

# RCE Setup with ICON LEM physics on a Torus

(using icon-2.3.00, following the [RCEMIP protocol](#))

## 1. Generation of a Torus Grid (100 km x 100 km with doubly-periodic lateral boundary conditions)

The GridGenerator ([git@git.mpimet.mpg.de:GridGenerator.git](https://git.mpimet.mpg.de/GridGenerator.git)) provides a script template `/run/grid_creators/grid.create_torus_grid`.

```
cp /run/grid_creators/grid.create_torus_grid into /run
```

```
make_runscript grid.create_torus_grid
```

Set appropriate values in the last line for number of rows and columns as well the edge length in meter, e.g.:

```
y_noOfRows = 22  
x_noOfColumns = 20  
edge_length = 5000 m
```

## 2. Generation of the Ozone input

a) Download

“profile\_analytical\_o3.nc”: [https://code.mpimet.mpg.de/attachments/19451/profile\\_analytical\\_o3.nc](https://code.mpimet.mpg.de/attachments/19451/profile_analytical_o3.nc). This file contains the 1-dimensional vertical profile of ozone concentration (in ppmv), as prescribed by the RCEMIP protocol.

b) expand to the full grid

```
cdo enlarge,Torus_Triangles_20x22_5000m.nc \  
    profile_analytical_o3.nc \  
    rcemip_analytical_o3_20x22_5000m.nc
```

When processed with newer versions of cdo, it may be required to rename cell to ncells for the NWP Setup:

```
ncrename -O -d cell,ncells rcemip_analytical_o3_20x22_5000m.nc \  
    rcemip_analytical_o3_20x22_5000m_nwp_ncells.nc  
mv rcemip_analytical_o3_20x22_5000m_ncells.nc  
rcemip_analytical_o3_20x22_5000m.nc
```

c) adapt the run script: set `irad_o3 = 4` in `radiation_nml` and link the file

```
add_link_file ${OZONEDIR}/rcemip_analytical_o3_20x22_5000m.nc  
o3_icon_DOM01.nc
```

### 3. Initial thermodynamic profile

To have the same initial thermodynamic profile as defined by the RCEMIP protocol, two new testcases need to be defined in the ICON source code, 'RCEMIP\_analytical' and 'RCEMIP\_ncdf'.

'RCEMIP\_analytical' initializes the base state using the analytical profiles of T,q defined by the RCEMIP protocol, and 'RCEMIP\_ncdf' reads in mean profiles of T,q from a netcdf file of a previous simulation, assuming an identical grid.

The new testcases are defined in `src/testcases/mo_nh_testcases.f90`. There the new subroutines 'init\_torus\_rcemip\_analytical\_sounding' and 'init\_torus\_rcemip\_ncdf\_sounding' are called, contained in `src/testcases/mo_nh_torus_exp.f90`.

In total, the following src files have been adapted:

[testcases/mo\\_nh\\_testcases.f90](#)

[testcases/mo\\_nh\\_torus\\_exp.f90](#)

### 4. Radiation settings

The solar constant, the zenith angle and the surface albedo need to be adapted. In the source code, the solar constant needs to be factorized into the individual bands for our newly defined testcases (in `mo_nwp_phy_init.f90`). Besides, the albedo needs to be defined as a namelist parameter (in `mo_nh_testcases_nml.f90`), and needs to be transferred to an ICON variable for the two testcases, replacing the albedos calculated as a function of soiltype and zenith angle (in `mo_nwp_rad_interface.f90`).

In total, the following src files have been adapted:

`atm_phy_nwp/mo_nwp_phy_init.f90`

`namelists/mo_nh_testcases_nml.f90`

`atm_phy_nwp/mo_nwp_rad_interface.f90`

In the run script, the `nh_testcase_nml` namelist then can be set like this (following the RCEMIP protocol):

```
&nh_testcase_nml
  nh_test_name      = 'RCEMIP_analytical'      ! test case identifier
  ape_sst_case      = 'sst_const'
  w_perturb         = 0.1                      ! m/s, default=0.05
  th_perturb        = 0.3                      ! K, default=0.2
  zenithang         = 42.05 ! degrees
  sol_const         = 551.58
  albedo_set        = 0.07 ! fixed surface albedo; default is 0.07
/
```

In the `radiation_nml`, `izenith = 5` needs to be set to have no diurnal cycle and the aerosol/gas mixing

ratios need to be set according to the RCEMIP protocol.

```
&radiation_nml
  irad_o3      = 4 ! read in external ozone file
  izenith      = 5 ! 5: prescribed zenith angle (no diurnal cycle); 6: equinox
diurnal cycle
  irad_aero    = 0 ! switch off aerosols
  irad_cfc11   = 0
  irad_cfc12   = 0
  irad_co2     = 2
  irad_ch4     = 2
  irad_n2o     = 2
  vmr_co2      = 348.e-06 ! RCEMIP values; modern-day values
  vmr_ch4      = 1650.e-09
  vmr_n2o      = 306.e-09
/
```

