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 write namelist files containing the specified Fortran namelists. Programs are stored in <icon home>/build/<architecture>/bin/.

Table 1: Namelist files

Namelist file	Purpose	Made by script	Used by program
NAMELIST_GRAPH	Generate graphs	create_global_grids.run	grid_command
$NAMELIST_GRID$	Generate grids	$create_global_grids.run$	$\operatorname{grid} _\operatorname{command}$
NAMELIST GRIDREF	Gen. nested domains	create global grids.run	grid command
NAMELIST OCEAN GRID	Gen. ocean grid	create ocean grid.run	grid command
NAMELIST TORUS GRID	Gen. torus grid	create torus grid.run	grid command
NAMELIST ICON	Run ICON models	exp. <name>.run</name>	$\frac{-}{\text{control}}$ model

1.2 Namelist parameters

The following subsections tabulate all available Fortran namelist parameters by name, type, default value, unit, description, and scope:

- Type refers to the type of the Fortran variable, in which the value is stored: I=INTEGER, L=LOGICAL, R=REAL, C=character string
- ullet 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.
- \bullet 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 to specific settings of other 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	Scope
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 grid.	
				The root level consists of 8 triangles.	

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

2.1.2 grid_ini (NAMELIST_GRID)

Parameter	Type	Default	Unit	Description	Scope
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 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	Scope
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	Scope
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	$i_type_optimize = 1 \text{ or } 4$
				applied	
beta_spring	R	0.90		tuning factor for target grid length	$i_type_optimize = 4$

Defined and used in: src/grid generator/mo io grid.f90

2.1.4 plane options (NAMELIST GRID)

Parameter	Type	Default	Unit	Description	Scope
tria_arc_km	R	10.0	km	length of triangle edge on plane	lplane=.TRUE.

The number of grid points is generated by root level section and further bisections. The double periodic root level consists of 8 triangles. The spatial coordinates are -1 <= x <= 1, and $-\sqrt{3}/2 <= y <= \sqrt{3}/2$. Currently the planar option can only be used as an f-plane. Defined and used in: $\text{src/grid_generator/mo_io_grid.f90}$

2.1.5 gridref ini (NAMELIST GRIDREF)

Parameter	Type	Default	Unit	Description	Scope
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Parameter	Type	Default	Unit	Description	Scope
grid_root	I	2		root subdivision of initial edges	
start_lev	I	4		number of edge bisections following the root	
				subdivision	
n_dom	I	2		number of logical model domains, including the	
				global one	
n_phys_dom	I	$n_{ m dom}$		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 of	
				more than one nested domain per grid level)	
logical_id	I(n_phys_	i+1		logical grid ID of parent domain (first entry refers	
	dom-1)			to first nested domain; needs to be specified only in	
				case of 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	lcirc=.FALSE.
				$l_circ = .FALSE.$	
write_hierarchy	I	1		0: Output only computational grids	
				1: Output in addition parent grid of global model	
				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	h I	\max_{rlcell}		Number of cell rows along the lateral boundary of a	
		(=8)		model domain for which the refin_ctrl fields	
				contain the distance from the lateral boundary;	
				needs to be enlarged when lateral boundary	
				nudging is required for one-way nesting	

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

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

$2.2 \quad \text{Namelist parameters defining the ocean grid (NAMELIST_OCEAN_GRID)}$

The ocean grids are created by the script run/create_ocen_grid.run

2.2.1 grid geometry conditions

Parameter	Type	Default	Unit	Description	Scope
$no_of_conditions$	I	0		Number of geometric conditions	
patch_shape	I(no_of_co	n@litions)		1=rectangle; 2=circle	
patch_center_x	R(no_of_o	o ûdi tions)	degrees	longitude of patch center	
patch_center_y	R(no_of_o	o ûdi tions)	degrees	latitude of patch center	
rectangle_xradious	R(no_of_o	o ûdi tions)	degrees	half meridional extension of a rectangular patch	${ m patch_shape}{=}1$
rectangle_yradious	R(no_of_o		$_{ m degrees}$	0 1	${ m patch_shape}{=}1$
circle_radious	R(no_of_o	$\operatorname{olditions})$	degrees	radius of a circular patch	${ m patch_shape}{=}2$

Defined in mo_grid_conditions.f90

2.2.2 create_ocean_grid

Parameter	Type	Default	Unit	Description	Scope
only_get_sea_land	таsk	.false.		.true.:returns the whole grid with a sea-land mask;	
				.false.:returns only the ocean grid	
smooth_ocean_bour	ıdlary	.true.		.true.:smooths the ocean boundaries so no triabgle	
				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	$no_of_conditions=0$
				the input_file	
elevation_field	С			name of the field containing the cell elevation values	$no_of_conditions=0$
\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_interp_n	$\operatorname{le}_{\operatorname{I}}\operatorname{hod}$	2		compute edge elevation from cells using: linear	
				interpolation=1; min value = 2	
output_refined_ocea	n <u>C</u> file			name of the output refined ocean grid file	

Defined in mo_create_ocean_grid.f90

${\bf 2.3}\quad {\bf Namelist~parameters~defining~the~torus~grid~(NAMELIST_TORUS_GRID)}$

2.3.1 torus grid parameters

Parameter	Type	Default	Unit	Description	Scope
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	C			the torus grid file name	
unfolded_torus_file	r G me			the unfolded torus grid file name (for plotting)	
ascii_filename	C			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 the experiment, and the properties of dynamics, transport, physics etc.

$3.1 \quad master_nml$

Parameter	Type	Default	Unit	Description	Scope
l_atmo_active	L	.FALSE.		If .TRUE., the atmosphere model is started.	
atmo_name	С			character string to identify the atmosphere	
				component.	
atmo_namelist_filer	a fi le			namelist containing the atmospere namelists.	
atmo_min_rank	I	0		Start MPI rank for the atmosphere.	
atmo_max_rank	I	-1		End MPI rank for the atmosphere.	
atmo_inc_rank	I	0		stride of MPI ranks.	
atmo_restart_info_	fil © name	restart.info		Name (including full path) of the restart info file for	
				the atm model	
l_ocean_active	L	.FALSE.		If .TRUE., the ocean model is started.	
ocean_name	C			character string to identify the ocean component.	
ocean_namelist_file	naGhe			namelist containing the ocean namelists.	

