

**Whole Atmosphere  
Community Climate Model:  
Going from v3.1.9 to v3.5.48**

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303-497-1469**

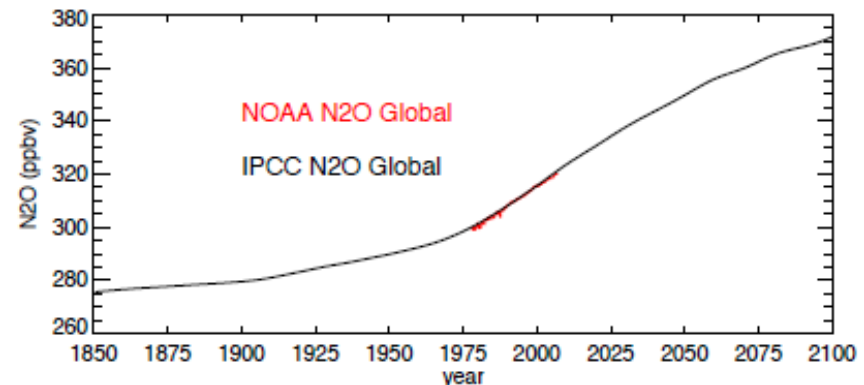
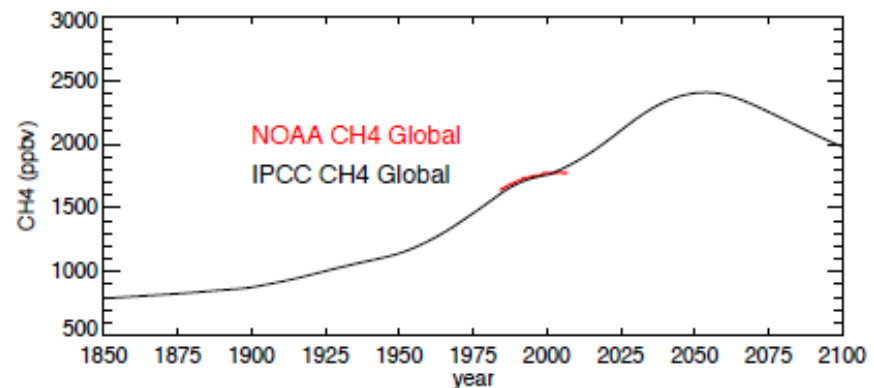
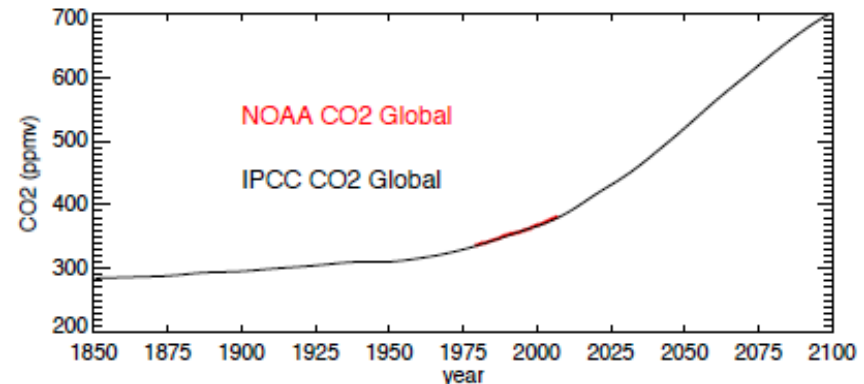


# Outline

- **Updated GW scheme [Richter / Garcia].**
- Parameterization of unresolved mountain drag [Sassi et al.].
- **Updated source gas boundary conditions [1850-2100].**
- Conservation of CLOY [and BrOY].
- **New Sulfate SAD time series [SPARC].**
- Volcanic heating module added.
- **Update PSCs module [NAT] – less denitrification [more SAD].**
- LEAN Extra Terrestrial Fluxes [replaced Woods/Rottman/Neckel]
- **QBO inclusion [not used extensively in 319]**
- Chemical Mechanism: JPL06 [+photolysis benchmark]
- **Specified Dynamics Option Developed [w/GEOS5.1].**
- Solar Proton Events [not used extensively in 319].
- **Unification of WACCM3 and CAM-CHEM modules [not complete].**

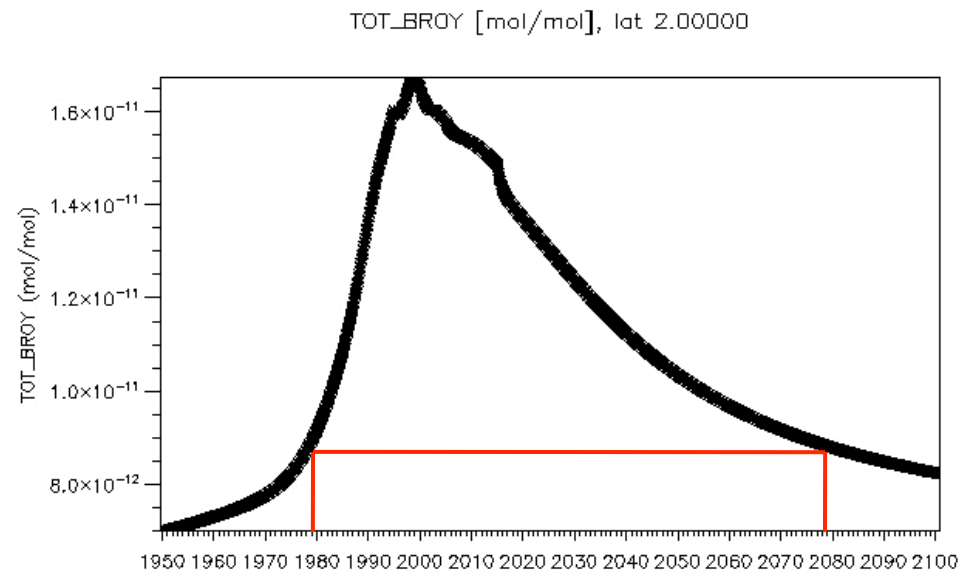
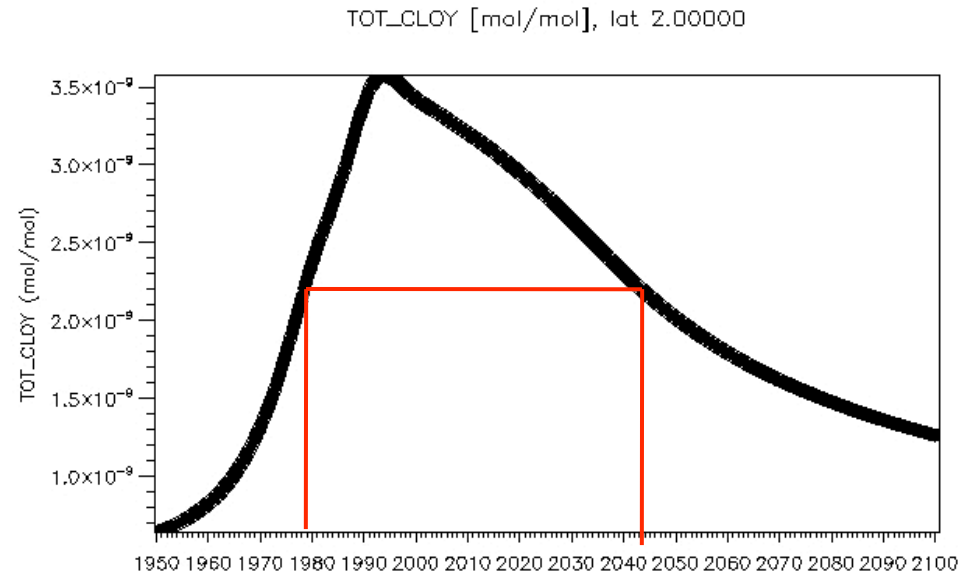
# Updated source gas boundary conditions [1850-2100]

- **CO<sub>2</sub>, N<sub>2</sub>O, and CH<sub>4</sub> from AR4**
  - A1B-medium scn
  - Not what CCSM used in AR4.
  - 1850 – 2100 [ready for WACCM3 / deep ocean sim's]
  - CO<sub>2</sub> and CH<sub>4</sub> have latitudinal and seasonal cycle included.
- **New Halocarbon time series [Guss Velders].**
  - 1951 - 2100
  - WMO 2006 [chpt8].
  - w/Montreal 2007 HCFC adj.
- **New CO, NO<sub>x</sub>, CH<sub>2</sub>O, NMHC Emissions for 1997-2008 [Emmons].**
  - Specific for each year.
  - 1999-2007 average
- **Coming soon...**
  - LBC for VSL Org BrOY and IOY spc.
  - Time series of NMHC's [IPCC AR5]



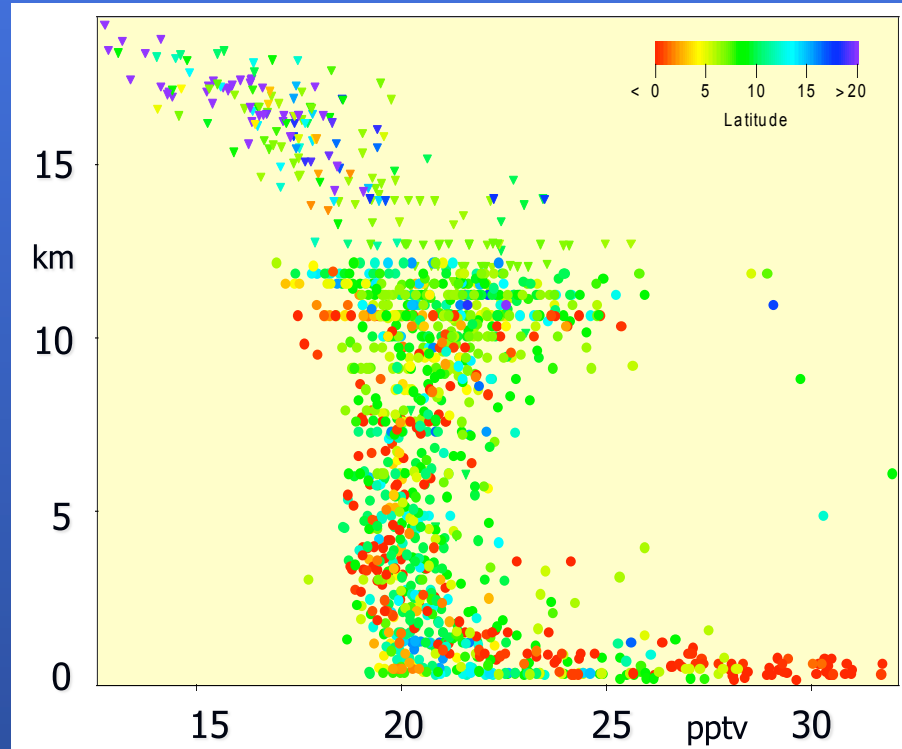
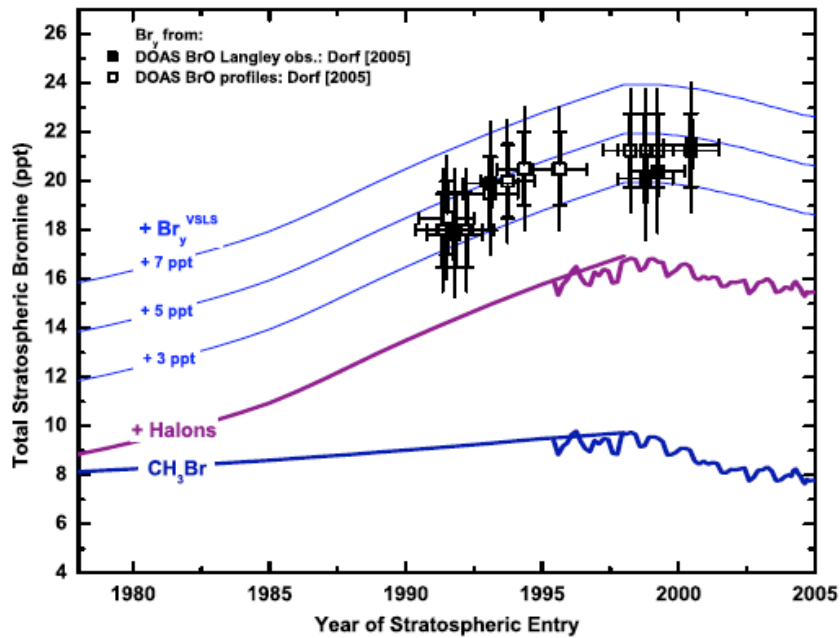
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# Comparison to WAS (NASA TC4 Campaign)



