ICON Namelist Overview

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1 ICON Namelists

1.1 Scripts, Namelist files and Programs

Run scripts starting the programs for the grid generation and the models are stored in run/. These scripts we the specified Fortran namelists. Programs are stored in <icon home>/build/<architecture>/bin/.

Table 1: Namelist files

N	Namelist file	Purpose	Made by script	Used b
N	NAMELIST_GRAPH	Generate graphs	create_global_grids.run	grid_c
N	$NAMELIST_GRID$	Generate grids	$create_global_grids.run$	$\operatorname{grid} \underline{} c$
N	$NAMELIST_GRIDREF$	Gen. nested domains	create_global_grids.run	$\operatorname{grid} \underline{} c$
N	NAMELIST_OCEAN_GRID	Gen. ocean grid	create_ocean_grid.run	$\operatorname{grid} \underline{} c$
N	NAMELIST_TORUS_GRID	Gen. torus grid	create_torus_grid.run	grid_c
N	NAMELIST ICON	Run ICON models	exp. <name>.run</name>	control

1.2 Namelist parameters

The following subsections tabulate all available Fortran namelist parameters by name, type, default value, u

- Type refers to the type of the Fortran variable, in which the value is stored: I=INTEGER, L=LOGICA string
- Default is the preset value, if defined, that is assigned to this parameter within the programs.
- *Unit* shows the unit of the control parameter, where applicable.
- Description explains in a few words the purpose of the parameter.
- Scope explains under which conditions the namelist parameter has any effect, if its scope is restricted namelist parameters.

Information on the file, where the namelist is defined and used, is given at the end of each table.

2 Namelist parameters for grid generation

2.1 Namelist parameters defining the atmosphere grid

2.1.1 graph_ini (NAMELIST_GRAPH)

Parameter	Type	Default	Unit	Description
nroot	I	2		root subdivision of initial edges
grid_levels	I	4		number of edge bisections following the root subdivision
lplane	L	.FALSE.		switch for generating a double periodic planar grant The root level consists of 8 triangles.
				The root level consists of a triangles.

Defined and used in: src/grid_generator/mo_io_graph.f90

2.1.2 grid_ini (NAMELIST_GRID)

Parameter	Type	Default	Unit	Description
nroot	I	2		root subdivision of initial edges
grid_levels	Ι	4		number of edge bisections following the root subdivision
lplane	L	.FALSE.		switch for generating planar grid. The root level consists of 8 triangles.

Defined and used in: $src/grid_generator/mo_io_grid.f90$

2.1.3 grid_options (NAMELIST_GRID)

Parameter	Type	Default	Unit	Description
x_rot_angle	R	0.0	deg	Rotation of the icosahedron about the x-axis
				(connecting the origin and [0°E, 0°N])

Parameter	Type	Default	Unit	Description
y_rot_angle	R	0.0	deg	Rotation of the icosahedron about the y-axis
				(connecting the origin and [90°E, 0°N), done after
				the rotation about the x-axis.
z_rot_angle	R	0.0	deg	rotation of the icosahedron about the z-axis
				(connecting the origin and [0°E, 90°N), done after
				the rotation about the y-axis.
itype_optimize	I	4		Grid optimization type
				0: no optimization
				1: Heikes Randall
				2: equal area
				3: c-grid small circle
				4: spring dynamics
l_c_grid	L	.FALSE.		C-grid constraint on last level
maxlev_optim	I	100		Maximum grid level where the optimization is
				applied
beta spring	R	0.90		tuning factor for target grid length

Defined and used in: $src/grid_generator/mo_io_grid.f90$

2.1.4 plane options (NAMELIST GRID)

Parameter	Type	Default	Unit	Description
tria arc km	R	10.0	km	length of triangle edge on plane

The number of grid points is generated by root level section and further bisections. The double periodic root The spatial coordinates are -1 <= x <= 1, and $-\sqrt{3}/2 <= y <= \sqrt{3}/2$. Currently the planar option can obefined and used in: $src/grid_generator/mo_io_grid.f90$

2.1.5 gridref ini (NAMELIST GRIDREF)

Parameter	Type	Default	Unit	Description

Parameter	Type	Default	Unit	Description
grid_root	Ι	2		root subdivision of initial edges
start_lev	I	4		number of edge bisections following the root
				subdivision
n_dom	Ι	2		number of logical model domains, including the
				global one
n_phys_dom	Ι	n_dom		number of physical model domains, may be larger
				than n_dom (in this case, domain merging is
				applied)
parent_id	I(n_phys_	i		ID of parent domain (first entry refers to first
	dom-1)			nested domain; needs to be specified only in case
				more than one nested domain per grid level)
logical_id	$I(n_{phys}$	i+1		logical grid ID of domain (first entry refers to firs
	dom-1)			nested domain; needs to be specified only in case
				domain merging, i.e. n_dom < n_phys_dom)
l_plot	L	.FALSE.		produces GMT plots showing the locations of the
				nested domains
l_circ	L	.TRUE.		Create circular (.T.) or rectangular (.F.) refined
				domains
l_rotate	L	.FALSE.		Rotates center point into the equator in case of 1 circ = .FALSE.
write hierarchy	Ι	1		0: Output only computational grids
write_merarchy	1	1		1: Output only computational grids 1: Output in addition parent grid of global mode
				domain (required for computing physics on a
				reduced grid)
				2: Output all grids back to level 0 (required for
				hierarchical search algorithms)
bdy_indexing_depth	, T	max rlcell		Number of cell rows along the lateral boundary o
bdy_indexing_depti	1 1	(=8)		model domain for which the refin ctrl fields
		(-0)		contain the distance from the lateral boundary;
				needs to be enlarged when lateral boundary
				nudging is required for one-way nesting
				madeling is required for one way nesting

	Parameter	Type	Default	Unit	Description
Ī	radius	R(n_dom-	30.	deg	radius of nested domain (first entry refers to first
		1)			nested domain; needs to be specified for each nest
					domain separately)
Ī	hwidth_lon	R(n_dom-	20.	deg	zonal half-width of refined domain (first entry ref
		1)			to first nested domain; needs to be specified for
					each nested domain separately)
Ī	hwidth_lat	R(n_dom-	20.	deg	meridional half-width of refined domain (first ent
		1)			refers to first nested domain; needs to be specified
					for each nested domain separately)
Ī	center_lon	R(n_dom-	90.	deg	center longitude of refined domain (first entry ref
		1)			to first nested domain; needs to be specified for
					each nested domain separately)
Ī	center_lat	R(n_dom-	30.	deg	center latitude of refined domain (first entry refer
		1)			to first nested domain; needs to be specified for
					each nested domain separately)

Defined and used in: $src/grid_generator/mo_gridrefinement.f90$

2.2 Namelist parameters defining the local grid generation

The ocean grids are created by the script ${\tt run/create_ocen_grid.run}$

${\bf 2.2.1 \quad grid_geometry_conditions}$

Parameter	Type	Default	Unit	Description
no_of_conditions	I	0		Number of geometric conditions
patch_shape	I(no_of_ condi-	0		1=rectangle; 2=circle
	tions)			
patch_center_x	R(no_of condi-	0.0	degrees	longitude of patch center
	tions)			

Parameter	Type	Default	Unit	Description
patch_center_y	R(no_of	0.0	degrees	latitude of patch center
	_ condi-			
	tions)			
rectangle_xradious	R(no_of_	0.0	degrees	half meridional extension of a rectangular patch
	condi-			
	tions)			
rectangle_yradious	R(no_of_	0.0	degrees	half zonal extension of a rectangular patch
	condi-			
	tions)			
circle_radious	R(no_of_	0.0	degrees	radius of a circular patch
	condi-			
	tions)			

Defined in mo_grid_conditions.f90

${\bf 2.2.2}\quad {\bf local_grid_optimization}$

Parameter	Type	Default	Unit	Description
$use_optimization$	L	.FALSE.		Apply, or not, optimization
use_edge_springs	L	.FALSE.		Use spring dynamics
prime_ref_length _coeff	R	1.0		Spring length coefficient
use_adaptive_ spring_length	L	.FALSE.		Use adaptive spring length
use_local_reference _length	L	.FALSE.		Use locally adaptive spring length
local_reference_ length_coeff	R	0.0		Coefficient of local vs global spring length
use_isotropy_force	L	.FALSE.		Use isotropy force, tends to create symmetric triangles
isotropy_rotation coeff	R	0.0		Coefficient of the rotational isotropy force

Parameter	Type	Default	Unit	Description
isotropy_stretch	R	0.0		Coefficient of the stretch isotropy force
_coeff				
optimize_vertex	I	1		For patches the min depth of the vertices that wi
_depth				be optimized. The boundary vertices have depth
				the next level 1, etc.