Parameter	Type	Default	Unit	Description	Scope
ocean_min_rank	I	0		Start MPI rank for the ocean.	
ocean_max_rank	I	-1		End MPI rank for the ocean.	
ocean_inc_rank	I	0		stride of MPI ranks.	
ocean_restart_info_	$ ext{fi}$ Cename	restart.info		Name (including full path) of the restart info file for	
				the atm model	
l_restart	L	.FALSE.		If .TRUE.: Current experiment is started form a	
				restart.	

$3.2 \quad time_nml$

Parameter	Type	Default	Unit	Description	Scope
dt_restart	R	86400.*30.	S	Length of restart cycle in seconds	
calendar	I	1		Calendar type:	
				0=Julian/Gregorian	
				1=proleptic Gregorian	
				$2{=}30{ m day/month}, 360{ m day/year}$	
ini_datatime_string	С	'2008-09-		Initial date and time of the simulation	
		01T00:00:0	DΖ'		
end_datatime_string	g C	2008-09-		End date and time of the simulation	
		01T01:40:0	DΖ'		
				Length of the run	
				If "nsteps" in run_nml (see below) is positive, then	
				nsteps*dtime is used to compute the end date and	
				time of the run.	
				Else the initial date and time, the end date and	
				time, dt_restart, as well as the time step are used	
				to compute "nsteps".	

$3.3 \quad parallel_nml$

Parameter	Type	Default	Unit	Description	Scope
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Parameter	Type	Default	Unit	Description	Scope
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	
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 only	
				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 summation	
iorder sendrecv	I	1		Sequence of send/receive calls: 1 = irecv/send; 2 =	
_				isend/recv; 3 = isend/irecv	
itype_comm	I	1		1: use local memory for exchange buffers	
				2: use global memory for exchange buffers	
				3: asynchronous halo communication for dynamical	
				core (NH tria only)	
num_io_procs	I	0		Number of I/O processors (running exclusively for	
				$\operatorname{doing} \operatorname{I/O})$	
pio_type	I	1		Type of parallel I/O. Only used if number of I/O	
				processors greater number of domains.	
				Experimental!	
nh_stepping_threa	ds 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!	

Parameter	Type	Default	Unit	Description	Scope
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!	

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

3.4 coupling_nml

Parameter	Type	Default	Unit	Description	Scope
name	С	blank		short name of the coupling field	
frequency	I	0	s	coupling frequency	
time_step	I	0	s	model time step	
lag	I	0		offset to coupling event in number of model time	
				m step	
l_time_average	L	.FALSE.		.TRUE.: time averaging between to coupling events	
l_time_accumulatio	n L	.FALSE.		.TRUE.: accumulation of coupling fields in time	
				between to coupling events	

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

$3.5 \quad run_nml$

Parameter	Type	Default	Unit	Description	Scope
ldump_states	L	.FALSE.		Dump patch/interpolation/grid refinement state of	
				every patch (after subdivision in case of a parallel	
				run) to a Netcdf file and exit program.	
lrestore_states	L	.FALSE.		Restore patch/interpolation/grid refinement states	
				from NetCDF dump files instead of calculating	
				them.	
nsteps	I	0		number of time steps of this run.	

Parameter	Type	Default	Unit	Description	Scope
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_nlev	I(max_do	m)31		Number of full levels (atm.) for each domain	$lvert_nest=.TRUE.$
nshift	I(max_do	m)0		vertical half level of parent domain which coincides	$lvert_nest=.TRUE.$
				with upper boundary of the current domain	
ltimer	L	.TRUE.		TRUE: Timer for monitoring thr runtime of specific	
				routines is on $(FALSE = off)$	
timers_level	I	1			
msg_level	I	10		controls how much printout is written during	
				runtime.	
				For values less than 5, only the time step is written.	

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

$3.6 \quad grid_nml$

Parameter	Type	Default	Unit	Description	Scope
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 geographical	lplane=.TRUE.
	_			latitude	
l_limited_area	L	.FALSE.			
lfeedback	L(n_dom)	.TRUE.		Specifies if feedback to parent grid is performed.	n_dom>1
				Setting lfeedback(1)=.false. turns off feedback for	
				all nested domains; to turn off feedback for selected	
				nested domains, set lfeedback(1)=.true. and set	
				".false." for the desired model domains	
patch_weight	R(n_dom)	0.		If patch_weight is set to a value > 0 for any of the	n_dom>1
				first level child patches, processor splitting will be	
				performed, i.e. every of the first level child patches	
				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	
				root patch and higher level childs, patch_weight is	
				not used. However, patch_weight must be set to 0	
1 1 1 1	т	DATCE		for these patches to avoid confusion.	
lredgrid_phys	L	.FALSE.		If set to .true. is calculated on a reduced grid (=	
1 ' '1	C			one grid level higher)	
dynamics_grid_				Array of the grid filenames to be used by the	
filename	T			dycore.	
dynamics_parent_	I			Array of the indexes of the parent grid filenames, as	
grid_id				described by the dynamics grid_filename array.	
no disting and	C			Indexes start at 1, an index of 0 indicates no parent.	
radiation_grid_				Array of the grid filenames to be used for the	
filename				radiation model. Filled only if the radiation grid is	
				different from the dycore grid.	

Parameter	Type	Default	Unit	Description	Scope
dynamics_radiation	I			Array of the indexes linking the dycore grids, as	
_grid_link				described by the dynamics_grid_filename array,	
				and the radiation_grid_filename array. It provides	
				the link index of the radiation_grid_filename, for	
				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. Only	
				needs to be filled when the	
				radiation_grid_filename is defined.	

