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	grid_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>	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 $[0^{\circ}E, 0^{\circ}N]$)	

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_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 domain (first entry refers to first	
	dom-1)			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	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	
I				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: $src/grid_generator/mo_gridrefinement.f90$

${\bf 2.1.6}\quad {\bf gridref_metadata~(NAMELIST_GRIDREF)}$

Parameter	Type	Default	Unit	Description	Scope
number_of_grid_used	$I(n_dom +$	1)0		sets the number of grid used in the netcdf header;	
				the number of entries must be n_dom+1 because	
				the first number refers to the radiation grid	
centre	I	0		centre running the grid generator: 78 - edzw	
				(DWD), 252 - MPIM	
subcentre	I	0		subcentre to be assigned by centre, usually 0	
outname_style	I	1		Output name style	
				1: Standard: $iconRXBXX_DOMXX.nc$	
				2: DWD: $icon_grid_XXXX_RXXBXX_X.nc$	

2.2 Namelist parameters defining the local grid generation

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

${\bf 2.2.1 \quad grid_geometry_conditions}$

Parameter	Type	Default	Unit	Description	Scope
no_of_conditions	I	0		Number of geometric conditions	
patch_shape	I(no_of_	0		1=rectangle; 2=circle	
	condi-				
	tions)				
patch_center_x	R(no_of	0.0	degrees	longitude of patch center	
	_ condi-				
	tions)				
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	patch_shape=1
	condi-				
	tions)				
rectangle_yradious	R(no_of_	0.0	degrees	half zonal extension of a rectangular patch	patch_shape=1
	condi-				
	tions)				
circle_radious	R(no_of_	0.0	degrees	radius of a circular patch	patch_shape=2
	condi-				
	tions)				

Defined in mo_grid_conditions.f90

2.2.2 local_grid_optimization

Parameter	Type	Default	Unit	Description	Scope
use_optimization	L	.FALSE.		Apply, or not, optimization	

Parameter	Type	Default	Unit	Description	Scope
use_edge_springs	L	.FALSE.		Use spring dynamics	
prime_ref_length _coeff	R	1.0		Spring length coefficient	
use_adaptive_	L	.FALSE.		Use adaptive spring length	
spring_length					
use_local_reference	L	.FALSE.		Use locally adaptive spring length	
_length					
local_reference_	R	0.0		Coefficient of local vs global spring length	
length_coeff					
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	
isotropy_stretch _coeff	R	0.0		Coefficient of the stretch isotropy force	
optimize_vertex _depth	I	1		For patches the min depth of the vertices that will	
				be optimized. The boundary vertices have depth 0,	
				the next level 1, etc.	

Defined in mo_local_grid_optimization.f90

${\bf 2.2.3 \quad create_ocean_grid}$

Parameter	Type	Default	Unit	Description	Scope
only_get_sea_	L	.false.		.true.:returns the whole grid with a sea-land mask;	
$land_mask$.false.:returns only the ocean grid	
smooth_ocean_	L	.true.		.true.:smooths the ocean boundaries so no triabgle	
boundary				has two boundary edges; .false.:no smoothing	
input_file	С			name of the input grid file	
elevation_file	C			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)		

Parameter	Type	Default	Unit	Description	Scope
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					

Defined in mo_create_ocean_grid.f90

${\bf 2.2.4}\quad torus_grid_parameters$

Parameter	Type	Default	Unit	Description	Scope
y_no_of_rows	I		4	number of triangle rows of the torus grid, >=2	
x_no_of_columns	I		8	number of triangle columns of the torus grid, >=2	
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	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 coupling_nml

Parameter	Type	Default	Unit	Description	Scope
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	
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	
$l_time_accumulation$	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

3.2 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	
lhdiff_w	L	.TRUE.		Diffusion on the vertical wind field	
hdiff_order	I	4 (hydro) 5 (NH)		Order of ∇ operator for diffusion:	
		<i>5</i> (NII)		 -1: no diffusion 2: ∇² diffusion (not available for NH model on triangles!) 3: Smagorinsky ∇² diffusion (includes frictional heating for the hexagonal model if lhdiff_temp=.TRUE.) 	

Parameter	Type	Default	Unit	Description	Scope
				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 hydrostatic	
				triangular model, 5 for nonhydrostatic triangular	
				NH model	24 1 421
				24 or 42: ∇2 diffusion from model top to a certain	24 and 42 currently
				level (cf. k2_pres_max and k2_klev_max below); ∇^4 for the lower levels.	allowed only in the
				V for the lower levels.	hydrostatic atm model (run nml:iequation = 1
					or 2).
itype vn diffu	I	1		Reconstruction method used for Smagorinsky	iequations=3,
V				diffusion:	hdiff order=3 or 5
				1: u/v reconstruction at vertices only	_
				2: u/v reconstruction at cells and vertices	
itype_t_diffu	I	2		Discretization of temperature diffusion:	iequations=3,
				1: $K_h \nabla^2 T$	hdiff_order=3 or 5
				$2: \nabla \cdot (K_h \nabla T)$	
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.
1 1:0	D	1.0		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	
		(hydro) 36.0		when using a 3 time level time stepping scheme)	
				(only for triangles currently; for triangular NH	
		(NH)		model, values above 30 are recommended when using hdiff order=5)	
hdiff w efdt ratio	R	15.0		ratio of e-folding time to time step for diffusion on	iequations=3
nam_w_erat_ratio	l n	15.0		vertical wind speed	requations—3
				vertical will speed	

Parameter	Type	Default	Unit	Description	Scope
hdiff_min_efdt_ratio	R	1.0		minimum value of hdiff_efdt_ratio near model top	iequations=3 .AND.
					hdiff_order=4
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
				coefficient for nested domains	
hdiff_smag_fac	R	0.15		Scaling factor for Smagorinsky diffusion	for triangles only with
		(hydro)			iequations=3, for
		0.015			hexagons with
		(NH)			hdiff_order=3

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

3.3 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:	
				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	$idiv_method=2$

Parameter	Type	Default	Unit	Description	Scope
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: $src/namelists/mo_dynamics_nml.f90$

$3.4 \quad echam_conv_nml$

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	
nevmicro	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]) across the top of cloud	iforcing = 2 .AND. lconv = .TRUE.
cprcon	R	1.0e-4		Coefficient for determining conversion from cloud water to rain.	iforcing = 2 .AND. lconv = .TRUE.
cminbuoy	R	0.025		Minimum excess buoyancy.	iforcing = 2 .AND. lconv = .TRUE.

Parameter	Type	Default	Unit	Description	Scope
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

$3.5 \quad echam_phy_nml$

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
icover	I	1		1 = diagnostic Sunquist cloud cover scheme,	iforcing = 2
				2 = prognostic Tompkins cloud cover scheme.	Note: icover $= 2$, has not
					been used (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

Parameter	Type	Default	Unit	Description	Scope
dt_rad	R	3600.	second	time interval of full radiation computation	$run_nml/iforcing =$
					iecham

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

$3.6 \quad \text{gribout_nml}$

Parameter	Type	Default	Unit	Description	Scope
backgroundProcess	I	0		Background process	filetype=2
				- GRIB2 code table backgroundProcess.table	
generatingCenter	I	-1		Output generating center. If this key is not set,	filetype=2
				center information is taken from the grid file	
				DWD: 78	
				MPIMET: 98	
				ECMWF: 98	
generatingProcessIdentifier	I(n_dom)	1		generating Process Identifier	filetype=2
				- GRIB2 code table	
				generatingProcessIdentifier.table	
generatingSubcenter	I	-1		Output generating Subcenter. If this key is not set,	filetype=2
				subcenter information is taken from the grid file	
				DWD: 255	
				MPIMET: 232	
				ECMWF: 0	
ldate_grib_act	L	.TRUE.		GRIB creation date	filetype=2
				.TRUE.: add creation date	
				.FALSE.: add dummy date	
lgribout_24bit	L	.FALSE.		If TRUE, write thermodynamic fields ρ , θ_v , T , p	filetype=2
				with 24bit precision instead of 16bit	
localDefinitionNumber	I	254		local Definition Number	filetype=2
				- GRIB2 code table	
				grib2LocalSectionNumber.78.table	
localNumberOfExperiment	I	1		local Number of Experiment	filetype=2

Parameter	Type	Default	Unit	Description	Scope
numberOfForecastsInEnser	nb l le	-1		Local definition for ensemble products, (only set if value changed from default)	filetype=2
perturbationNumber	I	-1		Local definition for ensemble products, (only set if value changed from default)	filetype=2
preset	С	"none"		Setting this different to "none" enables a couple of defaults for the other gribout_nml namelist parameters. If, additionally, the user tries to set any of these other parameters to a conflicting value, an error message is thrown. Possible values are "none", "deterministic", "ensemble".	filetype=2
productionStatusOfProcess	sedData	1		Production status of data - GRIB2 code table 1.3	filetype=2
significanceOfReferenceTin	ne I	1		Significance of reference time - GRIB2 code table 1.2	filetype=2
typeOfEnsembleForecast	I	-1		Local definition for ensemble products (only set if value changed from default)	filetype=2
type Of Generating Process	I	2		Type of generating process - GRIB2 code table 4.3	filetype=2
typeOfProcessedData	I	1		Type of data - GRIB2 code table 1.4	filetype=2

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

$3.7 \quad \mathrm{grid_nml}$

Parameter	Type	Default	Unit	Description	Scope
cell_type	I	3		Cell type: not used	
lplane	L	.FALSE.		planar option	
is_plane_torus	L	.FALSE.		f-plane approximation on triangular grid	
corio_lat	R	0.0	deg	Center of the f-plane is located at this geographical	lplane=.TRUE. and
				latitude	$is_plane_torus = .TRUE.$
grid_angular _velocity	R	Earth's	rad/sec	The angular velocity in rad per sec.	

Parameter	Type	Default	Unit	Description	Scope
l_limited_area	L	.FALSE.			
grid_rescale_factor	R	1.0		The geometry and the timestep will be multiplied by this factor. The angular velocity will be divided by this factor.	
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 selected nested domains, set lfeedback(1)=.true. and set ".false." for the desired model domains	n_dom>1
ifeedback_type	I	2		1: incremental feedback 2: relaxation-based feedback Note: vertical nesting requires option 2 to run numerically stable over longer time periods	n_dom>1
start_time	R(n_dom)	0.	s	Time when a nested domain starts to be active (namelist entry is ignored for the global domain)	n_dom>1
end_time	R(n_dom)	1.E30	S	Time when a nested domain terminates (namelist entry is ignored for the global domain)	n_dom>1
patch_weight	R(n_dom)			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 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 for these patches to avoid confusion.	n_dom>1
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. May contain the keyword <path> which will be substituted by model_base_dir.</path>	

Parameter	Type	Default	Unit	Description	Scope
dynamics_parent_	I			Array of the indexes of the parent grid filenames, as	
grid_id				described by the dynamics_grid_filename array.	
				Indexes start at 1, an index of 0 indicates no parent.	
radiation_grid_ filename	С			Array of the grid filenames to be used for the	
				radiation model. Filled only if the radiation grid is	
				different from the dycore grid. May contain the	
				keyword <path> which will be substituted by</path>	
				model_base_dir.	
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.	
create_vgrid	L	.FALSE.		.TRUE.: Write vertical grid files containing (vct_a,	
				vct_b, z_ifc, and z_ifv.	
vertical_grid_filename	С			Array of filenames. These files contain the vertical	
	(n_dom)			grid definition (vct_a, vct_b, z_ifc.	