Taken from Chapter 2, VLSL, WMO, 2006

TC4, Whole Air Sampler, Schauffler, Atlas, et al., AGU, 2007

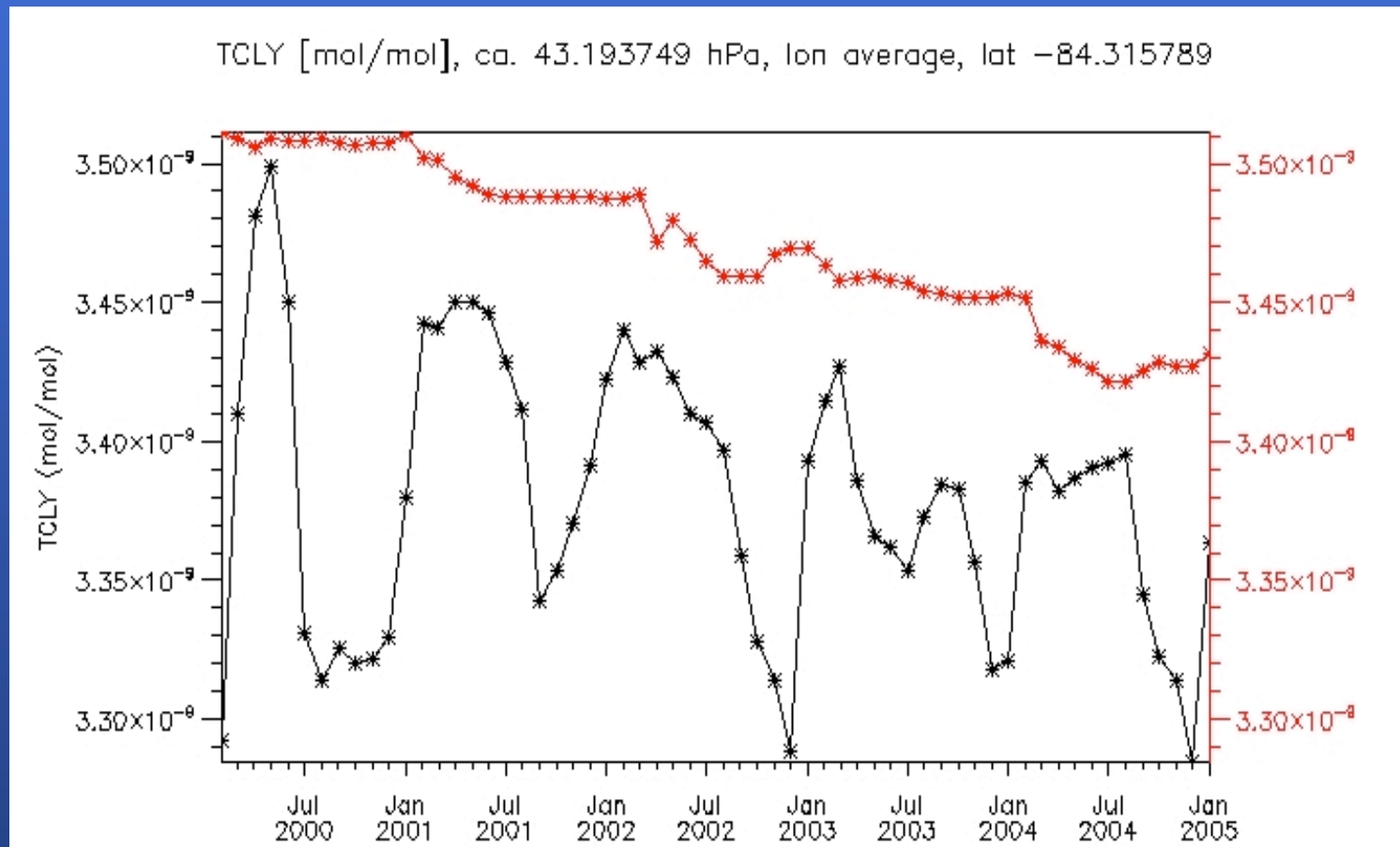
# Halogenated Very Short-lived Substances (VSLs) currently being added the WA3 UTLS mechanism

## VSL Organic Bromine Species

Source Gas	Formula	Local Lifetime (days)	Main Loss processes	WAS
Bromochloromethane	CH <sub>2</sub> BrCl	150	OH	√
Trichloromethane (chloroform)	CHCl <sub>3</sub>	150	OH	√
Methylene chloride	CH <sub>2</sub> Cl <sub>2</sub>	140	OH	√
Dibromomethane	CH <sub>2</sub> Br <sub>2</sub>	120	OH	√
Bromodichloromethane	CHBrCl <sub>2</sub>	78	OH, hv	√
Dibromochloromethane	CHBr <sub>2</sub> Cl	69	hv, OH	√
Tribromomethane (bromoform)	CHBr <sub>3</sub>	26	hv	√

# Conservation of CLOY [and BrOY].

- Conservation of inorganic chlorine:  $TCLY = org + inorganic$



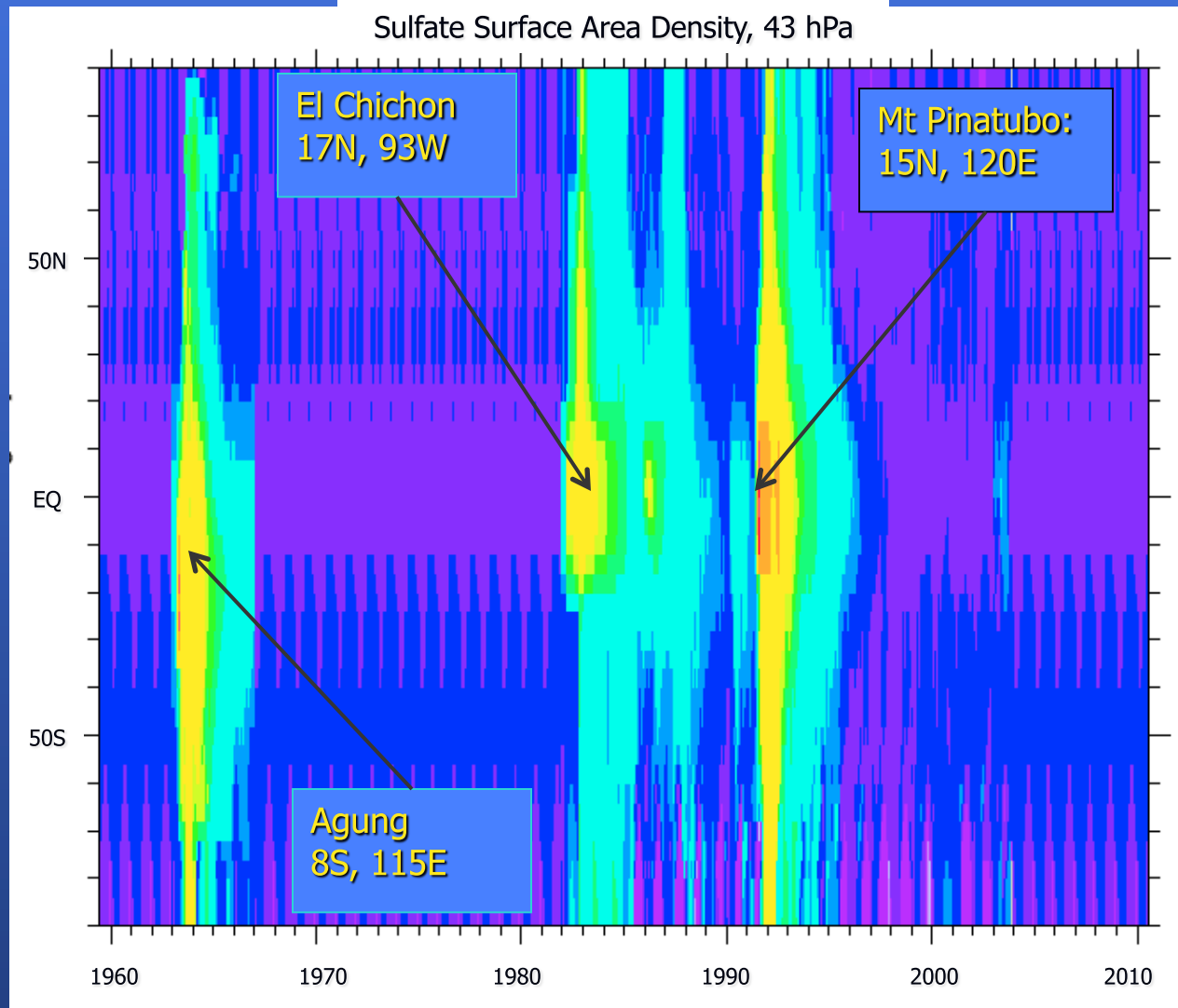
# New Sulfate SAD climatology [SPARC].

## SPARC [2006] Time Series – Modified by CCMVal

Observations used: SAGEI, SAGEII, SAM II, and SME instruments. See Thomason et al., [2007]. Range: 12-40 km and 80S-80N.

