables within each 2-D open boundary data file 1: depth-integrated currents and elevations 2: elevations only 3: depth-integrated currents only
<code>isur1dtype</code>	Identifies the variables within a 1-D surface forcing data file. 1: components of the pressure gradient and elevation 2: surface elevation 3: components of the pressure gradient
<code>ityp2dobu</code>	Type of open boundary conditions at the U-nodes (0/13)

	0 : clamped
	1 : zero slope
	2 : zero volume flux
	3 : specified elevation
	4 : specified transport
	5 : radiation condition using shallow water speed
	6 : Orlanski (1976) condition
	7 : Camerlengo & O'Brien (1980)
	8 : Flather (1976) with specified elevation and transport
	9 : Flather with specified elevation
	10: Røed & Smedstad (1984)
	11: characteristic method with specified elevation and transport
	12: characteristic method with specified elevation
	13: characteristic method using a zero normal gradient condition
ityp2dobv	Type of open boundary condition at the V-nodes. Meaning is the same as for ityp2dobu with U replaced by V and West/East by South/North.
no2dobc	Number of input open boundary locations for each 2-D open boundary data file
obcsalatu	Storage array for salinity S in case the open boundary conditions at the U-nodes require the solution of a differential equation in time
obcsalav	Storage array for salinity S in case the open boundary conditions at the V-nodes require the solution of a differential equation in time
obctmpatu	Storage array for temperature T in case the open boundary conditions at the U-nodes require the solution of a differential equation in time
obctmpav	Storage array for temperature T in case the open boundary conditions at the V-nodes require the solution of a differential equation in time

<code>obc2uvatv</code>	Storage array for the X-component of the transport U in case the open boundary conditions at the U-nodes require the solution of a differential equation in time
<code>obc2uvatv_old</code>	Value of <code>obc2uvatv</code> at the first (outer) iteration with the implicit scheme
<code>obc3uvatv</code>	Storage array for the Y-component of the transport V in case the open boundary conditions at the V-nodes require the solution of a differential equation in time
<code>obc3uvatv_old</code>	Value of <code>obc3uvatv</code> at the first (outer) iteration with the implicit scheme
<code>ud2obu</code>	X-component of the transport U at U-open boundaries as given by (4.354). The external term term is stored in elements (*,1), the full expression (including the harmonic expansion) in elements (*,2) [m ² /s]
<code>ud2obu_amp</code>	Amplitudes A_n in the harmonic expansion of the X-component of the transport U at U-open boundaries [m ² /s]
<code>ud2obu pha</code>	Phases φ in the harmonic expansion of the X-component of the transport U at U-open boundaries [rad]
<code>vd2obv</code>	Y-component of the transport V at V-open boundaries as given by (4.354). The external term term is stored in elements (*,1), the full expression (including the harmonic expansion) in elements (*,2) [m ² /s]
<code>vd2obv_amp</code>	Amplitudes A_n in the harmonic expansion of the Y-component of the transport V at V-open boundaries [m ² /s]
<code>vd2obv pha</code>	Phases φ in the harmonic expansion of the Y-component of the transport V at V-open boundaries [rad]
<code>zeta_amp</code>	Amplitudes of the surface elevation ζ in case of a 1-D application [m]
<code>zeta pha</code>	Phases of the surface elevation ζ in case of a 1-D application [rad]

<code>zetobu</code>	Surface elevation at U-open boundaries as given by (4.354). The external term term is stored in elements (*,1), the full expression (including the harmonic expansion) in elements (*,2) [m]
<code>zetobu_amp</code>	Amplitudes A_n in the harmonic expansion of the surface elevation ζ at U-open boundaries [m]
<code>zetobu pha</code>	Phases φ in the harmonic expansion of the surface elevation ζ at U-open boundaries [rad]
<code>zetobv</code>	Surface elevation at V-open boundaries as given by (4.354). The external term term is stored in elements (*,1), the full expression (including the harmonic expansion) in elements (*,2) [m]
<code>zetobv_amp</code>	Amplitudes A_n in the harmonic expansion of the surface elevation ζ at V-open boundaries [m]
<code>zetobv pha</code>	Phases φ in the harmonic expansion of the surface elevation ζ at V-open boundaries [rad]

33.14 Optical arrays

```

MODULE optics
REAL, DIMENSION(ncloc,nrloc) :: optattcoef2, qrad
REAL, DIMENSION(ncloc,nrloc,nz+1) :: radiance

```

File

optics.f90

Type

Module

Purpose

Optical arrays

Description

<code>optattcoef2</code>	Inverse optical attenuation depth for the short-wave spectrum [m ⁻¹]
<code>qrad</code>	Solar downward irradiance within the water column [W/m ²]
<code>radiance</code>	Solar irradiance incident on the surface [W/m ²]

33.15 Parameters for parallel processing

```

MODULE paralpars
LOGICAL :: parallel_set = .FALSE.
LOGICAL :: master, spare, worker
INTEGER :: idmaster = 0, nprocs = 1, nprocsx = 0, nprocsy = 0
INTEGER :: comm_work, icoordloc, idloc, iprocloc, jcoordloc, npworld
INTEGER, DIMENSION(2*nprocs) :: comprocs
INTEGER, DIMENSION(npworld) :: idprocs
INTEGER, DIMENSION(nprocs) :: icoordprocs, jcoordprocs, nprocs, nc1procs, &
& nc2procs, nobupprocs, nobvprocs, nobxprocs, &
& nobyprocs, nodisprocs, nosbupprocs, nosbvprocs, &
& nowbarupprocs, nowbarvprocs, nrprocs, nrvbupprocs, &
& nrvbvprocs, nr1procs, nr2procs
INTEGER, DIMENSION(0:nprocsx+1,0:nprocsy+1) :: iddomain
INTEGER, DIMENSION(nobu,nprocs) :: indexobupprocs
INTEGER, DIMENSION(nobv,nprocs) :: indexobvprocs
INTEGER, DIMENSION(nobx,nprocs) :: indexobxprocs
INTEGER, DIMENSION(noby,nprocs) :: indexobyprocs
INTEGER, DIMENSION(numdis,nprocs) :: indexdisprocs
INTEGER, DIMENSION(numwbaru,nprocs) :: indexwbaruprocs
INTEGER, DIMENSION(numwbarv,nprocs) :: indexwbarvprocs
TYPE(ExchComms), DIMENSION(MaxHaloComms) :: halocomms
!---MPI definitions
INTEGER :: bsend_overhead_MPI = 0, comm_null_MPI = 0, comm_world_MPI = 0, &
& proc_null_MPI = 0, undefined_MPI = 0

```

File

paralpars.f90

Type

Module

Purpose

Parameters for parallel applications

Description

<code>bsend_overhead_MPI</code>	Alias for <code>MPI_bsend_overhead</code>
<code>comm_null_MPI</code>	Alias for <code>MPI_comm_null</code>
<code>comm_work</code>	MPI communicator containing all worker (non-idle) processes

<code>comm_world_MPI</code>	Alias for <code>MPI_comm_world</code>
<code>comprocs</code>	Array defining the rank order for all-to-all communications (see Section 11.4.3.1)
<code>halocomms</code>	Parameters for exchange (send/receive) communications
<code>icoordloc</code>	Domain index in the X-direction of the local process on the parallel grid
<code>icoordprocs</code>	Array with the values of <code>icoordloc</code> for each process
<code>iddomain</code>	Process ranks as defined on the parallel domain grid
<code>idloc</code>	Rank of the local process
<code>idmaster</code>	Rank of the master process
<code>idprocs</code>	Array of process ranks (global)
<code>indexdisprocs</code>	Index mapping array at discharge locations. The element <code>indexdisprocs(lloc,iprocs)</code> maps the local discharge position index <code>lloc</code> on the local process <code>iprocs</code> to the corresponding global discharge position index.
<code>indexobuprocs</code>	Index mapping array at the U-open boundaries. The element <code>indexobuprocs(lloc,iprocs)</code> maps the local open boundary index <code>lloc</code> on the local process <code>iprocs</code> to the corresponding global open boundary index.
<code>indexobvprocs</code>	Index mapping array at the V-open boundaries. The element <code>indexobvprocs(lloc,iprocs)</code> maps the local open boundary index <code>lloc</code> on the local process <code>iprocs</code> to the corresponding global open boundary index.
<code>indexobxprocs</code>	Index mapping array at the X-open boundaries. The element <code>indexobxprocs(lloc,iprocs)</code> maps the local open boundary index <code>lloc</code> on the local process <code>iprocs</code> to the corresponding global open boundary index.
<code>indexobyprocs</code>	Index mapping array at the Y-open boundaries. The element <code>indexobyprocs(lloc,iprocs)</code> maps the local open boundary index <code>lloc</code> on the local process <code>iprocs</code> to the corresponding global open boundary index.

<code>indexwbaruprocs</code>	Index mapping array at U-node weir/barrier locations. The element <code>indexwbaruprocs(lloc,iproc)</code> maps the local U-node weir/barrier position index <code>lloc</code> on the local process <code>iproc</code> to the corresponding global U-node weir/barrier position index.
<code>indexwbarvprocs</code>	Index mapping array at V-node weir/barrier locations. The element <code>indexwbarvprocs(lloc,iproc)</code> maps the local U-node weir/barrier position index <code>lloc</code> on the local process <code>iproc</code> to the corresponding global V-node weir/barrier position index.
<code>iprocloc</code>	Local process number (<code>=idloc+1</code>)
<code>jcoordloc</code>	Domain index in the Y-direction of the local process on the parallel grid
<code>jcoordprocs</code>	Array with the values of <code>jcoordloc</code> for each process
<code>master</code>	<code>.TRUE.</code> on the master process, <code>.FALSE.</code> otherwise
<code>ncprocs</code>	Array with the X-dimension <code>ncloc</code> of each sub-domain (global)
<code>nc1procs</code>	Array with the values of <code>nc1loc</code> for each sub-domain (global)
<code>nc2procs</code>	Array with the values of <code>nc2loc</code> for each sub-domain (global)
<code>nobuprocs</code>	Array with the values of <code>nobuloc</code> for each process
<code>nobvprocs</code>	Array with the values of <code>nobvloc</code> for each process
<code>nobxprocs</code>	Array with the values of <code>nobxloc</code> for each process
<code>nobyprocs</code>	Array with the values of <code>nobyloc</code> for each process
<code>nodisprocs</code>	Array with the values of <code>numdisloc</code> for each process
<code>nosbuprocs</code>	Array with the values of <code>nosbuloc</code> for each process
<code>nosbvprocs</code>	Array with the values of <code>nosbvloc</code> for each process
<code>nowbaruprocs</code>	Array with the values of <code>numwbaruloc</code> for each process
<code>nowbarvprocs</code>	Array with the values of <code>numwbarvloc</code> for each process
<code>nprocs</code>	Total number of active (“worker”) processes used in the simulation
<code>nprocsx</code>	X-dimension of the parallel grid

<code>nprocsy</code>	Y-dimension of the parallel grid
<code>npworld</code>	Total number of processes in <code>MPI_comm_world</code> . This equals the number given in the script used to launch the program
<code>nrprocs</code>	Array with the Y-dimension <code>nrloc</code> of each sub-domain (global)
<code>nr1procs</code>	Array with the values of <code>nr1loc</code> for each sub-domain (global)
<code>nr2procs</code>	Array with the values of <code>nr2loc</code> for each sub-domain (global)
<code>nrvbuprocs</code>	Array with the values of <code>nrvbuloc</code> for each process
<code>nrvbvprocs</code>	Array with the values of <code>nrvbvloc</code> for each process
<code>parallel_set</code>	Enables/disables the parallel mode. The switch is automatically set by the program if the compiler option <code>-DMPI</code> is defined in <code>coherensflags.cmp</code> .
<code>proc_null_MPI</code>	Alias for <code>MPI_proc_null</code>
<code>spare</code>	.TRUE. if the local process is an idle (spare) process which is currently inactive
<code>undefined_MPI</code>	Alias for <code>MPI_undefined</code>
<code>worker</code>	.TRUE. for an active on (worker) process. Is the same as <code>.NOT.spare</code> .