 $Defined \ in \ {\tt mo_local_grid_optimization.f90}$

${\bf 2.2.3}\quad {\bf create_ocean_grid}$

Parameter	Type	Default	Unit	Description
only_get_sea_	L	.false.		.true.:returns the whole grid with a sea-land mas
land_mask				.false.:returns only the ocean grid
smooth_ocean_	L	.true.		.true.:smooths the ocean boundaries so no triabg
boundary				has two boundary edges; .false.:no smoothing
input_file	С			name of the input grid file
elevation_file	С			name of the file containing cell elevation values for
				the input_file
elevation_field	С			name of the field containing the cell elevation value
min_sea_depth	R	0.0	m	if cell elevation < min_sea_depth then the cell is
			(nega-	consider sea
			tive)	
set_sea_depth	R	0.0	m	if not 0, then sea cells are of set_sea_depth
			(nega-	elevation
			tive)	
set_min_sea_depth	R	0.0	m	if not 0, then sea cells have a maximum of
			(nega-	set_min_sea_depth elevation
			tive)	
edge_elev_	I	2		compute edge elevation from cells using: linear
interp_method				interpolation=1; min value = 2
output_refined_	С			name of the output refined ocean grid file
ocean file				

2.2.4 torus grid parameters

Parameter	Type	Default	Unit	Description
y_no_of_rows	I		4	number of triangle rows of the torus grid
x_no_of_columns	I		8	number of triangle columns of the torus grid
edge_length	R	m	1000.0	the triangle edge length
x_center	R	m	0.0	the x coordinate of the torus center
y_center	R	m	0.0	the y coordinate of the torus center
out_file_name	С			the torus grid file name
unfolded_torus_	С			the unfolded torus grid file name (for plotting)
file_name				
ascii_filename	С			the unfolded torus grid ascci file name (for plotting

Defined in mo_create_torus_grid.f90. See the run script run/create_torus_grid.run.

3 Namelist parameters defining the ICON model

Namelist parameters for the ICON models are organized in several thematic Fortran namelists controling properties of dynamics, transport, physics etc.

3.1 master nml

Parameter	Type	Default	Unit	Description
l_restart	L	.FALSE.		If .TRUE.: Current experiment is started from a
				restart.

3.2 master model nml (reapeated for each model)

Parameter	Type	Default	Unit	Description
model_name	С			Character string for naming this component.
model_namelist_	С	1		File name containing the model namelists.
filename	l'	[<u>'</u>	·	
model_type	I	0		Identifies which component to run. atmosphere=
	l'	[<u>'</u>	·	ocean=2, radiation=3, dummy_model=99
model_min_rank	I	0		Start MPI rank for this model.
model_max_rank	I	-1		End MPI rank for this model.
model_inc_rank	I	0		Stride of MPI ranks.
model_restart_info	С	restart.info		Name (including full path) of the restart info file
filename				this model

$3.3 \quad time_nml$

Parameter	Type	Default	Unit	Description
dt_restart	R	86400.*30.	S	Length of restart cycle in seconds. Note that the
				frequency of writing restart files is controlled by
				io nml:dt checkpoint. If the value of
				dt checkpoint resulting from model default or
				user's specification is longer than dt_restart, it w
				be reset (by the model) to dt_restart so that at
				least one restart file is generated during the resta
				cycle. If dt_restart is larger than but not a
				multiple of dt_checkpoint, restart file will NOT l
				generated at the end of the restart cycle.
calendar	I	1		Calendar type:
				0=Julian/Gregorian
				1=proleptic Gregorian
				$2{=}30 \mathrm{day/month}, 360 \mathrm{day/year}$
ini_datatime_string	С	'2008-09-		Initial date and time of the simulation
		01T00:00:0	þΖ'	
end_datatime_string	g C	2008-09-		End date and time of the simulation
		01T01:40:0	þΖ'	

Parameter	Type	Default	Unit	Description
				Length of the run
				If "nsteps" in run nml (see below) is positive, th
				nsteps*dtime is used to compute the end date an
				time of the run.
				Else the initial date and time, the end date and
				time, dt restart, as well as the time step are use
				to compute "nsteps".

3.4 parallel nml

Parameter	Type	Default	Unit	Description
nproma	I	1		chunk length
n_ghost_rows	I	1		number of halo cell rows
division_method	I	1		method of domain decomposition
				0: read in from file
				1: use built-in geometric subdivision
				2: use METIS
division_file_name	С			Name of division file
p_test_run	L	.FALSE.		.TRUE. means verification run for MPI
				parallelization (PE 0 processes full domain)
l_test_openmp	L	.FALSE.		if .TRUE. is combined with p_test_run=.TRUE
				and OpenMP parallelization, the test PE gets or
				1 thread in order to verify the OpenMP
				paralllelization
l_log_checks	L	.FALSE.		if .TRUE. messages are generated during each
				synchonization step (use for debugging only)
l_fast_sum	L	.FALSE.		if .TRUE., use fast (not
				processor-configuration-invariant) global summati
iorder_sendrecv	I	1		Sequence of send/receive calls: $1 = \text{irecv/send}$; 2
				isend/recv: $3 = isend/irecv$

Parameter	Type	Default	Unit	Description
itype_comm	I	1		1: use local memory for exchange buffers
				2: use global memory for exchange buffers
				3: asynchronous halo communication for dynamic
				core (NH tria only)
num_io_procs	I	0		Number of I/O processors (running exclusively fo
				doing I/O)
pio_type	I	1		Type of parallel I/O. Only used if number of I/O
				processors greater number of domains.
				Experimental!
nh_stepping_thread	s I	1		The number of OpenMP threads to be used by the
				non-hydrostatic dycore. Only used if the
				OMP_RADIATION flag is set during
				compilation. Experimental!
radiation_threads	I	1		The number of OpenMP threads to be used by the
				radiation. Only used if the
				OMP_RADIATION flag is set during
				compilation. Experimental!
use_icon_comm	L	.FALSE.		Enable the use of MPI bulk communication throu
				the icon_comm_lib
icon_comm_debug	L	.FALSE.		Enable debug mode for the icon_comm_lib
max_send_recv	I	131072		Size of the send/receive buffers for the
_buffer_size				icon_comm_lib.

Defined and used in: $src/namelists/mo_parallel_nml.f90$

3.5 coupling_nml

Parameter	Type	Default	Unit	Description
name	С	blank		short name of the coupling field
dt_coupling	I	0	s	coupling time step / coupling interval
dt model	I	0	s	model time step

Parameter	Type	Default	Unit	Description
lag	I	0		offset to coupling event in number of model time
				steps
l_time_average	L	.FALSE.		.TRUE.: time averaging between two coupling
]			events
l_time_accumulation	n L	.FALSE.		.TRUE.: accumulation of coupling fields in time
]			between two coupling events
l_diagnostic	L	.FALSE.		.TRUE.: simple diagnostics (min, max, avg) for
]			coupling fields is switched on
l_activated	L	.FALSE.		.TRUE.: activate the coupling of the respective
]			coupling field

Defined and used in: $src/namelists/mo_coupling_nml.f90$

$3.6 \quad run_nml$

Parameter	Type	Default	Unit	Description
ldump_states	L	.FALSE.		Dump patch/interpolation/grid refinement state
				every patch (after subdivision in case of a paralle
				run) to a Netcdf file and exit program.
lrestore_states	L	.FALSE.		Restore patch/interpolation/grid refinement state
				from NetCDF dump files instead of calculating
				them.
l_one_file_per_pat	chL	.FALSE.		Use one file per patch for all processors.
				This will decrease the amount of files used for
				dump/restore considerably, especially for massive
				parallel runs on hundreds or thousands of
				processors.
				Time for dumping will increase since the file has
				be written sequentially, the time for restore shoul
				stay roughly the same, however.