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

$3.7 \quad \mathbf{gridref_nml}$

Parameter	Type	Default	Unit	Description	Scope
$\operatorname{grf_intmethod_c}$	I	2		Interpolation method for grid refinement (cell-based	n_dom>1
				dynamical variables):	
				1: parent-to-child copying	
				2: gradient-based interpolation	
$grf_intmethod_ct$	I	2		Interpolation method for grid refinement (cell-based	$n_dom > 1$
				tracer variables):	
				1: parent-to-child copying	
				2: gradient-based interpolation	
$\operatorname{grf_intmethod_e}$	I	4		Interpolation method for grid refinement	n_dom>1
				(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:	n_dom>1
				1: average of child edges 1 and 2	
				2: 2nd-order method using RBF interpolation	

Parameter	Type	Default	Unit	Description	Scope
grf_scalfbk	I	2		Feedback method for dynamical scalar variables	n_dom>1
				(T, p_{sfc}) :	
				1: area-weighted averaging	
				2: bilinear interpolation	
$\operatorname{grf_tracfbk}$	I	2		Feedback method for tracer variables:	$n_{dom}>1$
				1: area-weighted averaging	
				2: bilinear interpolation	
$grf_idw_exp_e12$	R	1.2		exponent of generalized IDW function for child	n_dom>1
				edges 1/2	
$grf_idw_exp_e34$	R	1.7		exponent of generalized IDW function for child	n_dom>1
				edges 3/4	
rbf_vec_kern_grf_	eΙ	1		RBF kernel for grid refinement (edges):	n_dom>1
				1: Gaussian	
				$2: 1/(1+r^2)$	
				3: inverse multiquadric	
$rbf_scale_grf_e$	R	0.5		RBF scale factor for grid refinement (edges)	$n_{dom}>1$
$denom_diffu_t$	R	135		Deniminator for lateral boundary diffusion of	n_dom>1
				temperature	
denom_diffu_v	R	200		Deniminator for lateral boundary diffusion of	n_dom>1
				velocity	

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

3.8 prepicon nml

Remark: prepicon_nml contains switches controlling the real-data initialization functionality of ICON. There are currently two ways of 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 requested to do the horizontal and 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 equations needed by ICONAM, and
 - (c) perform the vertical interpolation to the ICON grid if IFS2ICON is requested to do only the horizontal interpolation step.

• If ICONAM (iequations=3) is combined with NWP physics (iforcing=3), setting ltestcase=.false. activates functionality (c) while running the ICON executable.

Parameter	Type	Default	Unit	Description	Scope
i_oper_mode	I	1		Operating mode if the prep_icon executable is run:	
				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	
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 input	
				(mandatory when inwp_surface >0)	
l_zp_in	L	.FALSE.		Logical switch for diagnostic output on pressure	prep_icon only
				and height levels	
l_extdata_out	L	.FALSE.		Logical switch to write extdata fields into output	prep_icon only

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

3.9 interpol_nml

Parameter	Type	Default	Unit	Description	Scope
llsq_high_consv	L	.TRUE.		conservative (T) or non-conservative (F)	
				least-squares reconstruction for high order transport	
lsq_high_ord	I	3		polynomial order for high order reconstruction	
				2: quadratic	
				30: cubic (no 3^{rd} order cross deriv.)	
				3: cubic	

Parameter	Type	Default	Unit	Description	Scope
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	
rbf vec scale c	R(n dom)	resolution-		Scale factor for RBF reconstruction at cell centres	
	\ _ /	dependent			
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	Ι	8		Total width (in units of cell rows) for lateral	
				boundary nudging zone	
i_cori_method	Ι	3		Selector for tangential wind reconstruction method	currently only for
					$cell_type=6$
				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	$i_cori_method{=}3$
				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

3.10 dynamics_nml

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

Parameter	Type	Default	Unit	Description	Scope
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:	$grid_nml:cell_type=3$
				1: Standard Gaussian integral. Hydrostatic	
				atm. model: for unaveraged normal components,	
				Non-hydrostatic atm. model: for averaged normal	
				components	
				2: bilinear averaging of divergence	
divavg_cntrwgt	R	0.5		Weight of central cell for divergence averaging	$idiv_method = 2$
sw_ref_height	R	0.9*2.94e4	$_{ m gm}$	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

$3.11 \quad ha_dyn_nml$

This namelist is relevant if run_nml:ldynamics=.TRUE. and dynamics_nml:iequations=IHS_ATM_TEMP or IHS_ATM_THETA.

Parameter Type Default Un	Description	Scope
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Parameter	Type	Default	Unit	Description	Scope
itime_scheme	I	4		Time integration scheme:	
				11: pure advection (no dynamics)	
				12: 2 time level semi implicit (not yet implemented)	
				13: 3 time level explicit	
				14: 3 time level with semi implicit correction	
				15: standard 4th-order Runge-Kutta method	
				(4-stage)	
				16: $SSPRK(5,4)$ scheme (5-stage)	
ileapfrog_startup	I	1		How to integrate the first time step when the	$itime_scheme = 13 \text{ or } 14$
				leapfrog scheme is chosen. $1 = \text{Euler forward}$; $2 = a$	
				series of sub-steps.	
asselin_coeff	R	0.1		Asselin filter coefficient	$itime_scheme = 13 \text{ or } 14$
si_2tls	R	0.6		weight of time step $n+1$. Valid range: $[0,1]$	$itime_scheme = 12$
si_expl_scheme	I	2		scheme for the explicit part used in the 2 time level	$itime_scheme = 12$
				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 with	itime_scheme=14 and
				speeds larger than si_cmin	$ m lsi_3d=.FALSE.$
si_coeff	R	1.0		weight of the semi implicit correction	$itime_scheme=14$
si_offctr	R	0.7			$itime_scheme=14$
si_rtol	R	1.0e-3		relative tolerance for GMRES solver	$itime_scheme=14$
lsi_3d	L	.FALSE.		3D GMRES solver or decomposistion into 2D	$lshallow_water=.FALSE.$
				problems	and itime_scheme= 14
ldry_dycore	L	.TRUE.		Assume dry atmosphere	iequations $\in \{1,2\}$
lref_temp	L	.FALSE.		Set a background temperature profile as base state	$iequations \in \{1,2\}$
				when computing the pressure graident force	

