

ecRad radiation scheme: User Guide (draft)

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Chapter 1

Introduction

1.1 What is *ecRad*?

ecRad is an atmospheric radiation scheme designed for computing profiles of solar (or *shortwave*) and thermal infrared (or *longwave*) irradiances from the surface up to the middle mesosphere. It is incorporated into the Integrated Forecasting System (IFS), the weather forecast model used operationally by the European Centre for Medium Range Weather Forecasts (ECMWF), in which it is used to compute radiative heating and cooling rates of the atmosphere and surface. An offline version of the scheme is available for educational and non-commercial research uses only.

A scientific overview of *ecRad* was provided by [Hogan and Bozzo \(2018\)](#). It incorporates the Rapid Radiative Transfer Model for GCMs (RRTMG; [Iacono et al., 2008](#)) for representing absorption by atmospheric gases and a flexible treatment of the optical properties of aerosol particles. Three different solvers capable of representing the effects of subgrid cloud structure: McICA ([Pincus et al., 2003](#)), Tripleclouds ([Shonk and Hogan, 2008](#)) and SPARTACUS ([Hogan et al., 2016](#)). It is coded in Fortran 2003 in a way that is efficient and flexible.

1.2 License

The Meteorological Services of ECMWF Member States are permitted to use *ecRad* (along with any IFS software) for any in-house purpose. Other researchers may apply for an institutional license to use the software, as described on the *ecRad* web site*. The license is the same as for the OpenIFS software and permits educational and non-commercial research use only, and prohibits redistribution. The full terms of the license are available on the *ecRad* web site.

1.3 Overview of this document

Chapter 2 describes how to compile and use the offline version of *ecRad*, which is essentially a Unix program that reads a configuration file and a NetCDF file containing a description of the atmospheric state, and outputs a NetCDF file containing the computed irradiance profiles. At some future date this document will be expanded to include chapters describing the API (for incorporating *ecRad* into a larger Fortran program), the internal architecture and the detailed scientific documentation.

*<https://confluence.ecmwf.int/display/ECRAD>

Chapter 2

Using the offline radiation scheme

2.1 Compiling the package

The offline version of *ecRad* is designed to be used on a Unix-like platform. You will need a Fortran compiler that supports the 2003 standard such as `gfortran`. As a prerequisite, you will need to install the NetCDF library, including the Fortran interface (packages to install on a Linux system are typically called `libnetcdf-dev` or `libnetcdf-devel`). To run some of the tests, you will also need to install the `nc` utilities for manipulating NetCDF Files.

First unpack the package and enter the subdirectory as follows:

```
tar xvfz ecrad-1.1.0.tar.gz
cd ecrad-1.1.0
```

On a non-GNU platform you may need to untar and unzip the package using the `tar` and `gunzip` commands separately. The `README` file contains concise instructions on compilation and testing, while the `NOTICE` file outlines the license conditions. The subdirectories are as follows:

radiation The *ecRad* source code for atmospheric radiation

radsurf The *ecRad* source code for surface radiation (under development)

ifsaux Source code providing a (sometimes dummy) IFS environment

ifsrrtm The IFS implementation of the RRTMG gas optics scheme

utilities Source code for useful utilities, such as reading NetCDF files

driver The source code for the offline driver program `ecrad`

ifs Source files from the IFS that are used to illustrate how *ecRad* can be incorporated into a large model, but note that these files are not used in the offline version

mod Where Fortran module files are written

lib Where the static libraries are written

bin Where the executable `ecrad` is written

data Contains configuration data files read at run-time

test Test cases including Matlab code to plot the outputs

include Automatically generated interface blocks for non-module routines

Compilation on different platforms using different compilers is facilitated by the various `Makefile.include.<prof>` files in the top-level directory: if you type

```
make PROFILE=gfortran
```

the code will be compiled using the `gfortran` compiler via the Makefile variables set in the `Makefile_include.gfortran` file. If everything goes to plan this should create the executable `bin/ecrad` and various static libraries in the `lib` directory. The other profiles are `pgi` (for the PGI compiler), `ecmwf` and `uor` (more specific configurations for the ECMWF and University of Reading computer systems). If you wish to add a new configuration for a different compiler, the cleanest way is to create a new `Makefile_include.<prof>` file modelled on one of the existing ones.