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

$3.8 \quad {\rm gridref_nml}$

Parameter	Type	Default	Unit	Description	Scope
$grf_intmethod_c$	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):	

Parameter	Type	Default	Unit	Description	Scope
				1: parent-to-child copying	
				2: gradient-based interpolation	
grf_intmethod_e	I	6		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	
				5/6: same as $3/4$, respectively, but direct	
				interpolation of mass fluxes along nest interface	
				edges	
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	
$\operatorname{grf}_\operatorname{scalfbk}$	I	2		Feedback method for dynamical scalar variables	n_dom>1
				(T, p_{sfc}) :	
				1: area-weighted averaging	
				2: bilinear interpolation	
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 $3/4$	
$rbf_vec_kern_grf_e$	I	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	

Parameter	Type	Default	Unit	Description	Scope
denom_diffu_v	R	200		Deniminator for lateral boundary diffusion of	n_dom>1
				velocity	
l_mass_consvcorr	L	.FALSE.		.TRUE.: Apply mass conservation correction in	n_dom>1
				feedback routine	
l_density_nudging	L	.FALSE.		.TRUE.: Apply density nudging near lateral nest	n_dom>1 .AND.
				boundary if grf_intmethod_e ≤ 4	led
fbk_relax_timescale	R	10800		Relaxation time scale for feedback	n_dom>1 .AND.
					led back = .TRUE.
					$AND. ifeedback_type =$
					2

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

$3.9 \quad gw_hines_nml \; (Scope: \; lgw_hines = .TRUE. \; in \; echam_phy_nml)$

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	$lrmscon_lat = .TRUE.$

Parameter	Type	Default	Unit	Description	Scope
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

$3.10 \quad ha_dyn_nml$

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

Parameter	Type	Default	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 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 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	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

Parameter	Type	Default	Unit	Description	Scope
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 gradient force	

$3.11 \quad initicon_nml$

Parameter	Type	Default	Unit	Description	Scope
init mode	I	1		1: start from DWD analysis	
				2: start from IFS analysis	
				3: combined mode: IFS atm + GME soil	
				4: start from COSMO-DE forecast	
				5: start from DWD analysis with incremental	
				analysis update	
dt iau	R	10800	s	Time interval during which an incremental analysis	init mode=5
_				update (IAU) is performed	_
type iau wgt	I	1		Weighting function for performing IAU	init mode=5
				1: Top-Hat	_
				2: SIN2	
nlevsoil_in	I	4		number of soil levels of input data	init_mode=2
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_sst_in	L	.TRUE.		Logical switch. If true, the surface temperature of	$init_mode=2$
				the water sea points is initialized with the SST	
				provided in the ifs2icon file. If false, it is initialized	
				with the skin temperature. If the SST is not	
				provided in the ifs2icon file,l_sst_in is reset to	
				false.	
lread_ana	L	.TRUE.		If .FALSE., ICON is started from first guess only.	init_mode=1,3
				Analysis field is not required, and skipped if	
				provided.	

Parameter	Type	Default	Unit	Description	Scope
l_coarse2fine_mode	L(max_do	m)FALSE.		If true, apply corrections for coarse-to-fine mesh	
				interpolation to wind and temperature	
ifs2icon_filename	С			Filename of IFS2ICON input file, default	$init_mode=2$
				" <path>ifs2icon_R<nroot>B<jlev>_DOM<idom>.nc"</idom></jlev></nroot></path>	
				May contain the keywords <path> which will be</path>	
				substituted by model_base_dir, as well as nroot,	
				jlev, and idom defining the current patch.	
dwdfg_filename	С			Filename of DWD first-guess input file, default	$init_mode=1,3$
				" <path>dwdFG_R<nroot>B<jlev>_DOM<idom>.nc".</idom></jlev></nroot></path>	
				May contain the keywords <path> which will be</path>	
				substituted by model_base_dir, as well as nroot,	
				jlev, and idom defining the current patch.	
dwdana_filename	С			Filename of DWD analysis input file, default	$init_mode=1$
				" <path>dwdana_R<nroot>B<jlev>_DOM<idom>.nc".</idom></jlev></nroot></path>	
				May contain the keywords <path> which will be</path>	
				substituted by model_base_dir, as well as nroot,	
				jlev, and idom defining the current patch.	
filetype	I	-1		One of CDI's FILETYPE_XXX constants.	
		(undef.)		Possible values: 2 (=FILETYPE_GRB2), 4	
				(=FILETYPE_NC2). If this parameter has not	
				been set, we try to determine the file type by its	
				extension "*.grb*" or ".nc".	
ana_varlist	C			List of mandatory analysis fields that must be	$init_mode=1$
				present in the analysis file. If these fields cannot be	
				found in the analysis file, the model aborts. For all	
				other analysis fields, the FG-fields will serve as	
				fallback position.	
ana_varnames_map_file	C			Dictionary file which maps internal variable names	
				onto GRIB2 shortnames or NetCDF var names.	
				This is a text file with two columns separated by	
				whitespace, where left column: ICON variable	
				name, right column: GRIB short name.	

$3.12 \quad interpol_nml$

Parameter	Type	Default	Unit	Description	Scope
l_intp_c2l	L	.TRUE.		If .TRUE. directly interpolate scalar variables from	
				cell centers to lon-lat points, otherwise do gradient	
				interpolation and reconstruction.	
l_mono_c2l	L	.TRUE.		Monotonicity can be enforced by demanding that	
				the interpolated value is not higher or lower than	
				the stencil point values.	
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	
<u>-</u> 0 =				1: linear	ihadv tracer=4
				2: quadratic	_
				30 : cubic (no 3^{rd} order cross deriv.)	
				3: cubic	
llsq lin consv	L	.FALSE.		conservative (T) or non-conservative (F)	
- -				least-squares reconstruction for 2nd order (linear)	
				transport	
nudge_efold_width	R	2.5		e-folding width (in units of cell rows) for lateral	
9				boundary nudging coefficient	
nudge max coeff	R	0.02		Maximum relaxation coefficient for lateral	
<u> </u>				boundary nudging	
nudge zone width	I	8		Total width (in units of cell rows) for lateral	
♥ _				boundary nudging zone. If < 0 the patch	
				boundary depth index is used.	
rbf dim c2l	I	10		stencil size for direct lon-lat interpolation: 4 =	
				nearest neighbor, $13 = \text{vertex stencil}$, $10 = \text{edge}$	
				stencil.	

Parameter	Type	Default	Unit	Description	Scope
rbf_scale_mode_ll	I	2		Specifies, how the RBF shape parameter is	
				determined for lon-lat interpolation. 1 : lookup	
				table based on grid level (default) 2: determine	
				automatically. So far, this routine only estimates	
				the smallest value for the shape parameter for	
				which the Cholesky decomposition is likely to	
				succeed in floating point arithmetic.	
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_ll	I	1		Kernel type for reconstruction at lon-lat-points:	
				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			

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

3.13 io_nml

Parameter	Type	Default	Unit	Description	Scope

Parameter	Type	Default	Unit	Description	Scope
lkeep_in_sync L	L	.FALSE.		Sync output stream with file on disk after each	
				timestep	
dt_{diag}	R	86400.		diagnostic integral output interval	
dt_checkpoint	R	2592000	s	Time interval for writing restart files. Note that if	output /= "none"
				the value of dt_checkpoint resulting from model	(run_nml)
				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 file	
				is generated during the restart cycle.	
inextra_2d	I	0		Number of extra 2D Fields for	iequations = 3 (to be
				diagnostic/debugging output.	done for $1, 2$
inextra_3d	I	0		Number of extra 3D Fields for	iequations = 3 (to be
				diagnostic/debugging output.	done for $1, 2$)
lflux_avg	L	.TRUE.		if .FALSE. the output fluxes are accumulated	iequations=3
				from the beginning of the run	iforcing=3
				if .TRUE. the output fluxes are average values	
				from the beginning of the run, except of	
				TOT_PREC that would be accumulated	
$itype_pres_msl$	I	1		Specifies method for computation of mean sea level	
				pressure (and geopotential at pressure levels below	
				the surface).	
				1: GME-type extrapolation,	
				2: stepwise analytical integration,	
				3: current IFS method,	
				4: IFS method with consistency correction	
itype_rh I	I	1		Specifies method for computation of relative	
				humidity	
				1: WMO-type: water only (e_s=e_s_water),	
				2: IFS-type: mixed phase (water and ice),	
				3: IFS-type with clipping (rh ≤ 100)	

Parameter	Type	Default	Unit	Description	Scope
output_nml_dict C	, ,		File containing the mapping of variable names to	output_nml namelists	
				the internal ICON names. May contain the	
				keyword <path> which will be substituted by</path>	
				model_base_dir.	
				The format of this file:	
				One mapping per line, first the name as given in	
				the ml_varlist, hl_varlist, pl_varlist or	
				il_varlist of the output_nml namelists, then the	
				internal ICON name, separated by an arbitrary	
				number of blanks. The line may also start and end	
				with an arbitrary number of blanks. Empty lines or	
				lines starting with $\#$ are treated as comments.	
				Names not covered by the mapping are used as they	
				are.	
netcdf_dict	С	, ,		File containing the mapping from internal names to	output_nml namelists,
				names written to NetCDF. May contain the	NetCDF output
				keyword <path> which will be substituted by</path>	
				model_base_dir.	
				The format of this file:	
				One mapping per line, first the name written to	
				NetCDF, then the internal name, separated by an	
				arbitrary number of blanks (inverse to the	
				definition of output_nml_dict). The line may also	
				start and end with an arbitrary number of blanks.	
				Empty lines or lines starting with $\#$ are treated as	
				comments.	
			Names not covered by the mapping are output as		
			they are.		
				Note that the specification of output variables, e.g.	
				in ml_varlist, is independent from this renaming,	
				see the namelist parameter varnames_map_file for	
				this.	