33.16 Physical and numerical model parameters

```

MODULE physpars
!---general
REAL :: Rearth = 6371000.0, rho_air = 1.2, specheat = 3987.5
!---reference and minimum values
REAL :: atmpres_ref = 101325.0, beta_sal_ref, beta_temp_ref, density_ref, &
      & dlat_ref = 0.0, dlon_ref = 0.0, dlon_ref_anal = 0.0, &
      & dlon_ref_obc = 0.0, gacc_mean, gacc_ref = real_undef, &
      & sal_ref = 33.0, sst_ref = 12.0, temp_min = 0.0, temp_ref = 12.0
!---model grid
REAL :: b_SH = 0.1, dl_BB = 1.5, du_BB = 1.5, hcrit_SH = 200.0, &
      & sigstar_DJ = 0.0, sig0_DJ = 0.1, theta_SH = 8.0
!---diffusion coefficients

```



```

REAL :: hdifmom_cst = 0.0, hdifscal_cst = 0.0, kinvisc_cst = 1.0E-06, &
      & smag_coef_mom = 0.1, smag_coef_scal = 0.1, vdifmom_cst = 1.0E-06, &
      & vdifscal_cst = 1.0E-06
!---water depths
REAL :: depmean_cst = 0.0, depmean_flag = 0.0
!---inundation schemes
INTEGER, PARAMETER :: nofldmasks = 11
REAL :: dcrit_fld = 0.1, dmin_fld = 0.02, dthd_fld = 0.1
INTEGER, DIMENSION(nofldmasks) :: fld_mask
!---bottom/surface fluxes
REAL :: bdragcoef_cst = 0.0, bdraglin = 0.0, ccharno = 0.014, &
      & cds_cst = 0.0013, ces_cst = 0.00113, chs_cst = 0.00113, ckar = 0.4, &
      & zbtoz0lim = 2.0, zref_atm = 10.0, zrough_cst = 0.0
!---relaxation distance for momentum advection
REAL :: distrlx_obc = 0.0
!---open boundary conditions
REAL :: cgravratio = 0.03
!---optical parameters
REAL :: optattcoef1_cst = 10.0, optattcoef2_cst = 0.067, opt_frac = 0.54
!---parameters for exchange coefficients in tabular form
INTEGER :: nrelhum, ntemp, nwind
REAL :: drelhum = 0.05, dtempdif = 1.0, dtempmax = 5.0, dtempmin = -5.0, &
      & dwind = 0.25, relhummax = 1.0, relhummin = 0.5, uwindmax = 50.0, &
      & uwindmin = 3.0
!---implicit code
INTEGER :: itsimp = 0, maxitsimp = 1, noitsimp
REAL :: dzetaresid, dzetaresid_conv = 1.0E-14, petsc_tol = 1.0E-07
!---numerical
REAL :: theta_cor = 0.5, theta_vadv = 0.501, theta_vdif = 1.0
REAL, PARAMETER :: eps_adv = 1.0E-12

```

File

physpars.f90

Type

Module

Purpose

Physical and numerical model parameters

Description

atmpres_ref Reference atmospheric pressure P_{ref} [Pa]

bdragcoef_cst	Constant bottom drag coefficient C_{db} when iopt_bstres_drag=1
bdraglin	Bottom friction velocity k_{lin} used in the linear bottom friction law if iopt_bstres_form=1 [m/s]
beta_sal_ref	Reference value for the salinity contraction coefficient β_S [PSU ⁻¹]
beta_temp_ref	Reference value for the temperature expansion coeffi- cient β_T [°C ⁻¹]
b_SH	Parameter b in the Song & Haidvogel (1994) vertical grid transformation
ccharno	Charnock's constant a used in Charnock's relation (4.291)
cds_cst	Constant surface drag coefficient C_{ds} when iopt_sflux_cds=0
ces_cst	Constant surface exchange coefficient C_e when iopt_sflux_cehs=0
cgravratio	Ratio of the internal to the external wave speed (used in the open boundary condition (4.382))
chs_cst	Constant surface exchange coefficient C_h when iopt_sflux_cehs=0
ckar	von Karman's constant κ
dcrit fld	Critical water depth d_{crit} used in the drying/wetting algorithm [m]
density_ref	Reference density ρ_0 [kg/m ³]
depmean_cst	Constant water depth used to set up a default bathy- metry [m]
depmean_flag	Data flag marking land points in the bathymetry [m]
distrx_abc	Maximum distance d_{max} (from the open boundaries) used in the relaxation factor (5.292) for momentum ad- vection
dlat_ref	Reference latitude to be used for the Coriolis frequency in case of a Cartesian grid [degrees]
dlon_ref	Reference longitude to be used in case of a Cartesian grid [degrees, positive East]
dlon_ref_anal	If iopt_astro_anal=1, the harmonically analysed phases are taken with respect to the astronomical argument for this reference longitude. [degrees, positive East]

dlon_ref_abc	If <code>iopt_astro_pars</code> >0, the phases at open boundaries are assumed to be taken with respect to the astronomical argument at this reference value. If zero, the reference longitude is taken at Greenwich. [degrees, positive East]
dl_BB	Parameter d_l in the Burchard & Bolding (2002) vertical grid transformation (4.26)
dmin fld	Minimum water depth d_{min} used in the drying/wetting algorithm [m]
drelhum	Interval of relative humidities taken to calculate the surface exchange coefficients using the Monin-Obukhov theory in tabular form
dtempdif	Interval of air minus sea temperature differences taken to calculate the surface exchange coefficients using the Monin-Obukhov theory in tabular form [$^{\circ}C$]
dtempmax	Maximum air minus sea surface temperature difference taken to calculate the surface exchange coefficients using the Monin-Obukhov theory in tabular form [$^{\circ}C$]
dtempmin	Minimum air minus sea surface temperature difference taken to calculate the surface exchange coefficients using the Monin-Obukhov theory in tabular form [$^{\circ}C$]
dthd fld	Threshold water depth d_{th} used in mask criteria for the inundation scheme [m]
dwind	Interval of wind speeds taken to calculate the surface exchange coefficients using the Monin-Obukhov theory in tabular form [m/s]
du_BB	Parameter d_u in the Burchard & Bolding (2002) vertical grid transformation (4.26)
dzetaresid	Value of $\ \zeta'\ _{\infty}$ (free surface corrector method). Its value is saved at the last iteration until the next time step.
dzetaresid_conv	Threshold value ϵ_{imp} used in the convergence criterium for the outer loop (free surface corrector method)
eps_adv	Tolerance factor for the calculation of the flux ratios with the TVD advection scheme

<code>fld_mask</code>	Enables (0) or disables (1) a specific mask criterium
<code>gacc_mean</code>	Mean acceleration of gravity <ul style="list-style-type: none"> • If <code>gacc_ref</code> is defined, <code>gacc_mean=gacc_ref</code>. • If <code>gacc_ref</code> is undefined and the grid is Cartesian, <code>gacc_mean</code> is defined by the geodetic formula (4.58) at the latitude given by <code>dlat_ref</code>. • If <code>gacc_ref</code> is undefined and the grid is spherical, <code>gacc_mean</code> is defined by the geodetic formula (4.58) applied at each C-node point and averaged over the physical domain.
<code>gacc_ref</code>	If different from <code>real_fill</code> , the acceleration of gravity is taken as horizontally uniform. Otherwise, g is evaluated as function of latitude using (4.58) [m/s ²].
<code>hcrit_SH</code>	Parameter h_{crit} in the Song & Haidvogel (1994) vertical grid transformation
<code>hdifmom_cst</code>	Constant coefficient for horizontal momentum diffusion ν_H when <code>iopt_hdif_coef=1</code> [m ² /s]
<code>hdifscal_cst</code>	Constant coefficient for horizontal scalar diffusion λ_H when <code>iopt_hdif_coef=1</code> [m ² /s]
<code>itsimp</code>	Current iteration number for the outer loop of the implicit scheme
<code>kinvisc_cst</code>	Constant value for the kinematic viscosity [m/s ²]
<code>maxitsimp</code>	Largest allowed iteration number for the outer loop (free surface corrector method)
<code>noitsimp</code>	Last iteration number of the outer loop (free surface corrector method).
<code>nofldmasks</code>	Number of available mask criteria
<code>nrelhum</code>	Number of relative humidities taken to calculate the surface exchange coefficients using the Monin-Obukhov theory in tabular form
<code>ntemp</code>	Number of air minus sea temperature differences taken to calculate the surface exchange coefficients using the Monin-Obukhov theory in tabular form
<code>nwind</code>	Number of wind speeds taken to calculate the surface exchange coefficients using the Monin-Obukhov theory in tabular form

optattcoef1_cst	Inverse optical attenuation depth (λ_1^{-1}) for the absorption of long-wave solar radiation as used in (4.59) [m^{-1}]
optattcoef2_cst	Inverse optical attenuation depth (λ_2^{-1}) for the absorption of short-wave solar radiation as used in (4.59) [m^{-1}]
opt_frac	Long-wave fraction R of the surface solar radiance as used in (4.59)
petsc_tol	Relative tolerance used by PETSc for solving the linear system. (The parameters <code>atol</code> , <code>dtol</code> , <code>maxits</code> used by PETSc in the solution procedure are set to the PETSc defaults).
Rearth	Mean radius of the Earth [m]
relhummax	Maximum relative humidity taken to calculate the surface exchange coefficients using the Monin-Obukhov theory in tabular form
relhummin	Minimum relative humidity taken to calculate the surface exchange coefficients using the Monin-Obukhov theory in tabular form
rho_air	Air mass density ρ_a [kg/m^3]
sal_ref	Reference salinity S_{ref} used if <code>iopt_sal=0</code> or in the linear equation of state (4.108) or as default initial condition [PSU]
sigstar_DJ	Parameter σ_* in the Davies & Jones (1991) vertical grid transformations (4.23) and (4.24)
sig0_DJ	Parameter σ_0 in the Davies & Jones (1991) vertical grid transformations (4.23) and (4.24)
smag_coef_mom	Smagorinsky coefficient C_m for horizontal diffusion of momentum
smag_coef_scal	Smagorinsky coefficient C_s for horizontal diffusion of scalars
speheat	Specific heat of seawater c_p at constant pressure [$\text{J}/\text{kg}/^\circ\text{C}$]
sst_ref	Reference sea surface temperature [$^\circ\text{C}$]
temp_min	Minimum temperature. If set to <code>real_fill</code> , the minimum is taken as the freezing point of sea water which is a function of salinity. [$^\circ\text{C}$]
temp_ref	Reference temperature T_{ref} used if <code>iopt_temp=0</code> or in