To compile in single precision, type

```
make PROFILE=gfortran SINGLE\_PRECISION=1
```

To compile with debugging options (no optimization, bounds checking and initializing real numbers with not-a-number), type

```
make PROFILE=gfortran DEBUG=1
```

Finer tuning may be achieved by specifying the optimization and debugging flags explicitly, for example

```
make PROFILE=gfortran OPTFLAGS="-O1" DEBUGFLAGS="-g1 -pg"
```

Remember that if you change the compile settings you will probably want to recompile everything, in which case you first need to remove all compiled files with

```
make clean
```

2.2 Running the offline radiation scheme

To test the code, type

```
make test
```

which runs `make` in each of the subdirectories of the `test` directory. The `README` files in these directories provide more information on what they are doing, and some Matlab scripts are provided to visualize the outputs.

You will see in the output of the tests the command line in each invocation of *ecRad*, which is of the form

```
ecrad config.nam input.nc output.nc
```

where `ecrad` needs to be the full path to the *ecRad* executable, `config.nam` is a Fortran namelist file configuring the code, `input.nc` contains the input atmospheric profiles and `output.nc` contains the output irradiance (flux) profiles. The namelist file contains a `radiation` namelist that configures the *ecRad* scheme itself; the variables available are described in section 2.3. The file also contains a `radiation_config` namelist that configures aspects of the offline package. Only the `radiation` namelist is used when *ecRad* is incorporated into an atmospheric model.

The input NetCDF file contains numerous floating-point variables listed in Table 2.1. The dimensions are shown in the order that they are listed by the `ncdump` utility, with the first dimension varying slowest in the file. Note that this is opposite to the internal Fortran ordering. Most variables are stored as a function of column and level (dimensions named `col` and `level` in Table 2.1, although the actual dimension names are ignored by *ecrad*). The `half_level` dimension corresponds to the mid-points of the levels, plus the top-of-atmosphere and surface, and so must be one more than `level`. The `level_interface` dimension excludes the top-of-atmosphere and surface so must be one less than `level`. The optional `sw_albedo_band` and `lw_emiss_band` dimensions allow for shortwave albedo and longwave emissivity to be specified in user-defined spectral intervals, but in the offline code these are ignored: the first element along these dimensions will be used for the entire shortwave and longwave spectrum.

Table 2.1: Main variables contained in the input NetCDF file to *ecRad*. Note that some variables are not required if they are not used by the particular solver selected, for example `iseed` is only needed if the McICA solver is specified. Also, only one of `o3_mmr` and `o3_vmr` should be provided. In addition to ozone, further gases can be specified in either mass mixing ratio (suffix `_mmr`) or volume mixing ratio (suffix `_vmr`) units, where the prefixes are `co2` (carbon dioxide), `n2o` (nitrous oxide), `co` (carbon monoxide), `ch4` (methane), `o2` (molecular oxygen), `cfc11` (CFC-11), `cfc12` (CFC-12), `hcfcc22` (HCFC-22), `cc14` (carbon tetrachloride) and `no2` (nitrogen dioxide). These further trace gases may either be specified as variable (dimensioned `col, level`) or constant (a scalar value in the file).

Variable	Dimensions	Description
<code>solar_irradiance</code>	–	Solar irradiance at Earth’s orbit (W m^{-2})
<code>skin_temperature</code>	<code>col</code>	Skin temperature (K)
<code>cos_solar_zenith_angle</code>	<code>col</code>	Cosine of solar zenith angle
<code>sw_albedo</code>	<code>col, sw_albedo_band</code>	Shortwave albedo (if 1D then assumed spectrally constant)
<code>lw_emissivity</code>	<code>col, lw_emiss_band</code>	Longwave emissivity (if 1D then assumed spectrally constant)
<code>iseed</code>	<code>col</code>	Seed for McICA random-number generator (double precision)
<code>pressure_hl</code>	<code>col, half_level</code>	Pressure at half levels (Pa)
<code>temperature_hl</code>	<code>col, half_level</code>	Temperature at half levels (K)
<code>q</code>	<code>col, level</code>	Specific humidity (kg kg^{-1})
<code>o3_mmr</code>	<code>col, level</code>	Ozone mass mixing ratio (kg kg^{-1})
<code>o3_vmr</code>	<code>col, level</code>	Ozone volume mixing ratio ($\text{m}^3 \text{m}^{-3}$), used only if <code>o3_mmr</code> not provided
<code>q_liquid</code>	<code>col, level</code>	Liquid cloud mass mixing ratio (kg kg^{-1})
<code>q_ice</code>	<code>col, level</code>	Ice cloud mass mixing ratio (kg kg^{-1})
<code>re_liquid</code>	<code>col, level</code>	Liquid cloud effective radius (m)
<code>re_ice</code>	<code>col, level</code>	Ice cloud effective radius (m)
<code>overlap_param</code>	<code>col, level, interface</code>	Cloud overlap parameter
<code>fractional_std</code>	<code>col, level</code>	Fractional standard deviation of cloud optical depth
<code>inv_cloud_effective_size</code>	<code>col, level</code>	Inverse of cloud effective horizontal size for SPARTACUS solver (m^{-1})
<code>aerosol_mmr</code>	<code>col, aerosol_type, level</code>	Aerosol mass mixing ratio (kg kg^{-1})

