

# CAM-chem update

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# CAM-chem GMDD paper

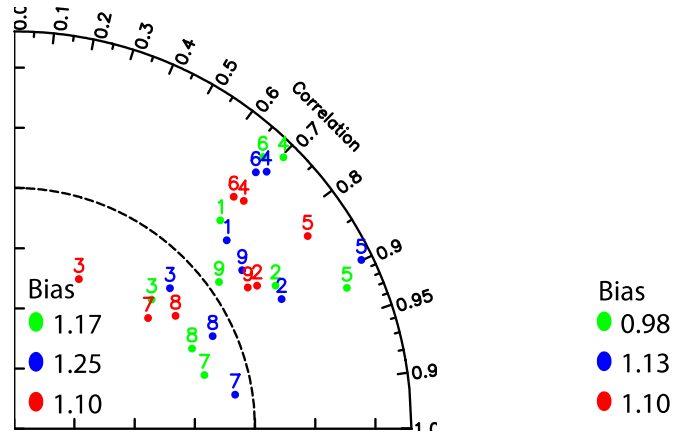
- Many thanks to all co-authors, especially Louisa, Simone, Peter and Francis.
- Should be available at GMD within a few weeks
- Led to many improvements in diagnostics

<http://www.geosci-model-dev-discuss.net/4/2199/2011/gmdd-4-2199-2011.html>

# CAM-chem GMDD paper

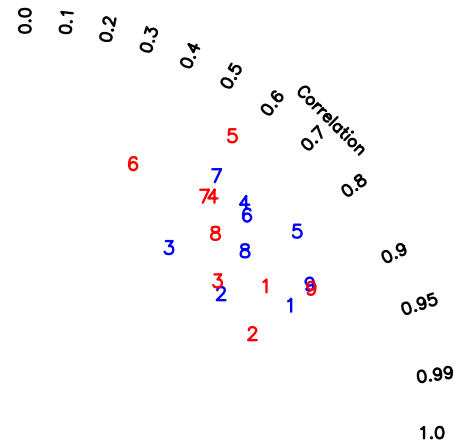
- Discusses all aspects of chemistry in CAM-chem
- Defines the MOZART-4 chemistry (not reduced or superfast) + BAM
- Compares simulations MERRA/GEOS-5/online with meteorological fields