	the linear equation of state (4.108) or as default initial condition	[°C]
theta_cor	Implicit factor θ_c for the Coriolis term (between 0.0 and 1.0)	
theta_SH	Parameter θ in the Song & Haidvogel (1994) vertical grid transformation	
theta_vadv	Implicit factor θ_a for vertical advection (between 0.0 and 1.0)	
theta_vdif	Implicit factor θ_d for vertical diffusion (between 0.0 and 1.0)	
uwindmax	Maximum wind speed taken to calculate the surface exchange coefficients using the Monin-Obukhov theory in tabular form	[m/s]
uwindmin	Minimum wind speed taken to calculate the surface exchange coefficients using the Monin-Obukhov theory in tabular form	[m/s]
vdifmom_cst	Constant coefficient for vertical diffusion of momentum used if <code>iopt_vdif_coef=1</code> or as background value if <code>iopt_turb_iwlim=0</code>	[m ² /s]
vdifscal_cst	Constant coefficient for vertical diffusion of scalars used if <code>iopt_vdif_coef=1</code> or as background value if <code>iopt_turb_iwlim=0</code>	[m ² /s]
zbtz0lim	Value of the limiting ratio ξ_{min} for z_b/z_0	
zref_atm	Reference height z_a for meteorological variables (taken by default at 10 m height)	[m]
zrough_cst	Constant bottom roughness length z_0 taken when <code>iopt_bstres_drag=3</code>	[m]
zbtz0lim	Value of the limiting ratio z_b/z_0	[]

33.17 Relaxation zones

```

MODULE relaxation
INTEGER :: norlxzones = 0
INTEGER, DIMENSION(2) :: inodesrlx = 0
INTEGER, DIMENSION(norlxzones) :: idirrlx, iposrlx, ityprlx, &
& jposrlx, ncrllx, nrllx
INTEGER, DIMENSION(ncloc,nrloc,2) :: indexrlxatc, indexrlxatuv

```

REAL, DIMENSION(ncloc,nrloc,2) :: rlxwghtatc, rlxwghtatuv

File

relaxation.f90

Type

Module

Purpose

Definitions for application of relaxation conditions near open boundaries

Description

- | | |
|--------------|--|
| idirrlx | Defines the orientation of each zone
1: western boundary
2: eastern boundary
3: southern boundary
4: northern boundary |
| indexrlxatc | Each C-node grid point within a zone must have an associated boundary location. The array stores both the corresponding zonal index and the (global) index of the corresponding open boundary index. If the grid point is outside all zones, its value is 0. The last array index refers to either U-node (1) or V-node (2) boundaries. |
| indexrlxatuv | Each velocity node grid point within a zone must have an associated boundary location. The array stores both the corresponding zonal index and the (global) index of the corresponding open boundary index. If the grid point is outside all zones, its value is 0. The last array index refers to either U-node (1) or V-node (2) boundaries. |
| inodesrlx | The first element disables (0) or enables (1) the application of relaxation condition for C-node (scalar quantities), the second element for quantities at velocity (U- or V-) nodes. |
| iposrlx | (Global) X-index of the southwest corner of each relaxation zone |
| ityprlx | Type of interpolation (weighting) scheme within the relaxation zones
1: linear |

	2: quadratic
	3: hyperbolic
jposrlx	(Global) Y-index of the southwest corner of each relaxation zone
ncrlx	Sizes of the zones (number of C- or velocity-node grid points) in the X-direction
norlxzones	Number of relaxation zones
nrrlx	Sizes of the zones (number of C- or velocity-node grid points) in the Y-direction
rlxwghtatc	Weight factors used in the interpolation for C-node quantities. If the last index is 1(2), weighting is performed with respect to U-node (V-node) boundaries.
rlxwghtatuv	Weight factors used in the interpolation for U-node or V-node quantities. If the last index is 1(2), weighting is performed with respect to U-node (V-node) boundaries.

33.18 Structures and discharges

```
!---dry cells
```

```
INTEGER :: numdry
```

```
INTEGER, ALLOCATABLE, DIMENSION(numdry) :: idry, jdry
```

```
!---thin dams
```

```
INTEGER :: numthinu, numthinuloc, numthinu, numthinu
```

```
INTEGER, ALLOCATABLE, DIMENSION(numthinu) :: ithinu, jthinu
```

```
INTEGER, ALLOCATABLE, DIMENSION(numthinu) :: ithinu, jthinu
```

```
INTEGER, ALLOCATABLE, DIMENSION(numthinuloc) :: ithinuloc, jthinuloc
```

```
INTEGER, ALLOCATABLE, DIMENSION(numthinu) :: ithinu, jthinu
```

```
!---weirs/barriers/orifices
```

```
INTEGER :: numwbaru, numwbaru, numwbaru, numwbaru
```

```
INTEGER, ALLOCATABLE, DIMENSION(numwbaru) :: iwbaru, jwbaru
```

```
INTEGER, ALLOCATABLE, DIMENSION(numwbaru) :: iwbaru, jwbaru
```

```
INTEGER, ALLOCATABLE, DIMENSION(numwbaru) :: iwbaru, jwbaru
```

```
INTEGER, ALLOCATABLE, DIMENSION(numwbaru) :: iwbaru, jwbaru
```

```
REAL :: wbarrlxu = 1.0, wbarrlxv = 1.0
```

```
REAL, ALLOCATABLE, DIMENSION(:) :: oricoefu, oricoefv, &
```

```
& oriheightu, oriheightv, &
```



```

& orisillu, orisillv
REAL, ALLOCATABLE, DIMENSION(:) :: wbarcoefu, wbarcoefv, wbarcrestu, &
& wbarcrestv, wbarmodlu, wbarmodlv
REAL, ALLOCATABLE, DIMENSION(:) :: wbarelossu, wbarelossv

!---discharges
INTEGER :: numdis, numdisloc, numdisloc_ext
LOGICAL, ALLOCATABLE, DIMENSION(:) : disflag
INTEGER, ALLOCATABLE, DIMENSION(:) :: idis, idisloc, indexdisloc, jdis, &
& jdisloc, kdis, kdistype, mdistype
REAL, ALLOCATABLE, DIMENSION(:) :: disarea, disdir, disspeed, disvol, &
& xdiscoord, ydiscoord, zdiscoord

```

File

structures.f90

Type

Module

Purpose

Model switches

Reference

Chapter 18

Description

disarea	Area over which the discharge takes place	[m ²]
disdir	Discharge direction	[radian]
disflag	.TRUE./FALSE if a discharge locations is located on a wet/dry cell	
disspeed	Discharge speed defined as the volume discharge rate divided by the cell area	[m ²]
disvol	Volume discharge rate	[m ³ /s]
idis	Global X-indices of discharge locations	
idisloc	Local X-indices of discharge locations	
idry	X-indices of dry cells	
indexdisloc	Global indices of the local discharge points	
indexwbaru	Global indices of the local U-node weirs/barriers	
indexwbarv	Global indices of the local V-node weirs/barriers	

<code>ithinu</code>	Global X-indices of thin dams at U-nodes
<code>ithinuloc</code>	Local X-indices of thin dams at U-nodes
<code>ithinv</code>	Global X-indices of thin dams at V-nodes
<code>ithinvloc</code>	Local X-indices of thin dams at V-nodes
<code>iwbaru</code>	Global X-indices of weirs/barriers at U-nodes
<code>iwbaruloc</code>	Local X-indices of weirs/barriers at U-nodes
<code>iwbarv</code>	Global X-indices of weirs/barriers at V-nodes
<code>iwbarvloc</code>	Local X-indices of weirs/barriers at V-nodes
<code>jdis</code>	Global Y-indices of discharge locations
<code>jdisloc</code>	Local Y-indices of discharge locations
<code>jdry</code>	Y-indices of dry cells
<code>jthinu</code>	Global Y-indices of thin dams at U-nodes
<code>jthinuloc</code>	Local Y-indices of thin dams at U-nodes
<code>jthinv</code>	Global Y-indices of thin dams at V-nodes
<code>jthinvloc</code>	Local Y-indices of thin dams at V-nodes
<code>jwbaru</code>	Global Y-indices of weirs/barriers at U-nodes
<code>jwbaruloc</code>	Local Y-indices of weirs/barriers at U-nodes
<code>jwbarv</code>	Global Y-indices of weirs/barriers at V-nodes
<code>jwbarvloc</code>	Local Y-indices of weirs/barriers at V-nodes
<code>kdis</code>	Vertical grid indices of the discharges locations. If zero, discharges are taken as homogeneous over the vertical
<code>kdistype</code>	Selects type of vertical location of the discharge 0: Uniformly distributed over the vertical 1: At the bottom 2: At the surface 3: At a fixed distance from the sea bed 4: At a fixed distance from the sea surface
<code>mdistype</code>	Selects method for flagging of discharge points located on dry cells 0: Locations on dry (C-node) cells are taken as invalid (default) 1: Locations on dry (C-node) cells are moved to the nearest neighbouring wet cell, provided such cell is available

	2: Locations on dry (C-node) cells are moved to the nearest wet cell	
numdis	Global number of discharge locations	
numdisloc	Local number of discharge locations	
numdisloc_ext	Local number of discharge locations including points within the first column of the western and first row of the southern halo	
numdry	Global number of dry cells	
numthinu	Global number of thin dams at U-nodes	
numthinuloc	Local number of thin dams at U-nodes	
numthinu	Global number of thin dams at V-nodes	
numthinu	Local number of thin dams at V-nodes	
numwbaru	Global number of weirs/barriers at U-nodes	
numwbaru	Local number of weirs/barriers at U-nodes	
numwbaru	Global number of weirs/barriers at V-nodes	
numwbaru	Local number of weirs/barriers at V-nodes	
oricoefu	Discharge coefficient for orifices at U-nodes	[m ^{1/2} /s]
oricoefv	Discharge coefficient for orifices at V-nodes	[m ^{1/2} /s]
oriheightu	Orifice width at U-nodes	[m]
oriheightv	Orifice width at V-nodes	[m]
orisillu	Orifice height at U-nodes	[m]
orisillv	Orifice height at V-nodes	[m]
wbarcoefu	Discharge coefficient for weirs/barriers at U-nodes	[m ^{1/2} /s]
wbarcoefv	Discharge coefficient for weirs/barriers at V-nodes	[m ^{1/2} /s]
wbarcrestu	Height of weir crest at U-nodes	[m]
wbarcrestv	Height of weir crest at V-nodes	[m]
wbarelossu	Energy loss sink term at U-node weirs/barriers	[1/s]
wbarelossv	Energy loss sink term at V-node weirs/barriers	[1/s]
wbarmodlu	Modular limit at U-node weirs/barriers	
wbarmodlv	Modular limit at V-node weirs/barriers	
wbarrlxu	Time relaxation coefficient at U-node weirs/barriers	
wbarrlxv	Time relaxation coefficient at V-node weirs/barriers	

xdiscoord	X-coordinates of discharge locations longitude]	[m or degrees
ydiscoord	Y-coordinates of discharge locations latitude]	[m or degrees
zdiscoord	Vertical coordinates (distance from sea bed or surface) of discharge location	[m]

