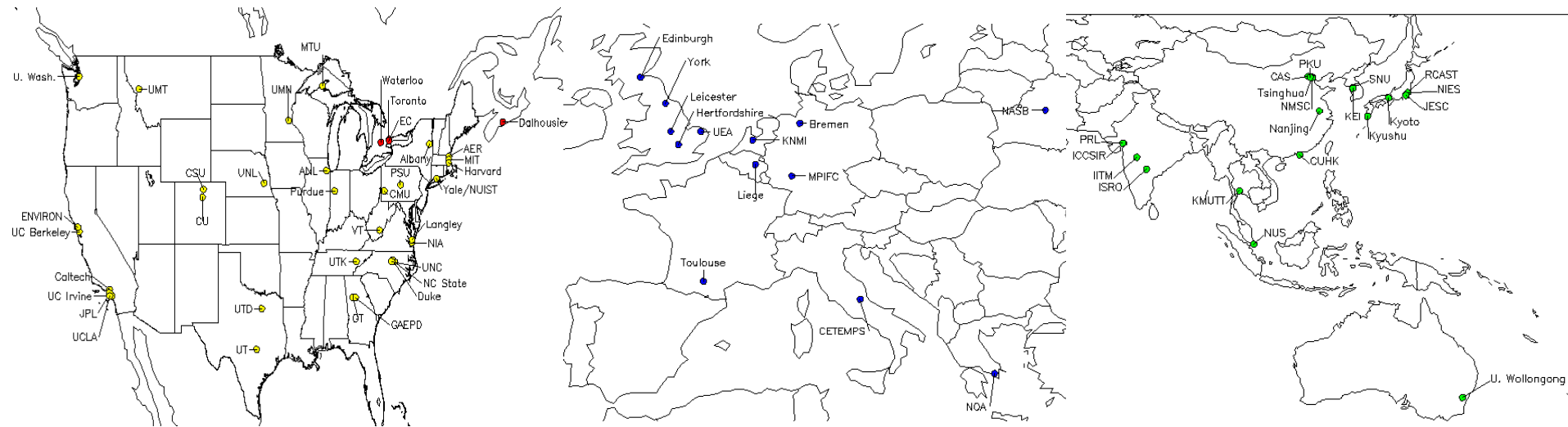


Welcome to the 6th International GEOS-Chem Meeting!

The GEOS-Chem user community: 70 institutions, 18 countries



Thanks to our sponsors:



NATIONAL ENVIRONMENT RESEARCH COUNCIL



ELECTRIC POWER
RESEARCH INSTITUTE



HARVARD

School of Engineering
and Applied Sciences



Harvard University
Center for the Environment

Thanks to Sandra Klemm, Brenda Mathieu, Harvard grad students for logistics

Meeting Objectives

Exchange information at the frontier of knowledge of atmospheric composition

- IGC6 is first and foremost a scientific meeting

Promote interactions both within and outside GEOS-Chem community

Sustain functioning of the GEOS-Chem community

- Model clinics
- Working Group meetings
- Business meeting: leadership/communications, model standards, long-term vision

Identify model development priorities for the next two years

- A major charge for the Working Group meetings

Full Agenda: 90 Talks, 70 Posters, Clinics, WG Meetings

Monday, May 6

Model Overview

GEOS-Chem Working Group Introductions

Aerosol Sources and Chemistry

Black Carbon and Related Processes

Model Clinics (Forward, Adjoint, Met & Grid Independent)

1st Poster Session (6:00 onward)

Tuesday, May 7

Aerosol Effects on Climate and Air Quality

Regional Air Quality (posters and introductions)

Mercury and POPs

1st Poster Session (12:40-1:40)

Climate-Chemistry

Working Group Meetings (discussion, development needs)

(Sources & Sinks, Chemistry-Climate, Hg & POPs)

(Oxidants and Chemistry, Carbon, Nested)

2nd Poster Session (6:00 onward)

**Interspersed
Scientific
Presentations
&
Users Meeting**



Full Agenda: 90 Talks, 70 Posters, Clinics, WG Meetings

Day 3: Wednesday, May 8

Sources and Sinks: Nitrogen

Tropospheric Ozone & Photochemistry (posters and intros)

Carbon Dioxide Fluxes

Sources of other Carbon Gases (posters and intros)

2nd Poster Session (1:10-2:10)

Sources of other Carbon Gases (oral)

Working Group Meetings (discussion, development needs)
(Adjoint, Aerosols, Software Engineering)

Day 4: Thursday, May 9

Tropospheric Ozone & Photochemistry (oral)

Regional Air Quality (oral)

GEOS-Chem Business Meeting (decide priorities, vision)

Awards for Poster Introductions

Live video cast organized by Patrick Kim and Alex Turner



Interspersed
Scientific
Presentations
&
Users Meeting



Some Logistics

Uploading Talks (Thanks to Patrick Kim and Alex Turner)

<http://www.fas.harvard.edu/~geoschem>

Username: GCMeeting

Password: GCUser

Directions on submission page

Presentations must be submitted **before** the break before your session

If you don't have web access, bring a thumb-drive to a member of the A/V team (Patrick Kim or Alex Turner)

in the front of the room at the **beginning** of the break

Oral Presentations

Max 7 minutes

Remaining 3 minutes for Q&A + changeover

Posters

Oral introductions: max 30 seconds

Session 1 posters must be taken down before Tuesday evening

GEOS-Chem Chemical Transport Model (CTM)

