

Yale



耶鲁大学-南京信息工程大学大气环境中心

Yale-NUIST Center on Atmospheric Environment

Testing of WRF-GHG on simulation of atmospheric methane in Yangtze River Delta region

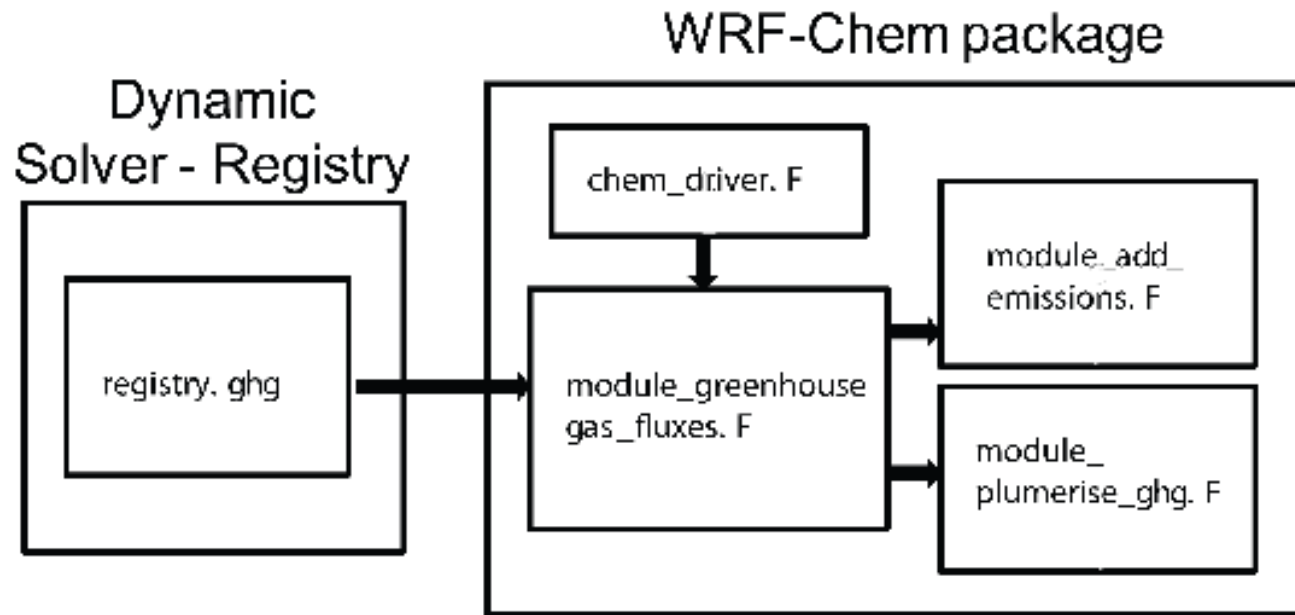
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2013-08-09

Outline

- Brief description of WRF-GHG
- A case study
- Next steps

Coupling GHG with WRF/Chem



Four additional modules are added to WRF-Chem and some minor modifications required by WRF-Chem itself.

Registry.ghg

CH4_1	total atmospheric CH ₄ concentration
CH4_2	changes in CH ₄ concentration from wetland emissions
CH4_3	changes in CH ₄ concentration from anthropogenic emissions
CH4_4	changes in CH ₄ concentration from biomass burning
CH4_5	changes in CH ₄ concentration from termite emissions
CH4_6	changes in CH ₄ concentration from soil uptake
CH4_7	changes in CH ₄ concentration from vegetation
CH4_B	atmospheric CH ₄ background concentration

CH₄ variables defined in
WRF/GHG

CH ₄		
fl_wet	ext. wetland	hourly [30]
fl_ant4	ext. anthropogenic	daily [1]
fl_antch4	ext. anthropogenic	hourly [30]
fl_bbch4	ext. biomass burning	daily [1]
bb_ch4	3d int. biomass burning	WRF timestep
fl_term	ext. termite emission	daily [1]
fl_soilu	ext. soil uptake	hourly [30]
fl_veg	ext. vegetation	hourly [30]
ch4_emiss	int. wetland	WRF timestep
ch4_term	int. termite emission	WRF timestep
ch4_soil	int. soil uptake	WRF timestep
ch4_veg	int. vegetation	WRF timestep

CH₄ flux variables
defined in WRF/GHG

Linkage of emission with CH₄ concentration

- `co2_surface_source_add`
- `co2_surface_source_ad2`
- `add_emis_anthro_ghg`
- `add_emis_burn_ghg`

Follow the same
principle

$$chem(i, 1, j) = chem_source(i, j) \left[\frac{kg}{m^2 s} \right] \times conv_rho(i, 1, j) \left[\frac{sm^2}{kg} \right]$$

$$conv_rho(i, 1, j) = \frac{1}{rho(i, 1, j) \left[\frac{kg}{m^3} \right]} \times \frac{dt[s]}{dz8(i, 1, j)[m]}.$$

Here $rho(i, 1, j)$ denotes the air density in the first model layer, dt the time step of the model and $dz8(i, 1, j)$ the thickness of the first model layer in [m].

