

# GEOS-Chem Nested Model Tutorial

## Nested Model Co-Scientist

Yuxuan Wang

Univ. of Houston



Lin Zhang

Peking Univ.

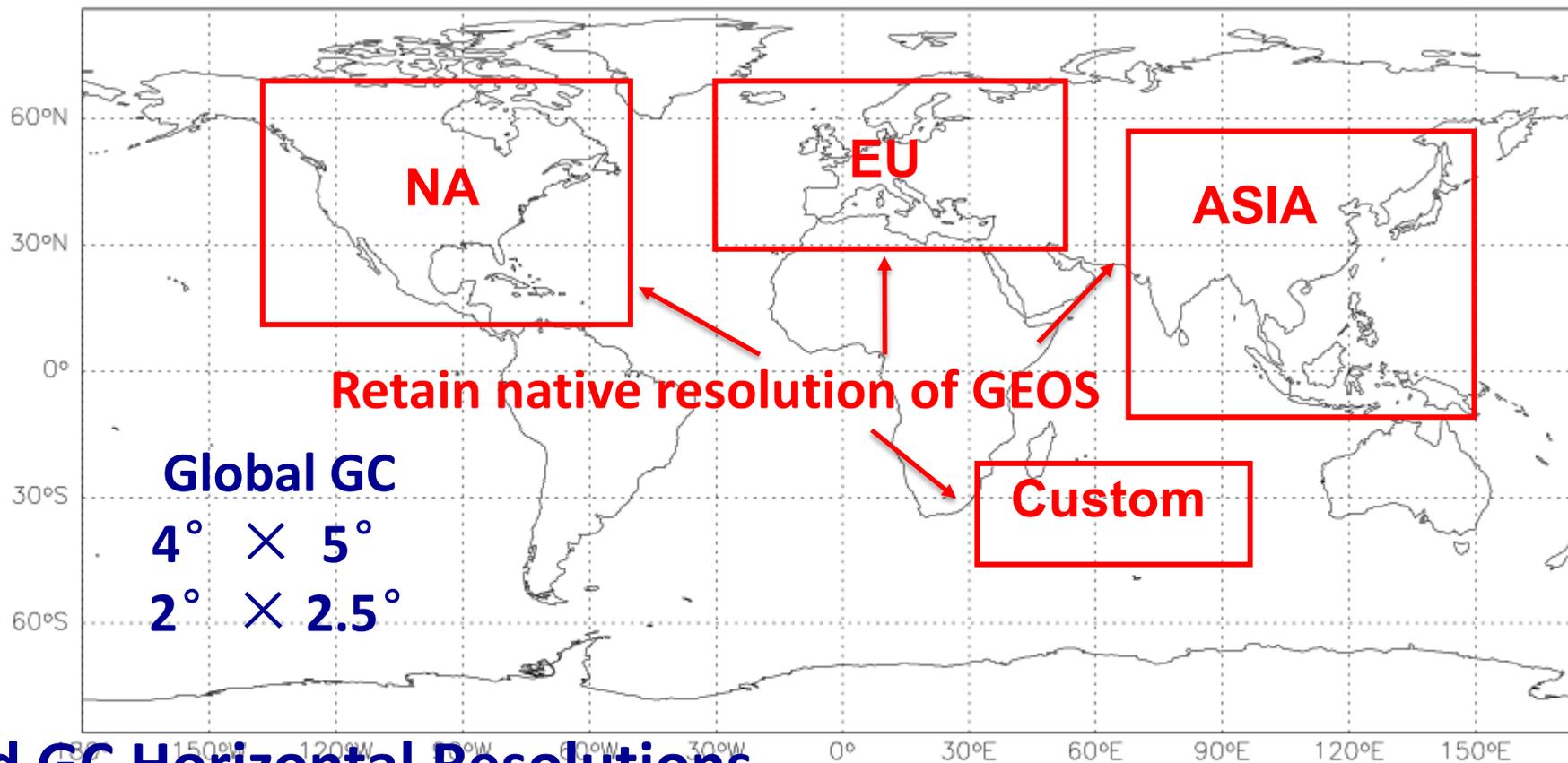


With contributions by: Tabitha Lee