All the test data store input fields in order of increasing pressure, i.e. starting at the top-of-atmosphere and working down to the surface. The output data are then provided using the same convention. If input data are provided in the opposite order then this should be automatically detected and under the bonnet the order is reversed before being passed to the radiation scheme. But if you use this convention then please test the results carefully.

The output NetCDF file contains the typical set of variables listed in Table 2.2. Clear-sky fluxes (i.e. computed on the same input profiles but in the absence of clouds) are provided if the `do_clear` namelist variable is set to `true` (see section 2.3) while surface spectral shortwave surface fluxes are provided if the `do_surface_sw_spectral_flux` is set to `true`. Note that if you want atmospheric heating rates then you will need to compute them yourself from the flux profiles.

Table 2.2: Variables contained in the output NetCDF file from *ecRad*, where all fluxes (or irradiances) have units of W m^{-2} . The `band_sw` dimension has the same size as the number of shortwave bands in the gas-optics scheme.

Variable	Dimensions	Description
<code>pressure_hl</code>	<code>col, half_level</code>	Pressure at half levels (Pa)
<code>flux_up_sw, flux_dn_sw</code>	<code>col, half_level</code>	Up- and downwelling shortwave fluxes
<code>flux_up_sw_clear, flux_dn_sw_clear</code>	<code>col, half_level</code>	Up- and downwelling clear-sky shortwave fluxes
<code>flux_dn_direct_sw</code>	<code>col, half_level</code>	Direct component of downwelling shortwave flux
<code>flux_dn_direct_sw_clear</code>	<code>col, half_level</code>	Direct component of downwelling clear-sky shortwave flux
<code>flux_up_lw, flux_dn_lw</code>	<code>col, half_level</code>	Up- and downwelling longwave fluxes
<code>flux_up_lw_clear, flux_dn_lw_clear</code>	<code>col, half_level</code>	Up- and downwelling clear-sky longwave fluxes
<code>lw_derivative</code>	<code>col, half_level</code>	Derivative of upwelling longwave flux with respect to surface value (Hogan and Bozzo, 2015)
<code>spectral_flux_dn_sw_surf</code>	<code>col, band_sw</code>	Downwelling surface shortwave flux in each band
<code>spectral_flux_dn_direct_sw_surf</code>	<code>col, band_sw</code>	Direct downwelling surface shortwave flux in each band

spectral_flux_dn_sw_surf_clear	col, band_sw	Clear-sky downwelling surface shortwave flux in each band
spectral_flux_dn_direct_sw_surf_clear	col, band_sw	Clear-sky direct downwelling surface shortwave flux in each band
cloud_cover_sw	col	Total cloud cover diagnosed by shortwave solver
cloud_cover_lw	col	Total cloud cover diagnosed by longwave solver

2.3 Configuring the radiation scheme

The detailed settings of *ecRad* are configured using the `radiation` namelist in the namelist file provided as the first command-line argument to the `ecrad` executable. The available namelist variables are listed in Table 2.3. One of the most important is `directory_name`, which provides the absolute or relative path to the directory containing all the configuration files. This is the `data` directory at the top level of the *ecRad* package.

Table 2.3: Options for the `radiation` namelist that configures the radiation scheme. The type of each variable can be inferred from its name: logicals begin with `do_` or `use_`, integers start with `i_` or `n_`, strings end with `_name`, and all other variables are real numbers.