Parameter	Type	Default	Unit	Description
ldump_dd	L	.FALSE.		Dump the domain decomposition (and a few relat
				fields). This can be done either in a parallel run
				in a single-CPU run. When done in a parallel run
				the domain decoposition is for the number of
				parallel processes in use. When done in a
				single-CPU run, nproc_dd (see below) determine
				the number of processes for the decomposition.
				Uses always only one file per patch,
lread_dd	L	.FALSE.		Read the domain decomposition when dumped w
				ldump_dd.
nproc_dd	I	1		Number of processors for the target domain
				decomposition (only relevant when running on a
				single processor).
nsteps	I	0		number of time steps of this run.
dtime	R	600.0	s	time step
ltestcase	L	.TRUE.		Idealized testcase runs
ldynamics	L	.TRUE.		Compute adiabatic dynamic tendencies
iforcing	I	0		Forcing of dynamics and transport by
				parameterized processes. Use positive indices for
				the atmosphere and negative indices for the ocean
				0: no forcing
				1: Held-Suarez forcing
				2: ECHAM forcing
				3: NWP forcing
				4: local diabatic forcing without physics
				5: local diabatic forcing with physics
				-1: MPIOM forcing (to be done)
ltransport	L	.FALSE.		Compute large-scale tracer transport
ntracer	I	0		Number of advected tracers handled by the
				large-scale transport scheme
lvert_nest	L	.FALSE.		If set to .true. vertical nesting is switched on (i.e.
				variable number of vertical levels)
num_lev	I(max_d	om)31		Number of full levels (atm.) for each domain

Parameter	Type	Default	Unit	Description
nshift	I(max_don	n)0		vertical half level of parent domain which coincid
				with upper boundary of the current domain
ltimer	L	.TRUE.		TRUE: Timer for monitoring thr runtime of spec
				routines is on $(FALSE = off)$
timers_level	Ι	1		
activate_sync_timer	sL	F		TRUE: Timer for monitoring runtime of
				communication routines (FALSE = off)
msg_level	Ι	10		controls how much printout is written during
				runtime.
				For values less than 5, only the time step is written

Defined and used in: $src/namelists/mo_run_nml.f90$

$3.7 \quad \mathrm{grid_nml}$

	Parameter	Type	Default	Unit	Description
	cell_type	I	3		Cell type
					3: triangular cells
					4: quadrilateral cells (to be done)
					6: pentagonal/hexagonal cells
	lplane	L	.FALSE.		planar option
	corio_lat	R	0.0	deg	Center of the f-plane is located at this geographic
					latitude
Ī	l_limited_area	L	.FALSE.		
Ī	lfeedback	L(n_dom)	.TRUE.		Specifies if feedback to parent grid is performed.
					Setting lfeedback(1)=.false. turns off feedback for
					all nested domains; to turn off feedback for select
					nested domains, set lfeedback(1)=.true. and set
					".false." for the desired model domains
Ī	ifeedback_type	I	1		1: incremental feedback
					2: relaxation-based feedback

Parameter	Type	Default	Unit	Description
patch_weight	R(n_dom)	0.		If patch_weight is set to a value > 0 for any of the first level child patches, processor splitting will be performed, i.e. every of the first level child patched gets a subset of the total number or processors corresponding to its patch_weight. A value of 0. corresponds to exactly 1 processor for this patch, regardless of the total number of processors. For the set of the set of the total number of processors.
	T	DALCE		root patch and higher level childs, patch_weight not used. However, patch_weight must be set to for these patches to avoid confusion.
lredgrid_phys	L	.FALSE.		If set to .true. is calculated on a reduced grid (= one grid level higher)
dynamics_grid_ filename	С			Array of the grid filenames to be used by the dycore.
dynamics_parent_ grid_id	Ι			Array of the indexes of the parent grid filenames, described by the dynamics_grid_filename array. Indexes start at 1, an index of 0 indicates no pare
radiation_grid_ filename	С			Array of the grid filenames to be used for the radiation model. Filled only if the radiation grid different from the dycore grid.
dynamics_radiation _grid_link	I			Array of the indexes linking the dycore grids, as described by the dynamics_grid_filename array, and the radiation_grid_filename array. It provid the link index of the radiation_grid_filename, fo each entry of the dynamics_grid_filename array. Indexes start at 1, an index of 0 indicates that the radiation grid is the same as the dycore grid. On needs to be filled when the radiation_grid_filename is defined.

Defined and used in: $src/namelists/mo_grid_nml.f90$

$3.8 \quad \mathrm{gridref_nml}$

Parameter	Type	Default	Unit	Description
grf_intmethod_c	I	2		Interpolation method for grid refinement (cell-base
				dynamical variables):
				1: parent-to-child copying
				2: gradient-based interpolation
grf_intmethod_ct	I	2		Interpolation method for grid refinement (cell-base
				tracer variables):
				1: parent-to-child copying
				2: gradient-based interpolation
grf_intmethod_e	I	4		Interpolation method for grid refinement
				(edge-based variables):
				1: inverse-distance weighting (IDW)
				2: RBF interpolation
				3: combination gradient-based / IDW
				4: combination gradient-based / RBF
grf_velfbk	I	1		Method of velocity feedback:
				1: average of child edges 1 and 2
				2: 2nd-order method using RBF interpolation
grf_scalfbk	I	2		Feedback method for dynamical scalar variables
				(T, p_{sfc}) :
				1: area-weighted averaging
				2: bilinear interpolation
grf_tracfbk	I	2		Feedback method for tracer variables:
_				1: area-weighted averaging
				2: bilinear interpolation
grf_idw_exp_e12	R	1.2		exponent of generalized IDW function for child
				edges $1/2$
grf_idw_exp_e34	R	1.7		exponent of generalized IDW function for child
				edges $3/4$
rbf vec kern grf e	ı I	1		RBF kernel for grid refinement (edges):
				1: Gaussian
				$2: 1/(1+r^2)$

Parameter	Type	Default	Unit	Description
				3: inverse multiquadric
rbf_scale_grf_e	R	0.5		RBF scale factor for grid refinement (edges)
denom_diffu_t	R	135		Deniminator for lateral boundary diffusion of
				temperature
denom_diffu_v	R	200		Deniminator for lateral boundary diffusion of
				velocity

Defined and used in: src/namelists/mo gridref nml.f90

3.9 prepicon nml

Remark: prepicon_nml contains switches controlling the real-data initialization functionality of ICON. The using it:

- Using the preprocessing tool prep_icon, it is possible to
 - (a) generate the three-dimensional coordinate fields needed by IFS2ICON if IFS2ICON is requeste vertical interpolation from the IFS grid to the ICON grid
 - (b) convert the hydrostatic set of variables provided by IFS2ICON to the nonhydrostatic set of equa and
 - (c) perform the vertical interpolation to the ICON grid if IFS2ICON is requested to do only the horiz
- If ICONAM (iequations=3) is combined with NWP physics (iforcing=3), setting ltestcase=.false. acti running the ICON executable.

Parameter	Type	Default	Unit	Description
i_oper_mode	I	1		Operating mode if the prep_icon executable is ru
				1: generate coordinate fields
				2: convert IFS2ICON output to NH prognostic
				variables
				3: do vertical interpolation
nlev_in	I	91		number of model levels of input data
nlevsoil_in	I	4		number of soil levels of input data
zpbl1	R	500.0	m	bottom height (AGL) of layer used for gradient
				computation

Parameter	Type	Default	Unit	Description
zpbl2	R	1000.0	m	top height (AGL) of layer used for gradient
				computation
l_w_in	L	.FALSE.		Logical switch if vertical wind is provided as input
l_sfc_in	L	.TRUE.		Logical switch if surface fields are provided as inp
				(mandatory when inwp_surface >0)
l_zp_in	L	.FALSE.		Logical switch for diagnostic output on pressure
				and height levels
l_extdata_out	L	.FALSE.		Logical switch to write extdata fields into output

Defined and used in: $src/namelists/mo_prepicon_nml.f90$

$3.10 \quad interpol_nml$

Parameter	Type	Default	Unit	Description
llsq_lin_consv	L	.FALSE.		conservative (T) or non-conservative (F)
				least-squares reconstruction for 2nd order (linear)
				transport
llsq_high_consv	L	.TRUE.		conservative (T) or non-conservative (F)
				least-squares reconstruction for high order transp
lsq_high_ord	I	3		polynomial order for high order reconstruction
				2: quadratic
				30: cubic (no 3^{rd} order cross deriv.)
				3: cubic
rbf_vec_kern_c	I	1		Kernel type for reconstruction at cell centres:
				1: Gaussian
				3: inverse multiquadric
rbf_vec_kern_e	I	3		Kernel type for reconstruction at edges:
				1: Gaussian
				3: inverse multiquadric
rbf_vec_kern_v	I	1		Kernel type for reconstruction at vertices:
				1: Gaussian
				3. inverse multiquadric

Parameter	Type	Default	Unit	Description
rbf_vec_scale_c	R(n_dom)	resolution- dependent		Scale factor for RBF reconstruction at cell centre
1.6	D(1)	-		
rbf_vec_scale_e	R(n_dom)	resolution-		Scale factor for RBF reconstruction at edges
		dependent		
rbf_vec_scale_v	R(n_dom)	resolution-		Scale factor for RBF reconstruction at vertices
		dependent		
nudge max coeff	R	0.02		Maximum relaxation coefficient for lateral
				boundary nudging
nudge_efold_width	R	2.5		e-folding width (in units of cell rows) for lateral
				boundary nudging coefficient
nudge_zone_width	I	8		Total width (in units of cell rows) for lateral
				boundary nudging zone
i_cori_method	I	3		Selector for tangential wind reconstruction method
				1: Almut's method for tangential wind, but PV
				usage as in TRSK
				2: method of Thuburn, Ringler, Skamarock and
				Klemp (TRSK)
				3: Almut's method for tangential wind and PV
				usage
l corner vort	L	.TRUE.		switch whether the rhombus averaged corner
				vorticity is averaged to the hexagon (.TRUE.) or
				the rhombi are directly averaged to the hexagon
				(.FALSE.)
				()

Defined and used in: $src/namelists/mo_interpol_nml.f90$

$3.11 \quad dynamics_nml$

This namelist is relevant if run_nml:ldynamics=.TRUE.