33.19 Model switches

MODULE switches

```

!---grid
INTEGER :: iopt_grid_htype = 1, iopt_grid_nodim = 3, iopt_grid_sph = 0, &
          & iopt_grid_vtype = 1, iopt_grid_vtype_transf = 0
!---interpolation
INTEGER :: iopt_arrint_hreg = 0, iopt_arrint_vreg = 0, iopt_arrint_3D = 0
!---hydrodynamics
INTEGER :: iopt_curr = 2, iopt_curr_wfall = 1, iopt_mode_2D, iopt_mode_3D
!---density
INTEGER :: iopt_dens = 0, iopt_dens_grad = 1, iopt_sal = 0, iopt_sal_sbc = 0, &
          & iopt_temp = 0, iopt_temp_optic = 1, iopt_temp_sbc = 1
!---external modules
INTEGER :: iopt_biolgy = 0, iopt_sed = 0
!---bottom stress
INTEGER :: iopt_bstres_drag = 3, iopt_bstres_form = 2, iopt_bstres_nodim = 3
!---transport
INTEGER :: iopt_transp_full = 0
!---advection
INTEGER :: iopt_adv_scal = 3, iopt_adv_turb = 0, iopt_adv_tvd = 1, &
          & iopt_adv_2D = 1, iopt_adv_3D = 1
!---diffusion
INTEGER :: iopt_hdif_coef = 0, iopt_hdif_scal = 0, iopt_hdif_turb = 0, &
          & iopt_hdif_2D = 0, iopt_hdif_3D = 0, iopt_kinvisc = 0,
          & iopt_vdif_coef = 3
!---turbulence
INTEGER :: iopt_turb_alg = 1, iopt_turb_dis_bbc = 2, iopt_turb_dis_sbc = 2, &
          & iopt_turb_iwlim = 0, iopt_turb_kinvisc = 0, iopt_turb_lmix = 4, &
          & iopt_turb_ntrans = 1, iopt_turb_param = 2, iopt_turb_stab_form = 3,
          & iopt_turb_stab_lev = 1, iopt_turb_stab_mod = 4, &
          & iopt_turb_stab_tke = 2, iopt_turb_tke_bbc = 2, &
          & iopt_turb_tke_sbc = 2

```

```
!---drying/wetting
INTEGER :: iopt_fld = 0
!---structures
INTEGER :: iopt_dischr = 0, iopt_drycel = 0, iopt_thndam = 0, iopt_weibar = 0
!---structures
!---explicit/implicit integration
INTEGER :: iopt_cor_impl = 1, iopt_hydro_impl = 0, iopt_scal_depos = 1, &
          & iopt_vadv_impl = 1, iopt_vdif_impl = 2
!---open boundary conditions
INTEGER :: iopt_obc_advflux = 1, iopt_obc_advrlx = 0, iopt_obc_invbar = 0, &
          & iopt_obc_relax = 0, iopt_obc_sal = 0, iopt_obc_sed = 0, &
          & iopt_obc_bio = 0, iopt_obc_temp = 0, iopt_obc_2D = 0, &
          & iopt_obc_3D = 0
!---astronomical tide
INTEGER :: iopt_astro_anal = 0, iopt_astro_pars = 0, iopt_astro_tide = 0
!---1-D applications
INTEGER:: iopt_sur_1D = 0
!---meteo surface forcing
INTEGER :: iopt_meteo = 0, iopt_meteo_heat = 0, iopt_meteo_salflx = 0, &
          & iopt_meteo_stres = 0
!---surface fluxes
INTEGER :: iopt_sflux_cds = 0, iopt_sflux_cehs = 0, iopt_sflux_strat = 0
!---surface waves
INTEGER :: iopt_waves = 0
!---nesting
INTEGER :: iopt_nests = 0
!---parallel processing (MPI)
INTEGER :: iopt_MPI, iopt_MPI_abort = 0, iopt_MPI_comm_all = 2, &
          & iopt_MPI_comm_coll = 0, iopt_MPI_comm_exch = 2, &
          & iopt_MPI_comm_full = 0, iopt_MPI_comm_gath = 2, &
          & iopt_MPI_comm_scatt = 2, iopt_MPI_partit = 1, iopt_MPI_sync = 0
!---output
INTEGER :: iopt_out_anal = 0, iopt_out_avrgd = 0, iopt_out_tsers = 1
!---netCDF
INTEGER :: iopt_CDF, iopt_CDF_abort = 0, iopt_CDF_fill = 0, &
          & iopt_CDF_format = 1
!---PETSC
INTEGER :: iopt_petsc, iopt_petsc_precond = 5, iopt_petsc_solver = 5
!---verification procedure
INTEGER :: iopt_verif
```

File

switches.f90

Type

Module

Purpose

Model switches

Reference

Section 14.4

Description

<code>iopt_adv_scal</code>	Type of scheme for the advection of scalar quantities 0: advection disabled 1: upwind scheme 2: Lax-Wendroff (explicit) in the horizontal, central (semi-implicit) in the vertical 3: TVD scheme
<code>iopt_adv_turb</code>	Type of scheme for the advection of turbulence quantities. 0: advection disabled 1: upwind scheme 2: Lax-Wendroff (explicit) in the horizontal, central (semi-implicit) in the vertical 3: TVD scheme
<code>iopt_adv_tvd</code>	Type of limiting function for the TVD scheme. 1: superbee limiter 2: monotone limiter
<code>iopt_adv_2D</code>	Type of scheme for the advection of 2-D transports. 0: advection disabled 1: upwind scheme 2: Lax-Wendroff (explicit) in the horizontal, central (semi-implicit) in the vertical 3: TVD scheme

<code>iopt_adv_3D</code>	Type of scheme for the advection of 3-D currents. 0: advection disabled 1: upwind scheme 2: Lax-Wendroff (explicit) in the horizontal, central (semi-implicit) in the vertical 3: TVD scheme
<code>iopt_arrint_hreg</code>	Disables/enables (0/1) the use of non-uniform weighted averages for interpolation in the horizontal of arrays on the model grid.
<code>iopt_arrint_vreg</code>	Disables/enables (0/1) the use of non-uniform weighted averages for interpolation in the vertical of arrays on the model grid.
<code>iopt_arrint_3D</code>	Selects dimension of mask or weight factor in some array interpolations 0: 2-D masks or weights 1: 3-D masks or weights
<code>iopt_astro_anal</code>	Disables/enables (0/1) the use of astronomical arguments for harmonic analysis if <code>iopt_astro_pars > 0</code> and <code>iopt_out_anal = 1</code> .
<code>iopt_astro_pars</code>	Enables or disables the inclusion of astronomical arguments and nodal corrections in the harmonic expansions (4.230) and (4.354). 0: astronomical argument set to zero, nodal factors set to 1, nodal phases set to zero 1: evaluate astronomical phases at a given time and reference longitude, nodal factors are set to 1, nodal phases set to zero 2: evaluate astronomical phases and nodal corrections at a given time and reference longitude
<code>iopt_astro_tide</code>	Disables/enables (0/1) the inclusion of the astronomical tidal force in the momentum equations. This requires that the model uses a spherical grid (<code>iopt_grid_sph=1</code>).
<code>iopt_biolgy</code>	Disables/enables (0/1) the biological module. Currently not implemented.

<code>iopt_bstres_drag</code>	Formulation for the bottom drag coefficient C_{db} . 0: not used 1: spatially uniform value 2: spatially non-uniform 3: using a spatially uniform roughness length 4: using a specified, spatially non-uniform roughness length
<code>iopt_bstres_form</code>	Type of formulation for the bottom stress. 0: bottom stress set to zero 1: linear bottom stress law (4.338) or (4.339) 2: quadratic bottom stress (4.340) or (4.339)
<code>iopt_bstres_nodim</code>	Type of currents in the (linear or quadratic) bottom stress formulation. 2: depth-mean currents 3: 3-D current taken at the bottom grid cell
<code>iopt_CDF</code>	Disables/enables (0/1) the use of <code>netCDF</code> format. The format of all user output is set by default to <code>netCDF</code> ('N') if switched on. This switch cannot be defined by the user, but is set automatically when <code>-DCDF</code> is defined as CPP compiler option.
<code>iopt_CDF_abort</code>	Disables/enables (0/1) automatic abortion of the program when an error is detected within a <code>netCDF</code> call.
<code>iopt_CDF_fill</code>	Disables/enables (0/1) the use of fill values in <code>netCDF</code> files.
<code>iopt_CDF_format</code>	Selects the type <code>netCDF</code> file format. 1: classic format 2: 64-bit offset format
<code>iopt_cor_impl</code>	Time-integration of the Coriolis term. 0: explicit 1: semi-implicit 2: implicit
<code>iopt_curr</code>	Type of current fields (2).