# Taylor diagrams

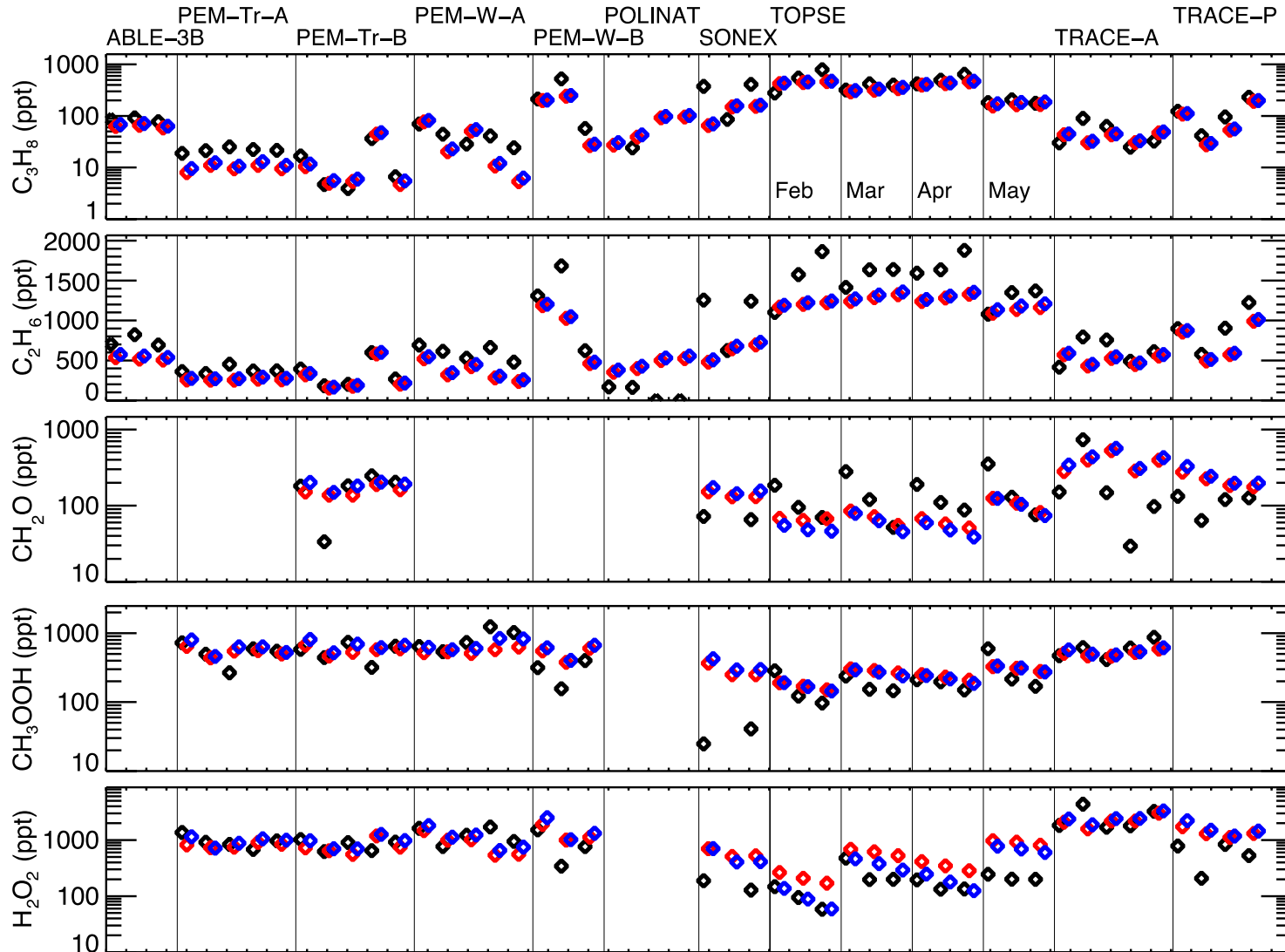


- GEOS5
- MERRA
- Online

- 1-Tropics
- 2-SH Midlat
- 3-SH Polar
- 4-North America
- 5-West. Europe
- 6-Japan
- 7-NH Polar East
- 8-NH Polar West
- 9-Canada