Parameter	Type	Default	Unit	Description

Parameter	Type	Default	Unit	Description
iequations	I	1		Equations and prognostic variables. Use positive
				indices for the atmosphere and negative indices for
				the ocean.
				0: shallow water model
				1: hydrostatic atmosphere, T
				2: hydrostatic atm., θ -dp
				3: non-hydrostatic atmosphere
				-1: hydrostatic ocean
idiv_method	I	1		Method for divergence computation:
				1: Standard Gaussian integral. Hydrostatic
				atm. model: for unaveraged normal components,
				Non-hydrostatic atm. model: for averaged norma
				components
				2: bilinear averaging of divergence
divavg_cntrwgt	R	0.5		Weight of central cell for divergence averaging
sw_ref_height	R	0.9*2.94e4	gm	Reference height of shallow water model used for
				linearization in the semi-implicit time stepping
				scheme
lcoriolis	L	.TRUE.		Coriolis force

Defined and used in: $src/namelists/mo_dynamics_nml.f90$

3.12 ha_dyn_nml

 $This \ name list \ is \ relevant \ if \ run_nml: ldynamics=. TRUE. \ and \ dynamics_nml: ie quations=IHS_ATM_TEMPLE and larger and leaves a support of the larger and larger a$

Parameter	Type	Default	Unit	Description
itime_scheme	I	4		Time integration scheme:
				11: pure advection (no dynamics)
				12: 2 time level semi implicit (not yet implemente
				13: 3 time level explicit
				14: 3 time level with semi implicit correction
				15: standard 4th-order Runge-Kutta method
				(4-stage)

Parameter	Type	Default	Unit	Description
				16: SSPRK(5,4) scheme (5-stage)
ileapfrog_startup	I	1		How to integrate the first time step when the
				leapfrog scheme is chosen. 1 = Euler forward; 2 =
				series of sub-steps.
asselin_coeff	R	0.1		Asselin filter coefficient
si_2tls	R	0.6		weight of time step $n+1$. Valid range: $[0,1]$
si_expl_scheme	I	2		scheme for the explicit part used in the 2 time levels
				semi-implicit time stepping scheme. $1 = \text{Euler}$
				forward; $2 = Adams$ -Bashforth 2nd order
si_cmin	R	30.0	m/s	semi implicit correction is done for eigenmodes w
				speeds larger than si_cmin
si_coeff	R	1.0		weight of the semi implicit correction
si_offctr	R	0.7		
si_rtol	R	1.0e-3		relative tolerance for GMRES solver
lsi_3d	L	.FALSE.		3D GMRES solver or decomposistion into 2D
				problems
ldry_dycore	L	.TRUE.		Assume dry atmosphere
lref_temp	L	.FALSE.		Set a background temperature profile as base star
				when computing the pressure graident force

3.13 nonhydrostatic_nml (relevant if run_nml:iequations=3)

Parameter	Type	Default	Unit	Description
itime_scheme	I	4		Time integration scheme:
				3: same as default, but computation of velocity
				tendencies in corrector step only
				4: Matsuno scheme
				6: same as default, but usage of velocity tendenci
				at $(nnow+nnew)/2$
rayleigh_coeff	R(n_dom)	0.05		Rayleigh damping coefficient (Klemp, Dudhia,
				Hassiotis: MWR136, pp.3987-4004)

Parameter	Type	Default	Unit	Description
damp_height	R(n dom)	30000	m	Height at which Rayleigh damping of vertical wir
	` - '			starts
htop_moist_proc	R	200000.0	m	Height above which moist physics and advection
				cloud and precipitation variables are turned off
htop_qvadv	R	250000.0	m	Height above which QV advection is turned off (
				not use except for debugging purposes)
hbot_qvsubstep	R	250000.0	m	Height above which QV is advected with
				substepping scheme (do not use except for
				debugging purposes)
k2_updamp_coeff	R	2.0e6		enhanced 2nd order diffusion coefficient in upper
				damping layer
vwind_offctr	R	0.05		Off-centering in vertical wind solver
ivctype	I	1		Type of vertical coordinate:
				1: Gal-Chen hybrid
				2: SLEVE (uses sleve_nml)
iadv_rcf	I	1		reduced calling frequency (rcf) for transport
				1: no rcf (every dynamics-step)
				2: transport every 2. step
				4:
l_nest_rcf	L	.TRUE.		Synchronize interpolation/feedback calls with
				advection (transport) time steps. l_nest_rcf is
				automatically reset to .FALSE. if iadv_rcf=1
l_masscorr_nest	L	.FALSE.		Apply mass conservation correction also in nested
				domain
iadv_rhotheta	I	2		Advection method for rho and rhotheta:
				1: centred differences horiz. + vert.
				2: 2nd order Miura horizontal
				3: 3rd order Miura horizontal (not recommended

Parameter	Type	Default	Unit	Description
igradp_method	I	1		Discretization of horizontal pressure gradient:
				1: conventional discretization with metric
				correction term
				2: Taylor-expansion-based reconstruction of
				pressure (advantageous at very high resolution)
				3: Similar discretization as option 2, but uses
				hydrostatic approximation for downward
				extrapolation over steep slopes
l_zdiffu_t	L	.FALSE.		.TRUE.: Compute Smagorinsky temperature
				diffusion truly horizontally over steep slopes
thslp_zdiffu	R	0.025		Slope threshold above which truly horizontal
msip_zumu	16	0.025		temperature diffusion is activated
				temperature diffusion is activated
thhgtd zdiffu	R	200	m	Threshold of height difference between neighboring
				grid points above which truly horizontal
				temperature diffusion is activated (alternative
				criterion to thslp_zdiffu)
exner_expol	R	0.5		Temporal extrapolation (fraction of dt) of Exner
				function for computation of horizontal pressure
				gradient
l_open_ubc	L	.FALSE.		.TRUE.: Use open upper boundary condition
				(rather than w=0) to better conserve sea-level
				pressure in the presence of diabatic heating
ltheta_up_hori	L	.FALSE.		upstream biased horizontal advection for theta (s
		1.0		also upstr_beta)
upstr_beta	R	1.0		Selection of order for horiz. theta advection: 3rd
		1.0.0		order=1.0, 4th order=0.0
gmres_rtol_nh	R	1.0e-6		relative tolerance for convergence in gmres solver

Defined and used in: $src/namelists/mo_nonhydrostatic_nml.f90$

$3.14 \quad sleve_nml \; (relevant \; if \; nonhydrostatic_nml:ivctype=2)$

Parameter	Type	Default	Unit	Description
min_lay_thckn	R	50	m	Layer thickness of lowermost layer
top_height	R	23500.0	m	Height of model top
stretch_fac	R	1.0		Stretching factor to vary distribution of model
				levels; values <1 increase the layer thickness near
				the model top
decay_scale_1	R	4000	m	Decay scale of large-scale topography component
decay_scale_2	R	2500	m	Decay scale of small-scale topography component
decay_exp	R	1.2		Exponent of decay function
flat_height	R	16000	m	Height above which the coordinate surfaces are fl

Defined and used in: src/namelists/mo_sleve_nml.f90

3.15 diffusion_nml

Parameter	Type	Default	Unit	Description
lhdiff_temp	L	.TRUE.		Diffusion on the temperature field
lhdiff_vn	L	.TRUE.		Diffusion on the horizontal wind field
hdiff_order	I	4		Order of ∇ operator for diffusion:
				-1: no diffusion
				2: ∇^2 diffusion
				3: Smagorinsky ∇^2 diffusion for the hexagonal
				model (includes frictional heating if
				lhdiff_temp=.TRUE.)
				4: ∇^4 diffusion
				5: Smagorinsky ∇^2 diffusion combined with ∇^4
				background diffusion as specified via
				hdiff_efdt_ratio
				defaults: 2 for hexagonal model, 4 for triangular
				model