Key	Default value, other values	Description
<i>General</i>		
<code>directory_name</code>	<code>.</code>	Directory containing NetCDF configuration files
<code>do_sw</code>	<code>true</code>	Compute shortwave fluxes?
<code>do_lw</code>	<code>true</code>	Compute longwave fluxes?
<code>do_sw_direct</code>	<code>true</code>	Do direct shortwave fluxes?
<code>do_clear</code>	<code>true</code>	Compute clear-sky fluxes?
<i>Gas and aerosol optics</i>		
<code>gas_model_name</code>	RRTMG-IFS , Monochromatic	Gas optics model
<code>use_aerosols</code>	<code>false</code>	Do we represent aerosols?
<code>do_lw_aerosol_scattering</code>	<code>true</code>	Do longwave aerosol scattering?
<code>n_aerosol_types</code>		Number of aerosol types
<code>i_aerosol_type_map</code>		Vector of integers that map from aerosol types to types in the NetCDF aerosol optics file, where positive indexes hydrophobic types, negative indexes hydrophilic types and zero indicates a type should be ignored
<i>Monochromatic scheme</i>		
<code>mono_lw_wavelength</code>	<code>-1.0</code>	Wavelength of longwave radiation, or if negative, a broadband calculation will be performed
<code>mono_lw_total_od</code>	<code>0.0</code>	Zenith optical depth of clear-sky atmosphere
<code>mono_sw_total_od</code>	<code>0.0</code>	Zenith optical depth of clear-sky atmosphere
<i>Cloud optics</i>		
<code>liquid_model_name</code>	SOCRATES , Slingo, Monochromatic	Liquid optics model, including the scheme in the SOCRATES radiation scheme and the older scheme of Slingo (1989)
<code>ice_model_name</code>	Fu-IFS , Baran2016, Yi, Monochromatic	Ice optics model, including the schemes of Fu (1996), Fu et al. (1998), Baran et al. (2016) and Yi et al. (2013)
<code>do_lw_cloud_scattering</code>	<code>true</code>	Do longwave cloud scattering?
<code>do_fu_lw_ice_optics_bug</code>	<code>false</code>	Reproduce bug in McRad implementation of Fu ice optics?
<i>Solver</i>		
<code>sw_solver_name</code>	Homogeneous, McICA , Tripleclouds, SPARTACUS	Shortwave solver
<code>lw_solver_name</code>	Homogeneous, McICA , Tripleclouds, SPARTACUS	Longwave solver
<code>overlap_scheme_name</code>	Max-Ran, Exp-Ran , Exp-Exp	Cloud overlap scheme
<code>use_beta_overlap</code>	<code>false</code>	Use Shonk et al. (2010) ‘ β ’ overlap parameter definition, rather than default ‘ α ’
<code>cloud_inhom_decorrr_scaling</code>	<code>0.5</code>	Ratio of overlap decorrelation lengths for cloud inhomogeneities and boundaries

cloud_fraction.threshold	10^{-6}	Ignore clouds with fraction below this
cloud_mixing_ratio.threshold	10^{-9}	Ignore clouds with total mixing ratio below this
cloud_pdf.shape.name	Gamma , Lognormal	Shape of cloud water PDF
cloud_pdf.override.file.name		Name of NetCDF file of alternative cloud PDF look-up table
do_sw.delta_scaling.with.gases	false	Apply delta-Eddington scaling to particle-gas mixture, rather than particles only
<i>SPARTACUS solver</i>		
do_3d.effects	true	Represent 3D effects when SPARTACUS solver selected
n_regions	2, 3	Number of regions
do_lw.side.emissivity	true	Represent effective emissivity of the side of clouds (Schäfer et al., 2016)
sw_encroachment.name	Minimum, Computed , Maximum	Encroachment (or ‘entrainment’) model
do_3d.lw.multilayer.effects	false	Maximum encroachment for longwave radiation?
max_3d.transfer.rate	10.0	Maximum rate of lateral exchange between regions in one layer
max_gas.od.3d	8.0	3D effects ignored for spectral intervals where gas optical depth of a layer exceeds this
max_cloud.od	18.0	Maximum in-cloud optical depth for stability
use_expm.everywhere	false	Use matrix-exponential method even when 3D effects not represented?
clear_to_thick.fraction	0.0	Fraction of cloud edge interfacing directly to the most optically thick cloudy region
overhead_sun.factor	0.0	Minimum tan-squared of solar zenith angle to allow some ‘direct’ radiation from overhead sun to pass through cloud sides (0.06 used by Hogan et al., 2016)
<i>Surface (under development)</i>		
use_canopy.full.spectrum.sw	false	Perform canopy shortwave radiative transfer at full atmospheric spectral resolution
use_canopy.full.spectrum.lw	false	Perform canopy longwave radiative transfer at full atmospheric spectral resolution
<i>Diagnostics</i>		
iverbose.setup	0, 1, 2, 3, 4, 5	Verbosity in setup, where 1=warning, 2=info, 3=progress, 4=detailed, 5=debug
iverbose	0, 1, 2, 3, 4, 5	Verbosity in execution
do_save.spectral.flux	false	Save flux profiles in each band?
do_save.gpoint.flux	false	Save flux profiles in each g-point?
do_surface.sw.spectral.flux	true	Save surface shortwave fluxes in each band for subsequent diagnostics?
do_lw.derivatives	false	Compute derivatives for Hogan and Bozzo (2015) approximate updates?
do_save.radiative.properties	false	Write intermediate NetCDF file(s) of properties sent to solver (radiative_properties*.nc)?