	0: Currents and elevations are set to their default (zero) values and are not updated.
	1: Currents and elevations are initialised but not updated in time.
	2: Currents and elevations are initialised and updated in time.
<code>iopt_curr_wfall</code>	Type of formulation for the settling of particulate matter.
	1: settling enabled without correction terms
	2: settling enabled with the correction terms (7.117)–(7.118) included
<code>iopt_dens</code>	Evaluation of the density and expansion coefficients.
	0: uniform density, zero expansion coefficients
	1: density from the linear equation of state (4.108), expansion coefficients are uniform
	2: from the McDougall <i>et al.</i> (2003) general equation of state (4.103)–(4.107) without pressure effects
	3: from the McDougall <i>et al.</i> (2003) general equation of state (4.103)–(4.107) with pressure effects included
<code>iopt_dens_grad</code>	Selects the numerical algorithm for the discretisation of the baroclinic pressure gradient.
	0: gradient set to zero
	1: traditional σ -coordinate (second order) method
	2: z -level method
	3: method of Shchepetkin & McWilliams (2003)
<code>iopt_disch</code>	Disables/enables (0/1) the discharge module.
<code>iopt_drycel</code>	Disables/enables (0/1) the dry cell module.
<code>iopt_fld</code>	Selects the type of drying/wetting scheme.
	0: Drying/wetting disabled
	1: Drying/wetting algorithm without dynamic masks
	2: Drying/wetting algorithm using dynamic masks

<code>iopt_grid_htype</code>	Type of horizontal grid. 1: uniform rectangular grid 2: non-uniform rectangular grid 3: curvilinear grid
<code>iopt_grid_nodim</code>	Grid dimension. 1: 1-dimensional grid (water column model) 2: 2-dimensional grid (depth-averaged model without vertical structure) 3: 3-dimensional grid
<code>iopt_grid_sph</code>	Type of coordinates. 0: Cartesian coordinates 1: spherical coordinates
<code>iopt_grid_vtype</code>	Type of vertical grid. 1: uniform σ -grid 2: non-uniform σ -grid in the vertical, uniform in the horizontal 3: non-uniform σ -coordinate grid in the horizontal and the vertical
<code>iopt_grid_vtype_transf</code>	Type of vertical grid transformation. 0 : uniform vertical grid (<code>iopt_grid_vtype=1</code>) or user-defined 11: log-transformation (4.23) at the bottom following Davies & Jones (1991) if <code>iopt_grid_vtype=2</code> 12: log-transformation (4.24) at the surface following Davies & Jones (1991) if <code>iopt_grid_vtype=2</code> 13: transformation with enhanced resolution near the bottom and/or the bottom as defined in Burchard & Bolding (2002) 21: Song & Haidvogel (1994) transformation given by (4.33) and (4.35) if <code>iopt_grid_vtype=3</code>
<code>iopt_hdif_coef</code>	Type of scheme for horizontal diffusion coefficients. 0: not used 1: spatially uniform

	2: Smagorinsky formulation (4.80) for momentum and (4.81) for scalars
<code>iopt_hdif_scal</code>	Disables/enables (0/1) horizontal diffusion in the scalar transport equations.
<code>iopt_hdif_turb</code>	Disables/enables (0/1) horizontal diffusion in the turbulence transport equations.
<code>iopt_hdif_2D</code>	Disables/enables (0/1) horizontal diffusion in the 2-D transport equations.
<code>iopt_hdif_3D</code>	Disables/enables (0/1) horizontal diffusion in the 3-D current transport equations.
<code>iopt_hydro_impl</code>	Disables/enables the implicit scheme. 0: The momentum equations are solved with the explicit (mode-splitting) scheme (default). 1: The momentum equations are solved using the implicit algorithm. The compiler option <code>-DPETSC</code> must be set.
<code>iopt_kinvisc</code>	Formulation for kinematic viscosity. 0: user-defined uniform value <code>kinvisc_cst</code> 1: ITTC (1978) relation (7.24)
<code>iopt_meteo</code>	Disables/enables (0/1) meteorological input and evaluation of all surface fluxes.
<code>iopt_meteo_heat</code>	Selects type of input data for the heat fluxes. 0: no input 1: air temperature T_a , relative humidity RH , cloud cover f_c 2: total (downward) non-solar surface heat flux, cloud cover f_c 3: total (downward) non-solar surface heat flux, surface solar radiance Q_{rad} 4: cloud cover f_c 5: surface solar radiance Q_{rad}
<code>iopt_meteo_salflx</code>	Selects type of input data for the salinity flux. 0: no input 1: evaporation minus precipitation rate $E_{vap} - P_{rc}$

	2: precipitation rate P_{rc}
<code>iopt_meteo_stres</code>	Selects type of input data for the barotropic mode, i.e. surface stress and pressure.
	0: no input
	1: components of the wind speed (U_{10}, V_{10}) and (unless <code>iopt_grid_nodim=1</code>) the atmospheric pressure P_a
	2: components of the surface stress (τ_s^u, τ_s^v) and (unless <code>iopt_grid_nodim=1</code>) the atmospheric pressure P_a
<code>iopt_mode_2D</code>	Status of the 2-D mode. Its value is set internally and cannot be changed by the user.
	0: The 2-D mode is disabled. Transports U , V and surface elevations ζ are set to their (zero) default values and are not updated.
	1: Transports and elevation are initialised, but not updated in time
	2: Transports and elevations are initialised and updated in time
<code>iopt_mode_3D</code>	Status of the 3-D mode. Its value is set internally and cannot be changed by the user.
	0: The 3-D current are set to their default (zero) values and are not updated.
	1: The 3-D current is initialised, but not updated in time.
	2: The 3-D current is initialised and updated in time.
<code>iopt_MPI</code>	Disables/enables (0/1) the use of parallel communications. This switch cannot be defined by the user, but is set automatically when <code>-DMPI</code> is defined as CPP compiler option in <code>coherens-flags.cmp</code> .
<code>iopt_MPI_abort</code>	Disables/enables (0/1) automatic abortion of the program when an error is detected within a MPI call.
<code>iopt_MPI_comm_all</code>	Communication type for “all to all” operations.

	1: blocking, standard send
	2: blocking, synchronous send
	3: non-blocking, standard send
	4: non-blocking, synchronous send
<code>iopt_MPI_comm_coll</code>	Disables/enables (0/1) the use of MPI collective calls.
<code>iopt_MPI_comm_exch</code>	Communication type for “exchange” operations.
	1: blocking, standard send
	2: blocking, synchronous send
	3: non-blocking, standard send
	4: non-blocking, synchronous send
	5: send-receive blocking calls
<code>iopt_MPI_comm_full</code>	Disables/enables (0/1) the exchange of 4-D arrays. Currently not implemented.
<code>iopt_MPI_comm_gath</code>	Communication type for “all to one” gather (combine) operations.
	1: blocking, standard send
	2: blocking, synchronous send
	3: non-blocking, standard send
	4: non-blocking, synchronous send
<code>iopt_MPI_comm_scatter</code>	Communication type for “one to all” scatter (distribute and copy) operations.
	1: blocking, standard send
	2: blocking, synchronous send
	3: non-blocking, standard send
	4: non-blocking, synchronous send
<code>iopt_MPI_partit</code>	Selects the method for domain decomposition.
	1: “simple” partition based on the values of <code>nprocsx</code> and <code>nprocsy</code>
	2: decomposition obtained from an external data file or defined in <code>usrdef_partition</code>
<code>iopt_MPI_sync</code>	Disables/enables (0/1) synchronisation calls at the end of a series of blocking or non-blocking operations.

<code>iopt_nests</code>	Disables/enables (0/1) the writing of open boundary data for nested sub-grids.
<code>iopt_obc_advflux</code>	Type of open boundary condition for the cross-stream (2-D and 3-D) advective fluxes (see Section 5.3.16.2) 1: zero gradient condition 2: quasi-upwind scheme
<code>iopt_obc_advrlx</code>	Disables/enables (0/1) the relaxation scheme for horizontal momentum advection (see Section 5.3.16.2) 0: relaxation scheme disabled (default) 1: relaxation scheme enabled. In that case the parameter <code>distrx_obc</code> (representing the parameter d_{max}) must be defined by the user in <code>usrdef_mod_params</code> or in the CIF.
<code>iopt_obc_bio</code>	(General) type of open boundary conditions for biological variables. Currently not implemented. 0: default conditions at all open boundaries 1: non-default conditions for at least one open boundary point
<code>iopt_obc_invbar</code>	Disables/enables (0/1) the inverse barometric effect at open boundaries.
<code>iopt_obc_relax</code>	Disables/enables (0/1) the open boundary relaxation as discussed in Section 4.10.3.
<code>iopt_obc_sal</code>	(General) type of open boundary conditions for salinity. 0: default conditions at all open boundaries 1: non-default conditions for at least one open boundary point
<code>iopt_obc_sed</code>	(General) type of open boundary conditions for sediments. 0: default conditions at all open boundaries 1: non-default conditions for at least one open boundary point
<code>iopt_obc_temp</code>	(General) type of open boundary conditions for temperature.

	0: default conditions at all open boundaries 1: non-default conditions for at least one open boundary point
<code>iopt_obc_2D</code>	(General) type of open boundary conditions for the 2-D mode. 0: default conditions at all open boundaries 1: non-default conditions for at least one open boundary point
<code>iopt_obc_3D</code>	(General) type of open boundary conditions for the 3-D baroclinic currents. 0: default conditions at all open boundaries 1: non-default conditions for at least one open boundary point
<code>iopt_out_anal</code>	Disables/enables (0/1) harmonic output.
<code>iopt_out_avrgd</code>	Disables/enables (0/1) time averaged output.
<code>iopt_out_tsers</code>	Disables/enables (0/1) time series output.
<code>iopt_petsc</code>	Disables/enables PETSC (0/1). This switch is only defined internally when <code>-DPETSC</code> is provided and cannot be reset by the user.
<code>iopt_petsc_precond</code>	Type of preconditioner used by PETSc. For details, see the PETSc User Manual. 1 : Jacobi (PCJACOBI) 2 : Block Jacobi (PCBJACOBI) 3 : SOR (and SSOR) (PCSOR) 4 : SOR with Eisenstat trick (PCEISENSTAT) 5 : Incomplete Cholesky (PCICC) 6 : Incomplete LU (PCILU) 7 : Additive Schwarz (PCASM) 8 : Linear solver (PCKSP) 9 : Combination of preconditioners (PCCOMPOSITE) 10: LU (PCLU) 11: Cholesky (PCCHOLESKY) 12: No preconditioning (PCNONE)

<code>iopt_petsc_solver</code>	Type of solver used by PETSc. For details, see the PETSc User Manual. 1 : Richardson (KSPRICHARDSON) 2 : Chebychev (KSPCHEBYCHEV) 3 : Conjugate Gradient (KSPCG) 4 : Biconjugate Gradient (KSPBICG) 5 : Generalised Minimal Residual (KSPGMRES) 6 : BiCGSTAB (KSPBCGS) 7 : Conjugate Gradient Squared (KSPCGS) 8 : Transpose-Free Quasi-Minimal Residual (1) (KSPTFQMR) 9 : Transpose-Free Quasi-Minimal Residual (2) (KSPTCQMR) 10: Conjugate Residual (KSPCR) 11: Least Squares Method (KSPLSQR) 12: Shell for no KSP method (KSPPREONLY)
<code>iopt_sal</code>	Salinity update. 0: uniform (space and time) salinity field 1: salinity field initialised but not updated in time 2: salinity field initialised and updated in time
<code>iopt_sal_sbc</code>	Type of surface boundary condition for salinity. 0: zero surface flux 1: surface flux given by (4.276)
<code>iopt_scal_depos</code>	Discretisation for the deposition (vertical advective flux at the sea bed) of particulate matter. 0: Deposition flux is set to zero. 1: first order (upwind) scheme 2: second order scheme using extrapolation
<code>iopt_sed</code>	Disables/enables (0/1) the activation of an external sediment module.
<code>iopt_sflux_cds</code>	Formulation for the neutral surface drag coefficient C_{ds} . 0: constant value as given by the parameter <code>cds_cst</code>

