


# Exercises for the NorESM user workshop 2024

# Hands-on session 1

Download the CMIP6 version of NorESM2:

1. `$ ssh -Y username@betzy.sigma2.no`
2. `$ mkdir NorESMworkshop2024`
3. `$ cd NorESMworkshop2024`
4. `$ mkdir cases`
5. `$ git clone https://github.com/NorESMhub/NorESM.git`
6. `$ cd NorESM`
7. `$ git checkout release-noresm2.0.9`
8. `$ ./manageExternals/checkoutExternals`

The “-Y” option enables X11 forwarding, e.g. for viewing model output on Betzy.



# Hands-on session 2

Create, setup, build and submit your (first?) NorESM2-LM piControl simulation running on 4 nodes by repeating these steps:

```
$ cd ~/NorESMworkshop2024/NorESM/cime/scripts/
```

```
$ ./create_newcase --case ~/NorESMworkshop2024/cases/N1850frc2_f19_tn14_test01_20241119  
--mach betzy --res f19_tn14 --compset N1850frc2 --project nn9039k --pecount=S
```

```
$ cd ~/NorESMworkshop2024/cases/N1850frc2_f19_tn14_test01_20241119/
```

```
$ ./case.setup
```

```
$ ./case.build
```

```
$ ./case.submit
```

Output will be available on `$USERWORK/archive/N1850frc2_f19_tn14_test01_20241119`

# Hands-on session 2 : checkout externals

In release-noresm2.0.9 there is a second externals file:

Externals\_continuous\_development.cfg

This (now outdated) file was used to build NorESM with alternative development branches of model components. We can use this file to illustrate how to use “status checker”:

1. Look at differences between externals files:

```
diff -u Externals.cfg Externals_continuous_development.cfg
```

2. Examine status for current and alternative externals file

```
./manageExternals/checkoutExternals -S ← for current externals
```

```
./manageExternals/checkoutExternals -e Externals_continuous_development.cfg -S  
← for alternative externals
```

Symbols used by status checker:

: all is fine	s : out-of-sync	o : optional source
e : empty	? : unknown (no .git or .svn)	M : modified source

# Hands-on session 2 : query\_config

Use query\_config to learn more about NorESM model options

1. Go to scripts directory: `cd cime/scripts`
2. Find details about compsets N1850frc, NorESM coupled, and blom
  - `./query_config --compsets | grep N1850frc2`
  - `./query_config --compsets | awk '$1 ~ /^N/'` ← NorESM compsets
  - `./query_config --compsets blom` ← OMIP compsets
  - `./query_config --compsets cam | awk '$1 ~ /^N/'` ← AMIP compsets
3. Find config options for a component or data replacement
  - `./query_config --components blom` ← Active ocean component
  - `./query_config --components docn` ← Data ocean component
4. Find details about grids
  - `./query_config --grids | grep -A 2 tn14` ← All grids using tn1v4 ocean grid

# Hands-on session 3 -

## A) piControl (N1850)

- create a new case with resolution f19\_tn14
- Set --pecount=S
- make it a branch run
- reference case: N1850\_f19\_tn14\_11062019
- reference date: 1600-01-01
- 1 month simulation time

## B) historical (NHIST)

- create a new case with resolution f19\_tn14
- Set --pecount=S
- make it a hybrid run
- reference case: N1850\_f19\_tn14\_11062019
- reference date: 1600-01-01
- Start date: 1850-01-01
- 1 month + 1 month simulation time i.e.  
run for 1 month, resubmit simulation once and  
run for 1 more month  
(set RESUBMIT=1)
- Dump restart files every 10 days

For A,B we need to change only env\_run.xml

[https://noresm-docs.readthedocs.io/en/noresm2/configurations/cmip6\\_compsets.html#reproduce-cmip6-picontrol-historical-and-ssp5-8-5-experiments](https://noresm-docs.readthedocs.io/en/noresm2/configurations/cmip6_compsets.html#reproduce-cmip6-picontrol-historical-and-ssp5-8-5-experiments)

- Check how xmlquery and xmlchange work together with the xml files:
  - Assuming you want to change the total run time for NorESM, but do not remember the exact keyword, find all keywords that include STOP or CLOCK:  
`./xmlquery --partial STOP`  
`./xmlquery --partial CLOCK`
  - STOP\_N is defined in env\_run.xml, view file with “less” or grep STOP\_N in env\_run.xml
  - Change STOP\_N value  
`./xmlchange STOP_N=10`
  - Confirm change has been made in env\_run.xml

- Create a clone case from (A), and add daily output for surface DMS flux from the ocean component
  - See “create\_clone” from “setting up a clone case”
  - Modify “user\_nl\_blom”: daily output for surface DMS flux  
     SRF\_DMSFLUX=4, 2, 2  
     (format for output is ‘daily, monthly, yearly’, default is ‘0, 2, 2’)
- Run case (A) on normal queue for 1 month with `--pecount=M` to run on 8 nodes. Compare model throughput in the log file with the previous run.
- Run NorESM with data component for either atmosphere or ocean for 1 month. Compare model throughput with fully coupled run.
- Run on development queue: Set `--pecount=128` when doing `create_newcase` ; In case:
  - `./xmlchange NTASKS_OCN=123`
  - `./xmlchange JOB_QUEUE="devel" --subgroup case.run --force`
  - `./xmlchange JOB_WALLCLOCK_TIME=0:30:00 --subgroup case.run`
  - `./xmlchange JOB_WALLCLOCK_TIME=0:30:00 --subgroup case.st_archive`



## Hands-on session on compsets

1. Create an ocean-sea ice only case (e.g. compset = NOICPLHISTOC)
2. Look at the README.case file in your case directory. How can you tell that only the ocean and sea ice components are active?
3. What can you tell about the atmosphere and land components? Are those active, data-components, stubb?
4. Try the same with `./xmlquery`. What additional information do you get?

1. `./xmlquery -p BLOM`
2. `./xmlquery -p CICE`
3. `./xmlquery -p CAM`
4. `./xmlquery -p CLM`

Repeat the steps with atmosphere-land only (e.g. compset = NF1850) and fully coupled (e.g. compset = N1850).

# Hands-on session make a case using bash script - part 1

There are two sample scripts on Betzy:

/cluster/shared/noesm/WORKSHOP/scripts/ReproExperimentScriptSimple.sh

/cluster/shared/noesm/WORKSHOP/scripts/ReproExperimentScript.sh

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1. make a copy and save it somewhere (e.g. in your NorESMWorkshop2024 directory)
2. make changes to suit your needs. In particular, review these four lines and at least replace xxUSERxx with your Betzy user ID.

```
COMPSET="N2000"
```

```
RES="f19_tn14"
```

```
SRCROOT="/cluster/projects/nn9039k/xxUSERxx/NorESM"
```

```
CASEDIR="/cluster/work/users/xxUSERxx/cases/${COMPSET}_${RES}"
```

# Hands-on session make a case using bash script - part 2

There are two sample scripts on Betzy:

`/cluster/shared/noresm/WORKSHOP/scripts/ReproExperimentScriptSimple.sh`

`/cluster/shared/noresm/WORKSHOP/scripts/ReproExperimentScript.sh`

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1. Make a copy and save it somewhere (e.g. in your NorESMWorkshop2024 directory)
2. Make a new script where you specify your run environment. E.g. Hands-on session 3B on slide 6