$3.12 \quad nonhydrostatic_nml \; (relevant \; if \; run_nml:iequations{=}3)$

Parameter	Type	Default	Unit	Description	Scope
$itime_scheme$	I	4		Time integration scheme:	

Parameter	Type	Default	Unit	Description	Scope
				3: same as default, but computation of velocity	iequations=3 and
				tendencies in corrector step only	$cell_type{=}3$
				4: Matsuno scheme	iequations=3 and
					$_{ m cell\ type=3}$
				6: same as default, but usage of velocity tendencies	iequations=3 and
				at $(nnow+nnew)/2$	$cell_type{=}3$
rayleigh coeff	R(n dom)	0.05		Rayleigh damping coefficient (Klemp, Dudhia,	cell_type=3
				Hassiotis: MWR136, pp.3987-4004)	
damp height	R(n dom)	30000	m	Height at which Rayleigh damping of vertical wind	
				starts	
damp height u	R	100000	m	Height at which Rayleigh damping of zonal wind	active only for inwp gwd
				starts	> 0
damp timescale u	R	259200	S	Shortest damping time scale (reached at model top)	
htop moist proc	R	200000.0	m	Height above which moist physics and advection of	
				cloud and precipitation variables are turned off	
htop qvadv	R	250000.0	m	Height above which QV advection is turned off (do	
				not use except for debugging purposes)	
k2_updamp_coeff	R	2.0e6		enhanced 2nd order diffusion coefficient in upper	cell type=6,
				damping layer	$hdiff_order=3$
					(Smagorinski)
vwind offetr	R	0.05		Off-centering in vertical wind solver	$cell_type=3$
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	cell_type=3
_				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	cell_type=3
				domain	

Parameter	Type	Default	Unit	Description	Scope
iadv_rhotheta	I	2		Advection method for rho and rhotheta:	$cell_type=3$
				1: centred differences horiz. + vert.	
				2: 2nd order Miura horizontal	
				3: 3rd order Miura horizontal (not recommended)	
igradp_method	I	1		Discretization of horizontal pressure gradient:	$cell_type=3$
				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	$cell_type=3$.AND.
				diffusion truly horizontally over steep slopes	$hdiff_order=5.AND.$
					$lhdiff_temp = .true.$
thslp_zdiffu	R	0.025		Slope threshold above which truly horizontal	$cell_type=3$.AND.
				temperature diffusion is activated	$hdiff_order=5$.AND.
					$lhdiff_temp=.true.$
					.AND. l_zdiffu_t=.true.
thhgtd_zdiffu	R	200	m	Threshold of height difference between neighboring	cell_type=3 .AND.
				grid points above which truly horizontal	$hdiff_order=5$.AND.
				temperature diffusion is activated (alternative	$lhdiff_temp=.true.$
				criterion to thslp_zdiffu)	.AND. l_zdiffu_t=.true.
$exner_expol$	R	0.5		Temporal extrapolation (fraction of dt) of Exner	$cell_type=3$
				function for computation of horizontal pressure	
				gradient	
l_open_ubc	L	.FALSE.		.TRUE.: Use open upper boundary condition	cell_type=3
				(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 (see	cell_type=6
_				also upstr_beta)	
upstr_beta	R	1.0		Selection of order for horiz, theta advection: 3rd	$cell_type=6$
				order=1.0, 4th order=0.0	

Parameter	Type	Default	Unit	Description	Scope
$gmres_rtol_nh$	R	1.0e-6		relative tolerance for convergence in gmres solver	$cell_type=6$

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

${\bf 3.13 \quad sleve_nml \ (relevant \ if \ nonhydrostatic_nml:ivctype=2)}$

Parameter	Type	Default	Unit	Description	Scope
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 flat	

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

$3.14 \quad diffusion_nml$

Parameter	Type	Default	Unit	Description	Scope
$lhdiff_temp$	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	

Parameter	Type	Default	Unit	Description	Scope
				5: Smagorinsky ∇^2 diffusion combined with ∇^4	
				background diffusion as specified via	
				hdiff_efdt_ratio	
				defaults: 2 for hexagonal model, 4 for triangular model	
				24 or 42: $\nabla 2$ diffusion from model top to a certain	24 and 42 currently
				level (cf. k2_pres_max and k2_klev_max below);	allowed only in the
				∇^4 for the lower levels.	hydrostatic atm model
					(run nml:iequation = 1)
					or 2).
k2_pres_max	R	-99.	Pa	Pressure level above which ∇^2 diffusion is applied.	$hdiff_order = 24 \text{ or } 42,$
					$and run_nml:iequation =$
					1 or 2.
k2_klev_max	I	0		Index of the vertical level till which (from the model	$hdiff_order = 24 \text{ or } 42,$
				top) ∇^2 diffusion is applied. If a positive value is	$and run_nml:iequation =$
				specified for k2_pres_max, k2_klev_max is reset	1 or 2.
				accordingly during the initialization of a model run.	
hdiff_efdt_ratio	R	1.0		ratio of e-folding time to time step (or 2* time step	
				when using a 3 time level time stepping scheme)	
				(only for triangles currently)	
hdiff_min_efdt_rat	ioR	1.0		minimum value of hdiff_efdt_ratio near model top	iequations = 3 . AND.
					cell_type=3
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	n_dom>1
	<u> </u>			coefficient for nested domains	
hdiff_smag_fac	R	0.15		Scaling factor for Smagorinsky diffusion	for triangles only with
					iequations=3, for
					hexagons with
					hdiff_order=3