Parameter	Type	Default	Unit	Description
				24 or 42: $\nabla 2$ diffusion from model top to a certai level (cf. k2_pres_max and k2_klev_max below
				∇^4 for the lower levels.
k2_pres_max	R	-99.	Pa	Pressure level above which $ abla^2$ diffusion is applied
k2_klev_max	I	0		Index of the vertical level till which (from the motop) ∇^2 diffusion is applied. If a positive value is specified for k2_pres_max, k2_klev_max is researched accordingly during the initialization of a model respectively.
hdiff_efdt_ratio	R	1.0		ratio of e-folding time to time step (or 2* time st when using a 3 time level time stepping scheme) (only for triangles currently)
hdiff_min_efdt_rati	ioR	1.0		minimum value of hdiff_efdt_ratio near model to
hdiff_tv_ratio	R	1.0		Ratio of diffusion coefficients for temperature and normal wind: $T: v_n$
hdiff_multfac	R	1.0		Multiplication factor of normalized diffusion coefficient for nested domains
hdiff_smag_fac	R	0.15		Scaling factor for Smagorinsky diffusion

Defined and used in: $src/namelists/mo_diffusion_nml.f90$

3.16 io_nml

Parameter	Type	Default	Unit	Description
no_output	L	.FALSE.		Main switch for turning off 'old' output

Parameter	Type	Default	Unit	Description
out_expname	С	'IIIEEEET	TTT'	Outfile basename
out_filetype	I	2		Type of output format:
				1: GRIB1 (not yet implemented)
				2: netCDF
lkeep_in_sync	L	.FALSE.		Sync output stream with file on disk after each
				timestep
dt_data	R	21600.0	S	Output time interval
dt_diag	R	86400.		diagnostic integral output interval
dt_file	R	2592000	S	Time interval of triggering new output file
dt_checkpoint	R	2592000	S	Time interval for writing restart files. Note that i
				the value of dt_checkpoint resulting from model
				default or user's specification is longer than
				time_nml:dt_restart, it will be reset (by the
				model) to dt_restart so that at least one restart :
				is generated during the restart cycle.
lwrite_vorticity	L	.TRUE.		write out averaged vorticity at vertices
lwrite_initial	L	.TRUE.		write out initial state
lwrite_dblprec	L	.FALSE.		write out double precision
lwrite_oce_timestep		.FALSE.		write out intermediate ocean vars
lwrite_divergence	L	.TRUE.		write out divergence at cells
lwrite_omega	L	.TRUE.		write out vertical velocity in pressure coords.
lwrite_pres	L	.TRUE.		write out full level pressure
lwrite_z3	L	.TRUE.		write out geopotential on full levels
lwrite_tracer	L(ntracer)	.TRUE.		write out tracer at cells
lwrite_tend_phy	L	.TRUE.		Physics induced tendencies.
		.FALSE.		
		(Scope)		
lwrite_radiation	L	.FALSE.		Radiation related fields.

Parameter	Type	Default	Unit	Description
lwrite_precip	L	.FALSE.		Precipitation
lwrite_cloud	L	.FALSE.		Cloud variables
lwrite_tke	L	.TRUE.		TKE
lwrite_surface	L	.FALSE.		surface variables
lwrite_extra	L	.FALSE.		debug fields
inextra_2d	I	0		Number of 2D Fields for diagnostic/debugging
				output.
inextra_3d	I	0		Number of 3D Fields for diagnostic/debugging output.
lwrite pzlev	L	.FALSE.	+	activate output on p- and/or z-levels
lflux avg	L	.FALSE.		if .FALSE. the output fluxes are accumulated
_ ~				from the beginning of the run
				if .TRUE. the output fluxes are average values
				from the beginning of the run, except of
				TOT PREC that would be accumulated

Defined and used in: $src/namelists/mo_io_nml.f90$

3.17 output_nml

Please note: There may be several instances of output_nml in the namelist file, every one defining a list attributes for output.

Parameter	Type	Default	Unit	Description
filetype	I	4		One of CDI's FILETYPE_XXX constants.
				Possible values: 2 (=FILETYPE_GRB2), 4
				(=FILETYPE_NC2), 5 (=FILETYPE_NC4)
namespace	С	, ,		'DWD' - DWD short names (or 'MPIM', 'CMIP'
				'ECMWF')
				Currently unused.
				RJ: For what exactly should that be used?
map_file	С	, ,		File containing the mapping from internal names t
				names written to NetCDF.
				The format of this file:
				One mapping per line, first the internal name, the
				the name written to NetCDF, separated by an
				arbitrary number of blanks. The line may also st
				and end with an arbitrary number of blanks.
				Empty lines or lines starting with $\#$ are treated a
				comments.
				Names not covered by the mapping are output as
				they are.
				Note that the specification of output variables, e.
				in ml_varlist, is independent from this renamin
mode	I	1		1 = forecast mode, $2 = $ climate mode
				Currently unused.
				RJ: For what exactly should that be used?
dom(:)	I	-1		Array of domains for which this name-list is used
				If not specified (or specified as -1 as the first arra
				member), this name-list will be used for all
				domains.
				Attention: Depending on the setting of the
				parameter l_output_phys_patch these are either
				logical or physical domain numbers!
output_time_unit	I	1		1 = second, 2=minute, 3=hour, 4=day, 5=month
				6=year

		T		
Parameter	Type	Default	Unit	Description
$output_bounds(3,:)$	R	None		post-processing times in units defined by
				output_time_unit: start, end, increment. There
				may be specified several triples (up to 100) which
				must be in increasing order.
steps_per_file	I	100		Max number of output steps in one output file. I
				this number is reached, a new output file will be
				opened.
include_last	L	.TRUE.		Flag whether to include the last time step
output_grid	L	.FALSE.		Flag whether grid information is output (in
				NetCDF output)
output_filename	С	None		Output filename prefix (which may include path)
				Domain number, level type, file number and
				extension will be added, so specifying 'XXX' for
				output_filename you will end up in a name like
				XXX_DOM01_ML_0001.nc
lwrite_ready	L	.FALSE.		Flag if a "ready file" (sentinel file) should be
				written at the end of each output stage.
				Not yet implemented.
ready_directory	С	None		Output directory for ready files.
				Not yet implemented.
ml_varlist(:)	С	None		Name of model level fields to be output.
pl_varlist(:)	С	None		Name of pressure level fields to be output.
p_levels(:)	R	None		pressure levels [hPa]
				Not yet implemented.
				The pressure levels are currently always taken fro
				array plevels in namelist nh_pzlev_nml.
hl_varlist(:)	С	None		Name of height level fields to be output.
h_levels(:)	R	None		height levels
				Not yet implemented.
				The height levels are currently always taken from
				array zlevels in namelist nh_pzlev_nml.
		1	1	

Parameter	Type	Default	Unit	Description
remap	I	0		interpolate horizontally, 0: none, 1: to regular
				lat-lon grid, 2: to Gaussian grids, (3:)
				Currently only 0 and 1 are implemented.
remap_internal	L	.FALSE.		do interpolations online in the model or external
				(including triggering)
				Currently unused, interpolations are always done
				internally.
$reg_lon_def(3)$	R	None		if remap=1: start, increment, end longitude in
				degrees
$reg_lat_def(3)$	R	None		if remap=1: start, increment, end latitude in
				degrees
gauss_tgrid_def	I	None		if remap=2: triangular truncation (e.g.63 for T63
				for which the Gauss grid should be used
				Currently unused since Gaussian grids are not
				implemented.
north_pole(2)	R	0,90		definition of north pole for rotated lon-lat grids.

Defined and used in: $src/namelists/mo_name_list_output.f90$

$3.18 \quad lonlat_intp_nml$

Parameter	Type	Default	Unit	Description
llonlat_enabled	L(n_dom)	.FALSE.		Flag. True, if lon-lat interpolation of output
				variables is desired.
lsupersede	L	.FALSE.		Flag. True, if standard grid variable is not writte
				for lon-lat vars.
lonlat_var_list	С	" 'PS',		List of variables for lon-lat interpolation or "all".
		'Q7',		
		'nor-		
		mal_veloci	ty'	
		"		
lon_delta	R(n_dom)	2.0	deg	Interpolation to lon-lat grid: resolution.

Parameter	Type	Default	Unit	Description
lat_delta	R(n_dom)	2.0	deg	Interpolation to lon-lat grid: resolution.
lon_corner1	R(n_dom)	-180.0	deg	South western corner of interpolation area (lon/la
lat_corner1	R(n_dom)	-90.0	deg	South western corner of interpolation area (lon/la
lon_corner2	R(n_dom)	180.0	deg	North eastern corner of interpolation area (lon/la
				Overrides corresponding "dimen" value.
lat_corner2	R(n_dom)	90.0	deg	North eastern corner of interpolation area (lon/la
				Overrides corresponding "dimen" value.
lon_poleN	R(n_dom)	0.	deg	Position of north pole for interpolation grid.
lat_poleN	R(n_dom)	90.	deg	Position of north pole for interpolation grid.
lon_dimen	I(n_dom)	-1		Dimensions of interpolation grid. Computed
				automatically when a second area corner is
				provided.
lat_dimen	I(n_dom)	-1		Dimensions of interpolation grid. Computed
				automatically when a second area corner is
				provided.

