### 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** 
  - + John Wong, (University od Colorado) + other **University contributions** + Saulo Freitas (**CPTEC**, **BRAZIL**)

### + 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

### 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

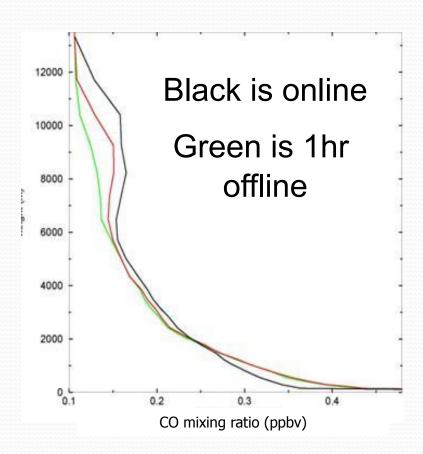
There are almost 50 chemistry options for the main gas phase chemistry and aerosol modules!

### 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
- Ideally suited 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



# What is needed for this type of modeling system?

- 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
- <u>Kinetic PreProcessor</u> (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)

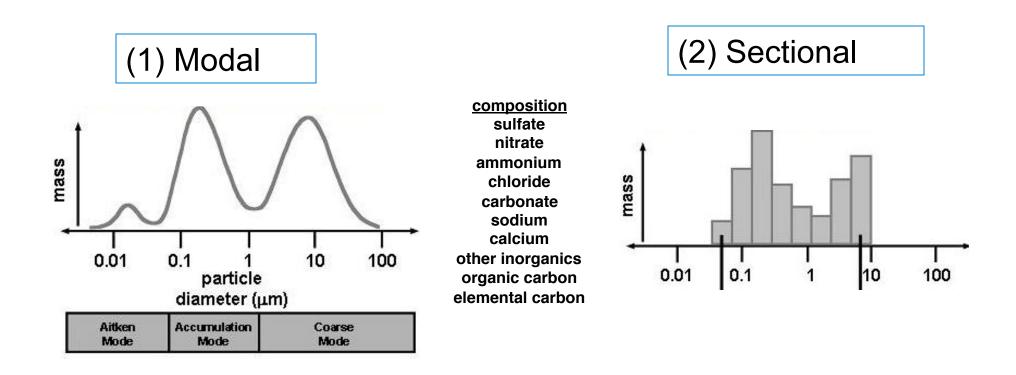
Talk by Ravan Ahmadov, Monday morning

# Photolysis Packages – all coupled to aerosols and hydrometeors

- Madronich Photolysis
- Madronich F-TUV
- Fast-j photolysis scheme

Talk by Sasha Madronich, Monday morning

#### Available aerosol modules



(3) Bulk: Sections for dust and sea salt, otherwise total mass only

Jan Kazil will give talk on aerosols in WRF-Chem Monday morning



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

Alma Hodzic will give talk on aerosol interactions on Monday



# 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)



## Selection of radiation parameterizations for aerosol "direct effect"

For 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"

For V3.5

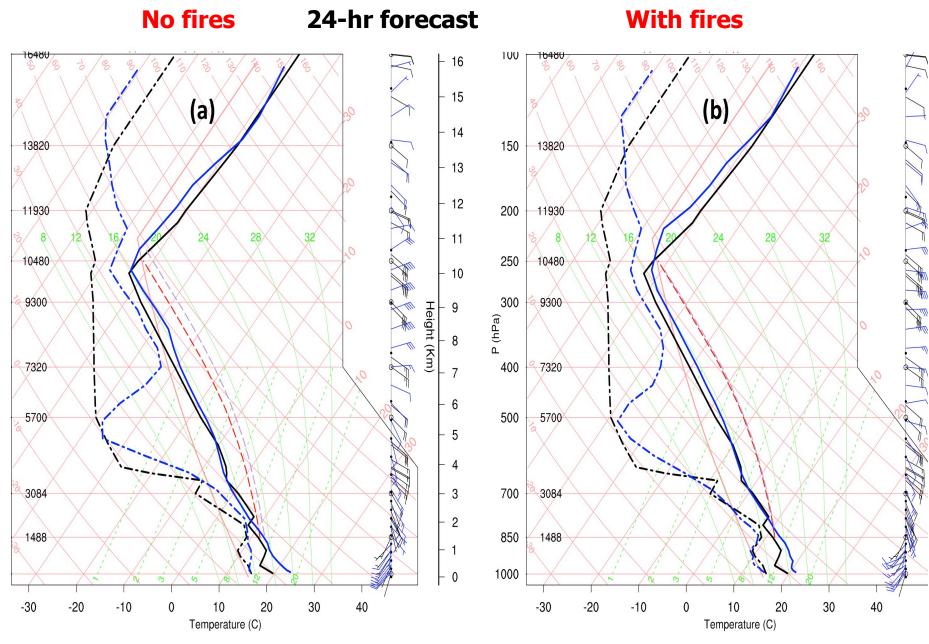
Modal and sectional schemes only can be used in combination with a version of the Lin et al. Microphysics scheme as well as the Morrison scheme

More to come in future

"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!!