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

3.15 io_nml

Parameter	Type	Default	Unit	Description	Scope
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	
lwrite_vorticity	L	.TRUE.		write out averaged vorticity at vertices	
lwrite_divergence	L	.TRUE.		write out divergence at cells	
lwrite_omega	L	.TRUE.		write out vertical velocity in pressure coords.	Always .FALSE. for
					nonhydrostatic and
					shallow water models
lwrite_pres	L	.TRUE.		write out full level pressure	$lshallow_water=.FALSE.$
lwrite_z3	L	.TRUE.		write out geopotential on full levels	$lshallow_water=.FALSE.$
lwrite_tracer	L(ntracer)	.TRUE.		write out tracer at cells	
lwrite_tend_phy	L	.TRUE.		Physics induced tendencies.	.TRUE. if
		.FALSE.			iforcing = iecham
		(Scope)			.FALSE. else
lwrite_radiation	L	.FALSE.		Radiation related fields.	Always .FALSE. if
					iforcing = inoforcing,
					iheldsuarez, ildf_dry
lwrite_precip	L	.FALSE.		Precipitation	Always .FALSE. if
					iforcing = inoforcing,
					iheldsuarez, ildf_dry
lwrite_cloud	L	.FALSE.		Cloud variables	Always .FALSE. if
					iforcing=inoforcing,
					iheldsuarez, ildf_dry

Parameter	Type	Default	Unit	Description	Scope
lwrite_tke	L	.TRUE.		TKE	.FALSE.
					Always .FALSE. if
					iforcing=inoforcing,
					iheldsuarez, ildf_dry
lwrite_surface	L	.FALSE.		surface variables	Always .FALSE. if
					iforcing=inoforcing,
					iheldsuarez, ildf_dry
lwrite_extra	L	.FALSE.		debug fields	.TRUE. if inextra_2d
					$/_3d > 0$
					.FALSE. else
inextra_2d	I	0		Number of 2D Fields for diagnostic/debugging	iequations = 3 (to be
				output.	done for $1, 2$)
inextra_3d	I	0		Number of 3D Fields for diagnostic/debugging	iequations = 3 (to be
				output.	done for $1, 2$)

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

$3.16 \quad transport_nml \; (used \; if \; run_nml/ltransport=.TRUE.)$

Parameter	Type	Default	Unit	Description	Scope
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:	
		4		0: no horiz. transport	
				1: upwind (1st order)	
				2: miura (2nd order, lin. reconstr.)	if cell_type=3
				3: miura3 (quadr. or cubic reconstr.)	$lsq_high_ord \in [2,3]$
				4: up3 (3rd or 4th order upstream)	if cell_type=6
ivadv_tracer	I(ntracer)	3		Tracer specific method to compute vertical	lvadv_tracer=TRUE
				advection:	
				0: no vert. transport	

Parameter	Type	Default	Unit	Description	Scope
				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	ihadv_tracer='miura'
				2: monotonous slope limiter	ihadv_tracer='miura'
				3: monotonous flux limiter	ihadv_tracer='miura[3]'
				4: positive definite flux limiter	ihadv_tracer='miura[3]', 'iup3[4]'
itype_vlimit	I(ntracer)	1		Type of limiter for vertical transport:	- 1
				0: no limiter	
				1: semi-monotone slope limiter	
				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	ihadv_tracer='miura'
				hardcoded)	
igrad_c_miura	I	1		Method for gradient reconstruction at cell center	
				for 2nd order miura	
				1: Least-squares (linear, non-consv)	$ihadv_tracer{=}2$
				2: Green-Gauss	
				3: gradient reconstruction (RBF) at cell center on	
				the basis of normal gradients at edges	
lclip_tracer	L	.FALSE.		Clipping negative values	

Parameter	Type	Default	Unit	Description	Scope
upstr_beta_adv	R	1.0		parameter to select 3rd order (=1) or 4th order	ihadv_tracer=iup3
				(=0) advection, or something inbetween (01)	
ivcfl_max	I	5		determines stability range of vertical PPM-scheme	ivadv_tracer=3
				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: $src/namelists/mo_advection_nml.f90$

$3.17 \quad nwp_phy_nml$

Parameter	Type	Default	Unit	Description	Scope
$inwp_gscp$	I	0		cloud microphysics and precipitation	$run_nml/iforcing = inwp$
				0: none	
				1: hydci (COSMO-EU microphysics)	
qi0	R	0.0	kg/kg	cloud ice threshold for autoconversion	$inwp_gscp=1$
qc0	R	0.0	kg/kg	cloud water threshold for autoconversion	inwp_gscp=1
inwp_convection	I	0		convection	$run_nml/iforcing = inwp$
				0: none	
				1: Tiedtke/Bechtold convection	
inwp_cldcover	I	1		cloud cover scheme for radiation	$run_nml/iforcing = inwp$
				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	0		radiation	${ m run_nml/iforcing} = { m inwp}$
				0: none	
				1: RRTM radiation	
				2: Ritter-Geleyn radiation	