# Comparison with Emmons' climatology



# Participation in ACCMIP

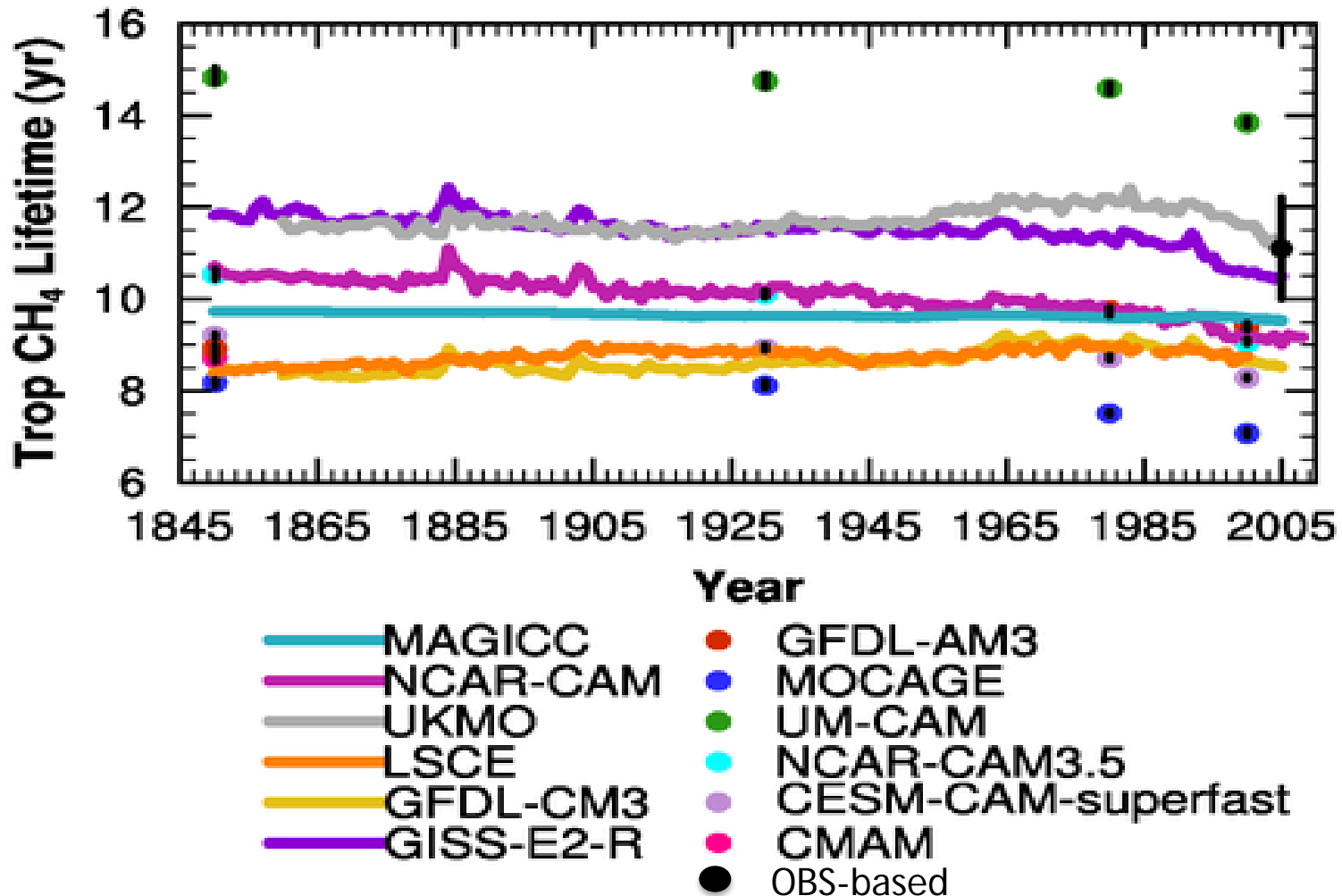
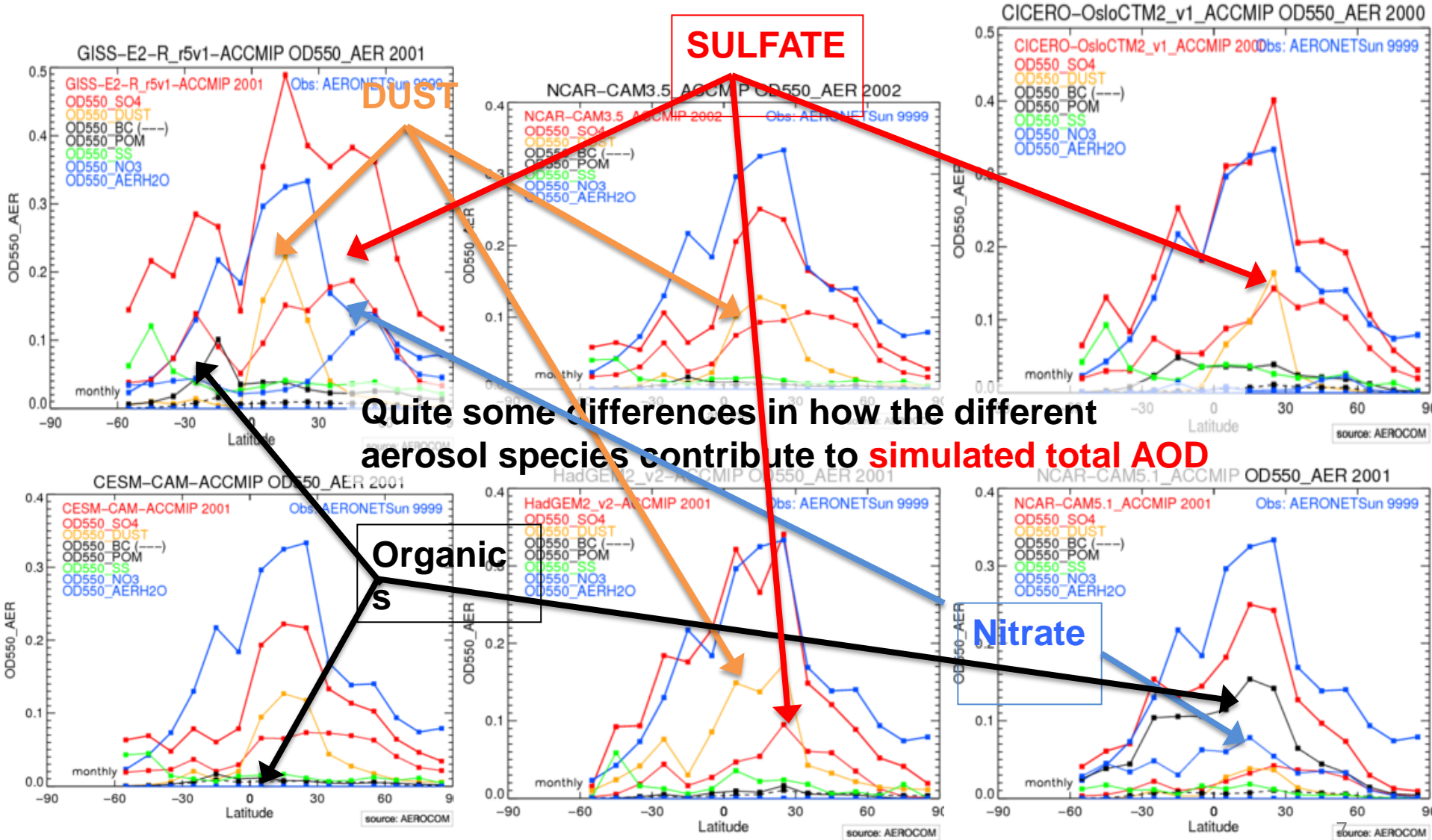


