Introduction to WRF-Chem

WMO GURME Regional Training Workshop on urban air quality modelling for ASEAN Countries

Malaysian Meteorological Department
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Introduction to WRF-Chem

Georg Grell

Steven E. Peckham, Stuart A. McKeen, Jan Kazil, R. Ahmadov + others from NOAA/ESRL

Jerome Fast, William Gustafson jr., P.L. Ma, B. Singh+ many others from PNNL

+ Alma Hodzic, Christine Wiedinmyer, Gabi Pfister, Mary Barth and many others from NCAR
  other University contributions
  + Saulo Freitas (CPTEC, BRAZIL)
  +new stuff from NCSU (Yang Zhang)

+ many more national and international collaborators

WRF-Chem web site - http://wrf-model.org/WG11
WRF-Chem

• Community effort

• Largest contributing groups: ESRL, PNNL, NCAR

• Other significant contributions from: National and international Universities, CPTEC Brazil, NASA, AFWA, NCSU
Structure of Talk

1. Brief description of only the *general features* of WRF-Chem
2. Some applications of what the model may be used for are mixed in

There are more than **50** chemistry options for the main gas phase chemistry and aerosol modules!
WRF-Chem

- Chemistry is online, completely embedded within WRF CI
- Consistent: all transport done by meteorological model
  - Same vertical and horizontal coordinates (no horizontal and vertical interpolation)
  - Same physics parameterization for subgrid scale transport
  - No interpolation in time
- Easy handling (Data management)
- Ideally suited:
  - to study feedbacks between chemistry and meteorology
  - for air quality forecasting on regional to cloud resolving scales
Why Online?

- Offline modeling introduces errors for air quality applications
  - Error for offline modeling is increasing with increasing horizontal resolution
  - Power spectrum analysis can show the amount of information that is lost in offline runs
- 2-way feedback in-between chemistry and meteorology
  - Process studies relevant for global climate change
  - Ultimately it should lead to improved data assimilation (meteorology) and improved weather forecasts

Grell and Baklanov, 2011, AE
What is needed for this type of modeling system?

1. Advection and diffusion (all done by WRF)
2. Sub-grid scale transport (WRF parameterizations, PBL, convection)
3. Some processes that are specific for chemical constituents, but need meteorology: emissions (biogenic, fire, sea salt, dust, volcanic, anthropogenic), dry deposition, wet scavenging
4. Treatment of chemical reactions, aqueous phase chemistry, gas phase species and aerosols
5. “Chemical” radiation routines (photolysis routines) that provide photolysis rates necessary for (4)
6. Capability of feedback from chemistry to meteorology (meteorological radiation and microphysics parameterizations, possibly also convective parameterizations)
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Gas Phase Chemistry Packages

Very complex part of the modeling system: Many additional species that are fully prognostic variables and require transport

- Hard coded: chemical mechanism from RADM2
- Hard coded: Carbon Bond (CBM-Z) based chemical mechanism
- *Kinetic PreProcessor* (KPP) – Many different equations files exist (also for RADM2 and CBM-Z). KPP will generate the modules from equation files. These generated modules will then be used by WRF-Chem
- IN V3.5.1/ V3.6: CRIMech gas phas scheme (U. of Manchester, 240 species, 652 reactions)
- V3.7: CB05
Photolysis Packages – all coupled to aerosols and hydrometeors

- Madronich Photolysis
- Madronich F-TUV
- Fast-j photolysis scheme
Available aerosol modules

(1) Modal

(2) Sectional

composition
sulfate
nitrate
ammonium
chloride
carbonate
sodium
calcium
other inorganics
organic carbon
elemental carbon

(3) Bulk: Sections for dust and sea salt, otherwise total mass only
Aerosols may have a significant impact on weather forecasts through interaction with radiation (sometimes also called “direct effect”) and microphysics (sometimes also called “indirect effect”)

Aerosols may also impact meteorological data assimilation
For NWP a bulk scheme is very attractive: GOCART (Currently used in real-time high resolution global (dx=30km) and regional modeling (up to dx=3km) at ESRL