	<ul style="list-style-type: none"> 1: equation (4.286) from Large & Pond (1981) 2: equation (4.287) from Smith & Banke (1975) 3: equation (4.288) from Geernaert <i>et al.</i> (1986) 4: equation (4.289) from Kondo (1975) 5: equation (4.290) from Wu (1980) 6: equation (4.291) from Charnock (1955)
<code>iopt_sflux_cehs</code>	<p>Formulation for the neutral surface (heat) exchange coefficients C_e, C_h.</p> <ul style="list-style-type: none"> 0: constant value as given by the parameter <code>ces_cst</code> or <code>chs_cst</code> 1: equation (4.292) from Large & Pond (1982) 2: equation (4.293) from Anderson & Smith (1981) 3: equation (4.294) from Kondo (1975) 4: equation (4.295) from Wu (1980)
<code>iopt_sflux_strat</code>	<p>Selects dependence of the surface drag and exchange coefficients on atmospheric stratification effects.</p> <ul style="list-style-type: none"> 0: no dependence 1: using the Kondo (1975) parameterisation (Section 4.8.2) 2: using the Monin-Obukhov similarity theory (Section 4.8.3)
<code>iopt_sur_1D</code>	<p>Disables/enables surface forcing (surface slopes and elevations) in case of a 1-D (<code>iopt_grid_nodim=1</code>) water column application.</p>
<code>iopt_temp</code>	<p>Temperature update.</p> <ul style="list-style-type: none"> 0: uniform (space and time) temperature field 1: temperature field initialised but not updated in time 2: temperature field initialised and updated in time
<code>iopt_temp_optic</code>	<p>Disables/enables (0/1) the optical module.</p> <ul style="list-style-type: none"> 0: all solar radiation is assumed to be absorbed at the surface, i.e. the water column is considered as opaque

	1: solar radiation is absorbed within the water column using specified values for the attenuation depths
<code>iopt_temp_sbc</code>	Type of surface boundary condition for temperature. <ul style="list-style-type: none"> 1: Neumann using the model's surface heat flux formulations 2: Dirichlet using prescribed surface temperatures taken at the first grid point below the surface 3: Dirichlet using prescribed surface temperature taken at the surface itself
<code>iopt_thndam</code>	Disables/enables (0/1) the thin dam module.
<code>iopt_transp_full</code>	Selects how a series of transport variables are updated in the transport routine. Currently not implemented. <ul style="list-style-type: none"> 0: Each equation is updated separately. 1: The equations are simultaneously updated.
<code>iopt_turb_alg</code>	Type of algebraic scheme if <code>iopt_vdif_coef=2</code> . <ul style="list-style-type: none"> 1: Pacanowski-Philander formulation (4.132)–(4.135) 2: Munk-Anderson formulation (4.136)–(4.140) 3: flow dependent formulation as described in Section 4.4.2.2 with α given by (4.148) 4: flow dependent formulation as described in Section 4.4.2.2 with α given by (4.149) 5: flow dependent formulation as described in Section 4.4.2.2 with α given by (4.150) 6: parabolic profile (4.154)
<code>iopt_turb_dis_bbc</code>	Type of the bottom boundary condition for the dissipation rate ε . <ul style="list-style-type: none"> 1: Neumann condition (4.353) 2: Dirichlet condition (4.351)
<code>iopt_turb_dis_sbc</code>	Type of the surface boundary condition for the dissipation rate ε . <ul style="list-style-type: none"> 1: Neumann condition (4.284) 2: Dirichlet condition (4.281)

<code>iopt_turb_iwlim</code>	Type of the background mixing scheme as described in Section 4.4.3.6. 0: using uniform background coefficients 1: using limiting conditions for turbulence parameters giving the background limits (4.225)–(4.226) 2: the Large <i>et al.</i> (1994) scheme given by (4.227)
<code>iopt_turb_kinvisc</code>	Selects type of background mixing. 0: user-defined constant value <code>vdifmom_cst</code> 1: kinematic viscosity as selected by <code>iopt_kinvisc</code>
<code>iopt_turb_lmix</code>	Mixing length formulation as described in Section 4.4.3.5. 1: parabolic law (4.213) 2: “modified” parabolic law (4.214) 3: “Xing” formulation (4.215) 4: “Blackadar” asymptotic formulation (4.216)
<code>iopt_turb_ntrans</code>	Number of transport equations as described in Section 4.4.3.4. 0: zero-equation model (equilibrium or Mellor-Yamada level 2 method) 1: turbulence energy equation with a mixing length selected by <code>iopt_turb_lmix</code> 2: k - ε of k - kl equation depending on the value of <code>iopt_turb_param</code>
<code>iopt_turb_param</code>	Selects the type of second turbulent variable. 1: mixing length l (k - l scheme) 2: dissipation rate ε (k - ε scheme)
<code>iopt_turb_stab_form</code>	Selects the type of stability function. 1: constant value (4.198) 2: Munk-Anderson form (4.199) 3: from a RANS model as explained in Section 4.4.3.3
<code>iopt_turb_stab_lev</code>	Selects the level for stability functions if <code>iopt_turb_stab_form=3</code> . 1: quasi-equilibrium method (Section 4.4.3.3)

	2: non-equilibrium method (Section 4.4.3.3)
<code>iopt_turb_stab_mod</code>	Selects the type of closure (RANS) model. 1: MY82-model (Mellor & Yamada, 1982) 2: KC94-model (Kantha & Clayson, 1994) 3: BB95-model (Burchard & Baumert, 1995) 4: HR82-model (Hossain & Rodi, 1982) 5: CA01-model (Canuto <i>et al.</i> , 2001) 6: CA02-model (Canuto <i>et al.</i> , 2001)
<code>iopt_turb_stab_tke</code>	Formulation for the turbulent diffusion coefficient ν_k (or stability coefficient S_k) of turbulent energy. 1: constant value for S_k as given by equation (4.200) 2: S_k is taken as proportional to the momentum stability function S_u as given by (4.201) 3: using the formulation of Daly & Harlow (1970) as given by (4.185) or (4.191)
<code>iopt_turb_tke_bcc</code>	Type of the bottom boundary condition for turbulence energy. 1: Neumann condition (4.352) 2: Dirichlet condition (4.351)
<code>iopt_turb_tke_sbc</code>	Type of the surface boundary condition for turbulence energy. 1: Neumann condition (4.283) 2: Dirichlet condition (4.281)
<code>iopt_vadv_impl</code>	Time-integration for vertical advection 0: explicit 1: semi-implicit 2: implicit
<code>iopt_vdif_coef</code>	Selects the (general) type of the vertical diffusion scheme. 0: vertical diffusion disabled 1: uniform diffusion coefficient 2: algebraic formulation as described in Section 4.4.2.2

	3: second order turbulence closure as described in Section 4.4.3
<code>iopt_vdif_impl</code>	Time-integration for vertical diffusion. 0: explicit 1: semi-implicit 2: implicit
<code>iopt_verif</code>	Disables/enables (0/1) the use of the verification procedure (not documented). This switch cannot be defined by the user, but is set automatically when <code>-DVERIF</code> is added as CPP compiler option in <code>coherensflags.cmp</code> . The only effect of the switch is that a different output is produced when running a pre-defined test case.
<code>iopt_waves</code>	Type of wave input wave input. 0: wave input disabled 1: wave height, period and wave direction 2: wave height, period, velocity, excursion and direction
<code>iopt_weibar</code>	Disables/enables (0/1) the weir/barrier module.

33.20 Constants and system parameters

```

MODULE syspars
!--kind parameters
INTEGER, PARAMETER :: knwchar = KIND('A'), knolog = KIND(.TRUE.), &
& kn dint = KIND(1), &
& kn dilong = MAX(4,SELECTED_INT_KIND(10)), &
& kn dreal = KIND(1.0), kn dlong = 8, kn dcplx = 8
!---data types
INTEGER, PARAMETER :: char_type = 1, log_type = 2, int_type = 3, &
& longint_type = 4, real_type = 5, long_type = 6, &
& cplx_type = 7
!---universal parameters
REAL, PARAMETER :: pi = 3.14159265, halfpi = 1.57079633, twopi = 6.28318531
REAL (KIND=kn dlong), PARAMETER :: pi_d = 3.1415926535897932_kn dlong
REAL (KIND=kn dlong), PARAMETER :: halfpi_d = 1.5707963267948966_kn dlong
REAL (KIND=kn dlong), PARAMETER :: twopi_d = 6.2831853071795865_kn dlong

```

```

REAL, PARAMETER :: enap = 2.718282
REAL, PARAMETER :: degtorad = pi/180.0, radtodeg = 180.0/pi
REAL, PARAMETER :: degtorad_d = pi_d/180.0_kndlong, &
    & radtodeg_d = 180.0_kndlong/pi_d
!---tidal parameters
INTEGER, PARAMETER :: MaxAstroTides = 56, MaxConstituents = 77
!---random generators
INTEGER, PARAMETER :: MaxGenerators = 32
!---MPI communications
INTEGER, PARAMETER :: MaxHaloComms = 8
!---model variables
INTEGER, PARAMETER :: MaxModArids = 454, MaxBioArids = 24, MaxSedArids = 51,
    & MaxTotArids = MaxModArids + MaxBioArids + MaxSedArids
!---model I/O
INTEGER, PARAMETER :: MaxCIFTypes = 4, MaxCIFVars = 50, MaxGridTypes = 5, &
    & MaxGridFiles = 2, MaxIOFiles = 33, MaxIOTypes = 39, &
    & MaxProgLevels = 20, MaxRestarts = 10
!---monitoring files and error coding
INTEGER, PARAMETER :: MaxErrCodes = 13, MaxErrMesgs = 50, MaxTimers = 36
!---character string lengths
INTEGER, PARAMETER :: lencifline = 300, lencifvar = 120, lendesc = 120, &
    & lenerrcode = 120, lenformat = 120, lenfreq = 7, &
    & leniofile = 120, lenname = 31, lennode = 3, &
    & lentime = 23, lentitle = 20, lenunit = 60, &
    & lenversion = 10

!---cif file
CHARACTER (LEN=1), PARAMETER :: cifcom = '!', cifend = '#', cifsep = ','

!---user output
LOGICAL, PARAMETER :: DegreesOut = .TRUE.
CHARACTER (LEN=lenversion), PARAMETER :: model_version = 'V2.6'
!---output formats
CHARACTER (LEN=lenformat), PARAMETER :: IntegerFormat='(50I11)',&
    & RealFormat='(50G16.7)'

!---undefined and zero values
LOGICAL, PARAMETER :: log_undef = .FALSE.
INTEGER, PARAMETER :: int_undef = -2147483647, izero = 0
INTEGER (KIND=kndilong), PARAMETER :: izero_d = 0_kndilong, &
    & longint_undef = -2147483647_kndilong
REAL, PARAMETER :: real_undef = -9.9692099683868690E+36, &

```

```

& real_flag = -9.9692099683868690E+35, rzero = 0.0
REAL (KIND=kndlong), PARAMETER :: rzero_d = 0.0_kndlong
CHARACTER (LEN=lentime), PARAMETER :: cdatetime_undef = &
& 'xxxx/xx/xx;00:00:00,000'

```

File

syspars.f90

Type

Module

Purpose

Constants and other system parameters. The **Max*** parameters are mostly intended for the dimensioning of arrays which cannot be declared as allocatable.