## 5. Further changes to the run script

To have a vertical grid similar to the RCEMIP protocol, `num_lev = 75` needs to be set in `run_nml`, and the `sleve_nml` needs to be adapted:

```
&sleve_nml
  top_height = 48000.0 ! these settings yield an actual model top of about
34 km
  min_lay_thckn = 75. ! thickness of lowest model layer
  max_lay_thckn  = 500. ! max thickness
  htop_thcknlimit = 40000.0 ! dont exceed max_lay_thckn below this height
  stretch_fac   = 1.2 ! values > 1 increase dz near model top
/
```

In the `dynamics_nml` namelist, `lcoriolis = .FALSE.` needs to be set.

In the `les_nml` namelist, `isrfc_type = 5` (fixed SST) and, e.g., `sst = 300` need to be set. In this namelist, a threshold value for the minimum surface wind speed can also be set, which is asked for by the RCEMIP protocol (`min_sfc_wind = 1.0`).

[A complete run script as it has been used for one of the RCEMIP simulations \(ICON LEM, small domain, 300K\) can be downloaded](#)

here

## 6. Further changes for ICON NWP

To switch from ICON LEM to ICON NWP, three modifications to the run script are necessary: In `nwp_phy_nml`, the Raschendorfer turbulence parameterization (`inwp_turb = 1`) and the diagnostic cloud cover parameterization by Martin Koehler (`inwp_cldcover = 1`) should be used, and in `nh_testcase_nml`, the fixed value for the SST has to be set as `ape_sst_val`. Note that `ape_sst_val` has

to be set in degrees Celsius!

WARNING: ape\_sst\_val is only used by the model if the new testcases are included in  
atm\_phy\_nwp/mo\_nwp\_phy\_init.f90:

```
IF (ltestcase .AND. (nh_test_name == 'APE_nwp' .OR. nh_test_name ==  
'dcmip_tc_52' .OR. nh_test_name == 'RCEMIP_analytical' .OR. nh_test_name ==  
'RCEMIP_ncdf')) THEN !Tobias Becker added new testcases for RCEMIP  
  ! t_g = ape_sst1  
  DO jc = i_startidx, i_endidx  
    zlat = p_patch%cells%center(jc,jb)%lat  
    p_prog_lnd_now%t_g(jc,jb) = ape_sst(ape_sst_case,zlat) ! set SST  
    p_prog_lnd_new%t_g(jc,jb) = ape_sst(ape_sst_case,zlat)  
    p_prog_lnd_now%t_g_t(jc,jb,1) = ape_sst(ape_sst_case,zlat)  
    p_prog_lnd_new%t_g_t(jc,jb,1) = ape_sst(ape_sst_case,zlat)  
    ! Humidity at water surface = humidity at saturation  
    p_diag_lnd%qv_s(jc,jb) = &  
    &  
    spec_humi(sat_pres_water(p_prog_lnd_now%t_g(jc,jb)),p_diag%pres_sfc(jc,jb))  
    p_diag_lnd%qv_s_t(jc,jb,1) = p_diag_lnd%qv_s(jc,jb)  
  END DO
```

## 7. Postprocessing - Remapping

The longitudes and latitudes on the Torus grid must not be interpreted. Instead we have to remap the data onto a regular grid ranging from 180°W to 180°E and 10°S to 10°N. Remapped data can then be displayed in the index space, not in the geographical domain. The grid description of the regular grid is stored in grid\_icon\_torus\_nx20ny22\_5km:

```
gridtype = lonlat  
xsize    = 20  
ysize    = 20  
xfirst   = -180.0 ! start longitude (fixed)  
xinc     = 18     ! 360 / xsize  
yfirst   = -10.0  ! start latitude (fixed)  
yinc     = 1      ! 20 / ysize
```

Generate a weight file once:

```
cdo gendis,grid_icon_torus_nx20ny22_5km <infile>  
remapweights_nx20ny22_5km.nc
```

Remap the data using the weight file:

```
cdo remap,grid_icon_torus_nx20ny22_5km,remapweights_nx20ny22_5km.nc  
<infile> <outfile>
```

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[https://wiki.mpimet.mpg.de/doku.php?id=models:pot-pourri:spec\\_configs:rce\\_icon\\_lem\\_torus](https://wiki.mpimet.mpg.de/doku.php?id=models:pot-pourri:spec_configs:rce_icon_lem_torus)

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