- Much simpler than the sectional and modal schemes
  - Calculates only with the total mass of the aerosol components
  - Provides no information on
    - Particle size
    - Particle concentration
  - E.g., when particles grow, the aerosol mass increases but we don't know how their size/number changes

- Numerically very efficient
- Coupled with radiation (Mie scattering and extinction calculations)
- Will be coupled to microphysics in future versions
For research on aerosol direct and indirect effects modal and sectional approaches are more attractive

Less assumptions are made when coupled to atmospheric radiation and/or microphysics

Interaction processes are very complex, they will not work for every radiation and microphysics scheme in WRF! (takes time to implement)
Examples of available Aerosol Modules

- **Bulk**: GOCART
- **Modal**: MADE/SORGAM (3 modes)
- **MAM** (Modal Aerosol Model from NCAR Climate model)
- **MOSAIC** (Sectional)
Selection of radiation parameterizations for aerosol “direct effect”

Since V3.5 all aerosol modules were hooked up to Goddard short wave radiation, and RRTMG short and long wave scheme, CAM radiation.
Selection of microphysics parameterizations for aerosol "indirect effect"

Since V3.6

Modal and sectional schemes only can be used in combination with a version of the Lin et al. Microphysics scheme as well as the Morrison scheme

Special physics choice available when using NCAR community climate model physics

"indirect effect" is a result of the interaction aerosols/microphysics
How is the meteorological forecast affected by aerosol?

- Large importance for climate simulations is recognized (when integrating models over 100’s of years, small differences in the earth’s energy budget are extremely important)
- Weather forecasting for only a few days?
  - Much research needed, but chemistry may positively influence forecasts when strong signals exist
  - Influence on meteorological data assimilation
Observed (black) and predicted (blue) sounding for Fairbanks, Alaska, on July 4, 0000UTC.
Biogenic emissions

• May be calculated “online” based on USGS landuse
  • Easy to use

• May be input

• BEISv3.13 (offline reference fields, online modified)
  • Good choice, but difficult to use

• **Use of MEGAN**
  • Best choice!!
Model of Emissions of Gases and Aerosols from Nature (MEGAN)

Global, high resolution biogenic emissions

Out of available biogenic emissions modules only BEIS and MEGAN are actively being worked on (developed)

Preprocessor for MEGAN exists and can be downloaded from NCAR
Fire Plumerise

1-D Cloud model used in WRF-Chem to determine injection height; wind shear effects are included

Satellite information (other aerial and ground observations may also be used) to determine fire location and fire properties

Emissions preprocessing may be done by (1) CPTEC preprocessor, or (2) NCAR’s FINN preprocessor.
Impact of Volcanoes

- Ash-fall near eruption
- Transport of fine ash in high concentrations for long distances
- Impact on weather, climate, and air quality

The plume of the 30 Sept/1 Oct 1994 eruption of Kliuchevskoi Volcano, Kamchatka taken from the space shuttle STS-68 mission (Russia)
10 size bins for prediction of ash-fall and transport of volcanic ash

<table>
<thead>
<tr>
<th>Particle Size Bin</th>
<th>Phi</th>
<th>Percentage of mass</th>
</tr>
</thead>
<tbody>
<tr>
<td>1 – 2mm</td>
<td>-1</td>
<td>2</td>
</tr>
<tr>
<td>0.5 – 1 mm</td>
<td>0</td>
<td>4</td>
</tr>
<tr>
<td>0.25 – 0.5 mm</td>
<td>1</td>
<td>11</td>
</tr>
<tr>
<td>125 – 250 µm</td>
<td>2</td>
<td>9</td>
</tr>
<tr>
<td>62.5 – 125 µm</td>
<td>3</td>
<td>9</td>
</tr>
<tr>
<td>31.25 – 62.5 µm</td>
<td>4</td>
<td>13</td>
</tr>
<tr>
<td>15.625 – 31.25 µm</td>
<td>5</td>
<td>16</td>
</tr>
<tr>
<td>7.8125 – 15.625 µm</td>
<td>6</td>
<td>16</td>
</tr>
<tr>
<td>3.9065 – 7.8125 µm</td>
<td>7</td>
<td>10</td>
</tr>
<tr>
<td>&lt; 3.9 µm</td>
<td>&gt; 8</td>
<td>10</td>
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• Options for transport only (4 bins or 10 bins +so2) – aerosol direct effect may be included