When calling *ecRad* from within a model, the variables listed in Table 2.3 are members of the `config_type` structure, and may be modified within the code at the appropriate place in the configuration stage. The exception is in the case of strings, which are prefixed by `_name` in the namelist. In the `config_type` structure there are equivalent integers to express these variables, which can be changed using the named constants listed in Table 2.4.

Table 2.4: Integers in the `config_type` structure that represents the strings in Table 2.3, where a namelist variable named `*_name` would be named `i_*` here.

Variable in <code>config_type</code>	Available named constants, default
<code>i_overlap.scheme</code>	<code>IOverlapMaximumRandom</code> , <code>IOverlapExponentialRandom</code> , <code>IOverlapExponential</code>
<code>i_solver.sw</code> , <code>i_solver.lw</code>	<code>ISolverHomogeneous</code> , <code>ISolverMcICA</code> , <code>ISolverSpartacus</code> , <code>ISolverTripleclouds</code>
<code>i_3d.sw.encroachment</code>	<code>IEncroachmentMinimum</code> , <code>IEncroachmentComputed</code> , <code>IEncroachmentMaximum</code>

i_gas_model	IGasModelMonochromatic, IGasModelIFSRRTMG
i_liq_model	ILiquidModelMonochromatic, ILiquidModelSOCRATES , ILiquidModelSlingo
i_ice_model	IIceModelMonochromatic, IIceModelFu , IIceModelBaran2016, IIceModelYi
i_cloud_pdf_shape	IPdfShapeGamma , IPdfShapeLognormal

When ecrad is run with the default verbosity settings, it outputs to the screen a summary of the configuration options, the files read and written and details of the aerosol mapping. The following is an example from the default test in the test/ifs directory.

```
----- OFFLINE ECRAD RADIATION SCHEME -----
Copyright (C) 2014-2018 European Centre for Medium-Range Weather Forecasts
Contact: Robin Hogan (r.j.hogan@ecmwf.int)
Floating-point precision: double
General settings:
  Data files expected in ".././data"
  Clear-sky calculations are ON                               (do_clear=T)
  Saving intermediate radiative properties OFF                 (do_save_radiative_properties=F)
  Saving spectral flux profiles OFF                             (do_save_spectral_flux=F)
  Saving surface shortwave spectral fluxes ON                  (do_surface_sw_spectral_flux=T)
  Gas model is RRTMG-IFS                                       (i_gas_model=1)
  Aerosols are ON                                              (use_aerosols=T)
  Longwave derivative calculation is ON                         (do_lw_derivatives=T)
Cloud settings:
  Cloud fraction threshold = .100E-05                         (cloud_fraction_threshold)
  Cloud mixing-ratio threshold = .100E-08                      (cloud_mixing_ratio_threshold)
  Liquid optics scheme is SOCRATES                             (i_liq_model=2)
  Ice optics scheme is Fu-IFS                                   (i_ice_model=2)
  Longwave ice optics bug in Fu scheme is OFF                  (do_fu_lw_ice_optics_bug=F)
  Cloud overlap scheme is Exp-Exp                              (i_overlap_scheme=2)
  Use "beta" overlap parameter is OFF                          (use_beta_overlap=F)
  Cloud PDF shape is Gamma                                     (i_cloud_pdf_shape=1)
  Cloud inhom decorrelation scaling = .500                     (cloud_inhom_decorr_scaling)
Solver settings:
  Shortwave solver is McICA                                    (i_solver_sw=1)
  Shortwave delta scaling after merge with gases OFF           (do_sw_delta_scaling_with_gases=F)
  Longwave solver is McICA                                     (i_solver_lw=1)
  Longwave cloud scattering is ON                              (do_lw_cloud_scattering=T)
  Longwave aerosol scattering is OFF                            (do_lw_aerosol_scattering=F)
Reading .././data/RADRRTM
Reading .././data/RADSRTM
Reading NetCDF file .././data/socrates_droplet_scattering_rrtm.nc
Reading NetCDF file .././data/fu_ice_scattering_rrtm.nc
Reading NetCDF file .././data/aerosol_ifs_rrtm_43R3.nc
Aerosol mapping:
  1 -> hydrophilic type 1: Sea salt, bin 1, 0.03-0.5 micron, OPAC
  2 -> hydrophilic type 2: Sea salt, bin 2, 0.50-5.0 micron, OPAC
  3 -> hydrophilic type 3: Sea salt, bin 3, 5.0-20.0 micron, OPAC
  4 -> hydrophobic type 1: Desert dust, bin 1, 0.03-0.55 micron, (SW) Dubovik et al 2002, (LW)...
  5 -> hydrophobic type 2: Desert dust, bin 2, 0.55-0.90 micron, (SW) Dubovik et al 2002, (LW)...
  6 -> hydrophobic type 3: Desert dust, bin 3, 0.90-20.0 micron, (SW) Dubovik et al 2002, (LW)...
  7 -> hydrophilic type 4: Hydrophilic organic matter, OPAC
  8 -> hydrophobic type 10: Hydrophobic organic matter, OPAC (hydrophilic at RH 20-30%)
  9 -> hydrophobic type 11: Black carbon, Olivier Boucher
  10 -> hydrophobic type 11: Black carbon, Olivier Boucher
  11 -> hydrophilic type 5: Ammonium sulfate, GACP
  12 -> hydrophobic type 14: Stratospheric sulfate (hydrophilic ammonium sulfate at RH 20-30%)
Reading NetCDF file .././data/mcica_gamma.nc
Reading NetCDF file ecrad_meridian.nc
Warning: variable co_vmr not found
Warning: variable no2_vmr not found
```

```

Writing NetCDF file inputs.nc
Thread 0 processing columns 1-32
Writing NetCDF file ecrad_meridian_default_out.nc
-----