### Data Gaps:

- El Chichon period, tropical region <16km. Not filled.
- >26km gaps were filled by  $0.01 \text{ um}^2/\text{cm}^3$
- Periods: 1850-1962; 1968-1979, 2005-2100 filled with monthly mean of 1998-2002 values.
- Agung (See Dameris, 2005)

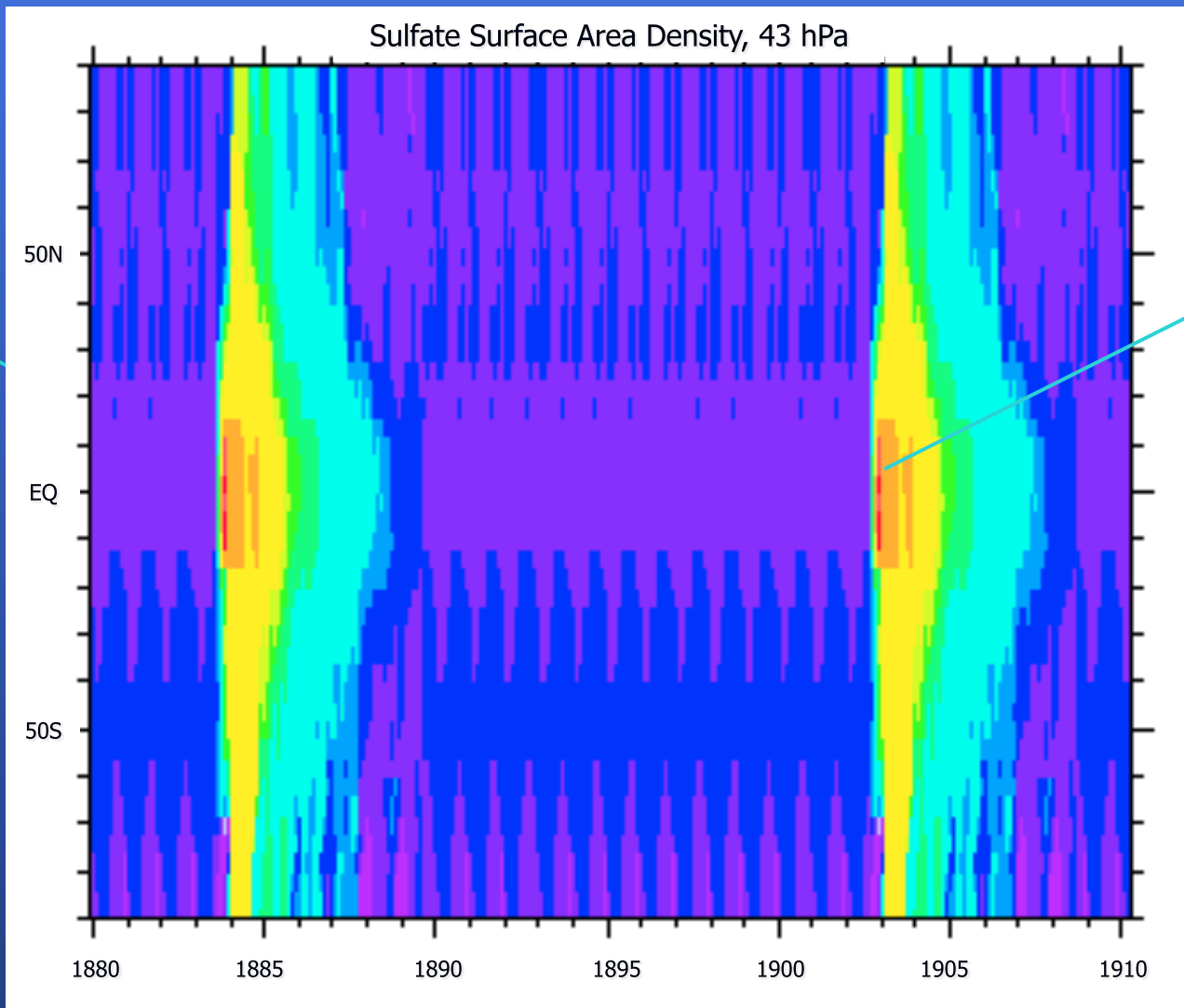




# Sulfate SAD before 1962

Used Mt Pinatubo for Krakatau and Santa Maria.

**Krakatou:**  
6S, 105E,  
27-28 Aug  
1883

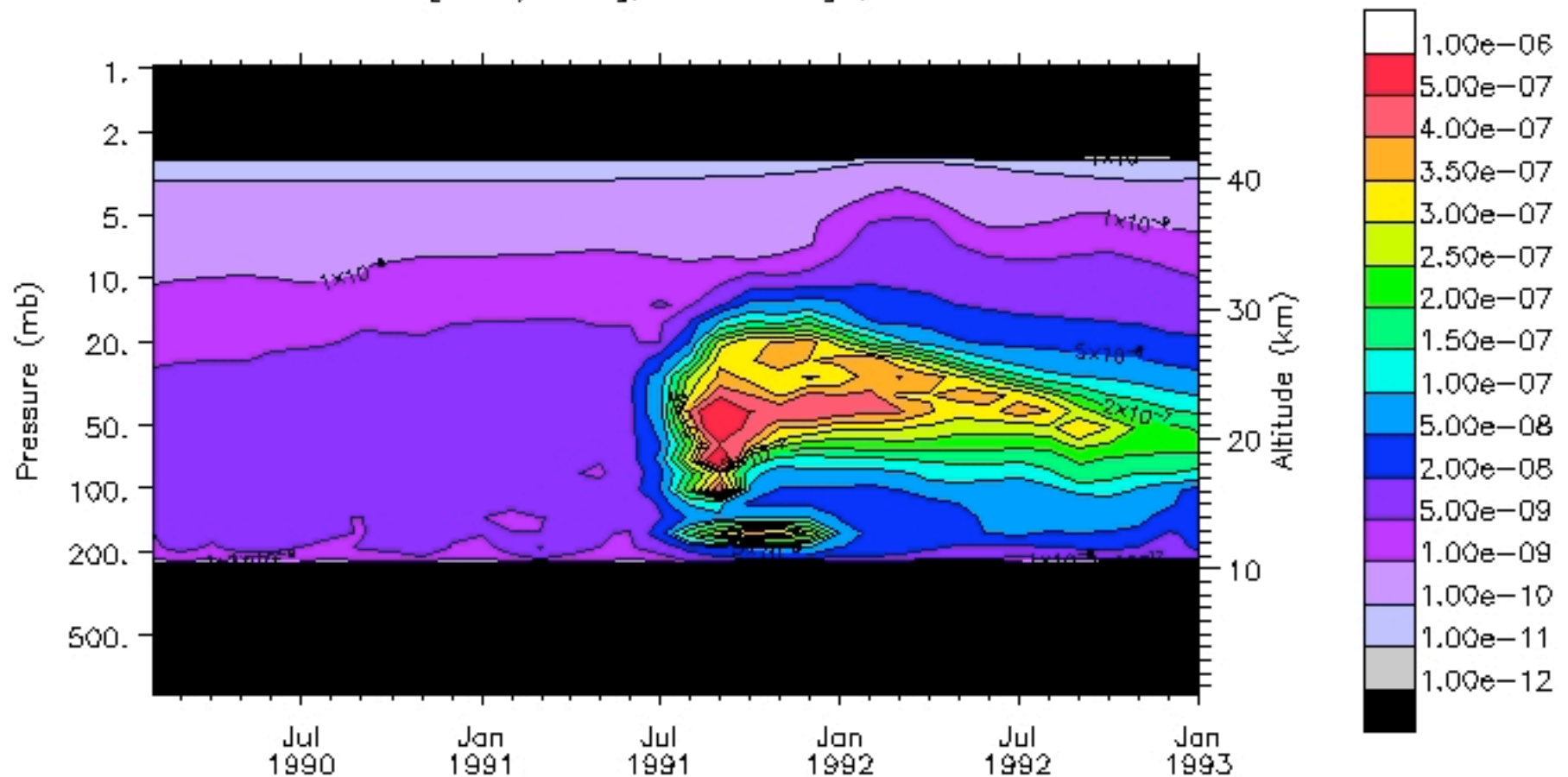


**Santa Maria:**  
14N, 91W,  
24 Oct. 1902

**Mt  
Pinatubo:**  
15N, 120E,  
15 June  
1991

# Volcanic Heating Module Added.

SAD\_SULFC [cm<sup>2</sup>/cm<sup>3</sup>], lon average, lat 2.0000000



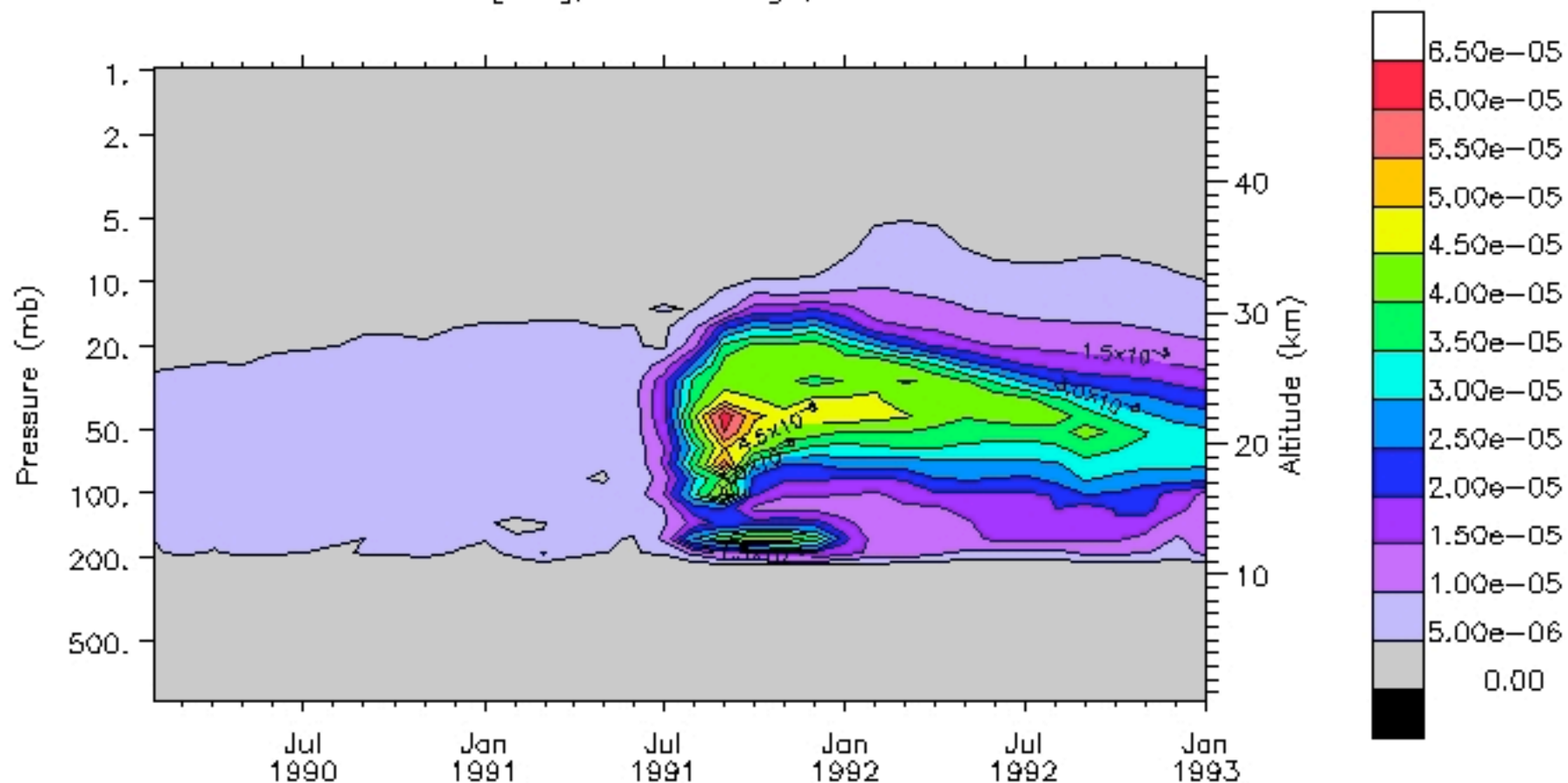
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dkln 26.09.2008 12:59



# Volcanic Heating Module Added.

RAD\_SULFC [cm], lon average, lat 2.0000000



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# Volcanic Heating Module Added.