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IGC10 @ WashU

# What is Nested GEOS-Chem?

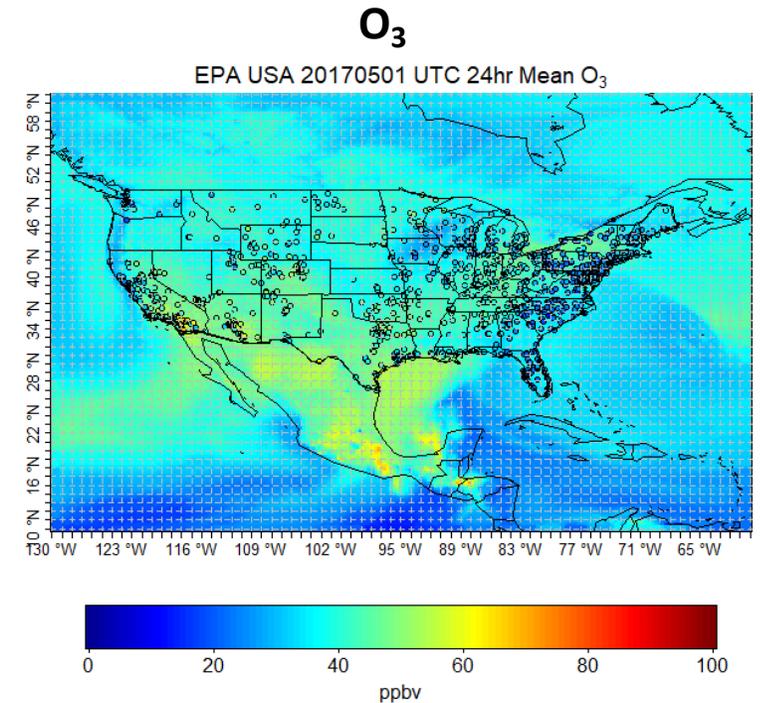
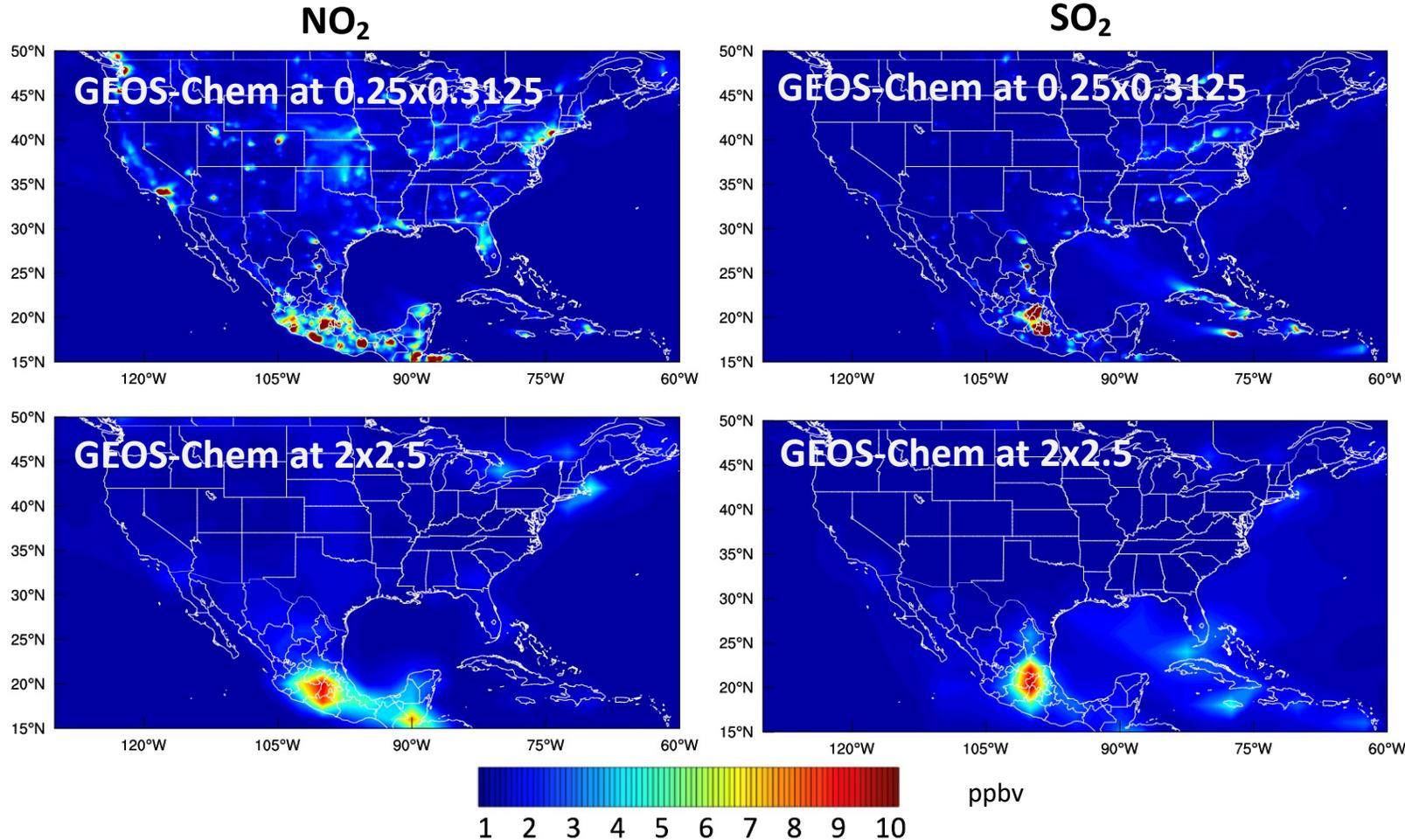