Calculation of CH₄ fluxes

- Online calculation of CH₄ wetland flux
- Flux from termite
- Flux from soil uptake
- Flux from vegetation
- Flux from biomass burning

Kaplan online calculation of CH₄ flux

● Inputs

- C_pool (external carbon pool, LPJ model [sitch et al.,2003])
- Wetland_map (Kaplan potential wetland map, wetland fraction/grid)
- Soil parameters (T_soil, Soil_type; Soil_M, etc.)

● Outputs

- CH₄_wet (CH₄ wetland emission)

$$k_r = \frac{\frac{1}{\tau_0} \cdot g(T) \cdot f_{SM}}{12 \cdot 24 \cdot 30}.$$

$$f_{SM} = 0.25 + 0.75 \frac{sm}{sm_{sat}}$$

$$g(T) = \exp \left(308.56 \cdot \left(\frac{1}{56.02} - \frac{1}{T + 46.01} \right) \right)$$

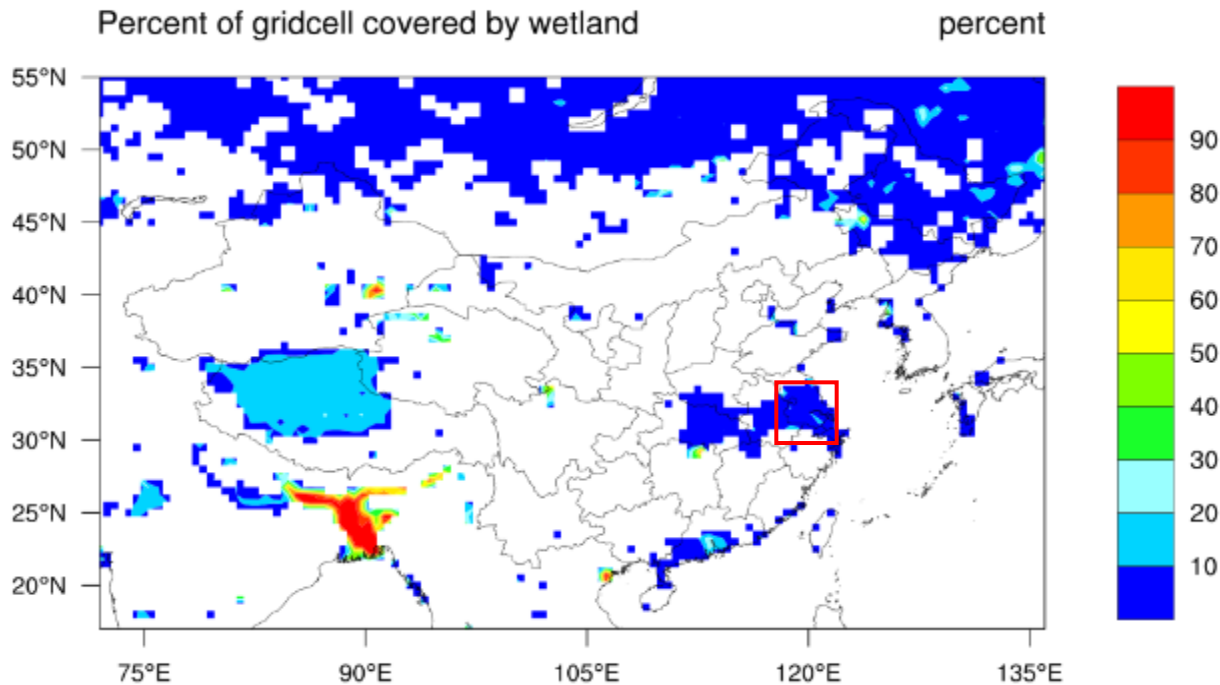
```

g_T(i,j) = EXP(308.56*(1.0/56.02 - 1.0/(T_2s(i,j) + 46.02)))
f_SM(i,j) = 0.25 + 0.75 * (sm_tot(i,j)/soil_sat(i,j))
k_r(i,j) = ((1/tau_10)*g_T(i,j)*f_SM(i,j))/(12.0*24.0*30.0)
Carbon(i,j) = C_pool(i,j)*(1.0 - EXP(-k_r(i,j)*1.0))
Carbon(i,j) = Carbon(i,j)/(3600.0)
HR(i,j) = 0.7*Carbon(i,j)
T_flood(i,j) = HR(i,j)*M_s*Wet_map(i,j)
T_peat(i,j) = HR(i,j)*E_f*Wet_map(i,j)
IF(wet_id == 2) then
  CH4_wet(i,j) = T_flood(i,j)
ELSE IF(wet_id == 3) then
  CH4_wet(i,j) = T_peat(i,j)
ELSE IF(wet_id == 1) then
  P_1(i,j) = EXP((T_a(i,j) - 303.0)/8.0)

