

**NORTHROP GRUMMAN**



1

# GMI Status

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**GODDARD SPACE FLIGHT CENTER**

**GMI Science Meeting  
June 13-15, 2007  
Greenbelt, MD**

**Software Integration and Visualization Office**

## Outline

- **SIVO GMI personnel**
- **SIVO support to GMI Community**
- **New GMI directory structure**
- **New GMI naming convention**
- **New GMI code features**
- **Automating/streamlining GMI runs**
- **Completed model runs**
- **Upcoming tasks**

## Personnel

- **Jules Kouatchou**
  - **AMTI, Programmer/developer**
- **Megan Damon**
  - **Northrop Grumman, Programmer/developer**
- **Tasks**
  - **Implement and document code changes**
  - **Manage the GMI CVS repository**
  - **Assist GMI community with building and running the code**
  - **Perform requested runs**
  - **Process, archive, and distribute data**



## **SIVO Support to GMI Community**

- **Bldg 33 GMI Scientists**
- **Georgia Tech**
- **Boeing**
- **AER**
- **GMAO**
- **University of Michigan**
- **NASA-Langley**
- **PNL**

## GMI Directory Structure

- Added new directory and directory structure on **dirac**
  - **/g8/anon/pub/gmidata2/input/**
    - output/**
    - .../output/gmic/aura2/2004/diagnostics**
    - docs/**
    - progs/**
    - users/**
- **/g8/anon/pub/gmidata** still exists, untouched
- Input data also available on **discover**
  - **/discover/nobackup/projects/gmi/gmidata2**
  - Met fields are on-line



## GMI File Naming Convention

- For input files
  - /g8/anon/pub/gmidata2/input/emissions/trop/  
emis**X\_yyyy\_Z\_Res\_Descriptor**.nc  
emist\_2004\_m\_2x2.5\_auraharvardwship.nc
- For output files
  - /g8/anon/pub/gmidata2/output/gmic/aura2/2004/  
**problemname.datatype.filetype**  
gmic\_aura2\_2004\_jul.monthly.nc



## New GMI Inputs

- **Emissions**
  - Added capability to use daily emissions
  - Emissions inputs now can include ship emissions
- **Chemistry**
  - **Photolysis**
    - Monthly rates can be used (GOCART aerosols)
    - Aerosol interaction with photolysis (fastjx53c) added
  - JPL06-2 mechanism is currently used
  - Default mechanisms
    - **124 species** for Combo Model
    - **85 species** for Troposphere Model

## New GMI Outputs

- **Variables added to all output files**
  - **Grid box height**
  - **Grid box area (mcor)**
  - **Grid info (ai, bi, am, bm)**
- **New outputs to some files**
  - **File description that is placed in header**
  - **Potential vorticity**
  - **Tropopause pressure**
  - **Relative humidity**
  - **Cloud optical depth**
  - **Overhead column O<sub>3</sub>**
  - **Surface emissions for aerosols**
  - **Grid box mass**





## New GMI Outputs

- **New output files**
  - **Set with freq settings**
    - **Column ozone (.columnoz.nc)**
    - **Hourly data (.hourly.nc; can select species/levels to output)**
    - **Instantaneous daily mixing ratios (.idaily.nc)**
  - **Four overpass times (formerly “noon\_species”; .overpass1.nc, etc.)**
- **Output file name changes**
  - **.const.nc is now .amonthly.nc**
  - **Station .column.nc is now .profile.nc**

## Namelist File Changes

- Namelist sections reorganized and restructured
- Input/output species changed from indices to names
  - **From:**
    - `flux_species(1:124) = 1, 1, 1, 0, 0, 1, 0...`
  - **To:**
    - `fluxSpeciesNames = 'CH2O, CO, H2O2, HNO3,...'`

# Namelist File Changes

- **Profile stations selection simplified**
  - **From:**
    - `col_diag_site(1:201) = 'SPO', 'MCM', 'HBA',...`
    - `col_diag_lat_lon(:,1) = -89.98, 335.20,`
    - `col_diag_lat_lon(:,2) = -77.83, 166.60,`
    - `col_diag_lat_lon(:,3) = -75.56, 333.50,`
  - **To:**
    - **stationsInputFileName**
      - Master list with station names, locations, and descriptions
      - File name set in namelist file
    - **colDiagStationsNames = 'SPO, MCM, HBA,...'**
      - User-selected list of stations to study

## Other Code Changes/Issues

- **New tropopause definition**
- **New lightning algorithm**
- **Tagged CO/age of air capability**
- **Species **names** replace species **indices****
- **Added flux diagnostic version to trunk**
- **Componentization of the code**
- **Assisting GMAO with integration of GEOS-5 and the Combo Model chemistry**

## Other Code Changes/Issues

- **Disclaimer/Notice of Release added**
- **Open-source process has begun**
- **Version numbering system established**
  - **Three digit system: X.Y.Z**
  - **Current version: 2.0.0**
- **Made GMI simulation date Y2K compliant**
- **New *gmi-users* mailing list**
- **Model runs transitioned to *discover***