Defined and used in: $src/namelists/mo_lonlat_intp_nml.f90$

$3.19 \quad meteogram_output_nml$

Parameter	Type	Default	Unit	Description
lmeteogram_enabled	L(n_dom)	.FALSE.		Flag. True, if meteogram of output variables is
				desired.
zprefix	C(n_dom)	"METEOG	RAM_"	string with file name prefix for output file
ldistributed	L(n_dom)	.TRUE.		Flag. Separate files for each PE.
n0_mtgrm	I(n_dom)	1		initial time step for meteogram output
ninc_mtgrm	I(n_dom)	1		output interval (in time steps)
stationlist_tot		53.633,		list of meteogram stations (triples with lat, lon,
		9.983,		name string)
		'Ham-		
		burg'		

Defined and used in: $src/namelists/mo_mtgrm_nml.f90$

$3.20 \quad nh_pzlev_nml$

			1	
Parameter	Type	Default	Unit	Description
lwrite_zlev	L	.TRUE.		Output on height levels
lwrite_plev	L	.TRUE.		Output on pressure levels
	-			
nzlev	I	10		number of height levels
nplev	I	10		number of pressure levels
zlevels	R	0,1000,	m	array of height levels
		2000,		
		,		
		10000		
plevels	R	100000,	Pa	array of pressure levels
		90000,		
		80000,		
		,		
		10000		

Defined and used in: $src/namelists/mo_nh_pzlev_nml.f90$

$3.21 \quad transport_nml \; (used \; if \; run_nml/ltransport = .TRUE.)$

Parameter	Type	Default	Unit	Description
lvadv_tracer	L	.TRUE.		TRUE : compute vertical tracer advection
				FALSE: do not compute vertical tracer advection
ihadv_tracer	I(ntracer)	2		Tracer specific method to compute horizontal
				advection:

Parameter	Type	Default	Unit	Description
		4		0: no horiz. transport
				1: upwind (1st order)
				2: miura (2nd order, lin. reconstr.)
				20: miura (2nd order, lin. reconstr.) with
				subcycling
				3: miura3 (quadr. or cubic reconstr.)
				22: combination of miura and miura with
				subcycling
				32: combination of miura and miura with
				subcycling
				4: up3 (3rd or 4th order upstream)
ivadv_tracer	I(ntracer)	3		Tracer specific method to compute vertical
				advection:
				0: no vert. transport
				1: upwind (1st order)
				2: $muscl_cfl$ (2nd order, handles $CFL > 1$)
				20: muscl (2nd order)
				3: ppm_cfl (3^{rd} order, handles CFL > 1)
				30: ppm (3rd order)
lstrang	L	.FALSE.		splitting into fractional steps
				- second order Strang splitting (.TRUE.)
				- first order Godunov splitting (.FALSE.)
ctracer_list	С	"		list of tracer names
itype_hlimit	I(ntracer)	3		Type of limiter for horizontal transport:
		4		0: no limiter
				1: semi-monotonous slope limiter
				2: monotonous slope limiter
				3: monotonous flux limiter
				4: positive definite flux limiter
itype_vlimit	I(ntracer)	1		Type of limiter for vertical transport:
				0: no limiter
				1: semi-monotone slope limiter

Parameter	Type	Default	Unit	Description
				2: monotonous slope limiter
				4: positive definite flux limiter
iord_backtraj	I	1		order of backward trajectory calculation:
				1: first order
				2: second order (iterative; currently 1 iteration
				hardcoded)
igrad_c_miura	I	1		Method for gradient reconstruction at cell center
				for 2nd order miura
				1: Least-squares (linear, non-consv)
				2: Green-Gauss
				3: gradient reconstruction (RBF) at cell center or
				the basis of normal gradients at edges
lclip_tracer	L	.FALSE.		Clipping negative values
upstr_beta_adv	R	1.0		parameter to select 3rd order (=1) or 4th order
				(=0) advection, or something inbetween (01)
ivcfl_max	I	5		determines stability range of vertical PPM-schem
				in terms of the maximum allowable CFL-number
llsq_svd	L	.FALSE.		use QR decomposition (FALSE) or SV
				decomposition (TRUE) for least squares design
				matrix A

Defined and used in: $src/namelists/mo_advection_nml.f90$

$3.22 \quad nwp_phy_nml$

Parameter	Type	Default	Unit	Description
inwp_gscp	I	1		cloud microphysics and precipitation
				0: none
				1: hydci (COSMO-EU microphysics)
qi0	R	0.0	kg/kg	cloud ice threshold for autoconversion
qc0	R	0.0	kg/kg	cloud water threshold for autoconversion

Parameter	Type	Default	Unit	Description
inwp_convection	I	1		convection
				0: none
				1: Tiedtke/Bechtold convection
inwp_cldcover	I	3		cloud cover scheme for radiation
				0: no clouds (only QV)
				1: grid-scale clouds and QV
				2: clouds from COSMO turbulence scheme
				3: clouds from COSMO SGS cloud scheme
inwp_radiation	I	1		radiation
				0: none
				1: RRTM radiation
				2: Ritter-Geleyn radiation
inwp_satad	I	1		saturation adjustment
				0: none
				1:
inwp_turb	I	1		vertical diffusion and transfer
				0: none
				1: COSMO diffusion and transfer
				2: ECHAM diffusion
				3: EDMF-DUALM (to be implemented)
inwp_sso	I	1		subgrid scale orographic drag
				0: none
				1: (COSMO) Lott and Miller scheme
inwp_gwd	I	1		non-orographic gravity wave drag
				0: none
				1:Orr-Ern-Bechtold-scheme(IFS)
inwp_surface	I	1		surface scheme
				0: none
				1: TERRA
ustart_raylfric	R	160.0	m/s	wind speed at which extra Rayleigh friction start
efdt_min_raylfric	R	10800.	s	minimum e-folding time of Rayleigh friction
				(effective for u $>$ ustart_raylfric $+$ 90 m/s)

Parameter	Type	Default	Unit	Description
latm_above_top	L	.FALSE.		.TRUE.: take into account atmosphere above mo
	(max_dom)		top for radiation computation
dt_conv	R	600.	seconds	time interval of convection call
	(max_dom)		currently each subdomain has
				the same value
dt _ccov	R	dt _conv	seconds	time interval of cloud cover call
	(max_dom)		currently each subdomain has
				the same value
dt_rad	R	1800.	seconds	time interval of radiation call
	(max_dom)		currently each subdomain has
				the same value
dt_sso	R	1200.	seconds	time interval of sso call
	(max_dom)		currently each subdomain has
				the same value
dt_gwd	R	1200.	seconds	time interval of gwd call
	(max_dom)		currently each subdomain has
				the same value

Defined and used in: $src/namelists/mo_atm_phy_nwp_nml.f90$

3.23 radiation_nml

Parameter	Type	Default	Unit	Description
ldiur	L	.TRUE.		switch for solar irradiation:
				.TRUE.:diurnal cycle,
				.FALSE.:zonally averaged irradiation
nmonth	I	0		0: Earth circles on orbit
				1-12: Earth orbit position fixed for specified mon
lyr_perp	L	.FALSE.		.FALSE.: transient Earth orbit following VSOP8'
				.TRUE.: Earth orbit of year yr_perp of the
				VSOP87 orbit is perpertuated
yr_perp	L	-99999		$year used for lyr_perp = .TRUE.$

defined in code.
from external file containing the
l insolation averaged over a yea
ted)
ll radiation computation
ngle formula for the radiative
ion.
verywhere
pends only on latitude
pends only on latitude. Local
at 07:14:15 for radiative transfer
time of day = 1/pi
anging with latitude and time
d irradiance changing with
nd time of day (iforcing=inwp
, ()
i i i i i i i i i i i i i i i i i i i

Parameter	Type	Default	Unit	Description
irad_h2o	I	1		Switches for the concentration of radiative agents
$irad_co2$		2		0: 0.
$irad_ch4$		3		1: prognostic variable
irad_n2o		3		2: global constant
irad_o3		3		3: externally specified
$irad_o2$		2		irad_aero = 5: Tanre aerosol climatology for
irad_cfc11		2		$run_nml/iforcing = 3 (NWP)$
irad_cfc12		2		irad_aero = 6: Tegen aerosol climatology for
irad_aero		2		$run_nml/iforcing = 3 (NWP) .AND. itopo = 1$
				irad_o3 = 2: ozone climatology from MPI
				irad_o3 = 4: ozone clim for Aqua Planet Exp
				$irad_o3 = 6$: ozone climatology with T5
				geographical distribution and Fourier series for
				seasonal cycle for $run_nml/iforcing = 3$ (NWP)
				irad_o3 = 7: GEMS ozone climatology (from IFS
				$for run_nml/iforcing = 3 (NWP)$
vmr_co2	R	353.9e-6		Volume mixing ratio of the radiative agents
vmr_ch4		1693.6e-9		
vmr_n2o		309.5e-9		
vmr_o2		0.20946		
vmr_cfc11		252.8e-12		
vmr_cfc12		466.2e-12		