```

Kaplan online calculation of CH₄ flux (cont.)

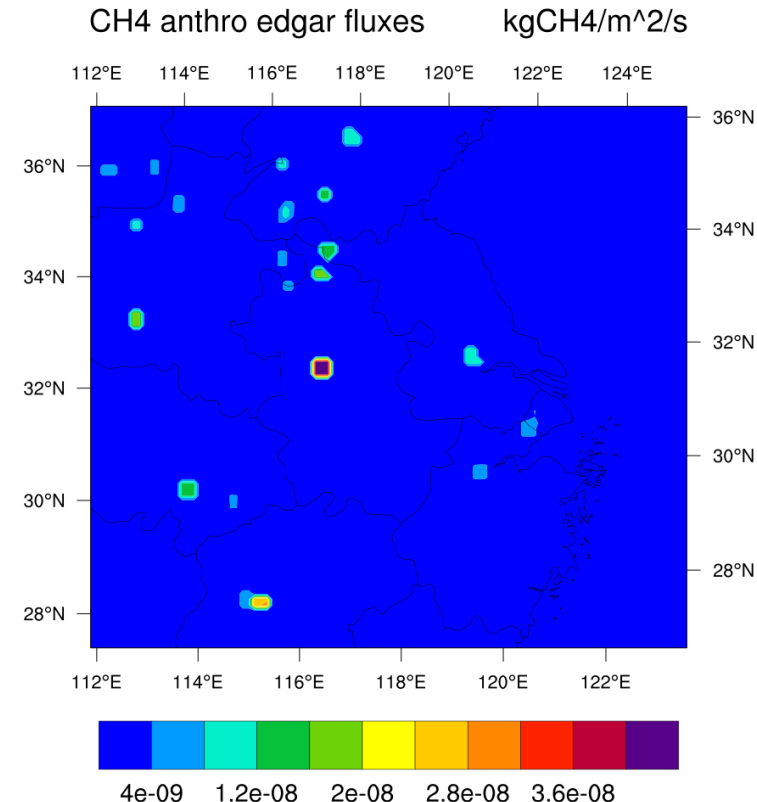
kaplan wetland map



From <http://arve.epfl.ch/pub/wetlands/index.html>

Anthropogenic emissions of CH₄

- Emission inventory: EDGAR V4 (<http://edgar.jrc.ec.europa.eu>)
- Resolution: $0.1^\circ \times 0.1^\circ$
- Base year: 2005
- Processing code: Prep_chem_source (a diurnal cycle peaking twice a day at 08:00 and 20:00 local time using a double Gaussian function are included here)



Kaplan module implemented in GHG

Take “CH4_2” for example

In the **namelist.input**

chem_opt 98

wetland_type

- 0 no CH₄ wetland emissions|
- 1 floodplain and peatland CH₄ wetland emissions (Kaplan)
- 2 only floodplain CH₄ wetland emissions (Kaplan)
- 3 only peatland CH₄ wetland emissions (Kaplan)
- 4 external CH₄ wetland emissions (Walter)

In **chem_driver.F**

Module_greenhouse_gases

SELECT CASE(wet_id)

CASE(1)

call KAPLAN

call co2_surface_source_add(...)

.....

CASE(4)

chem_sourcewet(:, :)=fl_wet(:, :)

.....