Parameter	Type	Default	Unit	Description	Scope
inwp_satad	I	1		saturation adjustment	$run_nml/iforcing = inwp$
				0: none	
				1:	
inwp_turb	I	0		vertical diffusion and transfer	$run_nml/iforcing = inwp$
				0: none	
				1: COSMO diffusion and transfer	
				2: ECHAM diffusion	
inwp_sso	I	0		subgrid scale orographic drag	$run_nml/iforcing = inwp$
				0: none	
				1: (COSMO) Lott and Miller scheme	
inwp_gwd	I	0		non-orographic gravity wave drag	$run_nml/iforcing = inwp$
				0: none	
				1:Orr-Ern-Bechtold-scheme(IFS)	
inwp_surface	I	0		surface scheme	$run_nml/iforcing = inwp$
				0: none	
				1:	
dt_conv	R	600.	seconds	time interval of convection call	$run_nml/iforcing = inwp$
	$ \max_{dom}$)		currently each subdomain has	
				the same value	
$\mathrm{dt}_\mathrm{ccov}$	R	${ m dt_conv}$	seconds	time interval of cloud cover call	$run_nml/iforcing = inwp$
	$(\max_{dom}$)		currently each subdomain has	currently is not used
				the same value	
dt_rad	R	1800.	seconds	time interval of radiation call	$run_nml/iforcing = inwp$
	$ \max_{dom}$)		currently each subdomain has	
				the same value	
dt_sso	R	3600.	seconds	time interval of sso call	$run_nml/iforcing = inwp$
	(max_dom)		currently each subdomain has	
				the same value	
dt_gwd	R	3600.	seconds	time interval of gwd call	$run_nml/iforcing = inwp$
	$(\max_{dom}$)		currently each subdomain has	
				the same value	

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

3.18 radiation_nml

Parameter	Type	Default	Unit	Description	Scope
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 month	
lyr_perp	L	.FALSE.		.FALSE.: transient Earth orbit following VSOP87	
				.TRUE.: Earth orbit of year yr_perp of the	
				VSOP87 orbit is perpertuated	
yr_perp	L	-99999		$year used for lyr_perp = .TRUE.$	
isolrad	I	0		Insolation scheme	
				0: Use insolation defined in code.	
				1: Use insolation from external file containing the	
				spectrally resolved insolation averaged over a year	
				(not yet implemented)	
dt_rad	R	7200.	second	time interval of full radiation computation	$run_nml/iforcing =$
					iecham
izenith	I	3		Choice of zenith angle formula for the radiative	
		4 (for		transfer computation.	
		iforcing		0: Sun in zenith everywhere	
		= inwp)		1: Zenith angle depends only on latitude	
				2: Zenith angle depends only on latitude. Local	
				time of day fixed at 07:14:15 for radiative transfer	
				= computation (sin(time of day) = 1/pi	
				3: Zenith angle changing with latitude and time of	
				day	
				4: Zenith angle and irradiance changing with	
				season, latitude, and time of day (iforcing=inwp	
				only)	

Parameter	Type	Default	Unit	Description	Scope
irad_h2o	I	1		Switches for the concentration of radiative agents	Note: until further notice,
$irad_co2$		2		0: 0.	please use
$irad_ch4$		3		1: prognostic variable	$irad_h2o = 1$
irad_n2o		3		2: global constant	$irad_co2 = 2$
irad_o3		3		3: externally specified	and 0 for all the other
irad_o2		2		irad_aero = 5: aerosol climatology for	agents for
irad_cfc11		2		$run_nml/iforcing = 3 (NWP) when$	$\operatorname{run_nml/iforcing} = 2$
irad_cfc12		2		$inwp_radiation = 2$	(ECHAM).
irad_aero		2		$irad_o3 = 6$: ozone climatology with T5	
				geographical distribution and Fourier series for	
				$seasonal\ cycle\ for\ run_nml/iforcing = 3\ (NWP)$	
${ m vmr_co2}$	R	353.9e-6		Volume mixing ratio of the radiative agents	
${ m vmr_ch4}$		1693.6e-9			
vmr_n2o		309.5e-9			
${ m vmr}_{ m o}2$		0.20946			
vmr_cfc11		252.8e-12			
${ m vmr_cfc}12$		466.2e-12			

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

$3.19 \quad nwp_lnd_nml$

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

$3.20 \quad echam_phy_nml$

Parameter	Type	Default	Unit	Description	Scope
lrad	L	.TRUE.		Switch on radiation.	iforcing = 2
lvdiff	L	.TRUE.		Switch on turbulent mixing (i.e. vertical diffusion).	iforcing = 2
lconv	L	.TRUE.		Switch on cumulus convection.	iforcing = 2
lcond	L	.TRUE.		Switch on large scale condensation.	iforcing = 2
lcover	L	.FALSE.		.TRUE. for prognostic cloud cover scheme, .FALSE.	iforcing = 2
				for diagnostic scheme.	Note: $lcover = .TRUE$.
					runs, but has not been
					evaluated (yet) in ICON.
lgw_hines	L	.FALSE.		.TRUE. for atmospheric gravity wave drag by the	iforcing = 2
				Hines scheme	
lssodrag	L	.FALSE.		.TRUE. for subgrid scale orographic drag	iforcing = 2
					Not implemeted yet
llandsurf	L	.FALSE.		.TRUE. for surface exchanges	iforcing = 2
					Not implemeted yet
lice	L	.FALSE.		.TRUE. for sea-ice temperature calculation	iforcing = 2
					Not implemeted yet
lmeltpond	L	.FALSE.		.TRUE. for calculation of meltponds	iforcing = 2
					Not implemeted yet
lhd	L	.FALSE.		.TRUE. for hydrologic discharge model	iforcing = 2
					Not implemeted yet
lmlo	L	.FALSE.		.TRUE. for mixed layer ocean	iforcing = 2
					Not implemeted yet