## Nested-grid GC Horizontal Resolutions

- ✓ 1° × 1° (GEOS-3) (Wang et al., 2004) (obsolete)
- ✓ 1/2° × 2/3° (GEOS-5) (Chen et al., 2009) (obsolete)
- ✓ 1/4° × 5/16° (GEOS-FP)
- ✓ 1/2° × 5/8° (MERRA-2)

# Nested-grid vs. coarse grid

## GEOS-Chem surface $\text{NO}_2$ , $\text{SO}_2$ and $\text{O}_3$ from May 2017 over the US



**Nested-grid provides better details of surface sources**

Plots generated by Lin Nan and Elizabeth Klovenski

# How to run GEOS-Chem nested model?

Tutorial based on GC version 13.4.0

**Step 1: Download the native-resolution met field**

**Step 2: Run the global simulation to generate boundary conditions**

*Either  $2^\circ \times 2.5^\circ$  or  $4^\circ \times 5^\circ$  boundary conditions*

**Step 3: Run the nested model (with boundary conditions from Step2 )**

# 1. Decide which met field to use

- Two options: GEOS-FP and MERRA2
- Check the horizontal grid first before you prepare the met field!
- [http://wiki.seas.harvard.edu/geos-chem/index.php/GEOS-Chem\\_horizontal\\_grids](http://wiki.seas.harvard.edu/geos-chem/index.php/GEOS-Chem_horizontal_grids)

Region	Met field	Resolution	Lon-lat coverage
US	MERRA2	0.5x0.625	140-40W, 10-70N
	GEOS-FP	0.25x0.3125	130-60W, 9.75-60N
Europe	MERRA2	0.5x0.625	30W-50E, 30-70N
	GEOS-FP	0.25x0.3125	15W-40E, 32.75-61.25N
Asia	MERRA2	0.5x0.625	60-150E, 11S-55N
China	GEOS-FP	0.25x0.3125	70-140E, 15-55N
<i>Custom (FlexGrid)</i>	<i>User defined</i>	<i>User defined</i>	<i>User defined</i>