## Unintended GMI “Features” Addressed

- **Soil NO<sub>x</sub>**
- **Truncated vertical ozone profile**
- **Mixing ratio blow-up due to convection**
- **Divide by zero potential in convection**
- **Advection code array bounds**
- **H<sub>2</sub> level not set properly**
- **Stratospheric chemistry divide by zero potential**
- **HO<sub>2</sub> heterogeneous chemistry reaction**



## Automating/Streamlining GMI Runs

- **Met field processing/regridding scripts**
  - Scripts automatically begin processing when GMAO releases new files
- **Model run submission/monitoring scripts**
  - Submits a series of model runs
  - Emails user when each phase is complete
- **Namelist/met file list generation scripts**

# Automating/Streamlining GMI Runs

- **GMI Workflow tool**
  - **GUI with “behind the scenes” scripts to do model runs “end to end”**
    - **Set up namelist file**
    - **Submit job**
    - **Graphically display run progress**
    - **View output data graphically**



# GMI Workflow Namelist GUI

Universal Experiment Designer: GMICombo-build.ued

6. do_wetchem	Boolean	<input type="checkbox"/>
7. loss_opt	Integer	0
8. do_AerDust_Calc	Boolean	<input checked="" type="checkbox"/>
9. AerDust_Effect_opt	Integer	0
10. AerDust_infile_name	String	._agcm_kg_per_m3_2%5x2x42.nc
11. h2oclim_opt	Integer	2
12. h2oclim_infile_name	String	2o_ch4_clim_2%5x2x42_fvccm.nc
13. sad_opt	Integer	2
14. lbssad_opt	Integer	3
15. lbssad_infile_name	String	ata/sad_dbc_2%5x2x42_2000.nc
16. forc_bc_opt	Integer	2
17. forc_bc_years	Integer	100
18. forc_bc_start_num	Integer	34
19. forc_bc_kmin	Integer	1
20. forc_bc_kmax	Integer	2
21. forcedBcSpeciesNames	ArrayList String	<input type="checkbox"/> CH2O <input checked="" type="checkbox"/> CH4 <input type="checkbox"/> CO <input type="checkbox"/> H <input type="checkbox"/> H2 <input type="checkbox"/> HCOOH <input type="checkbox"/> HNO2 <input type="checkbox"/> HNO3 <input type="checkbox"/> HN <input checked="" type="checkbox"/> N2O <input type="checkbox"/> NO <input type="checkbox"/> NO2 <input type="checkbox"/> NO3 <input type="checkbox"/> N2O5 <input type="checkbox"/> O <input type="checkbox"/> O1D <input type="checkbox"/> O3 <input type="checkbox"/> OH <input type="checkbox"/> Cl2 <input type="checkbox"/> ClO <input type="checkbox"/> Cl2O2 <input type="checkbox"/> ClONO2 <input type="checkbox"/> HCl <input type="checkbox"/> HOCl <input type="checkbox"/> OClO <input checked="" type="checkbox"/> CH3Br <input checked="" type="checkbox"/> CH <input checked="" type="checkbox"/> HCFC22 <input checked="" type="checkbox"/> HCFC141b <input checked="" type="checkbox"/> HCFC142b <input checked="" type="checkbox"/> CF2Br2 <input checked="" type="checkbox"/> CF2ClBr <input checked="" type="checkbox"/> CF3Br <input checked="" type="checkbox"/> H2402 <input type="checkbox"/> A302 <input type="checkbox"/> AC <input type="checkbox"/> ETO2 <input type="checkbox"/> ETP <input type="checkbox"/> GC03 <input type="checkbox"/> GLYC <input type="checkbox"/> GLYX <input type="checkbox"/> GP <input type="checkbox"/> GPAN <input type="checkbox"/> HAC <input type="checkbox"/> IAL <input type="checkbox"/> KO2 <input type="checkbox"/> MACR <input type="checkbox"/> MAN2 <input type="checkbox"/> MAO3 <input type="checkbox"/> MAOP <input type="checkbox"/> MAP <input type="checkbox"/> MCO3 <input type="checkbox"/> MEK <input type="checkbox"/> MG <input type="checkbox"/> PP <input type="checkbox"/> PPN <input type="checkbox"/> PRN1 <input type="checkbox"/> PRPE <input type="checkbox"/> PRPN <input type="checkbox"/> R4N1 <input type="checkbox"/> R4N2 <input type="checkbox"/> R4O2 <input type="checkbox"/> R4I <input type="checkbox"/> RIP <input type="checkbox"/> ROH <input type="checkbox"/> RP <input type="checkbox"/> VRO2 <input type="checkbox"/> VRP <input type="checkbox"/> ACET <input type="checkbox"/> N2 <input type="checkbox"/> O2 <input type="checkbox"/> Tot
22. forc_bc_infile_name	String	p2002_ch4latvar_1970_2069.asc