Parameter	Type	Default	Unit	Description	Scope
lzaxis_reference	L	.TRUE.		FALSE: use vertical axis ZAXIS_HYBRID for 3D	will be removed after
				atmospheric fields	some testing phase
				TRUE: use vertical axis ZAXIS_REFERENCE for	
				3D atmospheric fields	

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

$3.14 \quad les_nml \ (parameters \ for \ LES \ turbulence \ scheme; \ valid \ for \ inwp_turb=5)$

Parameter	Type	Default	Unit	Description	Scope
sst	R	300	K	sea surface temperature for idealized LES	nh_test_name=CBL,
				simulations	RICO
					$isrfc_type=5,4$
shflx	R	-999	Km/s	Kinematic sensible heat flux at surface	$isrfc_type = 2$
lhflx	R	-999	m/s	Kinematic latent heat flux at surface	$isrfc_type = 2$
isrfc_type	I	1		surface type	
				1 = TERRA land physics	
				2 = fixed surface fluxes	
				3 = fixed buoyancy fluxes	
				4 = RICO test case	
				5 = fixed SST	
ufric	R	-999	m/s	friction velocity for idealized LES simulations	
is_dry_cbl	L	.FALSE.		switch for dry convective boundary layer	
				simulations	
karman_constant	R	0.4		von Karman constant	
smag_constant	R	0.12		Smagorinsky constant	
turb_prandtl	R	0.333333		turbulent Prandtl number	
bflux	R	-999	$\mathrm{m^2/s^3}$	buoyancy flux for idealized LES simulations	isrfc_type=3
				(Stevens 2007)	
tran_coeff	R	-999	m/s	transfer coefficient near surface for idealized LES	isrfc_type=3
				simulation (Stevens 2007)	

Parameter	Type	Default	Unit	Description	Scope
vert_scheme_type	I	2		type of time integration scheme in vertical diffusion	
				1 = explicit	
				2 = fully implicit	
sampl_freq_sec	R	60	S	sampling frequency in seconds for statistical (1D	
				and 0D) output	
avg_interval_sec	R	900	S	(time) averaging interval in seconds for 1D	
				statistical output	
expname	С	ICOLES		expname to name the statistical output file	
ldiag_les_out	L	.TRUE.		Control for the statistical output in LES mode	

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

$3.15 \quad limarea_nml \; (Scope: \; l_limited_area=1 \; in \; grid_nml)$

Parameter	Type	Default	Unit	Description	Scope
itype_latbc	I	0		Type of lateral boundary nudging. Nudge from	
				0: the initial data,	
				1: IFS data analysis/forecast (if init_mode=4, we	
				take COSMO-DE data),	
				2: ICON output data (with the identical 3d grid)	
dtime_latbc	R	10800.0	S	Time difference between two consecutive boundary	$itype_latbc \ge 1$
				data.	
nlev_latbc	I	0	s	Number of vertical levels in boundary data.	itype_latbc ≥ 1
latbc_filename	С			Filename of boundary data input file, default:	$itype_latbc \ge 1$
				"prepiconR <nroot>B<jlev>_<y><m><d><h>.nc".</h></d></m></y></jlev></nroot>	
				<y>, <m>, <d>, and <h> will be automatically</h></d></m></y>	
				replaced during the run-time. In case the time span	
				between two consecutive boundary data is less than	
				1 hour, one can use <min> and <sec>. These files</sec></min>	
				must be located in the latbc_path directory.	
latbc_path	С	"./"		Absolute path to boundary data.	itype_latbc ≥ 1

$3.16 \quad lnd_nml$

Parameter	Type	Default	Unit	Description	Scope
nlev_snow	I	2		number of snow layers	lmulti_snow=.true.
				for lmulti_snow=.true.	
ntiles	I	1		number of tiles	
lsnowtile	L	.FALSE.		.TRUE.: consider snow-covered and snow-free tiles	ntiles>1
				separately	
frlnd_thrhld	R	0.05		fraction threshold for creating a land grid point	ntiles>1
frlake_thrhld	R	0.05		fraction threshold for creating a lake grid point	ntiles>1
frsea_thrhld	R	0.05		fraction threshold for creating a sea grid point	ntiles>1
frlndtile_thrhld	R	0.05		fraction threshold for retaining the respective tile	ntiles>1
				for a grid point	
lmulti_snow	L	.TRUE.		.TRUE. for use of multi-layer snow model	
max_toplaydepth	R	0.25	m	maximum depth of uppermost snow layer	lmulti_snow=.TRUE.
idiag_snowfrac	I	1		Type of snow-fraction diagnosis:	
				1 = based on SWE only	
				2-4 = more advanced experimental methods	
itype_lndtbl	I	1		Table values used for associating surface parameters	
				to land-cover classes:	
				1 = defaults from extpar (GLC2000 and	
				GLOBCOVER2009)	
				2 = Tuned version based on IFS values for	
				globcover classes (GLOBCOVER2009 only)	
				3 = even more tuned version (EXPERIMENTAL!!,	
				GLOBCOVER2009 only)	
itype_root	I	2		root density distribution:	
				1 = constant	
				2 = exponential	
lseaice	L	.TRUE.		.TRUE. for use of sea-ice model	
llake	L	.TRUE.		.TRUE. for use of lake model	

Parameter	Type	Default	Unit	Description	Scope
sstice_mode	I	1		1: SST and sea ice fraction are read from the	iequations=3
				analysis and kept constant. The sea ice fraction can	iforcing=3
				be modified by the seaice model.	
				2: SST and sea ice fraction are updated daily, based	
				on climatological monthly means	
				3: SST and sea ice fraction are updated daily, based	
				on actual monthly means	
				4: SST and sea ice fraction are updated daily, based	
				on actual daily means, not yet implemented	
sst_td_filename	С			Filename of SST input files for time dependent	$sstice_mode=2,3$
				SST. Default is	
				$"<\!path>\!SST_<\!year>_<\!month>_<\!gridfile>".$	
				May contain the keyword <path> which will be</path>	
				substituted by model_base_dir	
ci_td_filename	С			Filename of sea ice fraction input files for time	$sstice_mode=2,3$
				dependent sea ice fraction. Default is	
				$"<\!path>\!CI_<\!year>_<\!month>_<\!gridfile>".$	
				May contain the keyword <path> which will be</path>	
				substituted by model_base_dir	

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

$3.17 \quad ls_forcing_nml \ (parameters \ for \ large-scale \ forcing; \ valid \ for \ torus \ geometry)$

Parameter	Type	Default	Unit	Description	Scope
is_ls_forcing	L	.FALSE.		switch for enabling large-scale (LS) forcing on torus	is_plane_torus=.TRUE.
				grid	
is_subsidence_moment	L	.FALSE.		switch for enabling LS vertical advection due to	is_plane_torus=.TRUE.
				subsidence for momentum equations	
is_subsidence_heat	L	.FALSE.		switch for enabling LS vertical advection due to	is_plane_torus=.TRUE.
				subsidence for thermal equations	
is_advection	L	.FALSE.		switch for enabling LS horizontal advection	is_plane_torus=.TRUE.
				(currently only for thermal equations)	

Parameter	Type	Default	Unit	Description	Scope
is_geowind	L	.FALSE.		switch for enabling geostrophic wind	is_plane_torus=.TRUE.
is_rad_forcing	L	.FALSE.		switch for enabling radiative forcing	is_plane_torus=.TRUE.
					$inwp_rad=.FALSE.$
is_geowind	L	.FALSE.		switch for enabling geostrophic wind	is_plane_torus=.TRUE.
is_theta	L	.FALSE.		switch to indicate that the prescribed radiative	is_plane_torus=.TRUE.
				forcing is for potential temperature	is_rad_forcing=.TRUE.

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

${\bf 3.18}\quad {\bf master_model_nml}\ ({\bf repeated}\ {\bf for}\ {\bf each}\ {\bf model})$

Parameter	Type	Default	Unit	Description	Scope
model_name	С			Character string for naming this component.	
model_namelist_	С			File name containing the model namelists.	
filename					
model_type	I	0		Identifies which component to run. atmosphere=1,	
				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 for	
_filename				this model	

3.19 master_nml

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

Parameter	Type	Default	Unit	Description	Scope
model_base_dir	С	, ,		General path which may be used in file names of	
				other name lists: If a file name contains the	
				keyword " <path>", then this model_base_dir will</path>	
				be substituted.	

$3.20 \quad meteogram_output_nml$

Parameter	Type	Default	Unit	Description	Scope
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

$3.21 \quad nh_pzlev_nml$

Parameter	Type	Default	Unit	Description	Scope
nzlev	I	10		number of height levels	iequations=3
nplev	I	10		number of pressure levels	iequations=3
nilev	I	3		number of isentropes	iequations=3

Parameter	Type	Default	Unit	Description	Scope
zlevels	R	10000, 9000, , 1000, 0	m	array of height levels	iequations=3 level ordering from TOA to bottom
plevels	R	100000, 90000, 80000, , 10000	Pa	array of pressure levels	iequations=3 level ordering from TOA to bottom
ilevels	R	340, 320, 300	K	array of isentropic levels	iequations=3 level ordering from TOA to bottom

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

$3.22 \quad nonhydrostatic_nml \; (relevant \; if \; run_nml:iequations{=}3)$

Parameter	Type	Default	Unit	Description	Scope
Parameter itime_scheme	Type I	Default 4	Unit	Options for predictor-corrector time-stepping scheme: 4: Contravariant vertical velocity is computed in the predictor step only, velocity tendencies are computed in the corrector step only (most efficient option) 5: Contravariant vertical velocity is computed in both substeps (beneficial for numerical stability in very-high resolution setups with extremely steep slops, otherwise no significant impact)	Scope iequations=3
				6: As 5, but velocity tendencies are also computed in both substeps (no apparent benefit, but more expensive)	

Parameter	Type	Default	Unit	Description	Scope
rayleigh_type	I	2		Type of Rayleigh damping	
				1: CLASSICAL (requires velocity reference state!)	
				2: Klemp (2008) type	
rayleigh coeff	R(n dom)	0.05		Rayleigh damping coefficient $1/\tau_0$ (Klemp, Dudhia,	
	` - /			Hassiotis: MWR136, pp.3987-4004); higher values	
				are recommended for R2B6 or finer resolution	
damp_height	R(n_dom)	45000	m	Height at which Rayleigh damping of vertical wind	
				starts (needs to be adjusted to model top height;	
				the damping layer should have a depth of at least 20	
				km when the model top is above the stratopause)	
htop moist proc	R	22500.0	m	Height above which moist physics and advection of	
				cloud and precipitation variables are turned off	
hbot qvsubstep	R	22500.0	m	Height above which QV is advected with	ihadv tracer=22, 32, 42
_				substepping scheme (must be at least as large as	or 52
				htop moist proc)	
vwind_offctr	R	0.15		Off-centering in vertical wind solver. Higher values	
				may be needed for R2B5 or coarser grids when the	
				model top is above 50 km.	
rhotheta_offctr	R	-0.1		Off-centering of density and potential temperature	
				at interface level (may be set to 0.0 for R2B6 or	
				finer grids)	
veladv_offctr	R	0.25		Off-centering of velocity advection in corrector step	
ivctype	I	2		Type of vertical coordinate:	
				1: Gal-Chen hybrid	
				2: SLEVE (uses sleve_nml)	
iadv_rcf	I	5		reduced calling frequency (rcf) for	
				transport/diffusion/physics	
				1: no rcf (every dynamics-step)	
				n>1: transport every n-th step	
				Setting odd values (besides 1) requires l_nest_rcf	
				= .TRUE.	