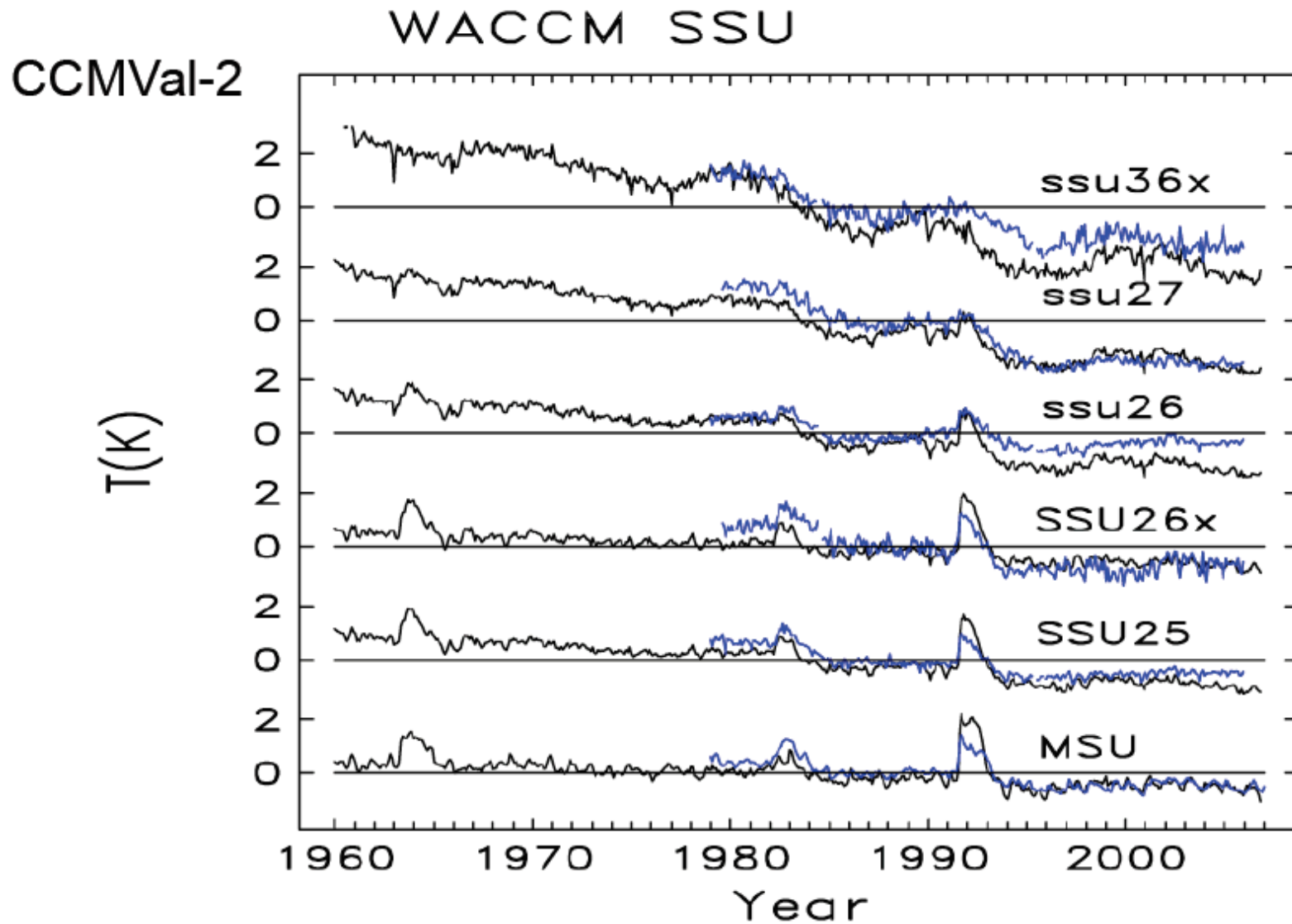
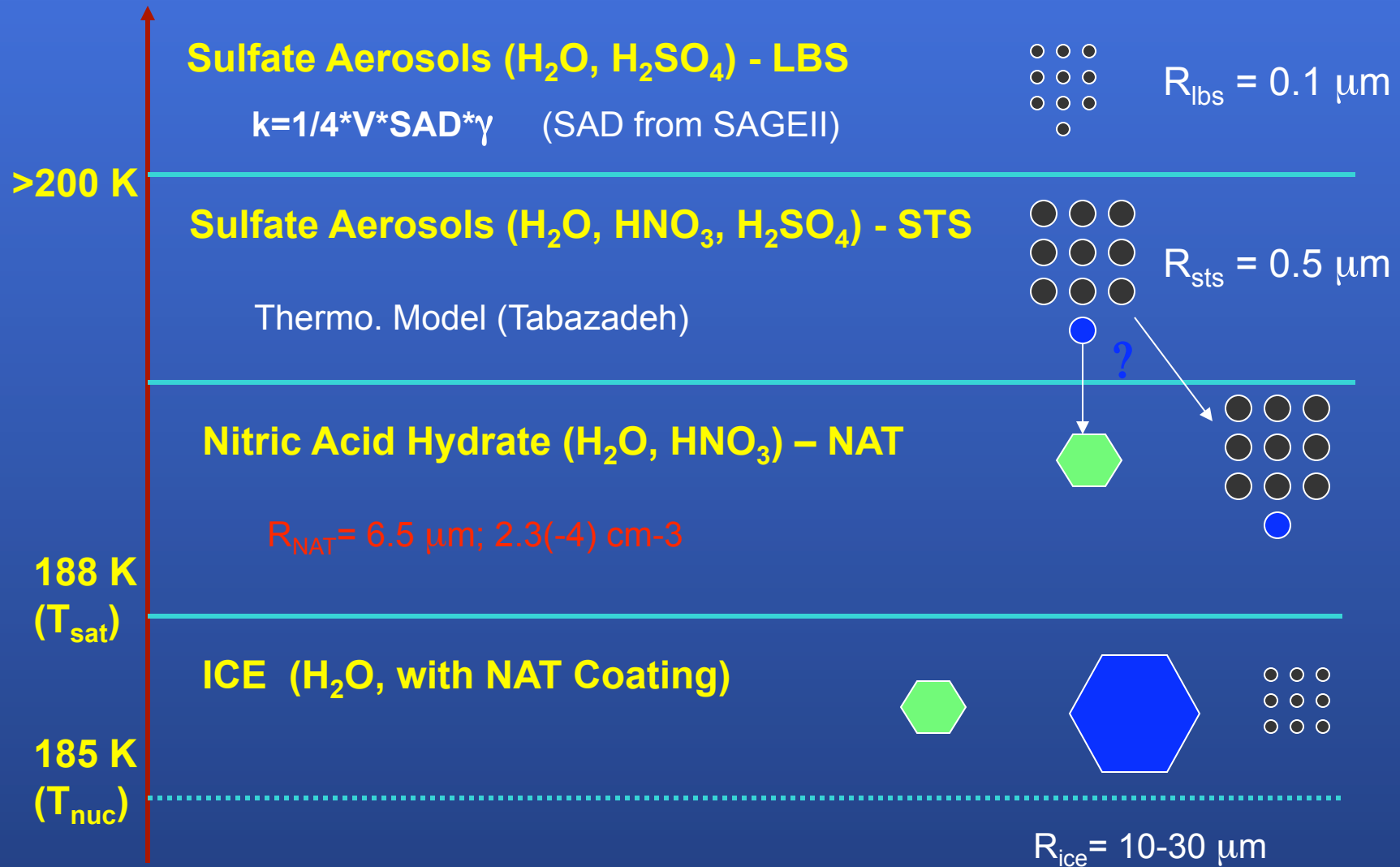
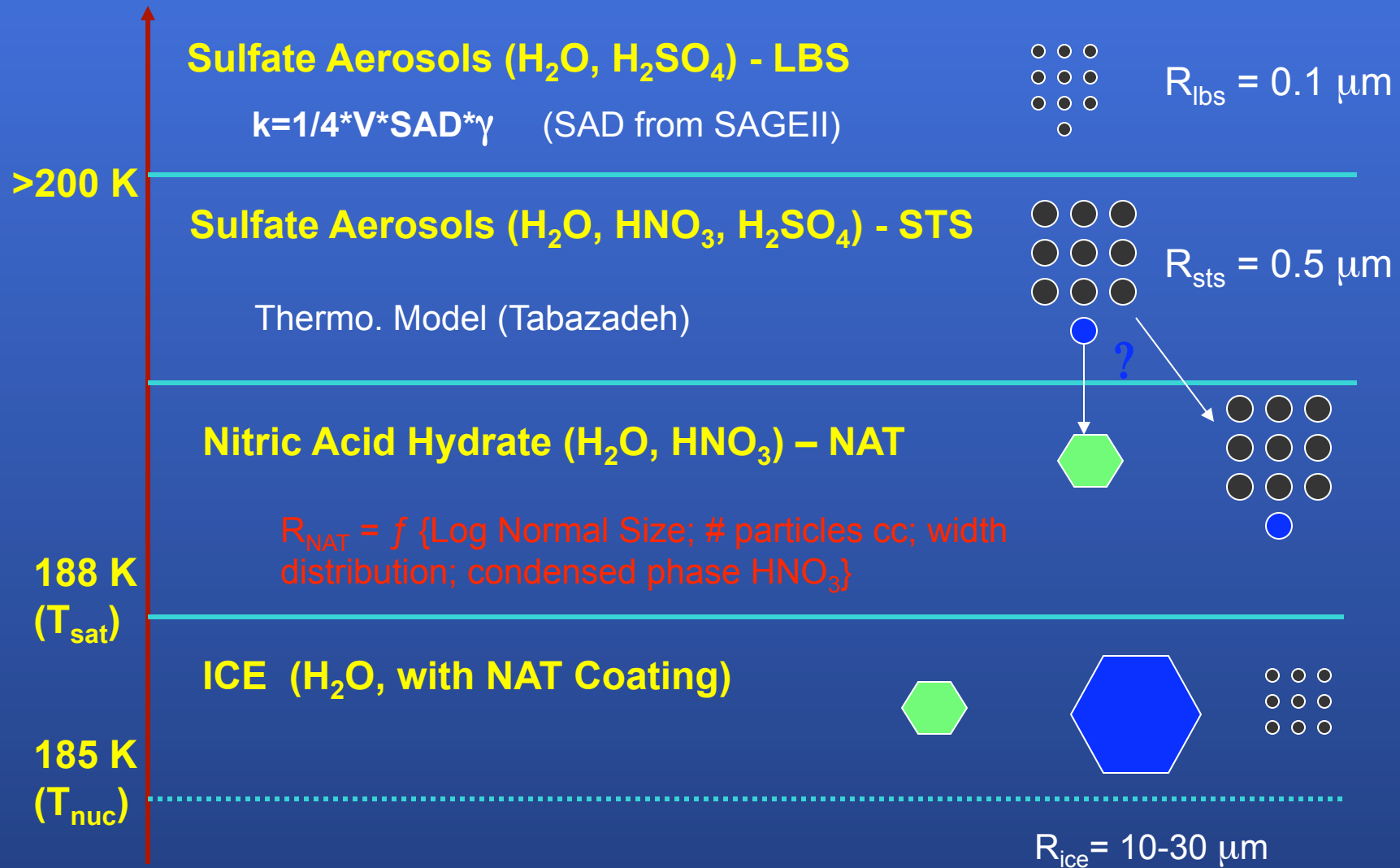