Defined and used in: $src/namelists/mo_radiation_nml.f90$

$3.24 \quad nwp_lnd_nml$

Parameter	Type	Default	Unit	Description
nlev_snow	I	1		number of snow layers
				for lmulti_snow=.true.
nsfc_subs	I	1		number of tiles

Parameter	Type	Default	Unit	Description
nsfc_snow	I	0		number of static surface types which can have sn
				as a tile
nztlev	I	2		used time integration scheme
lmulti_snow	L	.FALSE.		.TRUE. for use of multi-layer snow model
lseaice	L	.FALSE.		.TRUE. for use of sea-ice model
llake	L	.FALSE.		.TRUE. for use of lake model

Defined and used in: src/namelists/mo_nwp_lnd_nml.f90

$3.25 \quad echam_phy_nml$

Parameter	Type	Default	Unit	Description
lrad	L	.TRUE.		Switch on radiation.
lvdiff	L	.TRUE.		Switch on turbulent mixing (i.e. vertical diffusion
lconv	L	.TRUE.		Switch on cumulus convection.
lcond	L	.TRUE.		Switch on large scale condensation.
lcover	L	.FALSE.		.TRUE. for prognostic cloud cover scheme, .FALS
				for diagnostic scheme.
lgw_hines	L	.FALSE.		.TRUE. for atmospheric gravity wave drag by the
				Hines scheme
lssodrag	L	.FALSE.		.TRUE. for subgrid scale orographic drag
llandsurf	L	.FALSE.		.TRUE. for surface exchanges
lice	L	.FALSE.		.TRUE. for sea-ice temperature calculation
lmeltpond	L	.FALSE.		.TRUE. for calculation of meltponds
lhd	L	.FALSE.		.TRUE. for hydrologic discharge model

Parameter	Type	Default	Unit	Description
lmlo	L	.FALSE.		.TRUE. for mixed layer ocean

Defined and used in: $src/namelists/mo_echam_phy_nml.f90$

$3.26 \quad echam_conv_nml$

Parameter	Type	Default	Unit	Description
iconv	I	1		Choice of cumulus convection scheme.
				1: Nordeng scheme
				2: Tiedtke scheme
				3: hybrid scheme
ncvmicro	I	0		Choice of convective microphysics scheme.
lmfpen	L	.TRUE.		Switch on penetrative convection.
lmfmid	L	.TRUE.		Switch on midlevel convection.
lmfdd	L	.TRUE.		Switch on cumulus downdraft.
lmfdudv	L	.TRUE.		Switch on cumulus friction.
cmftau	R	10800.		Characteristic convective adjustment time scale.
cmfctop	R	0.3		Fractional convective mass flux (valid range [0,1]) across the top of cloud
cprcon	R	1.0e-4		Coefficient for determining conversion from cloud water to rain.
cminbuoy	R	0.025		Minimum excess buoyancy.
entrpen	R	1.0e-4		Entrainment rate for penetrative convection.

Parameter	Type	Default	Unit	Description
dlev	R	3.e4	Pa	Critical thickness necessary for the onset of
				convective precipitation.

Defined and used in: $src/namelists/mo_echam_conv_nml.f90$

3.27 vdiff nml

Parameter	Type	Default	Unit	Description
lsfc_mon_flux	L	.TRUE.		Switch on surface momentum flux.
lsfc heat flux	L	.TRUE.		Switch on surface sensible and latent heat flux.

Defined and used in: $src/namelists/mo_vdiff_nml.f90$

$3.28 \quad turbdiff_nml$

Parameter	Type	Default	Unit	Description
itype_tran	I	2		type of surface-atmosphere transfer
imode_tran	I	1		mode of surface-atmosphere transfer
icldm_tran	I	0		mode of cloud representation in transfer paramet
imode_turb	I	3		mode of turbulent diffusion parametrization
icldm_turb	I	2		mode of cloud representation in turbulence
				parametr
itype_sher	I	1		type of shear production for TKE
ltkesso	L	.FALSE.		calculation SSO-wake turbulence production for
				TKE
ltkecon	L	.FALSE.		consider convective buoyancy production for TKI
lexpcor	L	.FALSE.		explicit corrections of the implicit calculated turb
				diff.
ltmpcor	L	.FALSE.		consideration of thermal TKE-sources in the
				enthalpy budget

Parameter	Type	Default	Unit	Description
lprfcor	L	.FALSE.		using the profile values of the lowest main level
				instead of the mean value of the lowest layer for
				surface flux calulations
lnonloc	L	.FALSE.		nonlocal calculation of vertical gradients used for
				turbul. diff.
lcpfluc	L	.FALSE.		consideration of fluctuations of the heat capacity
				air
limpltkediff	L	.TRUE.		consideration of fluctuations of the heat capacity
				air
itype_wcld	I	2		type of water cloud diagnosis
itype_synd	I	2		type of diagnostics of synoptical near surface
				variables
lconst_z0	L	.FALSE.		TRUE: horizontally homogeneous roughness leng-
				z0
const_z0	R	0.001	m	value for horizontally homogeneous roughness
				lenght z0

Defined and used in: $src/namelists/mo_turbdiff_nml.f90$

$3.29 \quad gw_hines_nml \; (Scope: \; lgw_hines = .TRUE. \; in \; echam_phy_nml)$

Parameter	Type	Default	Unit	Description
lheatcal	L	.FALSE.		.TRUE.: compute drag, heating rate and diffusion
				coefficient from the dissipation of gravity waves
				.FALSE.: compute drag only
emiss_lev	I	10		Index of model level, counted from the surface,
				from which the gravity wave spectra are emitted
rmscon	R	1.0	m/s	Root mean square gravity wave wind at the
				emission level
kstar	R	5.0e-5	1/m	Typical gravity wave horizontal wavenumber
m_min	R	0.0	1/m	Minimum bound in vertical wavenumber
lrmscon lat	L	.FALSE.		.TRUE.: use latitude dependent rms wind

	Parameter	Type	Default	Unit	Description
					- $ latitude >= lat_rmscon$: use rmscon
					- latitude <= lat_rmscon_eq: use rmscon_eq
j					- $lat_rmscon_eq < latitude < lat_rmscon: use$
					linear interpolation between rmscon_eq and rmsc
					.FALSE.: use globally constant rms wind rmscon
Ì	lat_rmscon_eq	R	5.0	deg N	rmscon_eq is used equatorward of this latitude
Ì	lat_rmscon	R	10.0	deg N	rmscon is used polward of this latitude
Ī	rmscon_eq	R	1.2	m/s	is used equatorward of latitude lat_rmscon_eq

Defined and used in: $src/namelists/mo_gw_hines_nml.f90$

4 Namelist parameters for testcases (NAMELIST ICON)

The ICON model code includes several experiments, so-called test cases, for the shallow water model a atmosphere. Depending on the specified experiment, initial conditions and boundary conditions are computed to the computer of the specified experiment, initial conditions and boundary conditions are computed to the conditions are conditions and conditions are conditionally conditions are conditionally conditions.