Figure courtesy of V. Naik, GFDL, 2012

# AOD aggregated per latitude band Against *Aeronet AOD*



# ChemClim Development Plan (from Breckenridge 2011)

- Top Priority
  - Update to MEGAN/include maps when possible
  - Improvements to the dry deposition (better link with CLM)
  - Coupling chemistry with MAM and CAM5 physics
  - HOMME dynamical core
- Medium Priority
  - Update SOA mechanism: Colette Heald's additional SOA species
  - Implementation of FAST-J photolysis rate computation
- Low Priority
  - “Coarse resolution” FV
- Diagnostics:
  - Tools for model result differencing
  - Benchmark numbers: methyl chloroform lifetime, ozone budget terms, methane lifetime, mass-weighted tropospheric OH, lightning NO<sub>x</sub>, sf(co/nox/isoprene)

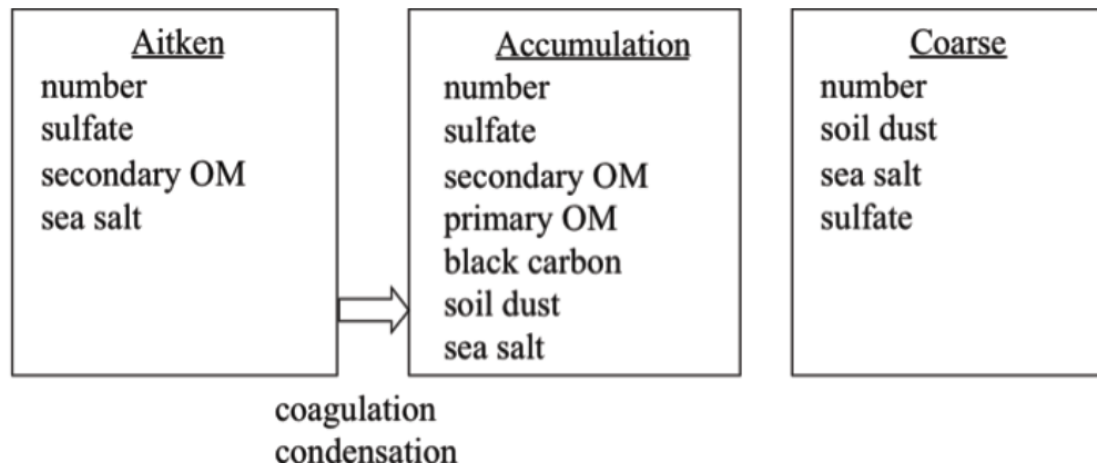


# ChemClim Development Plan (from Breckenridge 2011)

- Top Priority
  - Update to MEGAN/include maps when possible (see L. Emmons' presentation)
  - Improvements to the dry deposition (better link with CLM)
  - Coupling chemistry with MAM and CAM5 physics
  - HOMME dynamical core (DOE funding: A. Conley/P. Lauritzen)
  - kPP mechanism
  - Box Model or SCAM w/ chemistry
- Medium Priority
  - Update SOA mechanism: Colette Heald's additional SOA species (additional work by K. Barsanti)
  - Implementation of FAST-J photolysis rate computation (DOE funding: M. Prather/P. Cameron-Smith)
  - Conversion of preprocessor to KPP?
  - Vertical resolution and model top?
  - WACCM lite?
- Low Priority
  - "Coarse resolution" FV
- Diagnostics:
  - Tools for model result differencing
  - Benchmark numbers: methyl chloroform lifetime ozone budget terms, methane lifetime, mass-weighted tropospheric OH, lightning NO<sub>x</sub>, sf(co/NO<sub>x</sub>/isoprene)

# Coupling chemistry to MAM3 (1)

- MAM: Modal Aerosol Module (Liu et al., GMDD, 2011: [www.geosci-model-dev-discuss.net/4/3485/2011](http://www.geosci-model-dev-discuss.net/4/3485/2011))
- Very simplified chemistry (reads  $O_3$ , OH,  $NO_3$  and  $HO_2$  to compute DMS and  $SO_4$  oxidation to  $H_2SO_4$ )



# Coupling chemistry to MAM3 (2)

- Use MOZART mechanism (will also work with reduced NMHC) to provide  $O_3$ , OH,  $NO_3$ ,  $HO_2$  and  $H_2O_2$
- Modify scheme to include  $H_2SO_4$  (instead of directly making  $SO_4$ )
- Use surface area from MAM to compute rate for heterogeneous reactions on tropospheric aerosols

# Coupling chemistry to MAM3 (3)

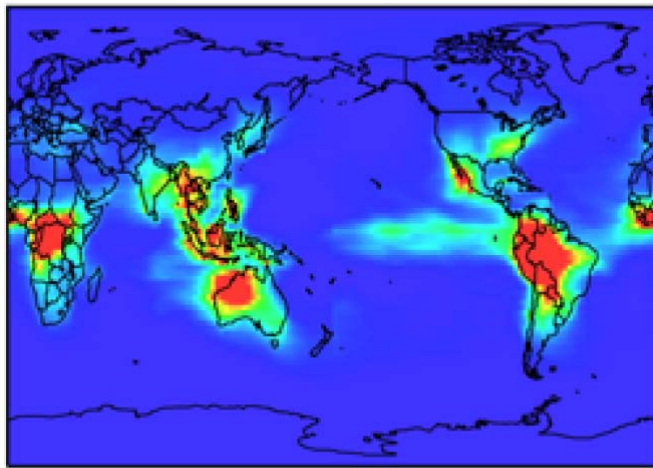
- Remaining issues (that we know 😊)
  - SOA in MAM not coupled with chemistry
  - Dry deposition is done differently (MAM needs size-dependent scheme)
  - SAD only from Aitken mode
- Status
  - Have performed several multi-year simulations for present-day, F case only
  - Initial evaluation (including clouds) indicates reasonable results
  - Code will be released in June 2012
  - Similar effort in WRF-chem
  - Evaluation in S. Tilmes' talk