WRITE(6,*) 'wetland walter was called'

```
IF ( config_flags%chem_opt==98 ) THEN
  call greenhouse_gases (  chem, num_3d_c, grid%alt, dz8w, grid%dt, grid%xtime,
                           ids, ide, jds, jde, kds, kde,
                           ims, ime, jms, jme, kms, kme,
                           its, ite, jts, jte, kts, kte,
                           config_flags%run hours, config_flags%co2 anthro.
                           &
                           &
                           &
                           &
                           &
```

Initial and boundary conditions

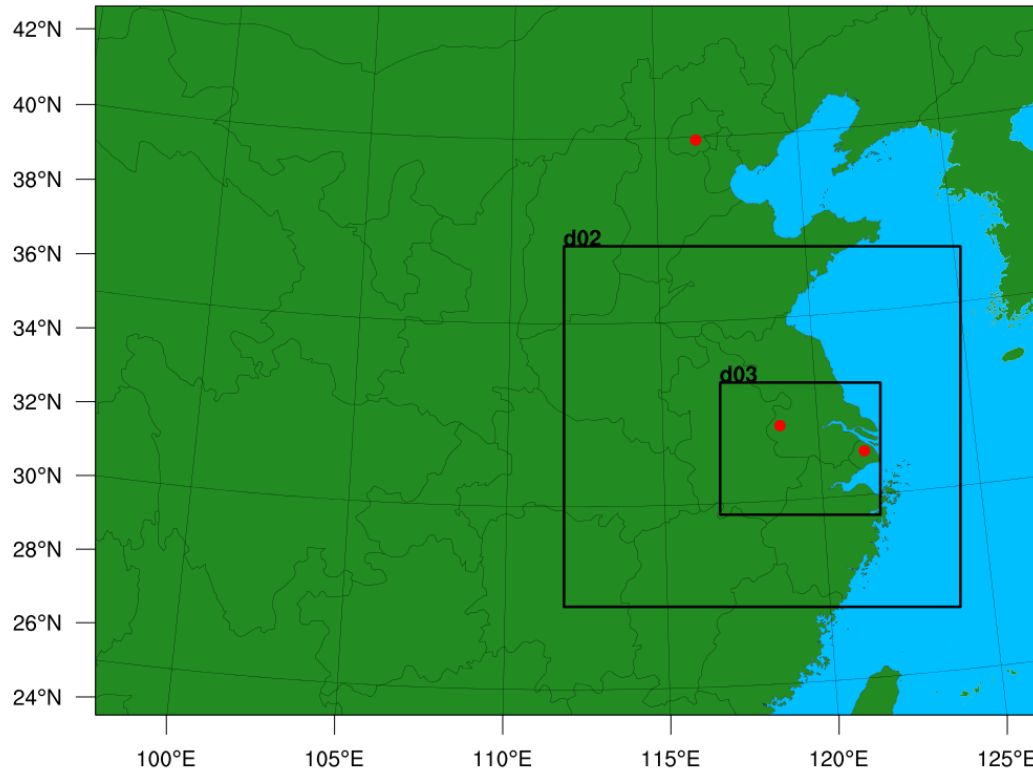
- **Meteorology:** NCEP FNL or ECMWF reanalysis data.
- **CH₄ and CO₂:** three-dimensional fields from global transport models TM5.
- The meteorology is initialized with ECMWF fields, while the tracer variables are initialized with outputs of global chemistry models like MOZART, TM5, or GOES-Chem on the first run and the simulations on the following runs are initialized with the previous day results to ensure continuity.

Steps to run WRF-GHG

- Compile source codes, WRF-GHG and WPS
- Run WPS to process all necessary fields on the predefined WRF grid
- Modify the namelist.input
- Run real.exe
- Use Matlab scripts process ICs/BCs of chemical species
- Run wrf.exe