$4.1 \quad \text{ha_testcase_nml (Scope: ltestcase=.TRUE. and iequations=[0,1,2] in run}$

Parameter	Type	Default	Unit	Description
ctest_name	C	'JWw'		Name of test case:
				'SW_GW': gravity wave 'USBR': unsteady solid body rotation 'Will_2': Williamson test 2 'Will_3': Williamson test 3 'Will_5': Williamson test 5
				'Will_6': Williamson test 6 'GW': gravity wave (nlev=20 only!) 'LDF': local diabatic forcing test without physics 'LDF-Moist': local diabatic forcing test with physics initalised with zonal wind field

Parameter	Type	Default	Unit	Description
				'HS': Held-Suarez test
				'JWs': Jablonowski-Will. steady state
				'JWw': Jablonowski-Will. wave test
				'JWw-Moist': Jablonowski-Will. wave test
				including moisture
				'APE': aqua planet experiment
				'MRW': mountain induced Rossby wave
				'MRW2': modified mountain induced Rossby way
				'PA': pure advection
				'SV': stationary vortex
				'DF1': deformational flow test 1
				'DF2': deformational flow test 2
				'DF3': deformational flow test 3
				'DF4': deformational flow test 4
				'RH': Rossby-Haurwitz wave test
rotate_axis_deg	R	0.0	deg	Earth's rotation axis pitch angle
gw_brunt_vais	R	0.01	1/s	Brunt Vaisala frequency
gw_u0	R	0.0	m/s	zonal wind parameter
gw_lon_deg	R	180.0	\deg	longitude of initial perturbation
gw_lat_deg	R	0.0	deg	latitude of initial perturbation
jw_uptb	R	1.0	m/s	amplitude of the wave pertubation
			(?)	
mountctr_lon_deg	R	90.0	deg	longitude of mountain peak
mountctr_lat_deg	R	30.0	deg	latitude of mountain peak
mountctr_height	R	2000.0	m	mountain height
mountctr_half_widt		1500000.0	m	mountain half width
mount_u0	R	20.0	m/s	wind speed for MRW cases
rh_wavenum	Ι	4		wave number
rh_init_shift_deg	R	0.0	deg	pattern shift

Parameter	Type	Default	Unit	Description
ihs_init_type	I	1		Choice of initial condition for the Held-Suarez tes
				1: the zonal state defined in the JWs test case;
				other integers: isothermal state (T=300 K,
				ps=1000 hPa, u=v=0.)
lhs_vn_ptb	L	.TRUE.		Add random noise to the initial wind field in the
				Held-Suarez test.
hs_vn_ptb_scale	R	1.	m/s	Magnitude of the random noise added to the init
				wind field in the Held-Suarez test.
lrh_linear_pres	L	.FALSE.		Initialize the relative humidity using a linear
				function of pressure.
		<u> </u>		
rh_at_1000hpa	R	0.75		relative humidity
				0,1
				at 1000 hPa
linit_tracer_fv	L	.TRUE.		Finite volume initialization for tracer fields
ape_sst_case	С	'sst1'		SST distribution selection
				'sst1': Control experiment
				'sst2': Peaked experiment
				'sst3': Flat experiment
				'sst4': Control-5N experiment
				'sst_qobs': Qobs SST distribution exp
				'sst ice': Control SST distribution with -1.8 C
				above $64 \text{ N/S}.$
ildf_init_type	I	0		Choice of initial condition for the Local diabatic
				forcing test. 1: the zonal state defined in the JWs
				test case; other: isothermal state (T=300 K,
				ps=1000 hPa, u=v=0.)
ldf symm	L	.TRUE.		Shape of local diabatic forcing:
_ •				.TRUE.: local diabatic forcing symmetric about t
				equator (at 0 N)
				.FALSE.: local diabatic forcing asym. about the
				equator (at 30 N)
L	- L	1	1	/

$4.2 \quad nh_testcase_nml~(Scope:~ltestcase=.TRUE.~and~iequations=3~in~run_nm$

Parameter	Type	Default	Unit	Description
nh_test_name	С	'jabw'		testcase selection
				'zero': no orography
				'bell': bell shaped mountain at 0E,0N
				'schaer': hilly mountain at 0E,0N
				'jabw': Initializes the full Jablonowski Williamson
				test case.
				'jabw_s': Initializes the Jablonowski Williamson
				steady state test case.
				'jabw_m': Initializes the Jablonowski Williamson
				test case with a mountain instead of the wind
				perturbation (specify mount_height).
				'mrw_nh': Initializes the full Mountain-induced
				Rossby wave test case.
				'mrw2_nh': Initializes the modified
				mountain-induced Rossby wave test case.
				'mwbr_const': Initializes the mountain wave with
				two layers test case. The lower layer is isotherma
				and the upper layer has constant brunt vaisala
				frequency. The interface has constant pressure.
				'PA': Initializes the pure advection test case.
				'HS_nh': Initializes the Held-Suarez test case. A
				the moment with an isothermal atmosphere at re
				(T=300K, ps=1000hPa, u=v=0, topography=0.0)
				'HS_jw': Initializes the Held-Suarez test case wit
				Jablonowski Williamson initial conditions and zer
				topography.
				'APE_nh': Initializes the APE experiments. Wit
ı				the jabw test case, including moisture.
i				'wk82': Initializes the Weisman Klemp test case

Parameter	Type	Default	Unit	Description
jw_up	R	1.0	m/s	amplitude of the u-perturbation in jabw test case
u0_mrw	R	20.0	m/s	wind speed for mrw(2) and mwbr_const cases
. 1 . 1 .	D	2222		(0)
mount_height_mrw	R	2000.0	m	maximum mount height in mrw(2) and
				mwbr_const
mount half width	R	1500000.0	m	half width of mountain in mrw(2), mwbr const
mount_nana.		1000000.0	""	and bell
mount_lonctr_mrw_	deg	90.	degrees	lon of mountain center in mrw(2) and mwbr_con
_				
mount_latctr_mrw_	d k g	30.	degrees	lat of mountain center in mrw(2) and mwbr_con
temp i mwbr cons	+ R	288.0	K	temp at isothermal lower layer for mwbr const ca
remb_i_mwoi_com	16 16	200.0	11	temp at isothermal lower rayer for mwbi_const c
p int mwbr const	R	70000.	Pa	pres at the interface of the two layers for
				mwbr_const case
bruntvais_u_mwbr_	cRnst	0.025	1/s	constant brunt vaissala frequency at upper layer:
				mwbr_const case
mount_height	R	100.0	m	peak height of mountain
layer_thickness	R	-999.0	m	thickness of vertical layers
n flat level	I	2	 	level number for which the layer is still flat and n
11_1100_10.01				terrain-following
nh u0	R	0.0	m/s	initial constant zonal wind speed
 nh_t0	R	300.0	K	initial temperature at lowest level
nh_brunt_vais	R	0.01	1/s	initial Brunt-Vaisala frequency

Parameter	Type	Default	Unit	Description
torus_domain_lengt	hR	100000.0	m	length of slice domain
rotate axis deg	R	0.0	deg	Earth's rotation axis pitch angle
lhs nh vn ptb	L	.TRUE.	0	Add random noise to the initial wind field in the
				Held-Suarez test.
lhs_fric_heat	L	.FALSE.		add frictional heating from Rayleigh friction in th
				Held-Suarez test.
hs_nh_vn_ptb_sca	m leR	1.	m/s	Magnitude of the random noise added to the init
				wind field in the Held-Suarez test.
rh_at_1000hpa	R	0.7	1	relative humidity at 1000 hPa
qv_max	R	20.e-3	kg/kg	specific humidity in the tropics
ape_sst_case	С	'sst1'		SST distribution selection
				'sst1': Control experiment
				'sst2': Peaked experiment
				'sst3': Flat experiment
				'sst4': Control-5N experiment
				'sst_qobs': Qobs SST distribution exp.
linit_tracer_fv	L	.TRUE.		Finite volume initialization for tracer fields
qv max wk	R	0.014	Kg/kg	maximum specific humidity near
				the surface, range $0.012 - 0.016$
				used to vary the buoyancy
u_infty_wk	R	20.	m/s	zonal wind at infinity height
				range 0 45.
				used to vary the wind shear
bub_amp	R	2.	K	maximum amplitud of the thermal perturbation
bubctr_lat	R	0.	deg	latitude of the center of the thermal perturbation
bubctr_lon	R	90.	deg	longitude of the center of the thermal perturbation
bubctr_z	R	1400.	m	height of the center of the thermal perturbation
bub_hor_width	R	10000.	m	horizontal radius of the thermal perturbation
bub_ver_width	R	1400.	m	vertical radius of the thermal perturbation

Defined and used in: $src/testcases/mo_nh_testcases.f90$

5 External data

5.1 ext_par_nml (Scope: itopo=1 in run_nml)

Parameter	Type	Default	Unit	Description
itopo	I	0		0: analytical topography/ext. data
				1: topography/ext. data read from file
n_iter_smooth_top	o I(n_dom)	0		iterations of topography smoother
fac_smooth_topo	R	0.015625		pre-factor of topography smoother
heightdiff_threshold	R(n_dom)	3000.	m	height difference between neighboring grid points
				above which additional local nabla2 diffusion is
				applied
l_emiss	L	.TRUE.		read and use external surface emissivity map

Defined and used in: $src/namelists/mo_extpar_nml.f90$

6 External packages

6.1 art nml

Parameter	Type	Default	Unit	Description
lart	L	.FALSE.		main switch for ART-package

Defined and used in: $src/namelists/mo_art_nml.f90$

7 Information on vertical level distribution

The hydrostatic and nonhydrostatic models need hybrid vertical level information to generate the terrain hybrid level specification is stored in <icon home>/hyb_params/HYB_PARAMS_<nlev>. The hydros pressure based coordinates, the nonhydrostatic model expects height based coordinates. For fur home>/hyb_params/README.