### <u>Model of Emissions of Gases and Aerosols</u> from <u>Nature (MEGAN)</u>

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

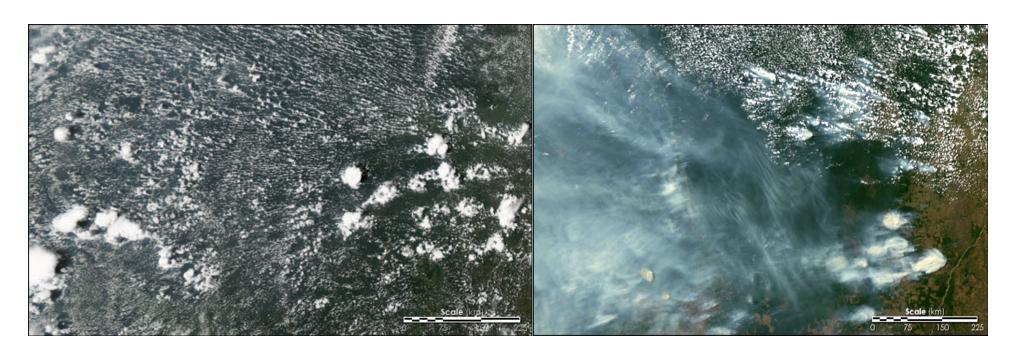
Alex Gunther will give talk on Monday

#### 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

Particle Size Bin	Phi	Percentage of mass
1 – 2mm	-1 - 0	2
0.5 – 1 mm	0 – 1	4
0.25 – 0.5 mm	1 – 2	11
125 – 250 μm	2 – 3	9
62.5 – 125 μm	3 – 4	9
31.25 – 62.5 μm	4 – 5	13
15.625 – 31.25 μm	5 – 6	16
7.8125 – 15.625 μm	6 – 7	16
3.9065 – 7.8125 μm	7 – 8	10
< 3.9 μm	> 8	10

- 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

Particle Size Bin	Phi	Percentage of mass
15.625 – 31.25 μm	5 – 6	16
7.8125 – 15.625 μm	6 – 7	16
3.9065 – 7.8125 μm	7 – 8	10
< 3.9 μm	> 8	10

3 size bins for coupling with other aerosol modules

## WRF-Chem Greenhouse Gas Packages (chem\_opt =17)-new in WRFChemV3.4

Online calculation of biospheric CH<sub>4</sub> fluxes wetland – Kaplan (2002) termite – Sanderson (1996) soil uptake – Ridgwell et al. (1999)

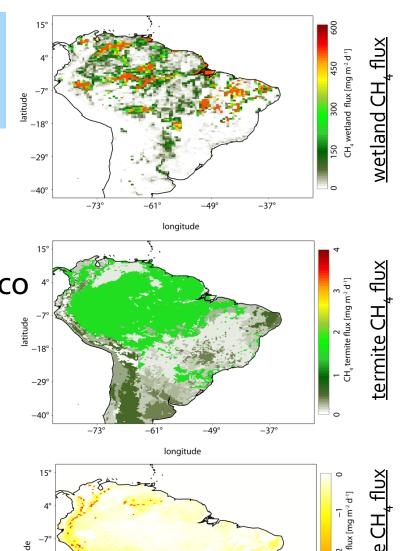
Passive tracer simulations for CO<sub>2</sub>, CH<sub>4</sub>, and CO (including all options of CO<sub>2</sub> tracer package, chem\_opt=16)

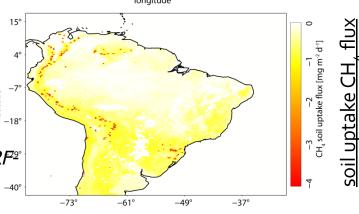
 Tuning of wetland fluxes through namelist options wpeat and wflood possible

 Separate biomass burning option for CO<sub>2</sub>, CH<sub>4</sub>, and CO including plumerise calculation (biomass\_burn\_opt = 5)