Parameter	Type	Default	Unit	Description	Scope
lhdiff_rcf	L	.TRUE.		.TRUE.: Compute diffusion only at advection time	
				steps (in this case, divergence damping is applied in	
				the dynamical core)	
lextra_diffu	L	.TRUE.		.TRUE.: Apply additional momentum diffusion at	
				grid points close to the stability limit for vertical	
				advection (becomes effective extremely rarely in	
				practice; this is mostly an emergency fix for	
				pathological cases with very large orographic	
				gravity waves)	
lbackward_integr	L	.FALSE.		.TRUE.: Integrate backward in time (preparation	
				for testing a digital filter initialization)	
divdamp_fac	R	0.0025		Scaling factor for divergence damping	$lhdiff_rcf = .TRUE.$
divdamp_order	I	4		Order of divergence damping:	$lhdiff_rcf = .TRUE.$
				2 = second-order divergence damping	
				4 = fourth-order divergence damping	
				24 = combined second-order and fourth-order	
				divergence damping and enhanced vertical wind	
				off-centering during the initial spinup phase (does	
				not allow checkpointing/restarting ealier than 2.5	
				hours of integration)	
divdamp_type	I	3		Type of divergence damping:	$lhdiff_rcf = .TRUE.$
				2 = divergence damping acting on 2D divergence	
				3 = divergence damping acting on 3D divergence	
l_nest_rcf	L	.TRUE.		Synchronize interpolation/feedback calls with	
				advection (transport) time steps. 1 nest rcf is	
				automatically reset to .FALSE. if iadv_rcf=1	
l masscorr nest	L	.FALSE.		.TRUE.: Apply mass conservation correction also in	
				nested domain	
iadv_rhotheta	I	2		Advection method for rho and rhotheta:	
_				1: simple second-order upwind-biased scheme	
				2: 2nd order Miura horizontal	
				3: 3rd order Miura horizontal (not recommended)	

Parameter	Type	Default	Unit	Description	Scope
igradp_method	I	3		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	
				4: Cubic/quadratic polynomial interpolation for	
				pressure reconstruction	
				5: Same as 4, but hydrostatic approximation for	
				downward extrapolation over steep slopes	
l_zdiffu_t	L	.TRUE.		.TRUE.: Compute Smagorinsky temperature	$hdiff_order=3/5$.AND.
				diffusion truly horizontally over steep slopes	$lhdiff_{temp} = .true.$
thslp_zdiffu	R	0.025		Slope threshold above which truly horizontal	$hdiff_order=3/5$.AND.
				temperature diffusion is activated	lhdiff_temp=.true.
111 1100		200			.AND. l_zdiffu_t=.true.
thhgtd_zdiffu	R	200	m	Threshold of height difference between neighboring	hdiff_order=3/5 .AND.
				grid points above which truly horizontal	lhdiff_temp=.true.
				temperature diffusion is activated (alternative	.AND. l_zdiffu_t=.true.
1		1 /0		criterion to thslp_zdiffu)	
exner_expol	R	1./3.		Temporal extrapolation (fraction of dt) of Exner	
				function for computation of horizontal pressure	
				gradient. This damps horizontally propagating	
				sound waves. For R2B5 or coarser grids, values	
1 anan uha	Т	EATCE		between 1/2 and 2/3 are recommended.	
l_open_ubc	L	.FALSE.		.TRUE.: Use open upper boundary condition	
				(rather than w=0) to allow vertical motions related	
				to diabatic heating to extend beyond the model top	

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

$3.23 \quad nwp_phy_nml$

The switches for the physics schemes and the time steps can be set for each model domain individually. If only one value is specified, it is copied to all child domains, implying that the same set of parameterizations and time steps is used in all domains. If the number of values given in the namelist is larger than 1 but less than the number of model domains, then the settings from the highest domain ID are used for the remaining model domains. If the time steps are not an integer multiple of the advective time step (dtime*iadv_rcf), then the time step of the respective physics parameterization is automatically rounded to the next higher integer multiple of the advective time step.

Parameter	Type	Default	Unit	Description	Scope
inwp_gscp	I	1		cloud microphysics and precipitation	$run_nml/iforcing = inwp$
	(max_dom)			
				0: none	
				1: hydci (COSMO-EU microphysics)	
				9: Kessler scheme	
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	1		convection	$run_nml/iforcing = inwp$
		$(\max_{dom_{dom_{dom_{dom_{dom_{dom_{dom_{dom$)		
				0: none	
				1: Tiedtke/Bechtold convection	
inwp_cldcover	I	1		cloud cover scheme for radiation	$run_nml/iforcing = inwp$
	(max_dom)			
				0: no clouds (only QV)	
				1: diagnostic cloud cover (by Martin Koehler)	
				2: prognostic total water variance (not yet started)	
				3: clouds from COSMO SGS cloud scheme	
				4: clouds as in turbulence (turbdiff)	
				5: grid scale clouds	
inwp_radiation	I	1		radiation	$run_nml/iforcing = inwp$
	(max_dom)			
				0: none	
				1: RRTM radiation	

Parameter	Type	Default	Unit	Description	Scope
				2: Ritter-Geleyn radiation	
inwp_satad	I	1		saturation adjustment	$run_nml/iforcing = inwp$
				0: none	
				1:	
inwp_turb	I	1		vertical diffusion and transfer	$run_nml/iforcing = inwp$
	(max_dom	վ)			
				0: none	
				1: COSMO diffusion and transfer	
				2: GME turbulence scheme (to be implemented)	
				3: EDMF-DUALM (work in progress)	
				5: Classical Smagorinsky diffusion	
$inwp_sso$	I	1		subgrid scale orographic drag	$ run_nml/iforcing = inwp $
	(max_dom	ų)			
				0: none	
				1: (COSMO) Lott and Miller scheme	
$inwp_gwd$	I	1		non-orographic gravity wave drag	$ run_nml/iforcing = inwp $
	(max_dom	ų)			
				0: none	
				1:Orr-Ern-Bechtold-scheme(IFS)	
inwp_surface	I	1		surface scheme	$ run_nml/iforcing = inwp $
	(max_dom	4)			
				0: none	
				1: TERRA	
ustart_raylfric	R	160.0	m/s	wind speed at which extra Rayleigh friction starts	$inwp_gwd > 0$
efdt_min_raylfric	R	10800.	S	minimum e-folding time of Rayleigh friction	$inwp_gwd > 0$
				(effective for u > ustart_raylfric + 90 m/s)	
latm_above_top	L	.FALSE.		.TRUE.: take into account atmosphere above model	$inwp_radiation > 0$
	(max_dom	4)		top for radiation computation	
itype_z0	I	1		Type of roughness length data used for turbulence	$inwp_turb > 0$
				scheme: $1 = \text{including contribution from sub-scale}$	
				orography, $2 = \text{land-cover-related roughness only}$	
dt_conv	R	600.	seconds	time interval of convection call	$run_nml/iforcing = inwp$
	(max_dom	4)		currently each subdomain has	

Parameter	Type	Default	Unit	Description	Scope
				the same value	
dt_ccov	R	dt _conv	seconds	time interval of cloud cover call	$run_nml/iforcing = inwp$
	(max_dom			currently each subdomain has	
				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	1200.	seconds	time interval of sso call	$run_nml/iforcing = inwp$
	(max_dom)		currently each subdomain has	
				the same value	
dt_gwd	R	1200.	seconds	time interval of gwd call	$run_nml/iforcing = inwp$
	(max_dom)		currently each subdomain has	
				the same value	
lrtm_filename	C(:)	"rrtmg_lw.	nc''	NetCDF file containing longwave absorption	
				coefficients and other data for RRTMG_LW	
				k-distribution model.	
cldopt_filename	C(:)	"ECHAM6	CldOptI	PrNpsGDF file with RRTM Cloud Optical Properties	
_				for ECHAM6.	

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

$3.24 \quad ocean_physics_nml$

Parameter	Type	Default	Unit	Description	Scope
i_sea_ice	I	1		0: No sea ice, 1: Include sea ice	
				.FALSE.: compute drag only	
richardson_factor_tracer	I	0.5e-5	m/s		
richardson_factor_veloc	I	0.5e-5	m/s		
l_constant_mixing	L	.FALSE.			

3.25 output_nml

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

Parameter	Type	Default	Unit	Description	Scope
dom(:)	I	-1		Array of domains for which this name-list is used.	
				If not specified (or specified as -1 as the first array	
				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!	
file_interval	С	5 5		Defines the length of a file in terms of an ISO-8601	
				duration string. An example for this time stamp	
				format is given below. This namelist parameter can	
				be set instead of steps_per_file.	
filename_format	C	see de-		Output filename format. Includes keywords path,	
		scription.		output_filename, physdom, etc. (see below).	
				Default is	
				<pre><output_filename>_DOM<physdom>_<levtype>_<jf< pre=""></jf<></levtype></physdom></output_filename></pre>	ile>
filetype	I	4		One of CDI's FILETYPE_XXX constants.	
				Possible values: 2 (=FILETYPE_GRB2), 4	
				(=FILETYPE_NC2), 5 (=FILETYPE_NC4)	
$h_{levels}(:)$	R	None	m	height levels	
				Not yet implemented.	
				The height levels are currently always taken from	
				array zlevels in namelist nh_pzlev_nml.	
hl_varlist(:)	C	None		Name of height level fields to be output.	
i_levels(:)	R	None	K	isentropic levels	
				Not yet implemented.	
				The isentropic levels are currently always taken	
				from array ilevels in namelist nh_pzlev_nml.	
il_varlist(:)	C	None		Name of isentropic level fields to be output.	
$include_last$	L	.TRUE.		Flag whether to include the last time step	

Parameter	Type	Default	Unit	Description	Scope
ml_varlist(:)	C	None		Name of model level fields to be output.	
mode	I	2		1 = forecast mode, $2 = $ climate mode	
				In climate mode the time axis of the output file is	
				set to TAXIS_ABSOLUTE. In forecast mode it is	
				set to TAXIS_RELATIVE. Till now the forecast	
				mode only works if the output is at multiples of 1	
				hour	
north_pole(2)	R	0,90		definition of north pole for rotated lon-lat grids.	
output_bounds(3)	R	None		Post-processing times (in seconds): start, end,	
				increment. We choose the advection time step	
				matching or following the requested output time,	
				therefore we require output_bounds(3) <	
				dtime*iadv_rcf. See namelist parameters	
				output_start, output_end, output_interval for	
				an alternative specification of output events.	
output_filename	C	None		Output filename prefix (which may include path).	
				Domain number, level type, file number and	
				extension will be added, according to the format	
				given in namelist parameter "filename_format".	
$output_grid$	L	.FALSE.		Flag whether grid information is added to output.	
$\operatorname{output_end}$	C	5 5		ISO8601 time stamp for end of output. An example	
				for this time stamp format is given below. See	
				namelist parameter output_bounds for an	
				alternative specification of output events.	
output_interval	C	5 5		ISO8601 time stamp for repeating output intervals.	
				We choose the advection time step matching or	
				following the requested output time, therefore we	
				require output_bounds(3) < dtime*iadv_rcf. An	
				example for this time stamp format is given below.	
				See namelist parameter output_bounds for an	
				alternative specification of output events.	