Figure courtesy Bill Randel

# New (old) NAT Approach Used

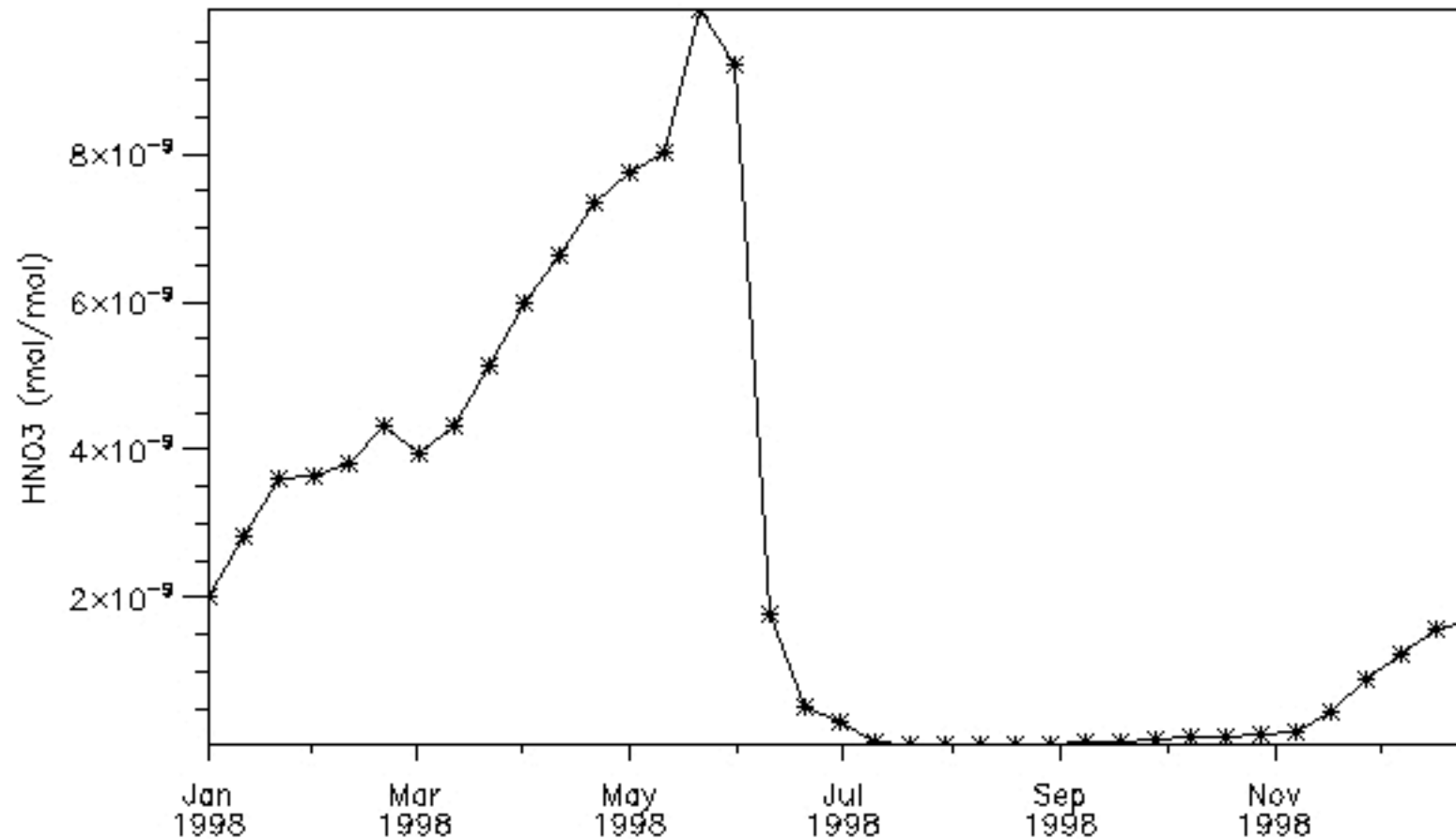


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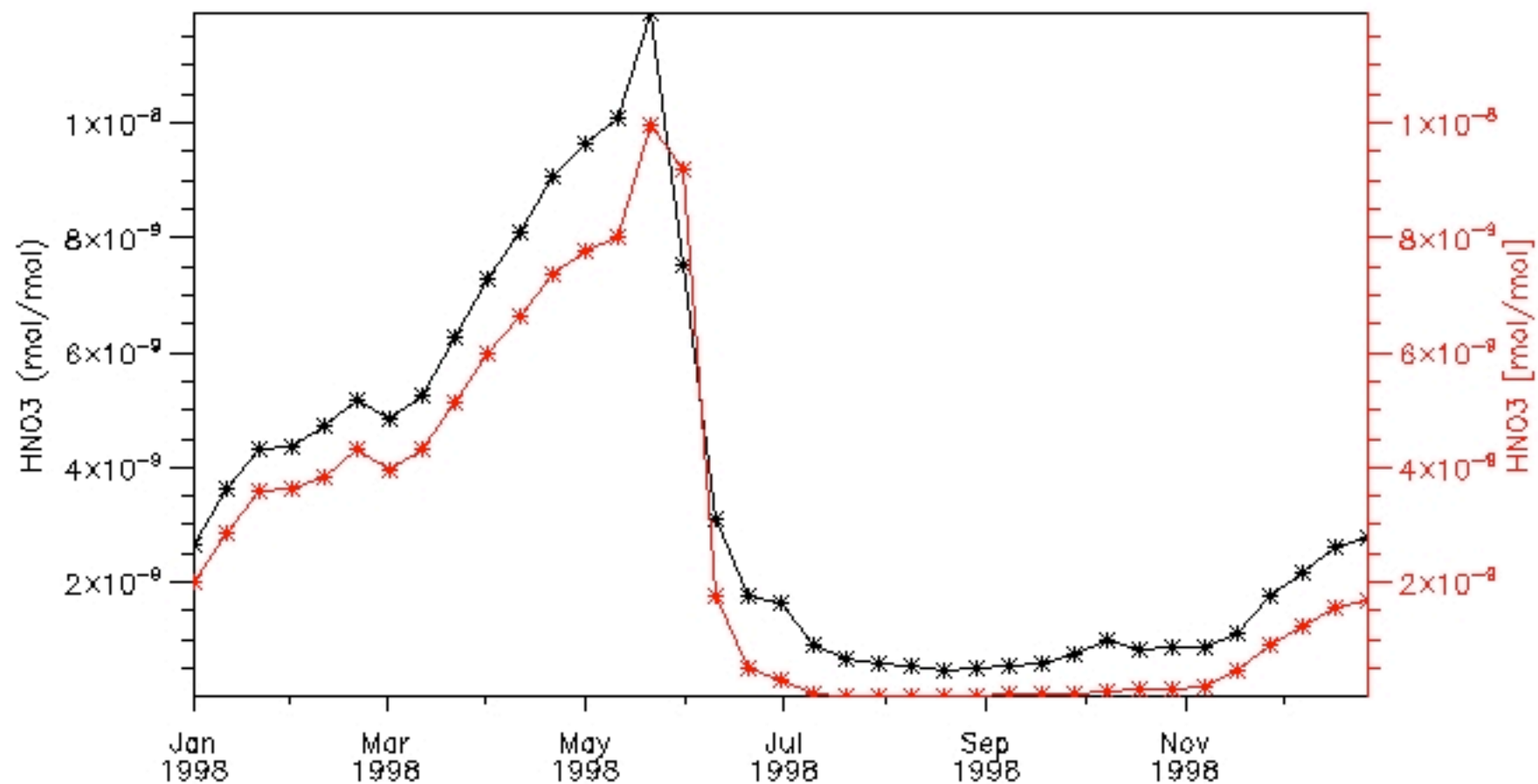
HNO<sub>3</sub> [mol/mol], ca. 51.677499 hPa, lon average, lat -78.000000