## 2. Run global simulation

Everything is the same except you need to modify the **Boundary Conditions** menu in **HISTORY.rc**

```
COLLECTIONS: 'Restart',
             'BoundaryConditions',
             ...

#-----
# %%%% THE BoundaryConditions COLLECTION %%%%
#
# GEOS-Chem boundary conditions for use in nested grid simulations
#
# Available for all simulations
#-----
BoundaryConditions.template: '%y4%m2%d2_%h2%n2z.nc4',
BoundaryConditions.format:   'CFIO',
BoundaryConditions.frequency: 00000000 030000
BoundaryConditions.duration: 00000001 000000
BoundaryConditions.mode:     'instantaneous'
→ BoundaryConditions.LON_RANGE: -130.0 -60.0,
→ BoundaryConditions.LAT_RANGE: 10.0 60.0,
BoundaryConditions.fields:   'SpeciesBC_?ADV?           ', 'GIGCchem',
::
```

**You need to add and set the BoundaryConditions.LON\_RANGE and BoundaryConditions.LAT\_RANGE region in HISTORY.rc ONLY IF you are running the simulation over a nested domain or custom domain.**

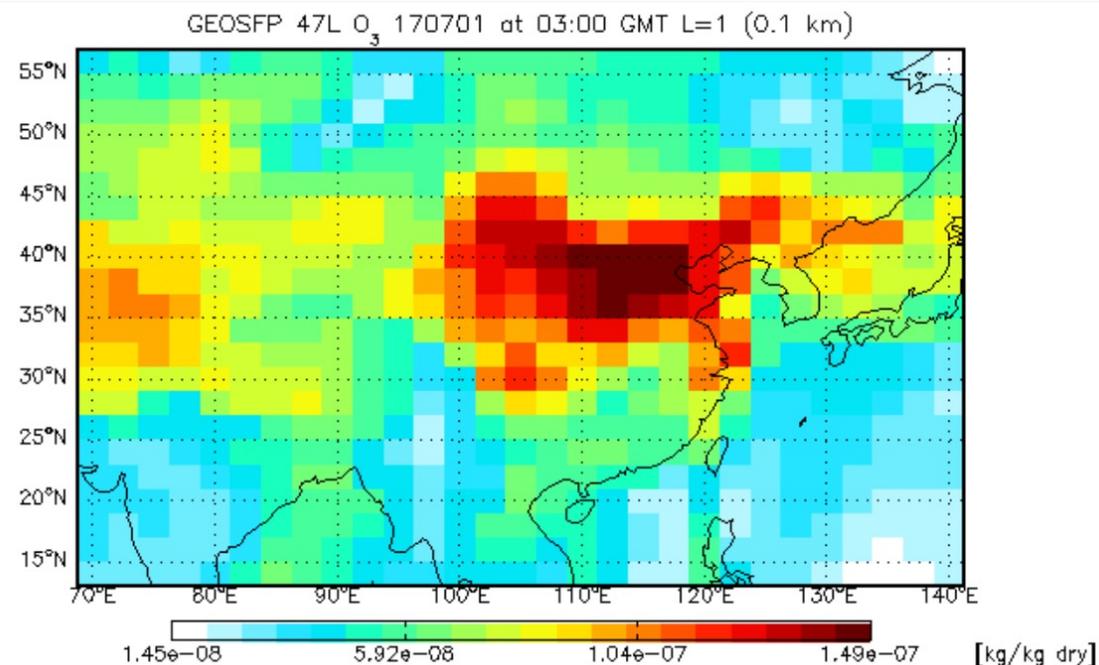
The first number in the LON\_ and LAT\_RANGE is the minimum value and the second number is the maximum value.