Task Properties

Task Name: nlGmiChemistry

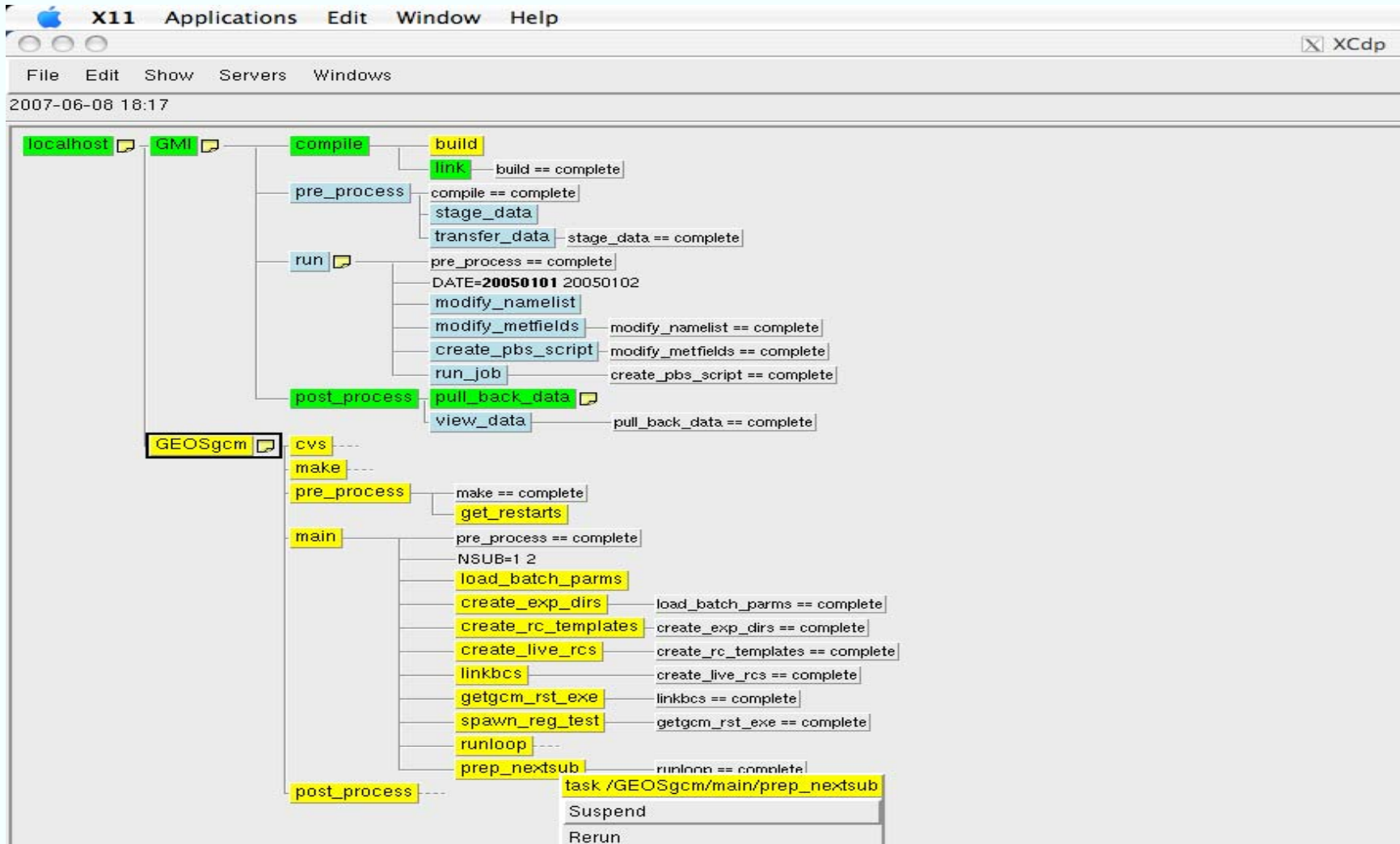
Task Description: GMI Chemistry section

Task Script:

Variable Script: ne/mrdamon/workflow/workflow/cfg/gmi/GMI/combo.in

Variable Script Type: FORTRAN Namelist

# GMI Workflow Run Monitoring



# Completed GMI Runs

- **Aura Runs**
  - **GEOS4-DAS met fields 2004-2006**
  - **Forecast met fields Jul 2004-Jun 2005**
    - 12 h met field forecasts
    - 24 h met field forecasts
- **Incremental Tests**
  - Tropopause definition
  - 124 species
  - JPL06-2
  - H<sub>2</sub> fix
  - HO<sub>2</sub> heterogeneous chemistry

## Completed GMI Runs

- **HTAP runs (with some post-processing)**
  - SR1 full chemistry and aerosol
  - SR6 aerosol
  - SR6 AEROCOM aerosol
- **Aerosol runs**
  - Present day (2000)
  - Pre-industrial (1750)
  - DAO, GISS, and FVGCM met fields
- **Boreal wildfire emissions runs**
  - Jun-Sep 2004
  - Forecast and GEOS4-DAS met fields



## Upcoming Tasks

- **Resolution doubling experiments**
- **Coupled aerosol/gas-phase chemistry**
- **Lightning experiments**
- **Support the integration of GEOS-5 with Combo chemistry**
- **Implement new TP Core**
- **GMI User's Guide**
- **HTAP runs/post-processing**
- **Evaluating fastjx53c and fastjx61**
- **Processing of GEOS5 met fields**