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

$3.21 \quad echam_conv_nml$

Parameter	Type	Default	Unit	Description	Scope
iconv	I	1		Choice of cumulus convection scheme.	iforcing = 2 .AND. $lconv$
				1: Nordeng scheme	= .TRUE.
				2: Tiedtke scheme	
				3: hybrid scheme	
nevmiero	I	0		Choice of convective microphysics scheme.	iforcing = 2 .AND. lconv = .TRUE.
lmfpen	L	.TRUE.		Switch on penetrative convection.	iforcing = 2 .AND. $lconv$
					= .TRUE.
lmfmid	L	.TRUE.		Switch on midlevel convection.	iforcing = 2 .AND. $lconv$
					= .TRUE.
lmfdd	L	.TRUE.		Switch on cumulus downdraft.	iforcing = 2 .AND. $lconv$
					= .TRUE.
lmfdudv	L	.TRUE.		Switch on cumulus friction.	iforcing = 2 .AND. $lconv$
					= .TRUE.
cmftau	R	10800.		Characteristic convective adjustment time scale.	iforcing = 2 .AND. $lconv$
					= .TRUE.
cmfctop	R	0.3		Fractional convective mass flux (valid range [0,1])	iforcing = 2 .AND. $lconv$
				across the top of cloud	= .TRUE.
cprcon	R	1.0e-4		Coefficient for determining conversion from cloud	iforcing = 2 .AND. $lconv$
				water to rain.	= .TRUE.
cminbuoy	R	0.025		Minimum excess buoyancy.	iforcing = 2 .AND. $lconv$
					= .TRUE.
entrpen	R	1.0e-4		Entrainment rate for penetrative convection.	iforcing = 2 .AND. $lconv$
					= .TRUE.
dlev	R	3.e4	Pa	Critical thickness necessary for the onset of	iforcing = 2 .AND. $lconv$
				convective precipitation.	= .TRUE.

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

$3.22 \quad vdiff_nml$

Parameter	Type	Default	Unit	Description	Scope

Parameter	Type	Default	Unit	Description	Scope
lsfc_mon_flux	L	.TRUE.		Switch on surface momentum flux.	lvdiff = .TRUE.
lsfc_heat_flux	L	.TRUE.		Switch on surface sensible and latent heat flux.	lvdiff = .TRUE.

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

$3.23 \quad gw_hines_nml \; (Scope: \; lgw_hines = .TRUE. \; in \; echam_phy_nml)$

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

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 as well as the 3-dimensional atmosphere. Depending on the specified experiment, initial conditions and boundary conditions are computed internally.

$4.1 \quad ha_testcase_nml \; (Scope: \; ltestcase=.TRUE. \; and \; iequations=[0,1,2] \; in \; run_nml)$

Parameter	Type	Default	Unit	Description	Scope
ctest_name	C	'JWw'		Name of test case:	
				CW CW,	l-b-ll TDIE
				'SW_GW': gravity wave	lshallow_water=.TRUE.
				'USBR': unsteady solid body rotation	lshallow_water=.TRUE.
				'Will_2': Williamson test 2	lshallow_water=.TRUE.
				'Will_3': Williamson test 3	$lshallow_water=.TRUE.$
				'Will_5': Williamson test 5	$lshallow_water=.TRUE.$
				'Will_6': Williamson test 6	$lshallow_water=.TRUE.$
				'GW': gravity wave (nlev=20 only!)	lshallow_water=.FALSE.
				'LDF': local diabatic forcing test without physics	lshallow_water=.FALSE.
					and iforcing=4
				'LDF-Moist': local diabatic forcing test with	lshallow water=.FALSE.,
				physics initalised with zonal wind field	and iforcing=5
				'HS': Held-Suarez test	lshallow water=.FALSE.
				'JWs': Jablonowski-Will. steady state	lshallow water=.FALSE.
				'JWw': Jablonowski-Will. wave test	lshallow water=.FALSE.
				'JWw-Moist': Jablonowski-Will. wave test	lshallow water=.FALSE.
				including moisture	_
				'APE': aqua planet experiment	lshallow_water=.FALSE.
				'MRW': mountain induced Rossby wave	lshallow water=.FALSE.
				'MRW2': modified mountain induced Rossby wave	lshallow water=.FALSE.
				'PA': pure advection	lshallow water=.FALSE.
				'SV': stationary vortex	lshallow water=.FALSE.,
				SV. Stationary volues	$\begin{array}{c} \text{Ishanow} _ \text{water} = .\text{FALSE.}, \\ \text{ntracer} = 2 \end{array}$
				'DF1': deformational flow test 1	
				'DF2': deformational flow test 2	

Parameter	Type	Default	Unit	Description	Scope
				'DF3': deformational flow test 3	
				'DF4': deformational flow test 4	
				'RH': Rossby-Haurwitz wave test	$lshallow_water=.FALSE.$
rotate_axis_deg	R	0.0	deg	Earth's rotation axis pitch angle	ctest_name= 'Will_2',
					'Will_3', 'JWs', 'JWw',
					'PA', 'DF1234'
gw_brunt_vais	R	0.01	1/s	Brunt Vaisala frequency	ctest_name= 'GW'
gw_u0	R	0.0	m/s	zonal wind parameter	ctest_name= 'GW'
gw_lon_deg	R	180.0	deg	longitude of initial perturbation	ctest_name= 'GW'
gw_lat_deg	R	0.0	deg	latitude of initial perturbation	ctest_name= 'GW'
jw_uptb	R	1.0	m/s	amplitude of the wave pertubation	ctest_name= 'JWw'
			(?)		
$mountctr_lon_deg$	R	90.0	deg	longitude of mountain peak	ctest_name= 'MRW(2)'
mountctr_lat_deg	R	30.0	deg	latitude of mountain peak	ctest_name= 'MRW(2)'
mountctr_height	R	2000.0	m	mountain height	ctest_name= 'MRW(2)'
mountctr_half_widt	$^{ m hR}$	1500000.0	m	mountain half width	ctest_name= 'MRW(2)'
mount_u0	R	20.0	m/s	wind speed for MRW cases	ctest_name= 'MRW(2)'
$\rm rh_wavenum$	I	4		wave number	ctest_name= 'RH'
$ ho_{init_shift_deg}$	R	0.0	deg	pattern shift	$ctest_name = 'RH'$
ihs_init_type	I	1		Choice of initial condition for the Held-Suarez test.	ctest_name= 'HS'
				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	ctest_name= 'HS'
				Held-Suarez test.	
hs_vn_ptb_scale	R	1.	m/s	Magnitude of the random noise added to the initial	ctest_name= 'HS'
				wind field in the Held-Suarez test.	
lrh_linear_pres	L	.FALSE.		Initialize the relative humidity using a linear	ctest_name=
				function of pressure.	'JWw-Moist','APE',
					'LDF-Moist'
		•			