## 2. Run global simulation

**BC will then be saved for daily output (3-hour frequency). Keep an eye on your storage!**

```
-rw----- 1 luxiao luxiao 133M Mar 30 08:17 BC.20171208
-rw----- 1 luxiao luxiao 133M Mar 30 09:27 BC.20171209
-rw----- 1 luxiao luxiao 133M Mar 30 10:39 BC.20171210
-rw----- 1 luxiao luxiao 133M Mar 30 11:51 BC.20171211
-rw----- 1 luxiao luxiao 133M Mar 30 13:04 BC.20171212
-rw----- 1 luxiao luxiao 133M Mar 30 14:18 BC.20171213
-rw----- 1 luxiao luxiao 133M Mar 30 15:31 BC.20171214
-rw----- 1 luxiao luxiao 133M Mar 30 16:44 BC.20171215
-rw----- 1 luxiao luxiao 133M Mar 30 17:55 BC.20171216
-rw----- 1 luxiao luxiao 133M Mar 30 19:08 BC.20171217
-rw----- 1 luxiao luxiao 133M Mar 30 20:22 BC.20171218
-rw----- 1 luxiao luxiao 133M Mar 30 21:36 BC.20171219
-rw----- 1 luxiao luxiao 133M Mar 30 22:48 BC.20171220
-rw----- 1 luxiao luxiao 133M Mar 31 00:02 BC.20171221
-rw----- 1 luxiao luxiao 133M Mar 31 01:14 BC.20171222
-rw----- 1 luxiao luxiao 133M Mar 31 02:27 BC.20171223
-rw----- 1 luxiao luxiao 133M Mar 31 03:40 BC.20171224
-rw----- 1 luxiao luxiao 133M Mar 31 04:54 BC.20171225
-rw----- 1 luxiao luxiao 133M Mar 31 06:07 BC.20171226
-rw----- 1 luxiao luxiao 133M Mar 31 07:21 BC.20171227
-rw----- 1 luxiao luxiao 133M Mar 31 08:35 BC.20171228
-rw----- 1 luxiao luxiao 133M Mar 31 09:48 BC.20171229
-rw----- 1 luxiao luxiao 133M Mar 31 11:01 BC.20171230
-rw----- 1 luxiao luxiao 117M Mar 31 12:06 BC.20171231
```

**You can also use IDL -> GAMAP to check the BC file.**



### 3. Create run directory and run the nested model

Choose horizontal resolution:

1. 4.0 x 5.0
2. 2.0 x 2.5
3. 0.5 x 0.625

Choose horizontal grid domain:

1. Global
2. Asia
3. Europe
4. North America
5. Custom

When creating a run directory from the GCClassic superproject folder, selections are the same as the Global simulation, but now you select **3** in **Choose horizontal resolution** for **Nested** or **Custom** simulations.

The **Nested-Grid Menu populates** where you can **Choose horizontal grid domain** matching your decided simulation domain (AS, CH, EU, NA, or Custom).

### 3. Create run directory and run the nested model

#### □ Make sure input.geos has correct Latitude and Longitude Bounds

```
%%% GRID MENU %%%           :
Grid resolution             : 0.25x0.3125
→ Longitude min/max        : -130.0 -60.0
→ Latitude min/max         : 9.75 60.0
Half-sized polar boxes?: F
Number of levels           : 47
Nested grid simulation?    : T
Buffer zone (N S E W )    : 3 3 3 3
```

**For existing domains:** Longitude min/max and Latitude min/max **are already set** when you create the Nested simulation run directory.

**For custom domains (FlexGrid):** Specify custom Latitude and Longitude ranges

# 3. Create run directory and run the nested model

## □ Edit HEMCO\_Config.rc and HEMCO\_Config.rc.gmao\_metfields

```
#####  
### BEGIN SECTION SETTINGS  
#####  
  
ROOT:                /n/holyufs/EXTERNAL_REPOS/GEOS-CHEM/gcgrid/data/ExtData/HEMCO  
METDIR:              /n/holyufs/EXTERNAL_REPOS/GEOS-CHEM/gcgrid/data/ExtData/GEOS_4x5/GEOS_FP  
...  
  
#####  
### NON-EMISSIONS DATA (subsection of BASE EMISSIONS SECTION)  
###  
### Non-emissions data. The following fields are read through HEMCO but do  
### not contain emissions data. The extension number is set to wildcard  
### character denoting that these fields will not be considered for emission  
### calculation. A given entry is only read if the assigned species name is  
### an HEMCO species.  
#####
```