Three-nested domains

WRF Domain YRD



d1:36km

d2:12km

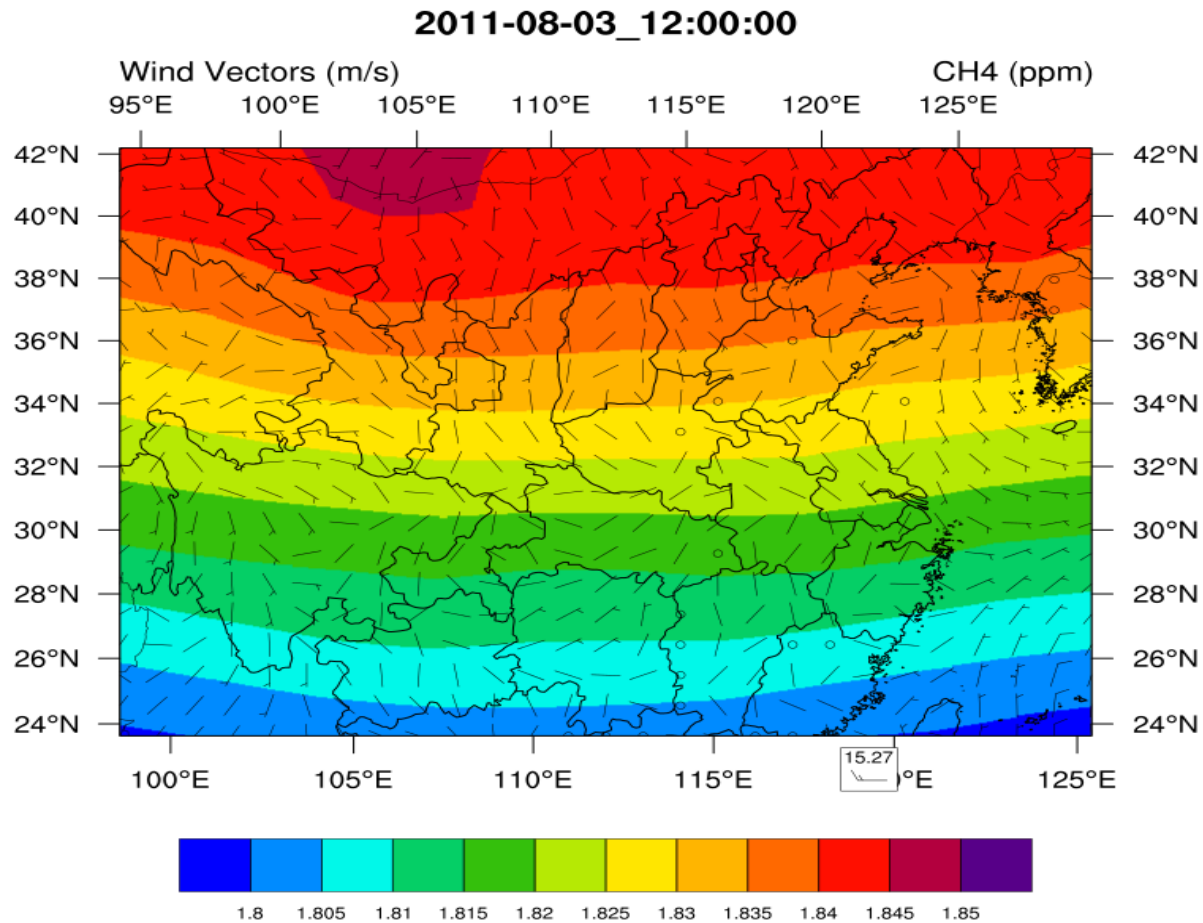
d3:4km

Vertical levels:40

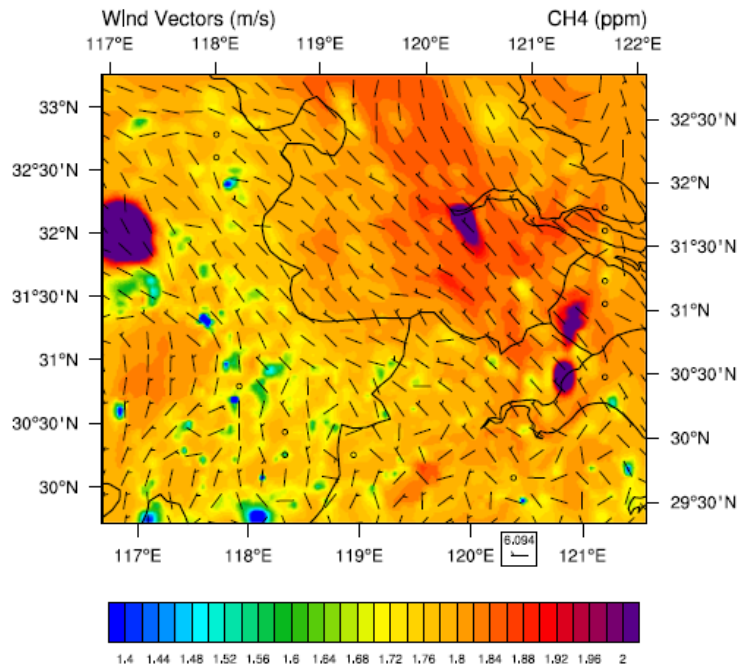
Namelist setting

vprm_class	= 8,	
vprm_par_file	= "VPRM_param_CERES2005_local.txt",	
co2_st_time	= 0,	
co2_anthro	= 2,	hourly anthropogenic emissions
wetland_type	= 0,	no CH4 wetland emissions
file_term	= "CH4_termite_NW.txt",	
term_id	= 0,	
bb_opt_ghg	= 0,	
plume_frq_ghg	= 180,	
soil_id	= 1,	Calculation of soil uptake fluxes
veg_id	= 1,	Calculation of CH4 emissions from vegetation
oce_id	= 0,	
chem_opt	98	
vertmix_onoff	1	to allow for vertical mixing of the tracer
have_bcs_chem	.true.	
chem_in_opt	0	
emiss_inpt_opt	0	
chem_conv_tr	1	to allow for subgrid convective tracer transport