Parameter	Type	Default	Unit	Description	Scope
output_start	С	5 5		ISO8601 time stamp for begin of output. An	
				example for this time stamp format is given below.	
				See namelist parameter output_bounds for an	
				alternative specification of output events.	
p_levels(:)	R	None	hPa	pressure levels	
				Not yet implemented.	
				The pressure levels are currently always taken from	
				array plevels in namelist nh_pzlev_nml.	
pl_varlist(:)	С	None		Name of pressure level fields to be output.	
ready_file	С	'default'		A ready file is a technique for handling	
				dependencies between the NWP processes. The	
				completion of the write process is signalled by	
				creating a small file with name ready_file.	
				Different output_nml's may be joined together to	
				form a single ready file event. The setting of	
				<pre>ready_file = "default" does not create a ready</pre>	
				file. The ready file name may contain string tokens	
				<pre><path>, <datetime>, <ddhhmmss> which are</ddhhmmss></datetime></path></pre>	
				substituted as described for the namelist parameter	
				filename_format.	
reg_def_mode	I	0		Specify if the "delta" value prescribes an interval	remap=1
				size or the total *number* of intervals: 0: switch	
				automatically between increment and no. of grid	
				points, 1: reg_lon/lat_def(2) specifies increment,	
				2: reg_lon/lat_def(2) specifies no. of grid points.	
$reg_lat_def(3)$	R	None		start, increment, end latitude in degrees.	remap=1
				Alternatively, the user may set the number of grid	
				points instead of an increment. Details for the	
				setting of regular grids is given below together with	
				an example.	

Parameter	Type	Default	Unit	Description	Scope
reg_lon_def(3)	R	None		The regular grid points are specified by three values: start, increment, end given in degrees. Alternatively, the user may set the number of grid points instead of an increment. Details for the setting of regular grids is given below together with an example.	remap=1
remap	I	0		interpolate horizontally, 0: none, 1: to regular lat-lon grid	
steps_per_file	I	-1		Max number of output steps in one output file. If this number is reached, a new output file will be opened.	
steps_per_file_inclfirst	L	see descr.		Defines if first step is counted wrt. steps_per_file files count. The default is .FALSE. for GRIB2 output, and .TRUE. otherwise.	
stream_partitions_hl	I	1		Splits height level output of this namelist into several concurrent alternating files. See namelist parameter stream_partitions_ml for details.	
stream_partitions_il	I	1		Splits isentropic level output of this namelist into several concurrent alternating files. See namelist parameter stream_partitions_ml for details.	
stream_partitions_ml	I	1		Splits model level output of this namelist into several concurrent alternating files. The output is split into N files, where the start date of part i gets an offset of $(i-1)*$ output_interval. The output interval is then replaced by $N*$ output_interval, the include_last flag is set to .FALSE., and the steps_per_file counter is set to 1.	
stream_partitions_pl	I	1		Splits pressure level output of this namelist into several concurrent alternating files. See namelist parameter stream_partitions_ml for details.	

Parameter	Type	Default	Unit	Description	Scope
taxis_tunit	I	2		$1 = TUNIT_SECOND$	mode=1
				$2 = TUNIT_MINUTE$	
				3 = TUNIT HOUR	
				Time unit of the TAXIS RELATIVE time axis.	
				For a complete list of possible values see cdi.inc	

Interpolation onto regular grids: Horizontal interpolation onto regular grids is possible through the namelist setting remap=1, where the mesh is defined by the parameters

- reg_lon_def: mesh latitudes in degrees,
- reg_lat_def: mesh longitudes in degrees,
- north_pole: definition of north pole for rotated lon-lat grids.

The regular grid points in reg_lon_def, reg_lat_def are each specified by three values, given in degrees: start, increment, end. The mesh then contains all grid points start + k * increment <= end, where k is an integer. Instead of defining an increment it is also possible to prescribe the number of grid points.

- Setting the namelist parameter reg_def_mode=0: Switch automatically from increment specification to no. of grid points, when the reg_lon/lat_def(2) value is larger than 5.0.
- 1: reg_lon/lat_def(2) specifies increment
- 2: reg_lon/lat_def(2) specifies no. of grid points

For longitude values the last grid point is omitted if the end point matches the start point, e.g. for 0 and 360 degrees.

Examples

local grid with 0.5 degree increment:

global grid with 720x361 grid points:

reg_lon_def = -30.,0.5,30.

 $reg_lat_def = 90.,-0.5, -90.$

 $reg_lon_def = 0.,720,360.$

 $reg_lat_def = -90.,360,90.$

Time stamp format: The namelist parameters output_start, output_end, output_interval allow the specification of time stamps according to ISO 8601. The general format for time stamps is YYYY-MM-DDThh:mm:ss where Y: year, M: month, D: day for dates, and hh: hour, mm: minute, ss: second for time strings. The general format for durations is PnYnMnDTnHnMnS. See, for example, http://en.wikipedia.org/wiki/ISO_8601 for details and further specifications.

NOTE: as the mtime library underlaying the output driver currently has some restrictions concerning the specification of durations:

- 1. Any number n in PnYnMnDTnHnMnS must have two digits. For instance use "PT06H" instead of "PT6H"
- 2. In a duration string PnyearYnmonMndayDTnhrHnminMnsecS the numbers nxyz must not pass the carry over number to the next larger time unit: 0<=nmon<=12, 0<=nhr<=23, 0<=nmin<=59, 0<=nsec<=59.999. For instance use "P01D" instead of "PT24H", or "PT01M" instead of "PT60S".

Soon the formatting problem will be resolved and the valid number ranges will be enlarged. (2013-12-16).

Examples

group:land_vars
group:multisnow_vars

```
date and time representation (output_start, output_end) 2013-10-27T13:41:00Z duration (output_interval) POODTO6H00M00S
```

Variable Groups: Using the "group:" keyword for the namelist parameters ml_varlist, hl_varlist, pl_varlist, sets of common variables can be added to the output:

```
group:all output of all variables (caution: do not combine with mixed vertical interpolation)

group:atmo_ml_vars basic atmospheric variables on model levels

group:nh_prog_vars same set as atmo_ml_vars, but except pres and height, respectively

additional prognostic variables of the nonhydrostatic model

group:rad_vars

group:precip_vars

group:pl_vars

group:pbl_vars

group:phys_tendencies
```

tile-averaged variables

Keyword substitution in output filename (filename_format):

path substituted by model_base_dir output_filename substituted by output_filename physdom substituted by physical patch ID

levtype substituted by level type "ML", "PL", "HL", "IL"

levtype_l like levtype, but in lower case substituted by output file counter

datetime substituted by ISO-8601 date-time stamp in format YYYY-MM-DDThh:mm:ss.sssZ

datetime2 substituted by ISO-8601 date-time stamp in format YYYYMMDDThhmmssZ substituted by ISO-8601 date-time stamp in format YYYYMMDDThhmmss.sssZ

ddhhmmss substituted by relative day-hour-minute-second string substituted by relative hour-minute-second string

npartitions If namelist is split into concurrent files: number of stream partitions.

ifile_partition If namelist is split into concurrent files: stream partition index of this file.

total_index If namelist is split into concurrent files: substituted by the file index,

which an "unsplit" namelist would have produced

Defined and used in: src/namelists/mo name list output.f90

3.26 parallel nml

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	
division_file_name	С			Name of division file	$\operatorname{division_method} = 0$

Parameter	Type	Default	Unit	Description	Scope
ldiv_phys_dom	L	.TRUE.		.TRUE.: split into physical domains before	$\operatorname{division_method} = 1$
				computing domain decomposition (in case of	
				merged domains)	
				(This reduces load imbalance; turning off this	
				option is not recommended except for very small	
				processor numbers)	
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.	$p_{test_run} = .TRUE.$
				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	
use_dycore_barrier	L	.FALSE.		if .TRUE., set an MPI barrier at the beginning of	
				the nonhydrostatic solver (do not use for	
				production runs!)	
itype_exch_barrier	I	0		1: set an MPI barrier at the beginning of each MPI	
				exchange call	
				2: set an MPI barrier after each MPI WAIT call	
				3: 1+2 (do not use for production runs!)	
iorder_sendrecv	I	1		Sequence of send/receive calls:	
				$1 = \mathrm{irecv/send}$	
				$2=\mathrm{isend/recv}$	
				$3 = \mathrm{isend/irecv}$	
itype_comm	I	1		1: use local memory for exchange buffers	
				3: asynchronous halo communication for dynamical	
				core (currently deactivated)	
num_io_procs	I	0		Number of I/O processors (running exclusively for	
				$\operatorname{doing} I/O)$	

Parameter	Type	Default	Unit	Description	Scope
num_restart_procs	I	0		Number of restart processors (running exclusively	
				for doing restart)	
pio_type	I	1		Type of parallel I/O. Only used if number of I/O	
				processors greater number of domains.	
				Experimental!	
use_icon_comm	L	.FALSE.		Enable the use of MPI bulk communication through	
				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.	
use_dp_mpi2io	L	.FALSE.		Enable this flag if output fields shall be gathered by	
				the output processes in DOUBLE PRECISION.	
restart_chunk_size	I	1		Number of levels to be buffered by the	
				asynchronous restart process. The (asynchronous)	
				restart is capable of writing and communicating	
				more than one 2D slice at once.	