Description

DegreesOut	Determines the unit of phases and angles in user-defined output files .TRUE. degrees .FALSE. radians
IntegerFormat	Format string used for reading/writing integer data from/to a ASCII ('A') file in standard COHERENS format
MaxAstroTides	Maximum allowed number of constituents for the astronomical forcing
MaxBioArids	Maximum available number of key ids for biological array variables.
MaxCIFTypes	Maximum number of CIF files
MaxCIFVars	Maximum number of data variables on a CIF line
MaxConstituents	Maximum allowed number of tidal constituents at open boundaries
MaxErrCodes	Maximum number of key ids for error messages
MaxErrMesgs	Default maximum number of error messages
MaxGenerators	Maximum available number of random generators which can be simultaneously used in the program
MaxGridFiles	Maximum number of surface grid files for each grid type. Current value is 1.

MaxGridTypes	Maximum number of surface grid types
MaxHaloComms	Maximum available number of exchange communications (send or receive)
MaxIOFiles	Maximum number of forcing files for a given file descriptor key id.
MaxIOTypes	Maximum number of file descriptor key ids.
MaxModArids	Maximum available number of key ids for physical model array variables
MaxProgLevels	Maximum number of subprogram levels
MaxRestarts	Maximum number of output times for the writing of restart conditions
MaxSedArids	Maximum available number of key ids for sediment array variables.
MaxTimers	Maximum available number of timers
MaxTotArids	Maximum available number of key ids for all model arrays
RealFormat	Format string used for reading/writing real data from/to a ASCII ('A') file in standard COHERENS format
cdatetime_undef	Flag for an undefined date/time string
char_type	Type parameter for character variables
cifcom	Comment character on a CIF line
cifend	Character marking the end of a data block in a CIF
cifsep	Data separator character within a CIF line
cmplx_type	Type parameter for complex variables
degtorad	Factor to convert degrees to radians
degtorad_d	Factor to convert degrees to radians in double precision
enap	Euler's number $e = 2.718282$
halfpi	The number $\pi/2$
halfpi_d	The number $\pi/2$ in double precision
int_type	Type parameter for integer variables
int_undef	Flag for undefined or invalid integer values
izero	The number zero in single precision
izero_d	The number zero in long integer format

<code>kndchar</code>	Kind parameter for character variables
<code>kndcmplx</code>	Kind parameter for complex variables
<code>kndilong</code>	Kind parameter for long integer variables
<code>kndint</code>	Kind parameter for integer variables
<code>kndlog</code>	Kind parameter for logical variables
<code>kndlong</code>	Kind parameter for double precision real variables
<code>kndreal</code>	Kind parameter for real variables
<code>lencifline</code>	Maximum length of a data line in a CIF
<code>lencifvar</code>	Maximum length of a CIF data value in string format
<code>lendesc</code>	Maximum length of a <code>long_name</code> attribute
<code>lenerrcode</code>	Maximum length of an error code message
<code>lenformat</code>	Maximum length of a string format specification
<code>lenfreq</code>	Maximum length of a frequency name
<code>leniofile</code>	Maximum length for the name of a file in standard COHERENS format
<code>lenname</code>	Maximum length of a <code>f90_name</code> attribute
<code>lennode</code>	Maximum length of a <code>node</code> attribute
<code>lentime</code>	Length of a date/time string
<code>lentitle</code>	Maximum length of a simulation title
<code>lenunit</code>	Maximum length of the <code>units</code> attribute
<code>lenversion</code>	Maximum length of the <code>model_version</code> string
<code>log_type</code>	Type parameter for logical variables
<code>log_undef</code>	Flag for undefined or invalid logical data
<code>longint_type</code>	Type parameter for a long integer variable
<code>longint_undef</code>	Flag for undefined or invalid long integer values
<code>long_type</code>	Type parameter for double precision real variables
<code>model_version</code>	Current COHERENS version number in string format
<code>pi</code>	The number π
<code>pi_d</code>	The number π in double precision
<code>radtodeg</code>	Factor to convert radians to degrees
<code>radtodeg_d</code>	Factor to convert radians to degrees in double precision

<code>real_flag</code>	Flag below which input data are considered as invalid (default value). Actual parameter used in the program is <code>real_min</code> .
<code>real_type</code>	Type parameter for real variables
<code>real_undef</code>	Default flag for undefined or invalid real values. Actual parameter used in the program is <code>real_fill</code> .
<code>rzero</code>	The number zero in single precision real format
<code>rzero_d</code>	The number zero in double precision real format
<code>twopi</code>	The number 2π
<code>twopi_d</code>	The number 2π in double precision

33.21 Tidal forcing

```

MODULE tide
!---number of tidal constituents
INTEGER :: nconastro = 0, nconobc = 0
!---tidal indices
INTEGER, DIMENSION(MaxAstroTides) :: index_astro = 0
INTEGER, DIMENSION(MaxConstituents) :: index_obc = 0
!---nodal factors
REAL, DIMENSION(nconastro) :: fnode_astro
REAL, DIMENSION(nconobc) :: fnode_obc
!---tidal phases
REAL, DIMENSION(nconastro) :: phase_astro
REAL, DIMENSION(nconobc) :: phase_obc
!---tidal force
REAL, DIMENSION(ncloc,nrloc) :: fxastro, fyastro
!---key ids for tidal constituents
INTEGER, PARAMETER :: &
    & icon_MS0 = 1, icon_Sa = 2, icon_Ssa = 3, icon_058 = 4, &
    ...
!---tidal species index
INTEGER, DIMENSION(MaxConstituents) :: ispec_tides
DATA ispec_tides /12*0, 21*1, 22*2, 1*3, 1*2, 4*3, 5*4, 6*6, 5*8/
!---amplitudes of the equilibrium tide
REAL, DIMENSION(MaxAstroTides) :: astro_ampl = &
    & (/0.198419, 0.003103, 0.019542, 0.001142, 0.004239, 0.022191, 0.003677, &
    ..., 0.003455/)
!---amplitude corrections for the earth tide

```

```

REAL, DIMENSION(MaxAstroTides) :: astro_earth
DATA astro_earth /15*0.693, 0.6946, 0.6948, 0.6950, 0.6956, 0.693, 0.6962, &
  ..., 24*0.693/
!---tidal frequencies
REAL, DIMENSION(MaxConstituents) :: tidal_spectrum = &
&( /0.0000000E+00, 1.9909688E-07, 3.9821277E-07, 5.9730965E-07, &
  & ..., 5.8177642E-04/)
!---names of tidal frequencies
CHARACTER (LEN=lenfreq), DIMENSION(MaxConstituents) :: tidal_freq_names = &
  & (/ 'MS0' , 'Sa' , 'Ssa' , '058' , 'Msm' , 'Mm' , &
  & ..., 'S8' '/')

```

File

tide.f90

Type

Module

Purpose

Parameters and arrays for tidal forcing

Description

<code>astro_amp</code>	Amplitudes of the astronomical equilibrium tide [m]
<code>astro_earth</code>	Elasticity factor α_{qn} representing the effect of the Earth tides on the tidal force
<code>fnode_astro</code>	Nodal factors of the constituents used for the astronomical tidal force
<code>fnode_obc</code>	Nodal factors of the tidal constituents used at open boundaries
<code>fxastro</code>	X-component of the astronomical tidal force [m ² /s]
<code>fyastro</code>	Y-component of the astronomical tidal force [m ² /s]
<code>icon_*</code>	Key ids of all tidal constituents
<code>index_astro</code>	Key ids of the constituents used in the calculation of the astronomical force
<code>index_obc</code>	Key ids of the constituents used at the open boundaries
<code>ispec_tides</code>	Tidal species indices q for each frequency
<code>nconastro</code>	Number of tidal constituents used in the calculation of the astronomical force

nconbc	Number of tidal constituents used at open boundaries
phase_astro	Astronomical phases of the constituents used for the astronomical tidal force, at Greenwich. Nodal corrections are included if <code>iopt_astro_pars=2</code> . [rad]
phase_obc	Astronomical phases of the constituents used at open boundaries with respect to the reference longitude <code>dlon_ref_obc</code> . Nodal corrections are included if <code>iopt_astro_pars=2</code> . [rad]
tidal_freq_names	Names of all pre-defined tidal frequencies
tidal_spectrum	Frequencies of all pre-defined tidal constituents [rad/s]

33.22 Time parameters

```

MODULE timepars
LOGICAL :: corrstep, metstepin, predstep, physinit, wavestep
CHARACTER (LEN=lentime) :: CDateTime, CEndDateTime = cdatetime_undef,&
                        & CStartDateTime = cdatetime_undef, ClockTime
INTEGER :: julianday, norestarts = 0, nstep = 0, nt = 0, ntobcrlx = 0
INTEGER :: ic3d = 1, icnodal = 0
INTEGER (KIND=kndilong) :: nosecsrun = 0
REAL :: delt2d, delt3d, time_zone = 0.0
INTEGER, DIMENSION(6) :: time_convert = (/1,60,3600,86400,2629800,31557600/)
INTEGER, DIMENSION(7) :: IDateTime, IEndDateTime, IStartDateTime
INTEGER, DIMENSION(12) :: days_in_month = &
                        & (/31,28,31,30,31,30,31,31,30,31,30,31/)
INTEGER, DIMENSION(13) :: monthdays = &
                        & (/0,31,59,90,120,151,181,212,243,273,304,334,365/)
INTEGER, DIMENSION(MaxRestarts) :: ntrestart = 0

```

File

timepars.f90

Type

Module

Purpose

Time parameters

Description

CDateTime Current date and time in string format

CEndDateTime	End date and time in string format
CStartDateTime	Start date and time in string format
ClockTime	Date and time from the machine's internal real-time clock at the start of the simulation
IDateTime	Current date and time in integer vector format
IEndDateTime	End date and time in integer vector format
IStartDateTime	Start date and time in integer vector format
corrstep	Set to <code>.TRUE.</code> at the corrector time steps
days_in_month	Number of days in each month (without correction for leap years)
delt2d	Barotropic (2-D) time step (mode-splitting scheme) or time step for all transport equations (implicit scheme) [s]
delt3d	Baroclinic (3-D) time step, equal to <code>delt2d</code> in case of an implicit scheme [s]
icnodal	Time step (measured in units of <code>delt2d</code>) for an update of the nodal tidal factors and astronomical arguments if <code>iopt_astro_pars > 0</code> . If zero, nodal corrections (amplitudes and phases) are evaluated at the initial time only.
ic3d	Ratio of the 3-D (baroclinic) and 2-D time steps
julianday	Julian day (year day between 1 and 365 or 366)
metstepin	Set to <code>.TRUE.</code> at time steps when an update of surface fluxes is required.
monthdays	Day number of the first day in each month (without correction for leap years)
norestarts	Number of times for the writing of re-start initial conditions
nosecsrun	Number of seconds since the start of the simulation
nstep	Number of 2-D time steps
nt	Time index (time since the start of the simulation divided by <code>delt2d</code>)
ntobcrlx	The relaxation period T_r , divided by the 2-D time step <code>delt2d</code> , (optionally) used to define the relaxation factor $\alpha_r(t)$ (see equation (4.356)) for the 2-D mode at open boundaries.

ntrestart	The times during a simulation, measured in time indices, when re-start conditions have to be written
predstep	Set to <code>.TRUE.</code> at the predictor time steps
physinit	Set to <code>.TRUE.</code> for (re)initialisation of physical conditions
time_convert	Factors for converting the time in seconds to another unit
time_zone	Time zone, i.e. the difference of the local time with respect to GMT. Difference is positive (negative) eastwards (westwards) from Greenwich. [hours]
wavestep	Set to <code>.TRUE.</code> at time steps when an update of surface wave data is required.