Detailed description

Beck et al., (2011): The WRF Greenhouse Gas Model (WRF<sup>2</sup> GHG) Technical Report No. 25, Max Planck Institute for Biogeochemistry, Jena, Germany, available online at http://www.bgc-jena.mpg.de/bgc-systems/index.shtml





longitude

#### 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

overview paper of CAM5 and coupling of these parameterizations (Rasch et al., 2013)

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

# University of Manchester: completed developments (Lowe et al.), and due for submission for inclusion in WRF-Chem 3.5.1/3.6

- Common Represetative Intermediate Mechanism (CRIMech) (CRIv2-R5; 240 species, 652 rxns) (Watson et al., 2008)
- N<sub>2</sub>O<sub>5</sub> 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

Developers are currently in the process of working with us to merge with the latest repository version

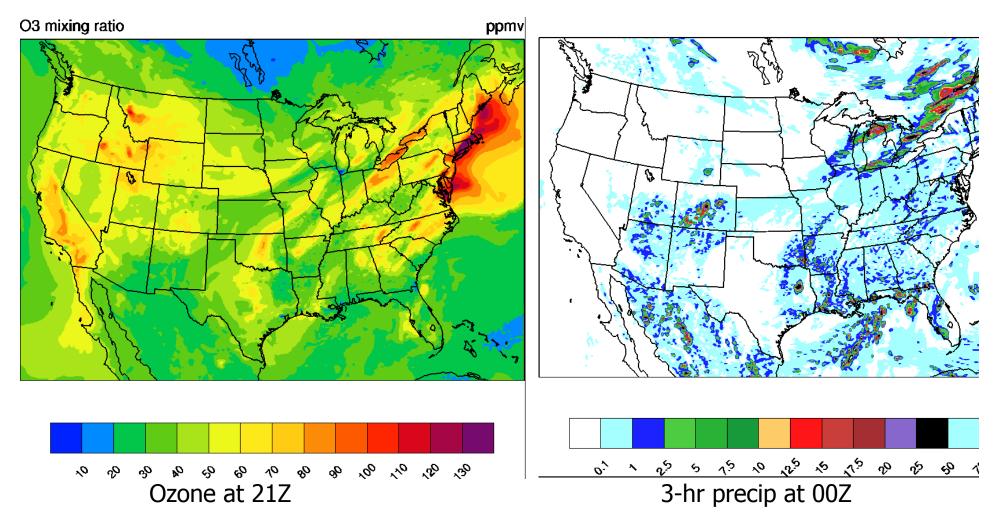
### Chemical data assimilation

- 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)

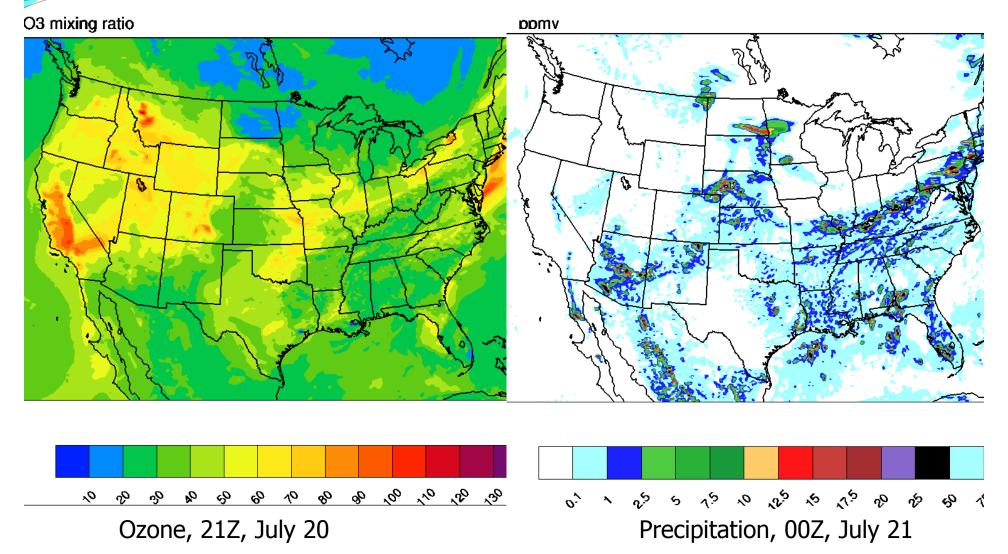
#### These approaches are not released to community yet

If you need chemical data assimilation to help develop or use, email wrfchemhelp for contact information