# SOA (from C. Heald)

- Current version: two-product method (Lack et al., 2004)
- From Colette Heald (Heald et al., JGR, 2008)
  - Two-product method with high- and low-NO<sub>x</sub> yields
  - Additional anthropogenic precursors (benzene, toluene and xylene) and isoprene
  - Split SOA into classes (one per precursor)
  - T-dependent partitioning between gas-phase and solid-phase (following Chung and Seinfeld, 2002)
  - Adds 13 species

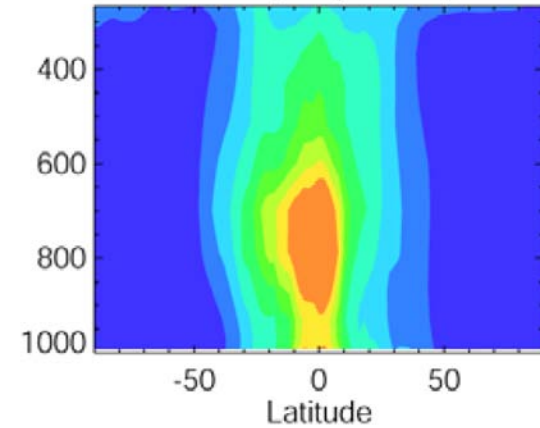
# Comparison w/ Heald et al. (2008)

SOA (isoprene)

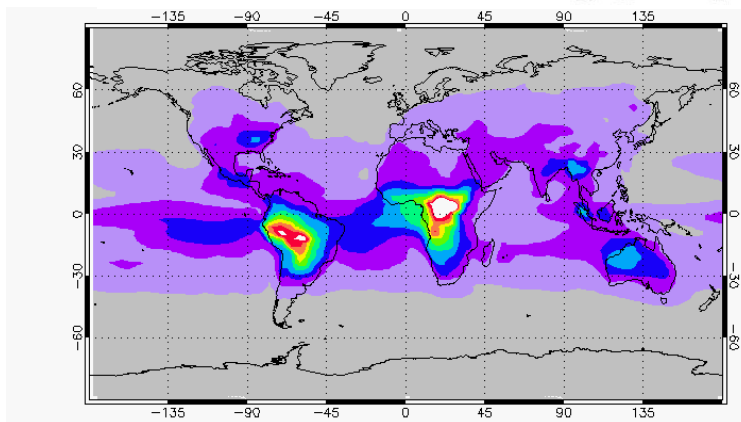


0.00 0.50 1.00 [ $\mu\text{gCm}^{-3}$ ]

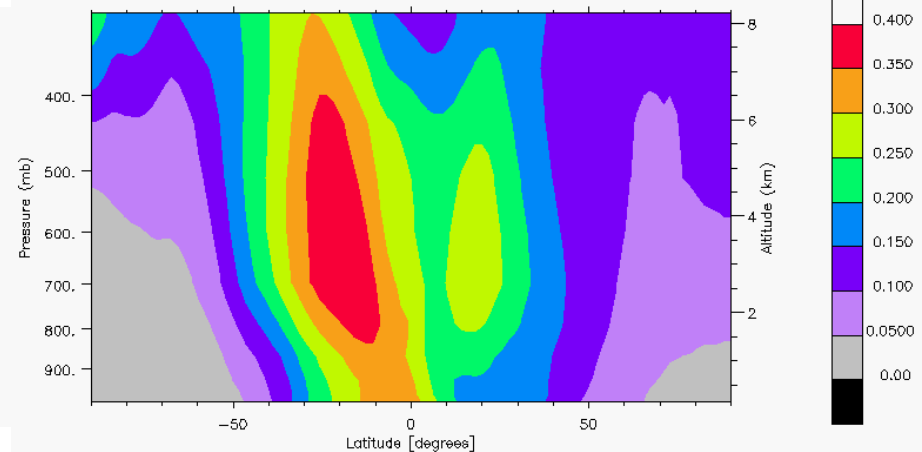
SOA (isoprene)