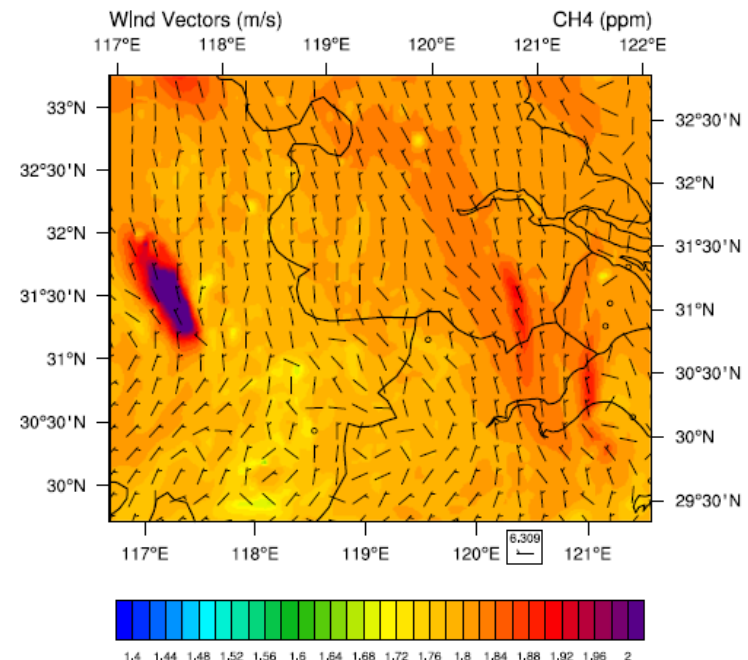
Case Study: Simulated CH₄



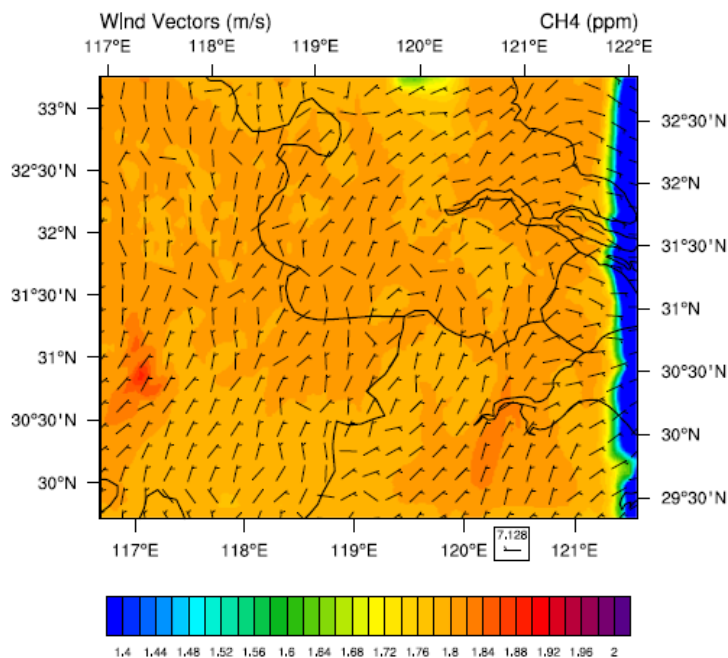
2011-08-03_18:00:00



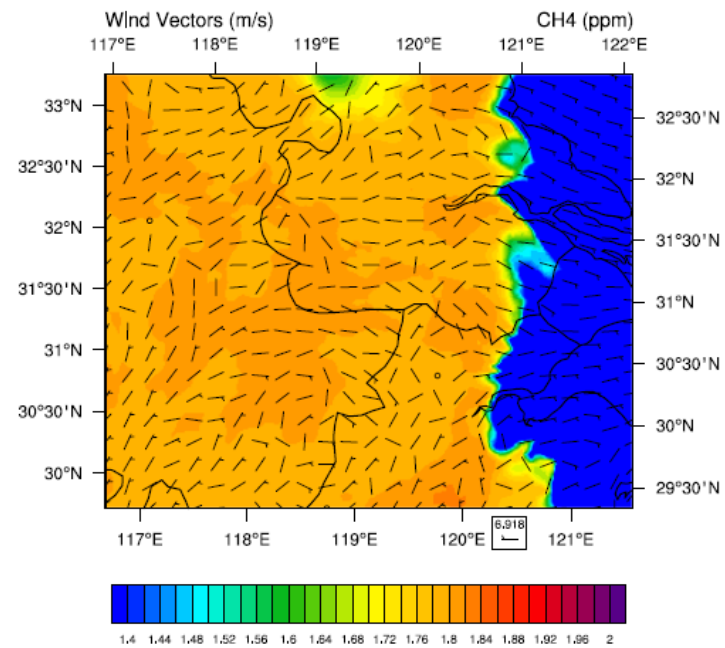
2011-08-04_00:00:00



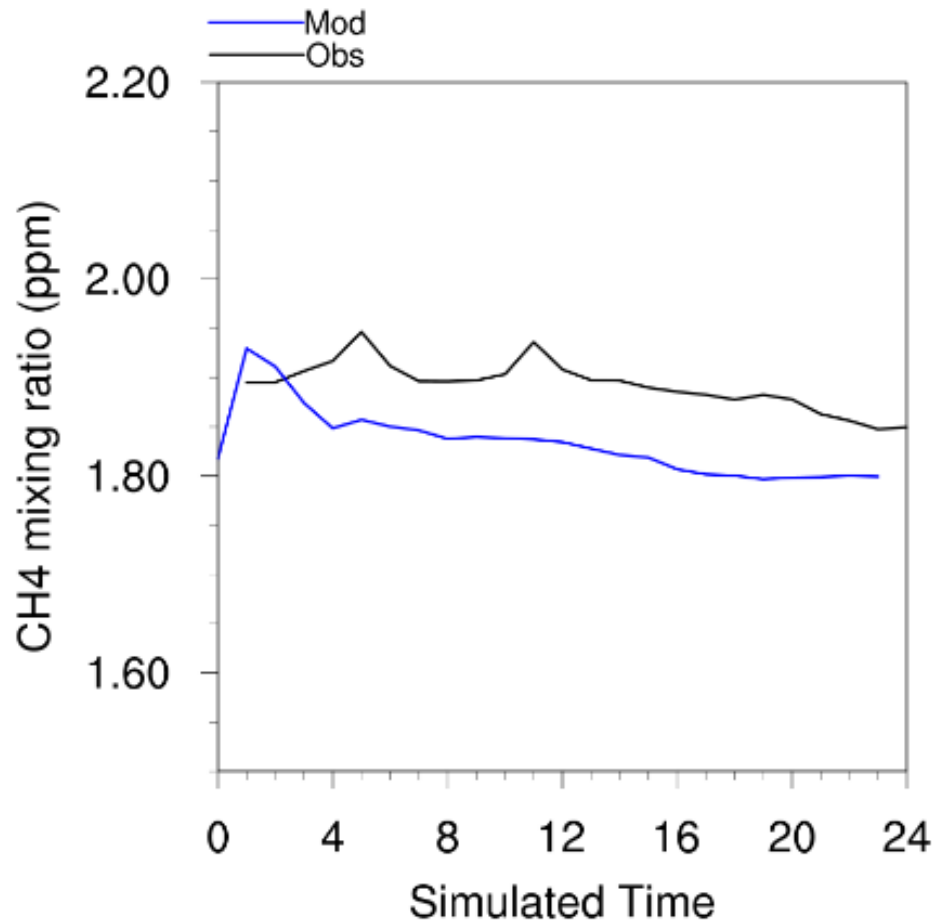
2011-08-04_06:00:00



2011-08-04_12:00:00



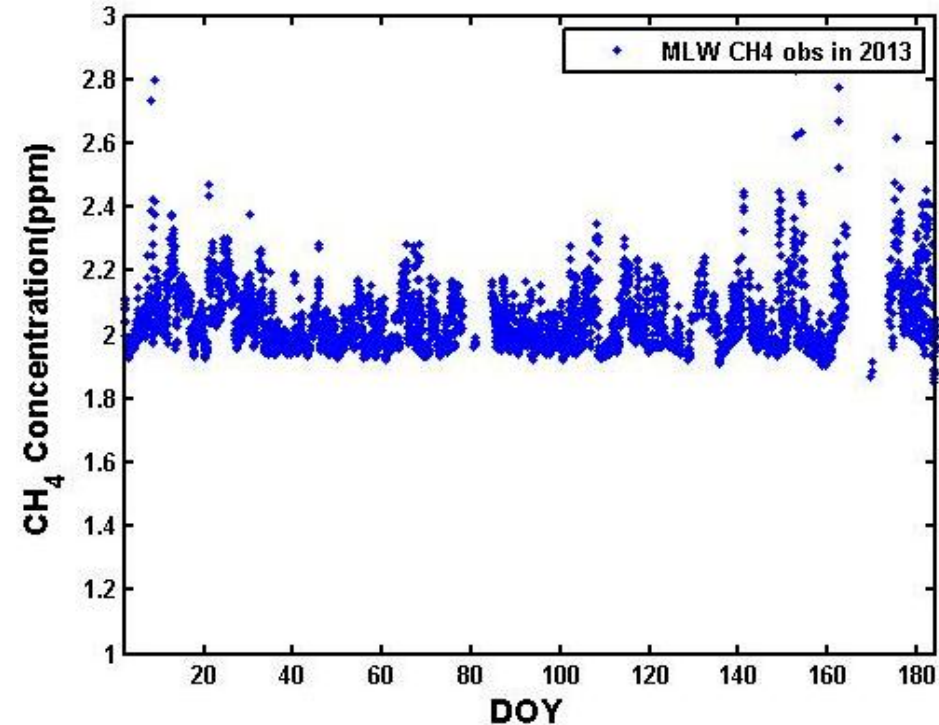
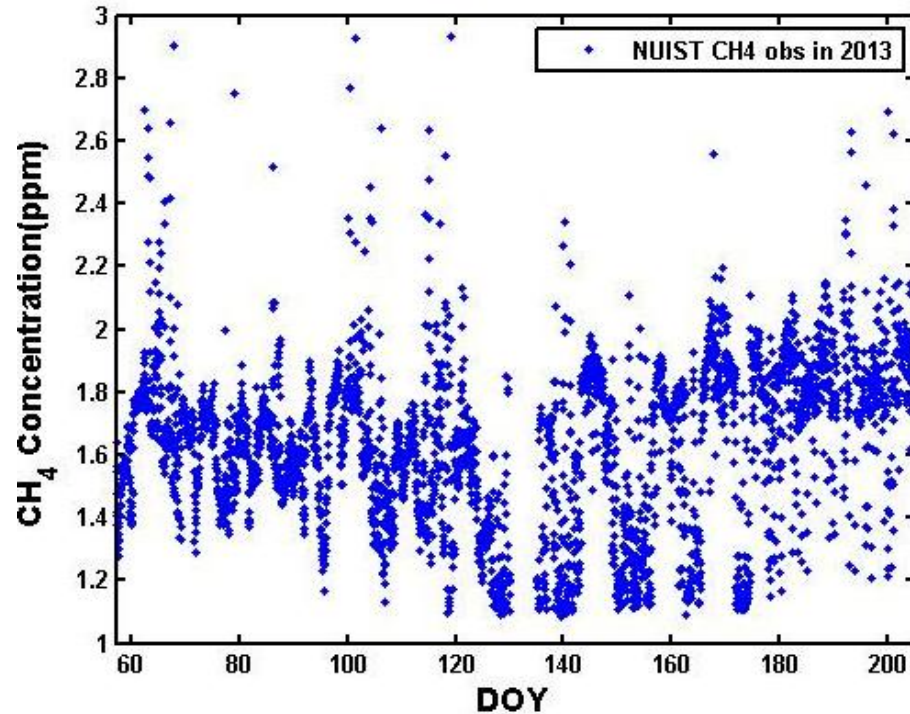
Comparison of simulated CH₄ with observed in MLW



Obs data is 2012.08.03-08.04

Mod data is 2011.08.03-08.04

CH₄ observed at NUIST and MLW



CH₄ concentration in NUIST is lower than that at MLW.

The diurnal pattern of CH₄ is not as obvious as CO₂.

Summary and future work

- Able to reproduce a case for WRF/GHG simulation on CH₄ using the available IC/BC and emission input files
- Will revisit all input parameters used in Kaplan to find a best scenario for application in Yangtze River Delta region (from a box model to a three dimensional model, WRF/GHG)
- Will learn how to generate emission, initial and boundary conditions of CH₄ and other chemical species for a new case.

Thank You!