• Coupled with chemistry/aerosol modules (only using up to three bins – depending on size)

4 size bins for prediction if transport only is of interest

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3 size bins for coupling with other aerosol modules
WRF-Chem Greenhouse Gas Packages (*chem_opt* =17)-new in WRF-ChemV3.4

- Online calculation of biospheric CH$_4$ fluxes
  - termite – Sanderson (1996)
  - soil uptake – Ridgwell et al. (1999)

- Passive tracer simulations for CO$_2$, CH$_4$, and CO
  (including all options of CO$_2$ tracer package, *chem_opt*=16)

- Tuning of wetland fluxes through namelist options `wppeat` and `wflood` possible

- Separate biomass burning option for CO$_2$, CH$_4$, and CO including plumerise calculation (biomass_burn_opt = 5)

- Detailed description
Direct connection to NCAR’s climate modeling system: Implementation of the Community Atmosphere Model version 5 (CAM5) Physics/Chemistry

- Includes different physics options for deep and shallow convection, microphysics, boundary layer

- **Aerosols:** Liu et al. (GMD, 2012), Modal Aerosol Model (MAM)

- **Gas-Phase Chemistry:** MOZART used by “CAM-Chem” already implemented in WRF-Chem by NCAR

- **PNNL** has coupled MAM with CBM-Z photochemistry in WRF-Chem

This Climate Model package also includes aerosol direct and indirect effect, but is limited on combinations with other packages
Several dust and sea-salt models, used for bulk, modal, and sectional approaches

Lightning parameterization for NOx emissions

Aerosol interaction with convective parameterization is currently evaluated in Grell-Freitas scheme (maybe released in V3.7.1)
University of Manchester: completed developments (Lowe et al.) was added in WRF-Chem 3.6

- Common Representative Intermediate Mechanism (CRIMech) (CRIv2-R5; 240 species, 652 rxns) (Watson et al., 2008)
- N$_2$O$_5$ heterogeneous chemistry in WRF-Chem sectional aerosol (Bertram & Thornton, 2009)
- Sea-spray emission scheme with organics (Fuentes et al., 2011)
- Organic Partial Derivative Fitted Taylor Expansion (PD-FiTE) added to MOSAIC sectional aerosol (Topping et al., 2009; 2012)

Douglas Lowe, Steven Utembe*, Scott Archer-Nicholls, David Topping, Mark Barley, Gordon McFiggans
Chemical data assimilation

- NCEP’s Grid Point Statistical Interpolation (GSI, 3DVAR) assimilation system can be used with surface chemical data as well as with AOD: Significant improvements in forecasts.
- EnKF assimilation system has been used for WRF-Chem
- Work is on-going with hybrid EnKF/GSI system (ESRL and NCAR)
- Work is also ongoing with WRF-Chem adjoint development (project lead by Greg Carmichael)

These approaches are not released to community yet, but one approach with WRF DART system maybe openly available

Otherwise, if you need chemical data assimilation to help develop or use, email wrfchemhelp for contact information
Real-time AQ forecasting with WRF-Chem

http://ruc.noaa.gov/wrf/WG11_RT/

2014_07_25_00  FORECAST 45  FIELD \(\text{O}_3\) AT 0

Ozone at 21Z, tomorrow

3-ht precip at 00Z

WRF-Chem using MADE/VBS/RACM on Rapid Refresh Domain, DX=13km
HRRR-Fire: 3km horizontal resolution, runs 4x day

PM2.5 from wildfires, forecast for midday today
Resources

- WRF project home page
  - http://www.wrf-model.org

- WRF users page (linked from above)
  - http://www.mmm.ucar.edu/wrf/users

- WRF users help desk
  - wrfhelp@ucar.edu

- WRF-Chem users help desk
  - wrfchemhelp.gsd@noaa.gov

- Publications (please send us yours)

Inter-journal special issue on WRF-Chem now also opened: ACP and GMD

Thank you for your interest, Questions?