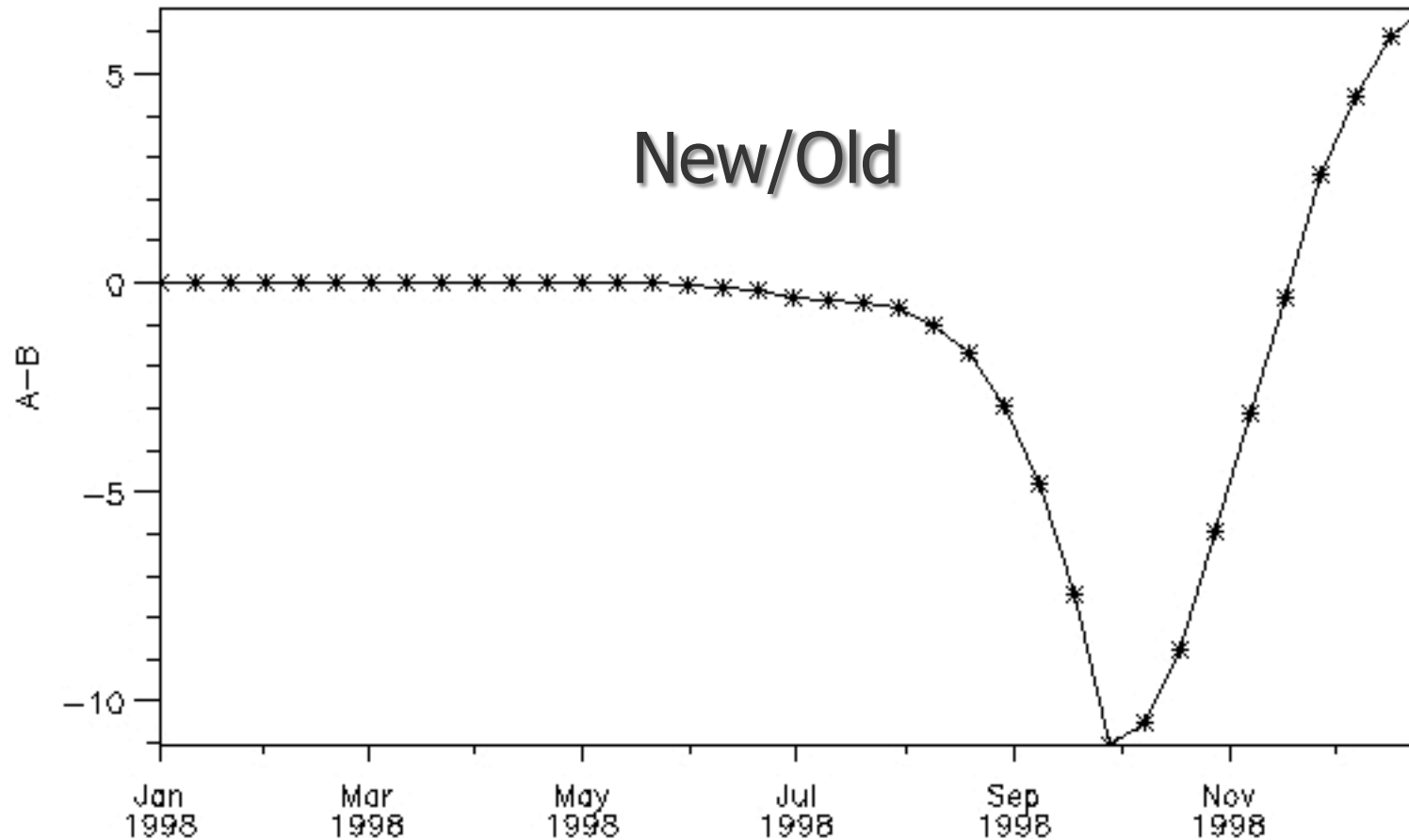
# New (old) NAT Approach Used

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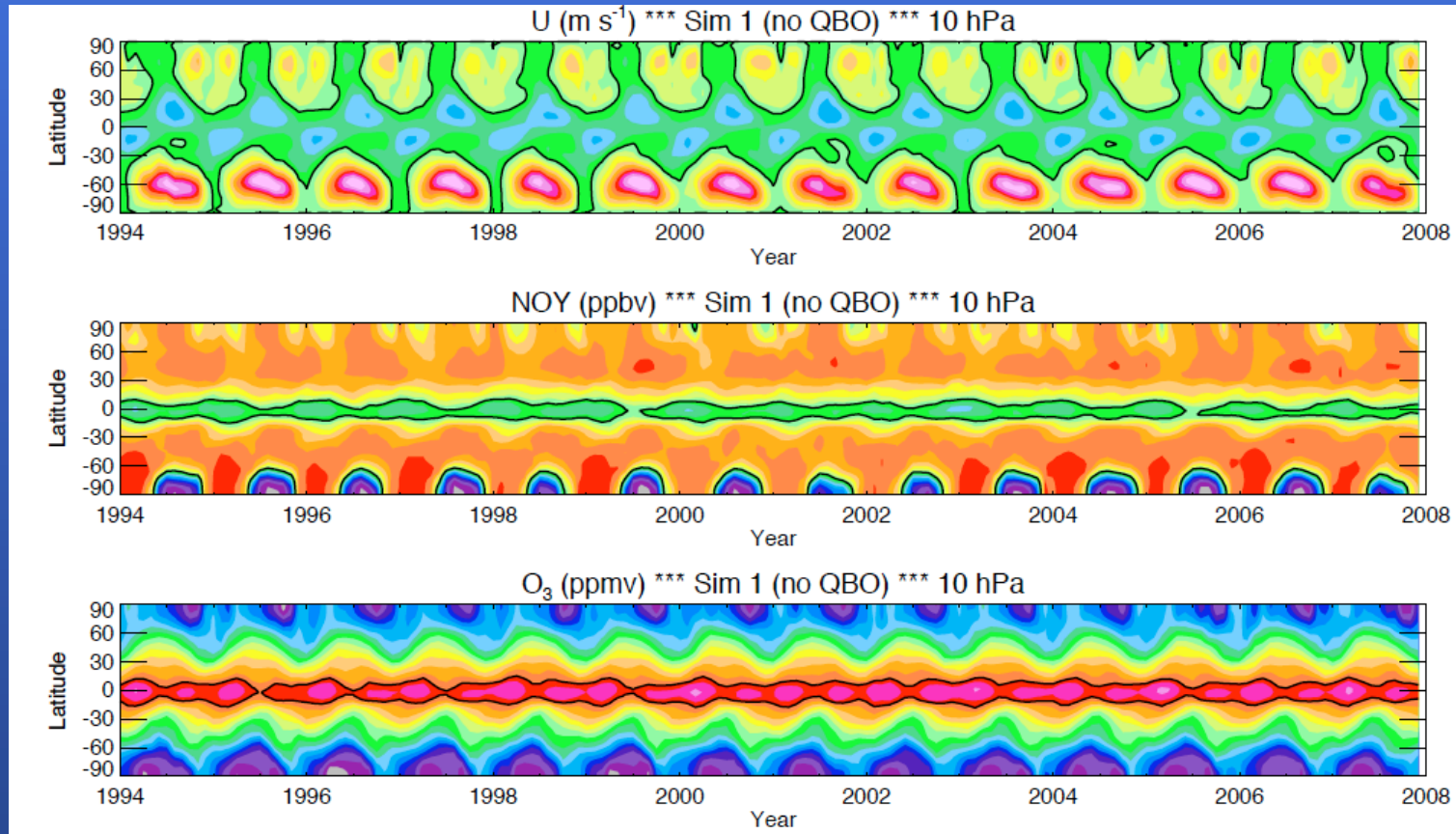


# New (old) NAT Approach Used

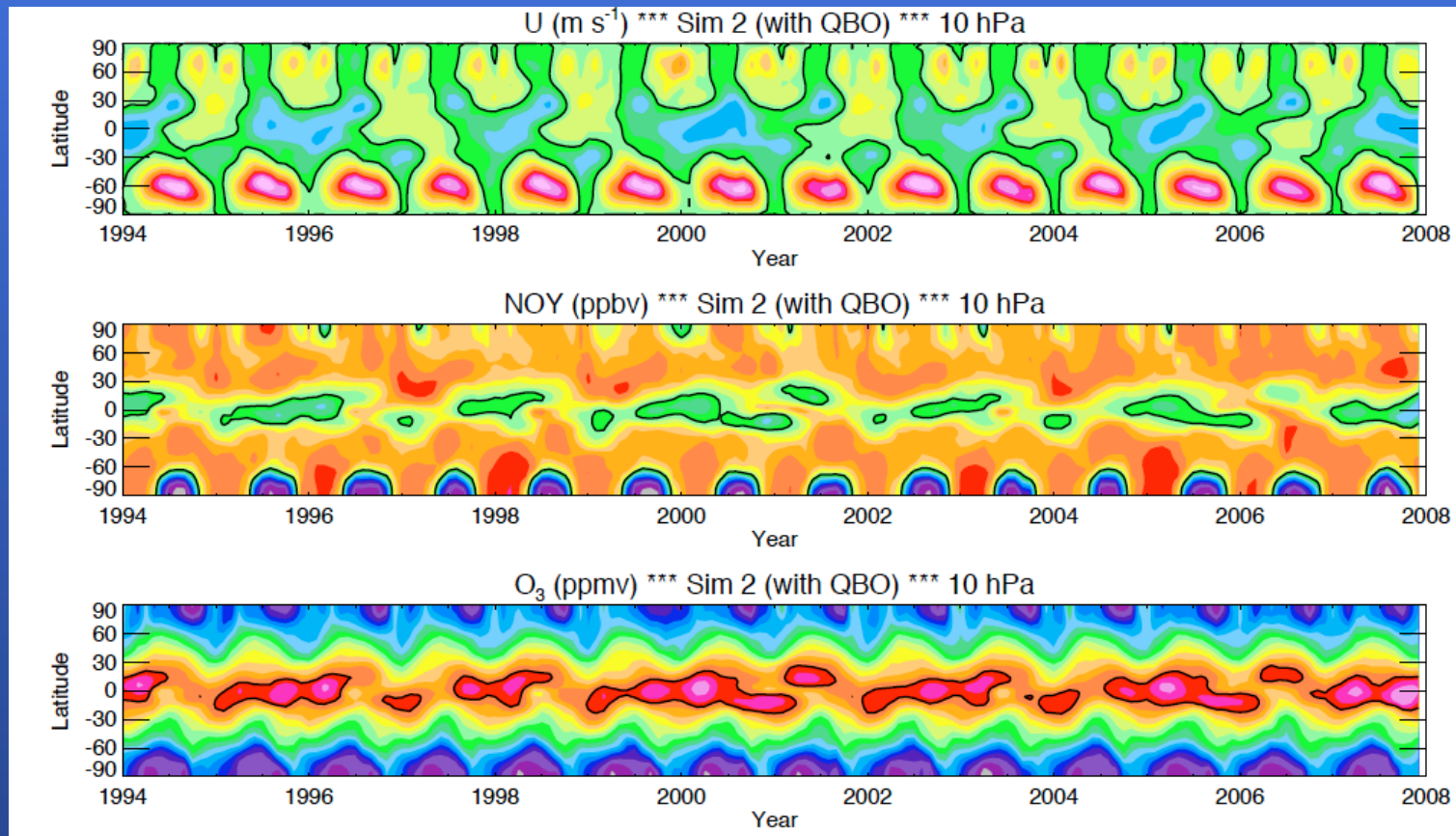
O3 Col Dens [DU], lon average, lat -78.000000



# QBO Included in 2009 CCMVal Simulations



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# Chemical Mechanism: ~60 species

- **Updated all reactions to JPL06**

- $O + O + M \Rightarrow O_2 + M$  [Smith and Robertson, 2008]
- $CO + OH \Rightarrow CO_2 + H$  [more complicated tree like expression]

- **Added:**

- $ClO + hv \Rightarrow Cl + O$
- $N + NO_2 \Rightarrow N_2O + O$  [Marsh recommendation]
- $HBr + O(1D) \Rightarrow Br + OH$
- $HBr + O \Rightarrow Br + OH$
- $HOBr + O \Rightarrow OH + BrO$
- $BrONO_2 + O \Rightarrow BrO + NO_2$
- $O(1D) + CH_3Br \Rightarrow Br$
- $O(1D) + H1211 \Rightarrow Br + Cl$
- $O(1D) + H1301 \Rightarrow Br$

- **Reactions: 159 total; 128 (gas-phase); 17 (het) ; 14 (ion / e)**

- **Photolysis: 73 total; 48 (120-750 nm); 25 (EUV)**

# Chemical Mechanism: 125 species

## Additional Surface Source Gases (16 additional) ...

### NHMCs:

CH<sub>3</sub>OH,  
C<sub>2</sub>H<sub>6</sub>, C<sub>2</sub>H<sub>4</sub>, C<sub>2</sub>H<sub>5</sub>OH, CH<sub>3</sub>CHO  
C<sub>3</sub>H<sub>8</sub>, C<sub>3</sub>H<sub>6</sub>, CH<sub>3</sub>COCH<sub>3</sub> (Acetone)  
C<sub>4</sub>H<sub>8</sub> (BIGENE), C<sub>4</sub>H<sub>8</sub>O (MEK)  
C<sub>5</sub>H<sub>8</sub> (Isoprene), C<sub>5</sub>H<sub>12</sub> (BIGALK)  
C<sub>7</sub>H<sub>8</sub> (Toluene)  
C<sub>10</sub>H<sub>16</sub> (Terpenes)