33.23 Turbulence model parameters

```

MODULE turbpars
!---stability functions
INTEGER :: ib22
REAL :: cfequil(5), cfstabtke(3), cfstab1(8), cfstab2(11)
REAL :: c_sk = 0.15, keps0, f0stabmom, f0stabscal, f0stabtke, &
      & skeps = 0.09, sq_my = 0.2
!---limiting conditions
REAL :: alphaM_max, alphaN_max, alphaN_min, alphaN_sl, &
      & dissipmin = 1.0E-012, tkelim = 1.0E-06, tkemin = 1.0E-014, &
      & zlmixmin = 1.7E-010
!---t.k.e.-equation
REAL :: sigma_k = 1.0, wfltke = 0.0
!---kl-equation
REAL :: e1_my = 1.8, e2_my = 1.33, e3_my = 1.0, sigma_kl
!---eps-equation
REAL :: c1_eps = 1.44, c2_eps = 1.92, c31_eps = 0.2, c32_eps = 1.0, sigma_eps
!---roughness lengths
REAL :: zrough_bot = 0.0, zrough_sur = 0.0
!---KPP scheme for internal waves
REAL :: riccrit_iw = 0.7, vdifmom_iw = 1.0E-04, vdifscal_iw = 5.0E-05, &
      & vdifshear_iw = 0.005
!---Pacanowski-Philander relations
REAL :: alpha_pp = 5.0, expmom_pp = 2.0, vbmom_pp = 1.0E-04, &
      & vbscal_pp = 1.0E-05, vmax_pp = 3.0, v0dif_pp = 0.01

```

```

!---Munk-Anderson relations
REAL :: alpha_ma = 10.0, beta_ma = 3.33, expmom_ma = 0.5, &
      & expscal_ma = 1.5, & vmaxmom_ma = 3.0, vmaxscal_ma = 4.0, v0dif_ma = 0.06
!---algebraic flow-dependent relations
REAL :: cnu_ad = 2.0, delta1_ad = 0.0, delta2_ad = 0.0, k1_ad = 0.0025, &
      & k2_ad = 2.0E-05, lambda_ad = 0.0, omega1_ad = 1.0E-04, &
      & r1_ad = 1.0, r2_ad = 1.0
!---algebraic mixing length formulations
REAL :: alpha_Black = 0.2, beta_Xing = 2.0

```

File

turbpars.f90

Type

Module

Purpose

Parameters for turbulence schemes

Description

<code>alpha_Black</code>	constant α_1 in the Blackadar (1962) mixing length formulation (4.217)
<code>alphaM_max</code>	maximum value for the stability parameter α_M in case of stable stratification when the the non-equilibrium method is applied
<code>alphaN_max</code>	maximum value for the stability parameter α_N in case of stable stratification
<code>alphaN_min</code>	minimum value for the stability parameter α_N in case of unstable stratification
<code>alphaN_sl</code>	slope of the limiting curve for α_N in the case of a stable stratification
<code>alpha_ma</code>	parameter α_m in the Munk & Anderson (1948) scheme (4.136)–(4.139)
<code>alpha_pp</code>	parameter α_p in the Pacanowski & Philander (1981) scheme (4.132)–(4.134)
<code>beta_ma</code>	parameter β_m in the Munk & Anderson (1948) scheme (4.136)–(4.139)
<code>beta_Xing</code>	attenuation factor β_1 in the Xing & Davies (1996) mixing length formulation (4.215)

cfequil	coefficients C_d used in the equilibrium formulation (4.194–4.195) for the stability functions
cfstabtke	coefficients used in the non-equilibrium (4.185) or the quasi-equilibrium (4.191) method for the stability function S_k
cfstab1	coefficients C_b or C_c used in the quasi-equilibrium formulation (4.187) or (4.190) for the stability functions
cfstab2	coefficients C_a used in the non-equilibrium formulation (4.182) for the stability functions
cnu_ad	parameter C_ν in equation (4.151)
c1_eps	constant $c_{1\varepsilon}$ in the shear production term of the ε -equation (4.205)
c2_eps	constant $c_{2\varepsilon}$ in the dissipation term of the ε -equation (4.205)
c31_eps	constant $c_{3\varepsilon}$ in the buoyance sink term of the ε -equation (4.205) in case of stable stratification ($N^2 > 0$)
c32_eps	constant $c_{3\varepsilon}$ in the buoyancy source term of the ε -equation (4.205) in case of unstable stratification ($N^2 < 0$)
c_sk	Daly-Harlow parameter c_{sk} in (4.177)
delta1_ad	parameter δ_1 in equation (4.143)
delta2_ad	parameter δ_2 in equation (4.143)
dissipmin	numerical lower limit ε_{min} for ε [W/kg]
eps0	parameter ϵ_0 as defined by (4.282)
expmom_ma	parameter n_1 in the Munk & Anderson (1948) scheme (4.136)–(4.139)
expmom_pp	parameter n_p in the Pacanowski & Philander (1981) scheme (4.132)–(4.134)
expscal_ma	parameter n_2 in the Munk & Anderson (1948) scheme (4.136)–(4.139)
e1_my	constant E_1 in the shear production term of the kl -equation (4.209)
e2_my	constant E_2 in the wall proximity term (4.210) of the kl -equation (4.209)
e3_my	constant E_3 in the buoyancy source/sink term of the kl -equation (4.209)

f0stabmom	neutral stability coefficient S_{u0} for momentum (quasi-equilibrium or equilibrium method)
f0stabsca	neutral stability coefficient S_{b0} for scalars (quasi-equilibrium or equilibrium method)
f0stabtke	constant stability coefficient S_{k0} for turbulent kinetic energy
ib22	Equals 1 when the parameter $c_{22\beta} \neq 0$, 0 otherwise
k1_ad	parameter K_1 in equations (4.148) and (4.150)
k2_ad	parameter K_2 in equation (4.149)
lambda_ad	parameter λ_* in equation (4.146) [m]
omega1_ad	parameter ω_1 in equation (4.151) [s ⁻¹]
riccrit_iw	critical Richardson number Ri_0 in the Large <i>et al.</i> (1994) background mixing scheme (4.227)
r1_ad	parameter r_1 in equation (4.143)
r2_ad	parameter r_2 in equation (4.143)
sigma_eps	the parameter σ_ϵ as obtained from (4.206)
sigma_k	parameter σ_k used to define S_k in (4.201)
sigma_kl	the parameter σ_{kl} obtained from (4.211)
skeps	neutral value S_{k0} of the stability coefficient S_k in the k - ϵ model (see equation (4.200))
sq_my	parameter S_q used to determine S_{k0} in the Mellor-Yamada model (see equation (4.202))
tkelim	background limit k_{lim} for k (see equation (4.226)) [J/kg]
tkemin	numerical lower limit k_{min} for k [J/kg]
vbmom_pp	parameter ν_{bp} in the Pacanowski & Philander (1981) scheme (4.132)–(4.134) [m ² /s]
vbscal_pp	parameter λ_{bp} in the Pacanowski & Philander (1981) scheme (4.132)–(4.134) [m ² /s]
vdifmom_iw	internal wave breaking diffusion coefficient ν_{T0} for momentum in the Large <i>et al.</i> (1994) background mixing scheme (4.227) [m ² /s]
vdifscal_iw	internal wave breaking diffusion coefficient λ_{T0} for scalars in the Large <i>et al.</i> (1994) background mixing scheme (4.227) [m ² /s]

<code>vdifshear_iw</code>	maximum mixing due to resolved vertical shear ν_0^s in the Large <i>et al.</i> (1994) background mixing scheme (4.227) [m ² /s]
<code>vmaxmom_ma</code>	parameter ν_{max} in the Munk & Anderson (1948) scheme (4.136)–(4.139)
<code>vmaxscal_ma</code>	parameter λ_{max} in the Munk & Anderson (1948) scheme (4.136)–(4.139)
<code>vmax_pp</code>	parameter ν_{max} in the Pacanowski & Philander (1981) scheme (4.132)–(4.134)
<code>v0dif_ma</code>	parameter ν_{0m} in the Munk & Anderson (1948) scheme (4.136)–(4.139) [m ² /s]
<code>v0dif_pp</code>	parameter ν_{0p} in the Pacanowski & Philander (1981) scheme (4.132)–(4.134) [m ² /s]
<code>wfltkc</code>	surface wave factor c_w used in the surface flux condition (4.283) for turbulent energy
<code>zlmixmin</code>	numerical lower limit l_{min} for l [m]
<code>zrough_bot</code>	bottom roughness length z_{0b} in the mixing length formulation (4.212) [m]
<code>zrough_sur</code>	surface roughness length z_{0s} in the mixing length formulation (4.212) [m]

33.24 Turbulence arrays

```

MODULE turbulence
REAL, DIMENSION(1-nhalo:ncloc+nhalo,1-nhalo:nrloc+nhalo,nz+1) :: &
    & dissip, tke, tke_old, zlmix
REAL, DIMENSION(ncloc,nrloc,nz+1) :: buofreq2, shearfreq2

```

File

turbulence.f90

Type

Module

Purpose

Turbulence arrays

Description

`buofreq2` Squared buoyancy frequency N^2 [1/s²]

dissip	Dissipation of turbulent kinetic energy	[W/kg]
tke	Turbulent kinetic energy	[J/kg]
tke_old	Turbulent kinetic energy at the old 3-D time step	[J/kg]
shearfreq2	Squared shear frequency M^2	[1/s ²]
zlmix	Turbulent mixing length	[m]

33.25 Surface wave arrays

MODULE waves

REAL, ALLOCATABLE, DIMENSION(ncloc,nrloc) :: wavedir, wavefreq

REAL, ALLOCATABLE, DIMENSION(0:ncloc,0:nrloc) :: waveexcurs, waveheight, &
& wavenum, waveperiod, waveuvel, wavevel, wavevvel

File

waves.f90

Type

Module

Purpose

Surface wave arrays

Description

wavedir	Wave direction	[rad]
wavexcurs	Near-bottom wave excursion amplitude	[m]
wavefreq	Peak wave frequency	[rad/s]
waveheight	Significant wave height	[m]
wavenum	Wave number	[1/m]
waveperiod	Peak wave period at	[s]
waveuvel	Near-bottom wave orbital velocity in the X-direction	[m/s]
wavevel	Near-bottom wave orbital velocity	[m/s]
wavevvel	Near-bottom wave orbital velocity in the Y-direction	[m/s]