Parameter	Type	Default	Unit	Description	Scope
rh_at_1000hpa	R	0.75		relative humidity	$ctest_name =$
				0,1	'JWw-Moist','APE',
				at 1000 hPa	'LDF-Moist'
linit_tracer_fv	L	.TRUE.		Finite volume initialization for tracer fields	$ctest_name='PA'$
ape_sst_case	C	'sst1'		SST distribution selection	$ctest_name='APE'$
				'sst1': Control experiment	
				'sst2': Peaked experiment	
				'sst3': Flat experiment	
				'sst4': Control-5N experiment	
				'sst_qobs': Qobs SST distribution exp.	
ildf_init_type	I	0		Choice of initial condition for the Local diabatic	ctest_name= 'LDF'
				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:	ctest_name=
				.TRUE.: local diabatic forcing symmetric about the	'LDF','LDF-Moist'
				equator (at 0 N)	
				FALSE: local diabatic forcing asym. about the	
				equator (at 30 N)	

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

$4.2 \quad nh_testcase_nml \; (Scope: \; ltestcase=.TRUE. \; and \; iequations=3 \; in \; run_nml)$

Parameter	Type	Default	Unit	Description	Scope
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.	

Parameter	Type	Default	Unit	Description	Scope
	-140			'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 isothermal 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. At the moment with an isothermal atmosphere at rest (T=300K, ps=1000hPa, u=v=0, topography=0.0). 'HS_jw': Initializes the Held-Suarez test case with Jablonowski Williamson initial conditions and zero topography. 'APE_nh': Initializes the APE experiments. At the moment with T=300K, ps=1013.25hPa,	
in up	R	1.0	m m/s	u=v=w=0). amplitude of the u-perturbation in jabw test case	nh test name='jabw'
jw_up u0_mrw	R	20.0	m/s	wind speed for mrw case	nh_test_name= jabw nh_test_name= 'mrw(2)_nh'
mount_height_mrw	R	2000.0	m	maximum mount height in mrw(2) and mwbr_const	nh_test_name= 'mrw(2)_nh' and 'mwbr_const'
mount_half_width	R	1500000.0	m	half width of mountain in mrw(2), mwbr_const and bell	nh_test_name= 'mrw(2)_nh', 'mwbr_const' and 'bell'
mount_lonctr_mrw_	_ de g	90.	degrees	lon of mountain center in mrw(2) and mwbr_const	nh_test_name= 'mrw(2)_nh' and 'mwbr_const'

Parameter	Type	Default	Unit	Description	Scope
mount_latctr_mrw	dReg	30.	degrees	lat of mountain center in mrw(2) and mwbr_const	$nh_test_name =$
					$'mrw(2)_nh'$ and
					'mwbr_const'
u0 mwbr const	R	20.0	m/s	wind speed for mwbr const case	nh test name=
					'mwbr_const'
temp_i_mwbr_con	st R	288.0	K	temp at isothermal lower layer for mwbr_const case	$nh_test_name =$
					'mwbr_const'
p int mwbr const	R	70000.	Pa	pres at the interface of the two layers for	nh test name=
				mwbr const case	'mwbr_const'
bruntvais u mwbr	cRnst	0.025	1/s	constant brunt vaissala frequency at upper layer for	nh test name=
	Ī		,	mwbr const case	'mwbr_const'
mount height	R	100.0	m	peak height of mountain	nh test name='bell'
layer thickness	R	-999.0	m	thickness of vertical layers	If layer thickness < 0 ,
· —					the vertical level
					distribution is read in
					from externally given
					HYB PARAMS XX.
n_flat_level	I	2		level number for which the layer is still flat and not terrain-following	${\text{layer_thickness}} > 0$
nh u0	R	0.0	m/s	initial constant zonal wind speed	nh test name = 'bell'
nh t0	R	300.0	K	initial temperature at lowest level	nh test name = 'bell'
nh brunt vais	R	0.01	1/s	initial Brunt-Vaisala frequency	nh test name = 'bell'
torus domain leng		100000.0	-	length of slice domain	nh test name = 'bell',
torus_domain_leng		100000.0	m	length of since domain	lplane=.TRUE.
rotate axis deg	R	0.0	deg	Earth's rotation axis pitch angle	nh test name='PA'
		.TRUE.	deg	Add random noise to the initial wind field in the	
$lhs_nh_vn_ptb$	L	I NUE.			nh_test_name= 'HS nh'
11. C: 1. /	т	DATCE		Held-Suarez test.	
lhs_fric_heat	L	.FALSE.		add frictional heating from Rayleigh friction in the	nh_test_name=
1. 1	l D	-	ļ.,	Held-Suarez test.	'HS_nh'
hs_nh_vn_ptb_sca	aleK	1.	m/s	Magnitude of the random noise added to the initial	nh_test_name=
1	<u></u>			wind field in the Held-Suarez test.	'HS_nh'
$rh_at_1000hpa$	R	0.7	1	relative humidity at 1000 hPa	nh_test_name= 'jabw',
					$nh_test_name = 'mrw'$

Parameter	Type	Default	Unit	Description	Scope
qv_max	R	20.e-3	kg/kg	specific humidity in the tropics	$nh_test_name = 'jabw',$
					nh_test_name= 'mrw'
ape_sst_case	C	'sst1'		SST distribution selection	nh_test_name='APE_nh'
				'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	ctest_name='PA'

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	Scope
itopo	I	0		0: analytical topography/ext. data	
				1: topography/ext. data read from file	
n_iter_smooth_top	o I	2		iterations of topography smoother	itopo = 1
fac_smooth_topo	R	0.015625		pre-factor of topography smoother	$n_iter_smooth_topo >$
					0

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

6 Information on vertical level distribution

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