May need to add abbreviation (AS, CH, EU, NA) to match nested domain selection

```
# --- CN fields ---  
* FRLAKE    $METDIR/$CNYR/01/$MET.$CNYR0101.CN.$RES.$NC    FRLAKE    */1/1/0    C xy 1 * - 1 1  
* FRLAND    $METDIR/$CNYR/01/$MET.$CNYR0101.CN.$RES.$NC    FRLAND    */1/1/0    C xy 1 * - 1 1  
* FRLANDIC  $METDIR/$CNYR/01/$MET.$CNYR0101.CN.$RES.$NC    FRLANDIC  */1/1/0    C xy 1 * - 1 1  
* PROCEAN  $METDIR/$CNYR/01/$MET.$CNYR0101.CN.$RES.$NC    PROCEAN   */1/1/0    C xy 1 * - 1 1  
* PHIS      $METDIR/$CNYR/01/$MET.$CNYR0101.CN.$RES.$NC    PHIS      */1/1/0    C xy 1 * - 1 1  
  
# --- A1 fields ---  
* ALBEDO    $METDIR/$YYYY/$MM/$MET.$YYYY$MM$DD.A1.$RES.$NC    ALBEDO    1980-2018/1-12/1-31/0-23/+30minute CY xy 1 * -  
* CLDTOT    $METDIR/$YYYY/$MM/$MET.$YYYY$MM$DD.A1.$RES.$NC    CLDTOT    1980-2018/1-12/1-31/0-23/+30minute CY xy 1 * -  
* EFLUX     $METDIR/$YYYY/$MM/$MET.$YYYY$MM$DD.A1  
* EVAP      $METDIR/$YYYY/$MM/$MET.$YYYY$MM$DD.A1  
...  
#####
```

May need to add abbreviation (AS, CH, EU, NA) to match nested domain selection

1: Specify **METDIR** in HEMCO\_Config.rc

2: Ensure that **meteorology file names** are correct in HEMCO\_Config.rc.gmao\_metfields

# 3. Create the run directory and run nested model

## □ Edit HEMCO\_Config.rc

```
# ExtNr ExtName          on/off Species
0      Base              : on      *
# ----- RESTART FIELDS -----
--> GC_RESTART           : true
--> GC_BCs              : true
--> HEMCO_RESTART        : true

...

#=====
# --- GEOS-Chem boundary condition file ---
#=====
(((GC_BCs
* BC_ $ROOT/SAMPLE_BCs/v2019-05/tropchem/GEOSChem.BoundaryConditions.$YYYY$MM$DD_$HH$MNz.nc4 SpeciesBC_?ADV? 1980-2019/1-12/1-31/0-23 RFY xyz 1 * - 1 1
)))GC_BCs
```

3: GC\_BC~~s~~ should be already set to **true** for nested simulations

4: **Specify path for boundary condition files** generated by your global simulation

# Nested Simulation Summary and References

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## Checklist

- 1 Step 1:** Have you decided and downloaded the met fields for your simulation?
  - 2 Step 2:** Have you turned on and correctly save the boundary conditions in your global simulation? (**HISTORY.rc in global simulation**)
  - 3 Step 3.1:** Have you created the nested model run directory from the GCCClassic superproject folder? (**~/GCCClassic.13+/run**)
  - 4 Step 3.2:** Have you set the Met and BC pathway correctly? (**HEMCO\_Config.rc in nested simulation**)
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[http://wiki.seas.harvard.edu/geos-chem/index.php/GEOS-Chem\\_Nested\\_Model](http://wiki.seas.harvard.edu/geos-chem/index.php/GEOS-Chem_Nested_Model)

[http://wiki.seas.harvard.edu/geos-chem/index.php/Setting\\_up\\_GEOS-Chem\\_nested\\_grid\\_simulations](http://wiki.seas.harvard.edu/geos-chem/index.php/Setting_up_GEOS-Chem_nested_grid_simulations)

<http://wiki.seas.harvard.edu/geos-chem/index.php/FlexGrid>