0.00 0.10 0.20 0.30 0.40 [ $\mu\text{gCm}^{-3}$ ]



1.00  
0.900  
0.800  
0.700  
0.600  
0.500  
0.400  
0.300  
0.200  
0.100  
0.00

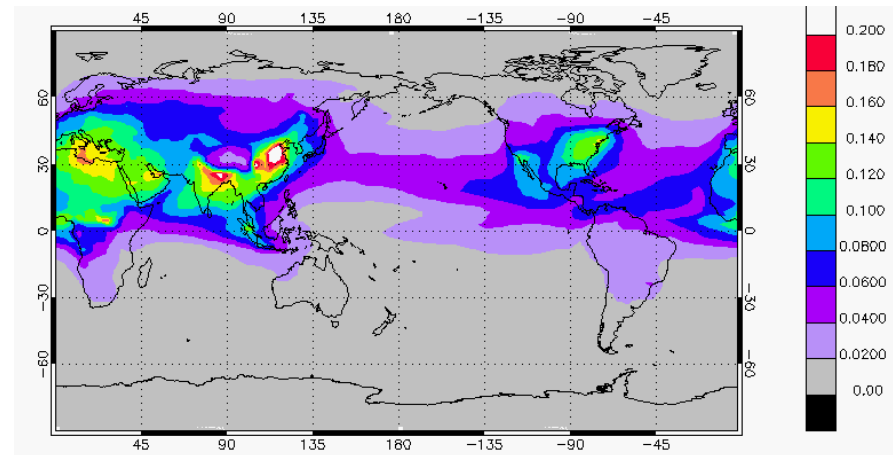
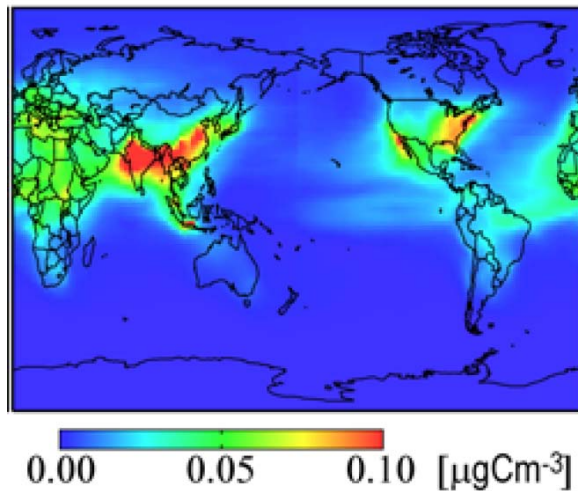


0.400  
0.350  
0.300  
0.250  
0.200  
0.150  
0.100  
0.0500  
0.00

Simulation is for one-year only

# Comparison w/ Heald et al. (2008)

SOA (aromatics)



Approximately 2x SOA from aromatics

- Bug?
- Emissions?

# Next steps: CSL allocation

- Development
  - New chemistry (halogen/CRI)
  - High-resolution (at least 0.5°)
  - Vertical resolution
  - SE dynamical core (ultimately with varying resolution!)
- Production
  - Chemistry-climate response to regional forcing
  - Benchmark simulations for CAM5-chem FV
    - PI control
    - 4xCO<sub>2</sub>



# Chemistry-climate coupling: BC