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

3.27 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 $lyr_perp = .TRUE$.	

Parameter	Type	Default	Unit	Description	Scope
isolrad	I	3		Insolation scheme 0: Use original SRTM insolation. 1: Use insolation from external file containing the spectrally resolved insolation (monthly means) 2: Use preindustrial insolation as in CMIP5 (average from 1844–1856) 3: Use insolation for AMIP-type CMIP5 simulation	
izenith	I	3 4 (for iforcing = inwp)		(average from 1979–1988) Choice of zenith angle formula for the radiative transfer computation. 0: Sun in zenith everywhere 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)	
albedo_type	I	1		Type of surface albedo 1: based on soil type specific tabulated values (dry soil) 2: MODIS albedo	iforcing=inwp

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	${ m irad_co2} = 2$
irad_o3		3		3: externally specified	and 0 for all the other
irad_o2		2		irad_aero = 5: Tanre aerosol climatology for	agents for
irad_cfc11		2		$run_nml/iforcing = 3 (NWP)$	$run_nml/iforcing = 2$
irad_cfc12		2		irad_aero = 6: Tegen aerosol climatology for	(ECHAM).
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			
${\rm vmr_cfc}12$		466.2e-12			

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

$3.28 \quad run_nml$

Parameter	Type	Default	Unit	Description	Scope
nsteps	I	0		number of time steps of this run.	
dtime	R	600.0	s	time step	
ltestcase	L	.TRUE.		Idealized testcase runs	

Parameter	Type	Default	Unit	Description	Scope
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_de			Number of full levels (atm.) for each domain	lvert_nest=.TRUE.
nshift	I(max_de	om)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			
activate_sync_timers	L	F		TRUE: Timer for monitoring runtime of	
				communication routines $(FALSE = off)$	
msg_level	I	10		controls how much printout is written during	
				runtime.	
				For values less than 5, only the time step is written.	
$msg_timestamp$	L	.FALSE.		If .TRUE., precede output messages by time stamp.	

Parameter	Type	Default	Unit	Description	Scope
test_mode	I	0		Setting a value larger than 0 activates a dummy	iequations = 3
				mode in which time stepping is changed into just	
				doing iterations, and MPI communication is	
				replaced by copying some value from the send	
				buffer into the receive buffer (does not work with	
				nesting and reduced radiation grid because the send	
				buffer may then be empty on some PEs)	
debug_check_level	I	0		Setting a value larger than 0 activates debug checks.	
output	C(:)	"nml","toti	nt"	Main switch for enabling/disabling components of	
				the model output. One or more choices can be set	
				(as an array of string constants). Possible choices	
				are:	
				• "none": switch off all output;	
				• "nml": new output mode (cf. output_nml);	
				• "totint": computation of total integrals.	
				If the output namelist parameter is not set	
				explicitly, the default setting "nml", "totint" is assumed.	
restart filename	C			File name for restart/checkpoint files (containg	
_				keyword substitution patterns <gridfile>, <idom>,</idom></gridfile>	
				<pre><rsttime>, <mtype>). default:</mtype></rsttime></pre>	
				" <gridfile>_restart_<mtype>_<rsttime>.nc".</rsttime></mtype></gridfile>	
profiling_output	I	1		controls how profiling printout is written:	
				TIMER_MODE_AGGREGATED=1,	
				$TIMER_MODE_DETAILED=2,$	
				TIMER_MODE_WRITE_FILES=3.	

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

$3.29 \quad sea_ice_nml$

Parameter	Type	Default	Unit	Description	Scope
i_ice_therm	I	2		Switch for thermodynamic model:	In an ocean run
				1: Zero-layer model	$i_{sea_ice must be} = 1.$
				2: Two layer Winton (2000) model	In an atmospheric run
				3: Zero-layer model with analytical forcing (for	the ice surface type must
				diagnostics)	be defined.
				4: Zero-layer model for atmosphere-only runs (for	
				diagnostics)	
i_ice_dyn	I	0		Switch for sea-ice dynamics:	
				0: No dynamics	
				1: FEM dynamics (from AWI)	
i_ice_albedo	I	1		Switch for albedo model. Only one is implemented	
				so far.	
i_Qio_type	I	2		Switch for ice-ocean heat-flux calculation method:	Defaults to 1 when
				1: Proportional to ocean cell thickness (like	i_ice_dyn=0 and 2
				MPI-OM)	otherwise.
				2: Proportional to speed difference between ice and	
				ocean	
kice	I	1		Number of ice classes (must be one for now)	
hnull	R	0.5	m	Hibler's h_0 parameter for new-ice growth.	
hmin	R	0.05	m	Minimum sea-ice thickness allowed.	
ramp_wind	R	10	days	Number of days it takes the wind to reach correct	
				strength. Only used at the start of an	
				OMIP/NCEP simulation (not after restart).	

$3.30 \quad sleve_nml \; (relevant \; if \; nonhydrostatic_nml:ivctype=2)$

Parameter	Type	Default	Unit	Description	Scope
min_lay_thckn	R	50	m	Layer thickness of lowermost layer; specifying zero	
				or a negative value leads to constant layer	
				thicknesses determined by top_height and nlev	

Parameter	Type	Default	Unit	Description	Scope
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	
lread_smt	L	.FALSE.		read smoothed topography from file (TRUE) or	
				compute internally (FALSE)	

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

$3.31 \quad time_nml$

Parameter	Type	Default	Unit	Description	Scope
calendar	I	1		Calendar type:	
				0=Julian/Gregorian	
				1=proleptic Gregorian	
				2=30 day/month, 360 day/year	
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 will	
				be reset (by the model) to dt_restart so that at	
				least one restart file is generated during the restart	
				cycle. If dt_restart is larger than but not a	
				multiple of dt_checkpoint, restart file will NOT be	
				generated at the end of the restart cycle.	
end_datetime_string	С	2008-09-		End date and time of the simulation	
		01T01:40:0	0Z'		

Parameter	Type	Default	Unit	Description	Scope
ini_datetime_string	С	'2008-09-		Initial date and time of the simulation	
		01T00:00:0	0 Z '		
is_relative_time	L	.FALSE.		.TRUE., if time loop shall start with step 0	
				regardless whether we are in a standard run or in a	
				restarted run (which means re-initialized run).	

Length of the run If "nsteps" in run_nml 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".

$3.32 \quad transport_nml \; (used \; if \; run_nml/ltransport=.TRUE.)$

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:	
		5		0: no horiz. transport (note that tracer mass ρq	
				instead of the specific tracer quantity q is kept	
				constant)	
				1: upwind (1st order)	
				2: Miura (2nd order, lin. reconstr.)	
				3: Miura3 (quadr. or cubic reconstr.)	$lsq_high_ord \in [2,3]$
				4: FFSL (quadr. or cubic reconstr.)	$lsq_high_ord \in [2,3]$
				5: hybrid Miura3/FFSL (quadr. or cubic reconstr.)	$lsq_high_ord \in [2,3]$
				20: miura (2nd order, lin. reconstr.) with	
				subcycling	
				22: combination of miura and miura with	
				subcycling	
				32: combination of miura and miura with	
				subcycling	
				42: combination of FFSL and miura with	
				subcycling	

Parameter	Type	Default	Unit	Description	Scope
				52: combination of hybrid FFSL/Miura3 with	
				subcycling	
ivadv_tracer	I(ntracer)	3		Tracer specific method to compute vertical	lvadv_tracer=TRUE
				advection:	
				0: no vert. transport (note that tracer mass ρq	
				instead of the specific tracer quantity q is kept	
				constant)	
				1: upwind (1st 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	
				3: monotonous flux limiter	$ihadv_tracer \neq 'iup3[4]'$
				4: positive definite flux limiter	
itype_vlimit	I(ntracer)	1		Type of limiter for vertical transport:	
				0: no limiter	
				1: semi-monotone slope limiter	
				2: monotonous slope limiter	
				4: positive definite flux limiter	
niter_fct	I	1		number of iterations of monotone flux correction	$ihadv_tracer = 3, 32, 4$
				procedure	$itype_hlimit = 3$
beta_fct	R	1.005		factor of allowed over-/undershooting in	$ihadv_tracer = 3, 32, 4,$
				monotonous limiter	42, 5, 52
					$itype_hlimit = 3$
iord_backtraj	I	1		order of backward trajectory calculation:	
				1: first order	
				2: second order (iterative; currently 1 iteration	ihadv_tracer='miura'
				hardcoded)	

Parameter	Type	Default	Unit	Description	Scope
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	
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	
lclip_tracer	L	.FALSE.		Clipping of negative values	

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

$3.33 \quad turbdiff_nml$

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

Parameter	Type	Default	Unit	Description	Scope
lprfcor	L	.FALSE.		using the profile values of the lowest main level	$inwp_turb = 1$
				instead of the mean value of the lowest layer for	
				surface flux calulations	
lnonloc	L	.FALSE.		nonlocal calculation of vertical gradients used for	$inwp_turb = 1$
				turbul. diff.	
lcpfluc	L	.FALSE.		consideration of fluctuations of the heat capacity of	$inwp_turb = 1$
				air	
limpltkediff	L	.TRUE.		consideration of fluctuations of the heat capacity of	$inwp_turb = 1$
				air	
itype_wcld	I	2		type of water cloud diagnosis	$inwp_turb = 1$
itype_synd	I	2		type of diagnostics of synoptical near surface	$inwp_turb = 1$
				variables	
lconst_z0	L	.FALSE.		TRUE: horizontally homogeneous roughness lenght	$inwp_turb = 1$
				z0	
const_z0	R	0.001	m	value for horizontally homogeneous roughness	$inwp_turb = 1$
				lenght z0	lconst_z0=.TRUE.

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

3.34 vdiff nml

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

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 ha_testcase_nml (Scope: ltestcase=.TRUE. and iequations=[0,1,2] in run_nml)

Parameter	Type	Default	Unit	Description	Scope
ctest_name	С	'JWw'		Name of test case:	
				'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.,
					ntracer = 2
				'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	lshallow_water=.FALSE.