```

2.4 Configuring the offline package

In addition to the namelist variables described in section 2.3 an additional set of variables are available in the `radiation_config` namelist that are specific to the offline version of *ecRad* and are listed in Table 2.5.

Table 2.5: Options for the `radiation_config` namelist that configures additional aspects of the offline radiation scheme. All entries must be scalars. If an override variable is present then it need not be included in the input file. The cloud effective sizes (used by the SPARTACUS solver) may be specified for low, middle and high clouds according to the cloud layer pressure p and the surface pressure p_0 .

Key	Description
<i>Execution control</i>	
<code>nrepeat</code>	Number of times to repeat, for benchmarking
<code>istartcol</code>	Start at specified input column (1 based)
<code>iendcol</code>	End at specified input column (1 based)
<code>iverbose</code>	Verbosity in offline setup (default 2)
<code>do_parallel</code>	Use OpenMP parallelism? (default <code>true</code>)
<code>nblocksize</code>	Number of columns per block when using OpenMP
<code>do_save_inputs</code>	Sanity check: save input variables in <code>inputs.nc</code>
<i>Override input variables</i>	
<code>solar_irradiance_override</code>	Override solar irradiance (W m^{-2})
<code>skin_temperature</code>	Override skin temperature (K)
<code>cos_solar_zenith_angle</code>	Override cosine of solar zenith angle
<code>sw_albedo</code>	Override shortwave albedo
<code>lw_emissivity</code>	Override longwave emissivity
<code>fractional_std</code>	Override cloud optical depth fractional standard deviation
<code>overlap_decorr_length</code>	Override cloud overlap decorrelation length (m)
<code>inv_effective_size</code>	Override inverse of cloud effective size (m^{-1})
<code>low_inv_effective_size</code>	...for low clouds ($p > 0.8p_0$)
<code>middle_inv_effective_size</code>	...for mid-level clouds ($0.45p_0 < p \leq 0.8p_0$)
<code>high_inv_effective_size</code>	...for high clouds ($p \leq 0.45p_0$)
<i>Scale input variables</i>	
<code>q_liquid_scaling</code>	Scaling for liquid water mixing ratio
<code>q_ice_scaling</code>	Scaling for ice water mixing ratio
<code>cloud_fraction_scaling</code>	Scaling for cloud fraction (capped at 1)
<code>overlap_decorr_length_scaling</code>	Scaling for cloud overlap decorrelation length
<code>effective_size_scaling</code>	Scaling for cloud effective size

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