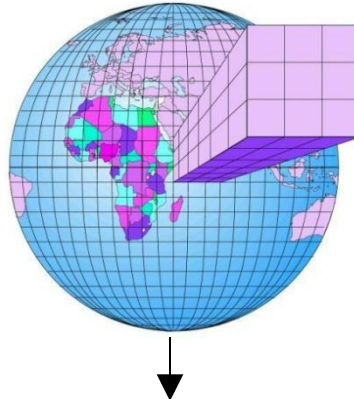
Input data

- NASA GEOS meteorological fields
- other

Model solves 3-D chemical continuity equations
on global or nested Eulerian grid

Modules

- emissions
- transport
- chemistry
- aerosols
- deposition
- sub-surface



Model adjoint

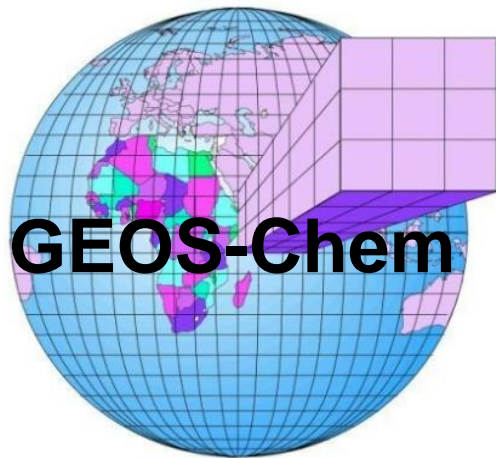
Applications

- chemical processes, transport, budgets
- inverse analyses
- radiative forcing
- air quality
- biogeochemistry
- ...

Model capabilities:

- Tropospheric aerosol-oxidant chemistry, aerosol microphysics, CO₂, methane, mercury, various tracers
- 1980-present GEOS meteorological data, future and paleoclimates (GCMs)
- Horizontal resolution: 0.25°x0.31° (native GEOS-5), 1/2°x2/3°, 2°x2.5°, 4°x5°

GEOS-Chem pushes the frontier of atmospheric composition modeling while also contributing to broader Earth science & applications



Observing systems

- **chemical data assimilation**

Chemistry-climate interactions

- **interface with GCMs, ESMs**

Satellite retrievals

- ***a priori* information**

Local air quality models

- **outer nest**

Functioning of the GEOS-Chem community

- **GEOS-Chem is a grass-roots community model**
 - Code is open-access, extensively documented and supported
 - All model development is initiated by users to serve their research needs
 - New developments are promptly implemented in the standard model
- **Strong model management maintains model unity, traceability, state of science**
 - GEOS-Chem Support Team (Harvard/Dalhousie/CU) maintains standard and adjoint model: developments, benchmarking, version control, user support
 - Support Team is funded by NASA (1.5 FTE) and NSERC (0.5 FTE); strong leverage from community volunteerism is essential
 - GEOS-Chem Steering Committee monitors model development, promotes good practices and communication, leads Working Groups
- **The model is owned by its users, and with ownership comes responsibilities:**
 - Keep up with the model (newsletters, wiki, working groups, IGC meetings)
 - Contribute to the community: help with requests, report bugs...
 - Share mature developments for incorporation into standard model
 - Update regularly to latest standard version of model
 - Be generous in credit to developers

The success of GEOS-Chem relies on the community spirit of its users!

Model development flow

Users

- develop modules
- report problems

Working Groups

- set model development priorities (IGCx)

Steering Committee

Scientists, WG chairs,
Engineer, GMAO reps

- implements and updates priorities
- designs successive model versions
- evaluates benchmarks

GEOS-Chem Support Team

New Model Versions

Future versions to have
hard releases every 9 months

Version (v9-1-3) released in September 2012;
Version (v9-2) to be released in summer 2013

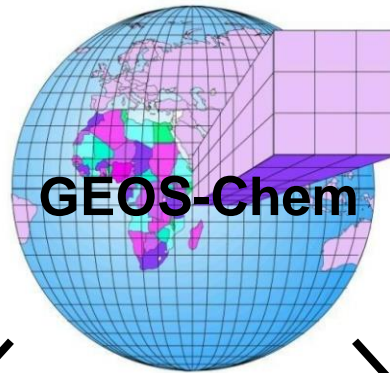
Working Groups:

1. Adjoint model & Data Assimilation
2. Aerosols
3. Carbon gases & organics
4. Chemistry-climate
5. Hg and POPs
6. Oxidants and chemistry
7. Nested model
8. Sources and sinks

Major GEOS-Chem developments over the horizon

- **0.25°x0.31° nested capability (Tsinghua, Dalhousie)**
 - Take it for a spin! Will be in standard model v9-2
- **Grid-independent GEOS-Chem (Harvard, NASA GMAO)**
 - Allows model to run on any grid, with any dynamical core, interface with any model through ESMF
 - Includes new emission libraries module
 - In advanced development stage; will eventually be GEOS-Chem
- **Increased computational performance (Harvard, York, GMAO)**
 - New implementation of KPP chemical solver
 - MPI parallelization through ESMF
- **RRTMG radiative transfer code (AER, MIT, SUNY-Albany)**
 - Allows radiative forcing calculations from UV to IR
- **Stratospheric chemistry (MIT)**
 - will substitute “climatological chemistry” presently in GEOS-Chem

Dual driving forces in GEOS-Chem development: A positive force!



**GEOS-Chem as a tool
for data interpretation**

**GEOS-Chem as a component
of Earth System Models**

**Push for improved processes,
input data, resolution**

**Push for improved performance,
flexibility, software engineering**