Parameter	Type	Default	Unit	Description	Scope
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_width	R	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)'
rh_wavenum	I	4		wave number	ctest_name= 'RH'
rh_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. 1: the zonal state defined in the JWs test case; other integers: isothermal state (T=300 K, ps=1000 hPa, u=v=0.)	ctest_name= 'HS'
lhs_vn_ptb	L	.TRUE.		Add random noise to the initial wind field in the Held-Suarez test.	ctest_name= 'HS'
hs_vn_ptb_scale	R	1.	m/s	Magnitude of the random noise added to the initial wind field in the Held-Suarez test.	ctest_name= 'HS'
lrh_linear_pres	L	.FALSE.		Initialize the relative humidity using a linear function of pressure.	ctest_name= 'JWw-Moist','APE', 'LDF-Moist'
rh_at_1000hpa	R	0.75		relative humidity $0,1$ at 1000 hPa	ctest_name= 'JWw-Moist','APE', 'LDF-Moist'
linit_tracer_fv	L	.TRUE.		Finite volume initialization for tracer fields	ctest_name='PA'

Parameter	Type	Default	Unit	Description	Scope
ape_sst_case	С	'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	
				'sst_ice': Control SST distribution with -1.8 C	
				above 64 N/S .	
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	C	'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
				'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. With	
				the jabw test case, including moisture.	
				'wk82': Initializes the Weisman Klemp test case	$l_limited_area = .TRUE.$
				'g lim area': Initializes a series of general	
				limited area test cases: itype_atmos_ana	
				determines the atmospheric profile,	
				itype_anaprof_uv determines the wind profile and	
				itype_topo_ana determines the topography	
				'dcmip_pa_12': Initializes Hadley-like	
				meridional circulation pure advection test case.	
				'dcmip_rest_200': atmosphere at rest test	lcoriolis = .FALSE.
				(Schaer-type mountain)	
				'dcmip_mw_2x': nonhydrostatic mountain	lcoriolis = .FALSE.
				waves triggered by Schaer-type mountain	
				'dcmip_gw_31': nonhydrostatic gravity waves	
				triggered by a localized perturbation (nonlinear)	

Parameter	Type	Default	Unit	Description	Scope
				'dcmip_gw_32': nonhydrostatic gravity waves	$l_limited_area = .TRUE.$
				triggered by a localized perturbation (linear)	and lcoriolis $=$.FALSE.
				'dcmip_tc_51': tropical cyclone test case with	lcoriolis = .TRUE.
				'simple physics' parameterizations (not yet	
				implemented)	
				'dcmip_tc_52': tropical cyclone test case with	lcoriolis = .TRUE.
				with full physics in Aqua-planet mode	
				'CBL': convective boundary layer simulations for	is_plane_torus= .TRUE.
				LES package on torus (doubly periodic) grid	
jw_up	R	1.0	m/s	amplitude of the u-perturbation in jabw test case	nh_test_name='jabw'
u0_mrw	R	20.0	m/s	wind speed for mrw(2) and mwbr_const cases	$nh_test_name =$
					'mrw(2)_nh' and
					'mwbr_const'
mount_height_mrw	R	2000.0	m	maximum mount height in $mrw(2)$ and	$nh_test_name =$
				$mwbr_const$	'mrw(2)_nh' and
					'mwbr_const'
mount_half_width	R	1500000.0	m	half width of mountain in mrw(2), mwbr_const	$nh_test_name =$
				and bell	'mrw(2)_nh',
					'mwbr_const' and 'bell'
mount_lonctr_mrw_deg	R	90.	degrees	lon of mountain center in mrw(2) and mwbr_const	$nh_test_name =$
					'mrw(2)_nh' and
					'mwbr_const'
mount_latctr_mrw_deg	R	30.	degrees	lat of mountain center in mrw(2) and mwbr_const	$nh_test_name =$
					'mrw(2)_nh' and
					'mwbr_const'
temp_i_mwbr_const	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_const	R	0.025	1/s	constant brunt vaissala frequency at upper layer for	nh_test_name=
				mwbr_const case	'mwbr_const'
mount_height	R	100.0	m	peak height of mountain	nh_test_name= 'bell'

Parameter	Type	Default	Unit	Description	Scope
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	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_length	R	100000.0	m	length of slice domain	nh_test_name = 'bell', lplane=.TRUE.
rotate_axis_deg	R	0.0	deg	Earth's rotation axis pitch angle	nh_test_name= 'PA'
lhs_nh_vn_ptb	L	.TRUE.		Add random noise to the initial wind field in the Held-Suarez test.	nh_test_name= 'HS_nh'
lhs_fric_heat	L	.FALSE.		add frictional heating from Rayleigh friction in the Held-Suarez test.	nh_test_name= 'HS nh'
hs_nh_vn_ptb_scale	R	1.	m/s	Magnitude of the random noise added to the initial wind field in the Held-Suarez test.	nh_test_name= 'HS nh'
rh_at_1000hpa	R	0.7	1	relative humidity at 1000 hPa	nh_test_name= 'jabw', nh_test_name= 'mrw'
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	pure advection tests, only
lcoupled_rho	L	.FALSE.		Integrate density equation 'offline'	pure advection tests, only

Parameter	Type	Default	Unit	Description	Scope
qv_max_wk	R	0.014	Kg/kg	maximum specific humidity near	nh_test_name='wk82'
				the surface, range 0.012 - 0.016	
				used to vary the buoyancy	
u_infty_wk	R	20.	m/s	zonal wind at infinity height	nh_test_name='wk82'
				range 0 45.	
				used to vary the wind shear	
bub_amp	R	2.	K	maximum amplitud of the thermal perturbation	nh_test_name='wk82'
bubctr_lat	R	0.	deg	latitude of the center of the thermal perturbation	nh_test_name='wk82'
bubctr_lon	R	90.	deg	longitude of the center of the thermal perturbation	nh_test_name='wk82'
bubctr_z	R	1400.	m	height of the center of the thermal perturbation	nh_test_name='wk82'
bub_hor_width	R	10000.	m	horizontal radius of the thermal perturbation	nh_test_name='wk82'
bub ver width	R	1400.	m	vertical radius of the thermal perturbation	nh test name='wk82'
itype_atmo_ana	I	1		kind of atmospheric profile:	nh test name=
·				1 piecewise N constant layers	'g lim area'
				2 piecewise polytropic layers	
itype anaprof uv	I	1		kind of wind profile:	nh test name=
				1 piecewise linear wind layers	'g lim area'
				2 constant zonal wind	
				3 constant meridional wind	
itype_topo_ana	I	1		kind of orography:	nh_test_name=
				1 schaer test case mountain	'g_lim_area'
				2 gaussian_2d mountain	
				3 gaussian_3d mountain	
				any other no orography	
nlayers_nconst	I	1		Number of the desired layers with a constant	nh_test_name=
				Brunt-Vaisala-frequency	'g_lim_area' and
					itype_atmo_ana=1
p_base_nconst	R	100000.	Pa	pressure at the base of the first N constant layer	nh_test_name=
					'g_lim_area' and
					itype_atmo_ana=1
theta0_base_nconst	R	288.	K	potential temperature at the base of the first N	nh_test_name=
				constant layer	'g_lim_area' and
					itype_atmo_ana=1

Parameter	Type	Default	Unit	Description	Scope
h_nconst	R(nlayers_	n00,n\$5)00.,	m	height of the base of each of the N constant layers	nh_test_name=
		12000.			'g_lim_area' and
					itype_atmo_ana=1
N_nconst	R(nlayers_	notonst)	1/s	Brunt-Vaisala-frequency at each of the N constant	nh_test_name=
				layers	'g_lim_area' and
					itype_atmo_ana=1
rh_nconst	R(nlayers_	n 0o nst)	%	relative humidity at the base of each N constant	nh_test_name=
				layers	'g_lim_area' and
					itype_atmo_ana=1
rhgr_nconst	R(nlayers_	n0onst)	%	relative humidity gradient at each of the N constant	nh_test_name=
				layers	'g_lim_area' and
					itype_atmo_ana=1
nlayers_poly	I	2		Number of the desired layers with constant gradient	nh_test_name=
				temperature	'g_lim_area' and
					itype_atmo_ana=2
p_base_poly	R	100000.	Pa	pressure at the base of the first polytropic layer	$nh_test_name =$
					'g_lim_area' and
					itype_atmo_ana=2
h_poly	R(nlayers_	p 0 ly)12000.	m	height of the base of each of the polytropic layers	nh_test_name=
					'g_lim_area' and
					itype_atmo_ana=2
t_poly	R(nlayers_	p 28 8., 213.	K	temperature at the base of each of the polytropic	$nh_test_name =$
				layers	'g_lim_area' and
					itype_atmo_ana=2
rh_poly	R(nlayers_	p ∂l§) 0.2	%	relative humidity at the base of each of the	$nh_test_name =$
				polytropic layers	'g_lim_area' and
					itype_atmo_ana=2
rhgr_poly	R(nlayers_	p 5 l y -)5, 0.	%	relative humidity gradient at each of the polytropic	nh_test_name=
				layers	'g_lim_area' and
					itype_atmo_ana=2
nlayers_linwind	I	2		Number of the desired layers with constant U	nh_test_name=
				gradient	'g_lim_area' and
					itype_anaprof_uv=1

Parameter	Type	Default	Unit	Description	Scope
h_linwind	R(nlayers_	li :0wi2x50 0.	m	height of the base of each of the linear wind layers	nh_test_name=
					'g_lim_area' and
					itype_anaprof_uv=1
u_linwind	R(nlayers_	li 5 w 10 d)	m/s	zonal wind at the base of each of the linear wind	nh_test_name=
				layers	'g_lim_area' and
					itype_anaprof_uv=1
ugr_linwind	R(nlayers_	lingwind)	1/s	zonal wind gradient at each of the linear wind layers	nh_test_name=
					'g_lim_area' and
					itype_anaprof_uv=1
vel_const	R	20.	m/s	constant zonal/meridional wind	nh_test_name=
				$(itype_anaprof_uv=2,3)$	'g_lim_area' and
					itype_anaprof_uv=2,3
$mount_lonc_deg$	R	90.	deg	longitud of the center of the mountain	nh_test_name=
					'g_lim_area'
$mount_latc_deg$	R	0.	deg	latitud of the center of the mountain	nh_test_name=
					'g_lim_area'
schaer_h0	R	250.	m	h0 parameter for the schaer mountain	nh_test_name=
					'g_lim_area' and
					itype_topo_ana=1
schaer_a	R	5000.	m	-a- parameter for the schaer mountain,	nh_test_name=
				also half width in the north and south side of the	'g_lim_area' and
				finite ridge to round the sharp edges	itype_topo_ana=1,2
schaer_lambda	R	4000.	m	lambda parameter for the schaer mountain	nh_test_name=
					'g_lim_area' and
					itype_topo_ana=1
lshear_dcmip	L	FALSE		run dcmip_mw_2x with/without vertical wind	nh_test_name=
				shear	'dcmip_mw_2x'
				FALSE: dcmip_mw_21: non-sheared	
				TRUE : dcmip_mw_22: sheared	
halfwidth_2d	R	10000.	m	half length of the finite ridge in the north-south	nh_test_name=
				direction	'g_lim_area' and
					itype_topo_ana=1,2