### Plus:

HCN, CH<sub>3</sub>CN

### Reactions:

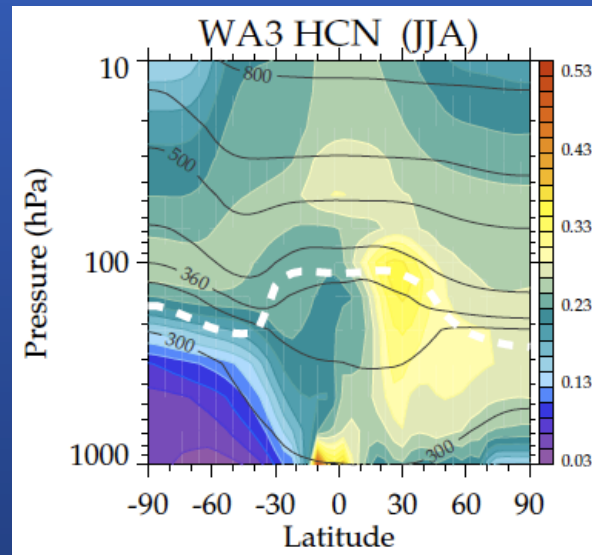
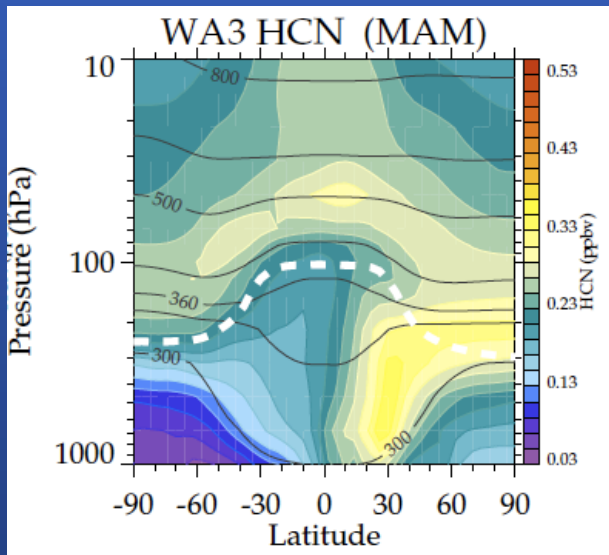
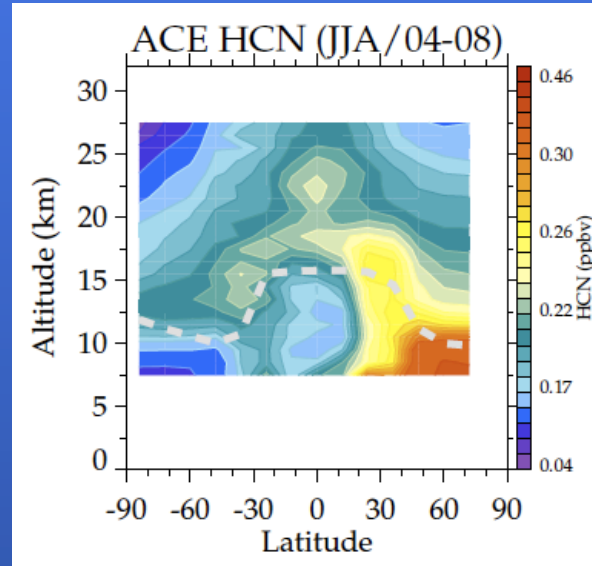
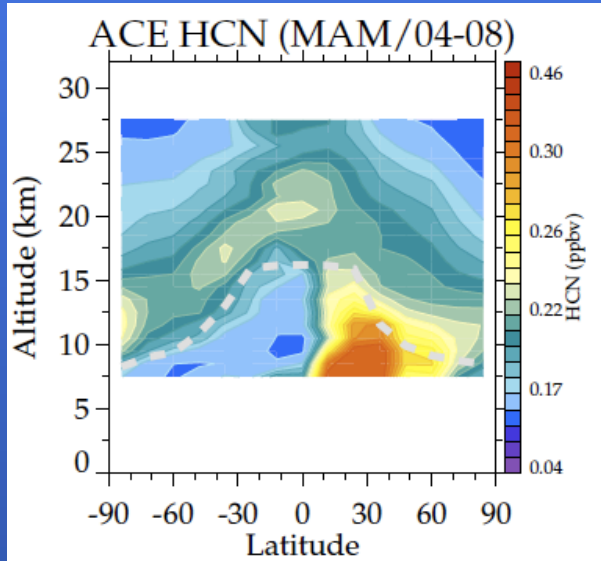
**178 total:** 147 (gas-phase); 17 (het); 14 (ions/  
electron)

### Photolysis:

**97 total:** 72 (120-750nm); 25 (EUV)

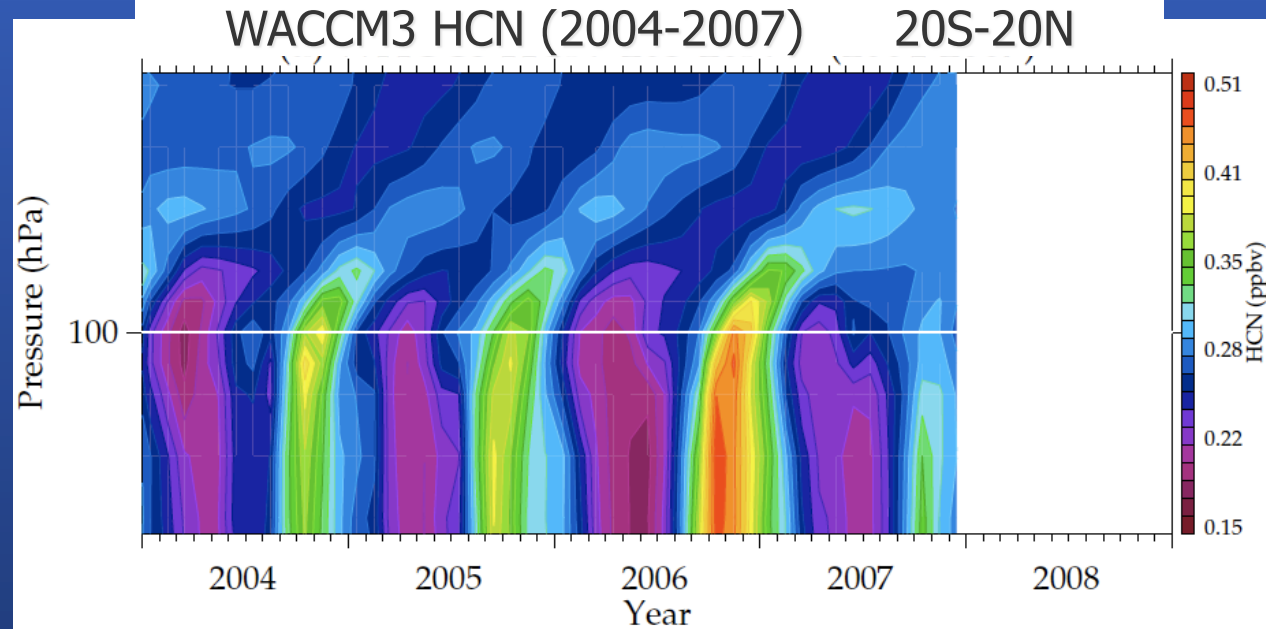
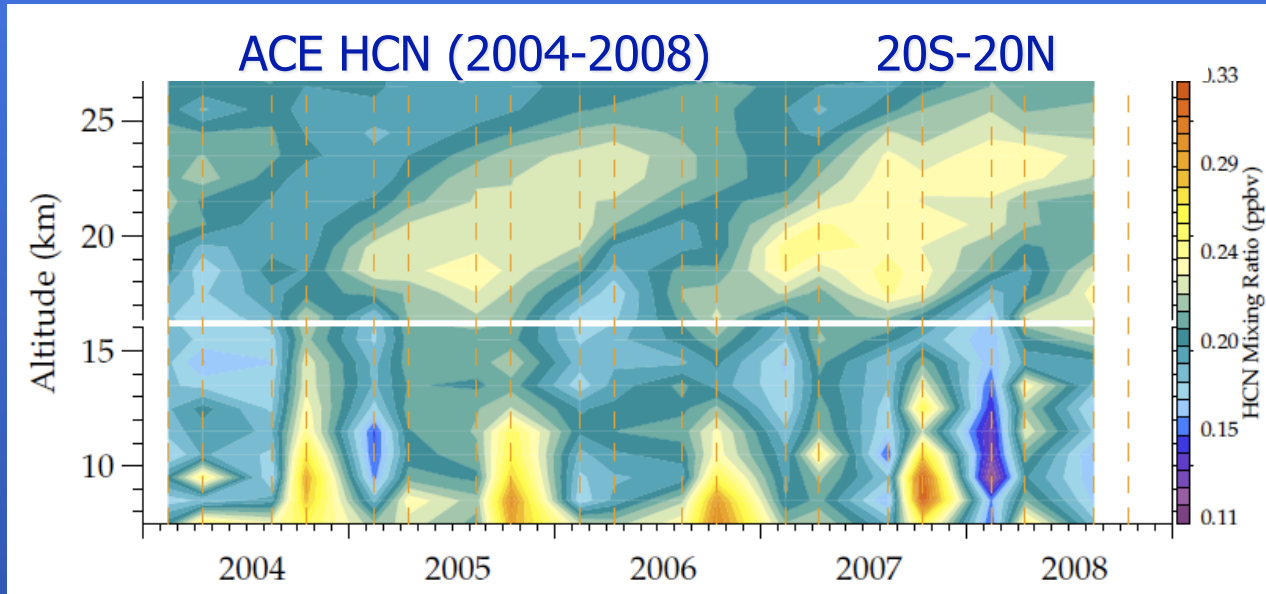
This version has been checked in: /wacm/fvitt/wacm08\_cam3\_5\_48\_b1.01

# HCN – Fully Interactive WACCM3 vs ACE-FTS



Courtesy of Mijeong Park

# HCN – Fully Interactive WACCM3 vs ACE-FTS



# New ETF's

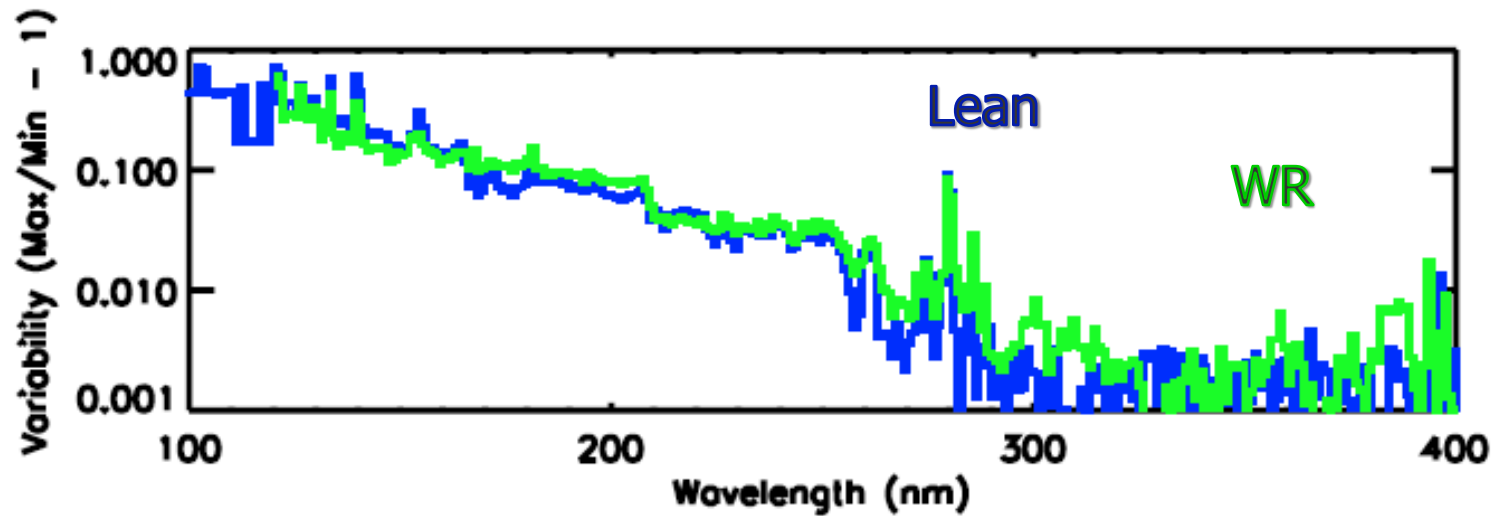
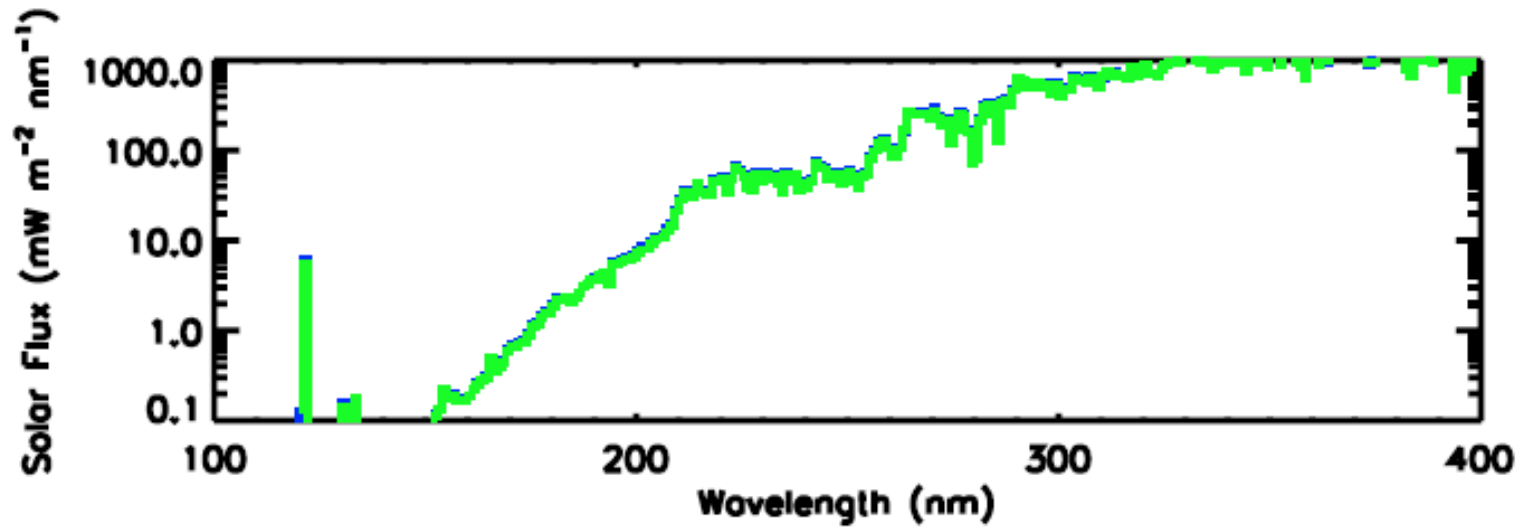
- **Changed to Lean Extraterrestrial Fluxes**