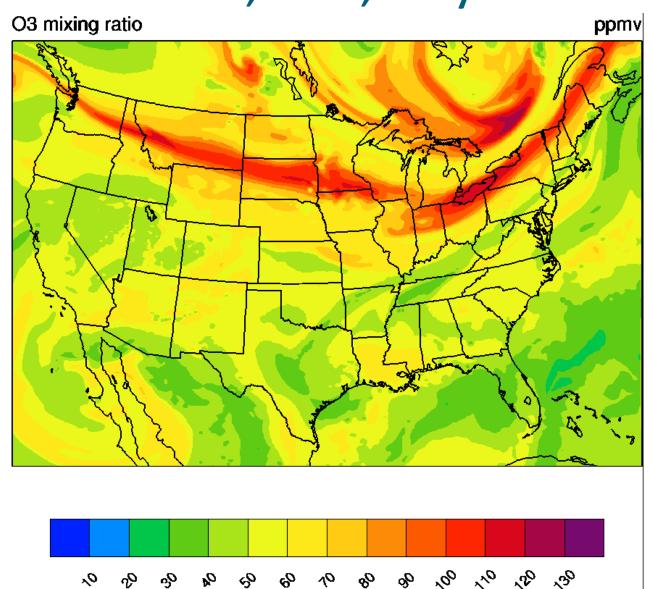
# Your real-time AQ forecast for Today http://ruc.noaa.gov/wrf/WG11\_RT/

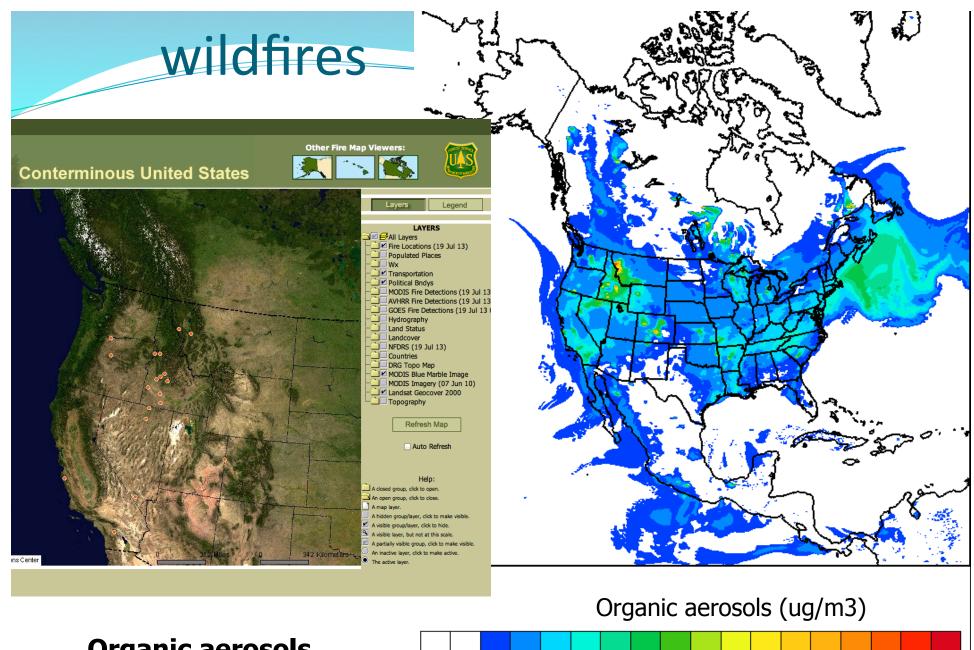


### Your real-time AQ forecast for Saturday

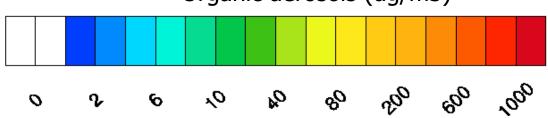


# Stratospheric Intrusion.... Ozone at 500mb, 21Z, July 20





Organic aerosols, 36hr forecast, 12Z, July 20



#### Resources

- WRF project home page
  - <a href="http://www.wrf-model.org">http://www.wrf-model.org</a>
- WRF users page (linked from above)
  - <a href="http://www.mmm.ucar.edu/wrf/users">http://www.mmm.ucar.edu/wrf/users</a>
- WRF users help desk
  - wrfhelp@ucar.edu
- WRF-Chem users help desk
  - wrfchemhelp.gsd@noaa.gov
- Publications (please send us yours)
  - <a href="http://ruc.noaa.gov/wrf/WG11/References/WRF-Chem.references.htm">http://ruc.noaa.gov/wrf/WG11/References/WRF-Chem.references.htm</a>

#### Thank you for your interest, Questions?