Parameter	Type	Default	Unit	Description	Scope
m_height	R	1000.	m	height of the mountain	nh_test_name=
					'g_lim_area' and itype topo ana=2,3
m_width_x	R	5000.	m	half width of the gaussian mountain in the	nh_test_name=
				east-west direction	'g_lim_area' and
				half width in the north-south direction in the rounding of the finite ridge (gaussian 2d)	itype_topo_ana=2,3
m_width_y	R	5000.	m	half width of the gaussian mountain in the	nh_test_name=
				north-south direction	'g_lim_area' and
					itype_topo_ana=2,3
gw_u0	R	0.	m/s	maximum amplitude of the zonal wind	nh_test_name=
					'dcmip_gw_3X'
gw_clat	R	90.	deg	Lat of perturbation center	nh_test_name=
					'dcmip_gw_3X'
gw_delta_temp	R	0.01	K	maximum temperature perturbation	nh_test_name=
					'dcmip_gw_32'
u_cbl(2)	R	0:0	m/s	to prescribe initial zonal velocity profile for	nh_test_name=CBL
			and	convective boundary layer simulations where	
			1/s	u_cbl(1) sets the constant and u_cbl(2) sets the	
			,	vertical gradient	
$v_{cbl}(2)$	R	0:0	m/s	to prescribe initial meridional velocity profile for	nh_test_name=CBL
			and	convective boundary layer simulations where	
			1/s	v_cbl(1) sets the constant and v_cbl(2) sets the	
1 11(0)	D	200 0 000	T7 1	vertical gradient	CDI
$th_cbl(2)$	R	290:0.006	K and	to prescribe initial potential temperature profile for	nh_test_name=CBL
			K/m	convective boundary layer simulations where	
				th_cbl(1) sets the constant and th_cbl(2) sets the	
				gradient	

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

5 External data

$5.1 \quad extpar_nml \ (Scope: itopo=1 \ in \ run_nml)$

Parameter	Type	Default	Unit	Description	Scope
itopo	I	0		0: analytical topography/ext. data	
				1: topography/ext. data read from file	
n_iter_smooth_topo	I(n_dom)	0		iterations of topography smoother	itopo = 1
fac_smooth_topo	R	0.015625		pre-factor of topography smoother	n_iter_smooth_topo >
					0
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	itopo = 1
extpar_filename	С			Filename of external parameter input file, default:	
				" <path>extpar_<gridfile>". May contain the</gridfile></path>	
				keyword <path> which will be substituted by</path>	
				model_base_dir.	
extpar_varnames_map_fi	le C	, ,		Filename of external parameter dictionary, This is a	
				text file with two columns separated by whitespace,	
				where left column: NetCDF name, right column:	
				GRIB short name.	

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

6 External packages

$6.1 \quad art_nml$

Parameter	Type	Default	Unit	Description	Scope
lart	L	.FALSE.		main switch for ART-package	
lemi_volc	L	.FALSE.		Emission of volcanic ash	

Parameter	Type	Default	Unit	Description	Scope
lconv_tracer	L	.FALSE.		Convection of tracers	
lwash_tracer	L	.FALSE.		Washout of tracers	
lrad_volc	L	.FALSE.		Radiative impact of volcanic ash	
lcld_tracer	L	.FALSE.		Impact on clouds	

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

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.

Changes incompatible with former versions of the model code

 $\begin{array}{ll} \textit{Change:} & \text{var_names_map_file, out_varnames_map_file} \\ \textit{Date of Change:} & 2013\text{-}04\text{-}25 \end{array}$

12016

- $\bullet \ {\rm Renamed} \ {\bf var} \quad {\bf names} \quad {\bf map} \quad {\bf file} \rightarrow {\bf output} \quad {\bf nml} \quad {\bf dict}.$
- ullet Renamed out varnames map file \to netcdf dict.
- The dictionary in netcdf dict is now reversed, s.t. the same map file as in output nml dict can be used to translate variable names to the ICON internal names and back.

output nml: namespace

2013-04-26 12051

• Removed obsolete namelist variable namespace from output_nml.

 $\begin{array}{ll} \textit{Change:} & \text{gribout_nml: generatingCenter, generatingSubcenter} \\ \textit{Date of Change:} & 2013\text{-}04\text{-}26 \\ \textit{Revision:} & 12051 \end{array}$

- Introduced new namelist variables generatingCenter and generatingSubcenter.
- If not set explicitly, center and subcenter information is copied from the input grid file

Change:radiation_nml: albedo_typeDate of Change:2013-05-03Revision:12118

- $\bullet\,$ Introduced new namelist variable ${\bf albedo_type}$
- If set to 2, the surface albedo will be based on the MODIS data set.

Change:initicon_nml: dwdinc_filenameDate of Change:2013-05-24Revision:12266

 \bullet Renamed dwdinc_filename to dwdana_filename

Change:initicon_nml: l_ana_sfcDate of Change:2013-06-25Revision:12582

 \bullet Introduced new namelist flag <code>l_ana_sfc</code>

• If true, soil/surface analysis fields are read from the analysis fiel dwdfg filename. If false, surface analysis fields are not read. Soil and surface are initialized with the first guess instead.

Change: new_nwp_phy_tend_list: output names consistent with variable names

Date of Change: 2013-06-25

12590

- temp tend radlw \rightarrow ddt temp radlw
- temp tend turb \rightarrow ddt temp turb
- temp tend $drag \rightarrow ddt$ temp drag

 $\begin{array}{ll} \textit{Change:} & \text{prepicon_nml, remap_nml, input_field_nml} \\ \textit{Date of Change:} & 2013\text{-}06\text{-}25 \\ \textit{Revision:} & 12597 \end{array}$

- Removed the sources for the "prepicon" binary!
- The "prepicon" functionality (and most of its code) has become part of the ICON tools.

 \bullet The number of vertical input levels is now read from file. The namelist parameter ${\bf nlev_in}$ has become obsolete in r12700 and has been removed.

14160

• The namelist parameter exch msgsize has been removed together with the option iorder sendrecv=4.

 $\begin{array}{ll} {\it Change:} & {\it parallel_nml} \\ {\it Date of Change:} & {\it 2013-08-14} \\ {\it Revision:} & {\it 14164} \end{array}$

• The namelist parameter **use_sp_output** has been replaced by an equivalent switch **use_dp_mpi2io** (with an inverse meaning, i.e. we have use dp mpi2io = .NOT. use sp output).

 $\begin{array}{ll} \textit{Change:} & \text{parallel_nml} \\ \textit{Date of Change:} & \textbf{2013-08-15} \\ \textit{Revision:} & \textbf{14175} \end{array}$

• The above-mentioned namelist parameter use dp mpi2io got the default .FALSE. By this, the output data are sent now in single precision to the output processes.

Change:initicon_nml: l_ana_sfcDate of Change:2013-10-21Revision:14280

• The above-mentioned namelist parameter l ana sfc has been replaced by lread ana. The default is set to .TRUE., meaning that analysis fields are required and read on default. With Iread ana=.FALSE. ICON is able to start from first guess fields only.

Change:output_nml:lwrite_ready, ready_directoryDate of Change:2013-10-25Revision:14391

• The namelist parameters lwrite_ready and ready_directory have been replaced by a single namelist parameter ready_file, where ready_file /= 'default' enables writing ready files.

• Different output_nml's may be joined together to form a single ready file event - they share the same ready_file.

Change:output_nml: output_boundsDate of Change:2013-10-25Revision:14391

 $\bullet \ \ \text{The namelist parameter } \mathbf{output_bounds} \ \text{specifies a start}, \ \text{end}, \ \text{and increment of output invervals}. \ \ \text{It does no longer allow multiple}$ triples.

Change:output_nml: steps_per_fileDate of Change:2013-10-30Revision:14422

 \bullet The default value of the namelist parameter ${\bf steps_per_file}$ has been changed to -1.

 $\begin{array}{ll} \textit{Change:} & \text{run_nml} \\ \textit{Date of Change:} & \textbf{2013-11-13} \\ \textit{Revision:} & \textbf{14759} \end{array}$

- The dump/restore functionality for domain decompositions and interpolation coefficients has been removed from the model code. This means, that the parameters
 - ldump_states,
 - lrestore_states,
 - ldump_dd,
 - lread_dd,
 - nproc_dd,
 - dd_filename,
 - dump_filename,

```
- l_one_file_per_patch
```

have been removed together with the corresponding functionality from the ICON model code.

 $\begin{array}{ll} \textit{Change:} & \text{output_nml: filename_format} \\ \textit{Date of Change:} & \textbf{2013-12-02} \\ \textit{Revision:} & \textbf{15068} \end{array}$

• The string token <ddhhmmss> is now substituted by the relative day-hour-minute-second string, whereas the absolute date-time stamp can be inserted using <datetime>.

 $\begin{array}{ll} {\it Change:} & {\it output_nml: ready_file} \\ {\it Date of Change:} & {\it 2013-12-03} \\ {\it Revision:} & {\it 15081} \end{array}$

• The ready file name has been changed and may now contain string tokens cpath>, <datetime>, <ddhmmss> which are substituted as described for the namelist parameter filename_format.

 $\begin{array}{ll} \textit{Change:} & \text{interpl_nml: rbf_vec_scale_ll} \\ \textit{Date of Change:} & \textbf{2013-12-06} \end{array}$

15156

- The real-valued namelist parameter rbf_vec_scale_11 has been removed.
- Now, there exists a new integer-valued namelist parameter, rbf_scale_mode_11 which specifies the mode, how the RBF shape parameter is determined for lon-lat interpolation.

- Removed remaining vlist-related namelist parameter. This means that the parameters
 - out filetype
 - out expname
 - dt_data
 - dt file
 - lwrite_dblprec, lwrite_decomposition, lwrite_vorticity, lwrite_divergence, lwrite_pres, lwrite_z3, lwrite_tracer, lwrite_tend_phy, lwrite radiation, lwrite precip, lwrite cloud, lwrite tke, lwrite surface, lwrite omega, lwrite initial, lwrite oce timestepping

are no longer available.

gridref nml Date of Change: gridrei_nm
2014-01-07 15436

• Changed namelist defaults for nesting: grf_intmethod_e, 1_mass_consvcorr, 1_density_nudging.

 $rac{\mathrm{interpol_nml}}{2014\text{-}02\text{-}10}$ Change:
Date of Change: 16047

• Changed namelist default for rbf_scale_mode_11: The RBF scale factor for lat-lon interpolation is now determined automatically by default.

echam_phy_nml 2014-02-27

Revision:16313

• Replace the logical switch lcover by the integer switch icover that is used in ECHAM-6.2. Values are transferred as follows: .FALSE. = 1 (=default), .TRUE. = 2.