- Previous: Woods / Rottman [120 – 350 nm]
- Neckel [>350nm no solar variability]

- **Why?**

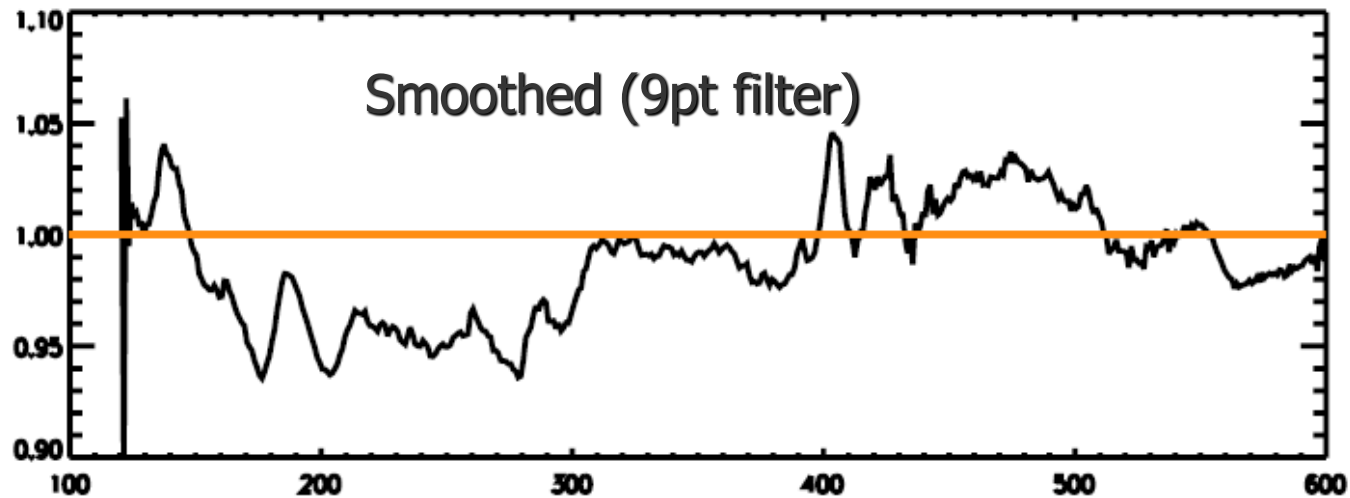
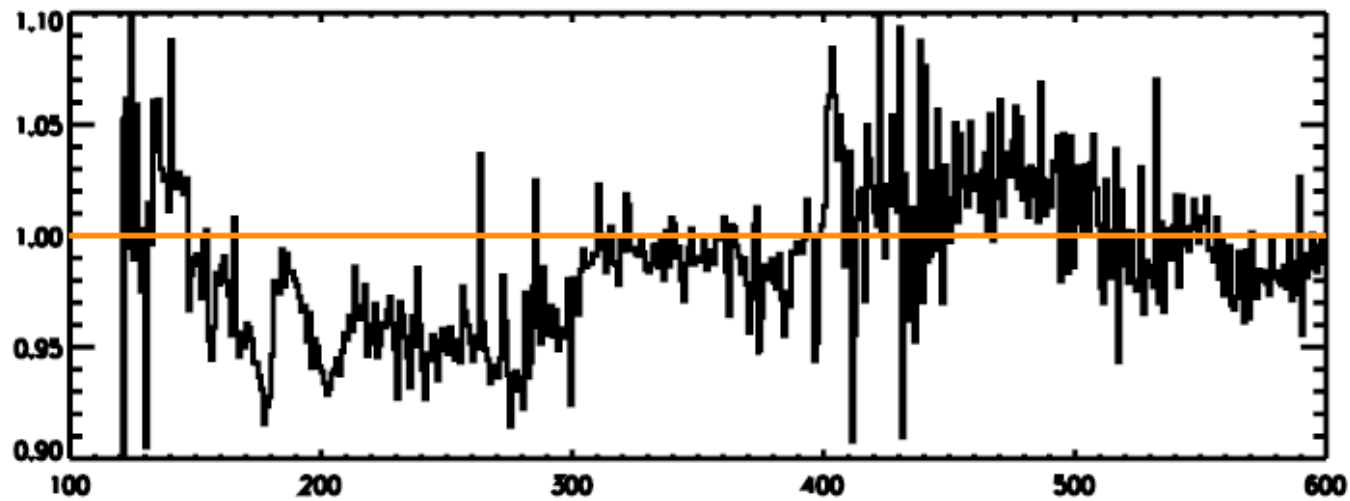
- Recommended by CCMVal-2
- Variability Extends long ward of 350 nm.
- Have wavelength dependent variability for pre-industrial simulations.

# Max to Min variability: Lean vs WR



Courtesy of Stan Solomon

# Ratio at low Solar Activity: Lean / WRN



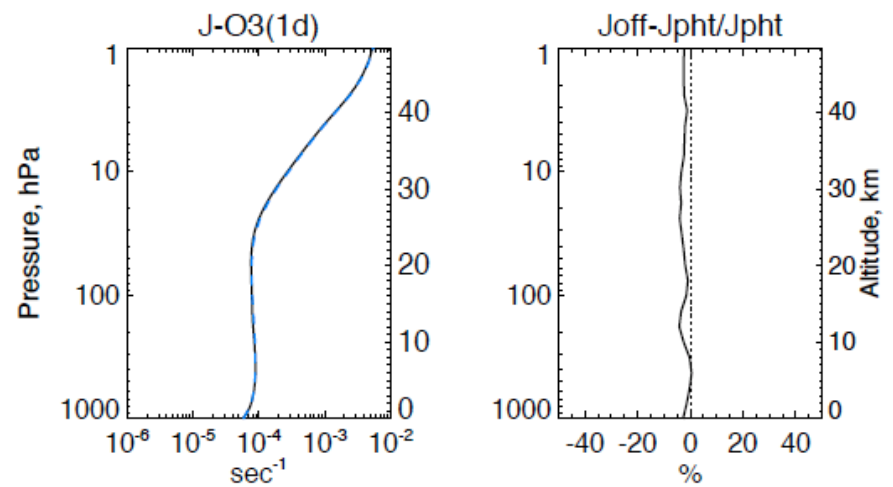
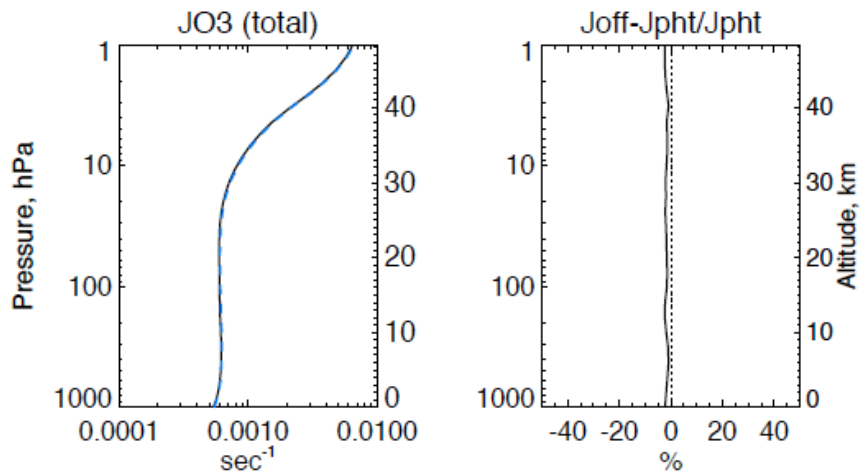
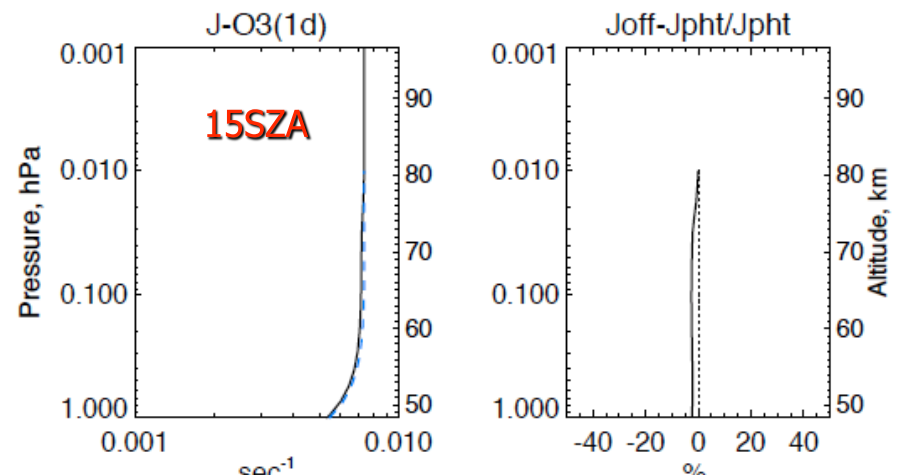
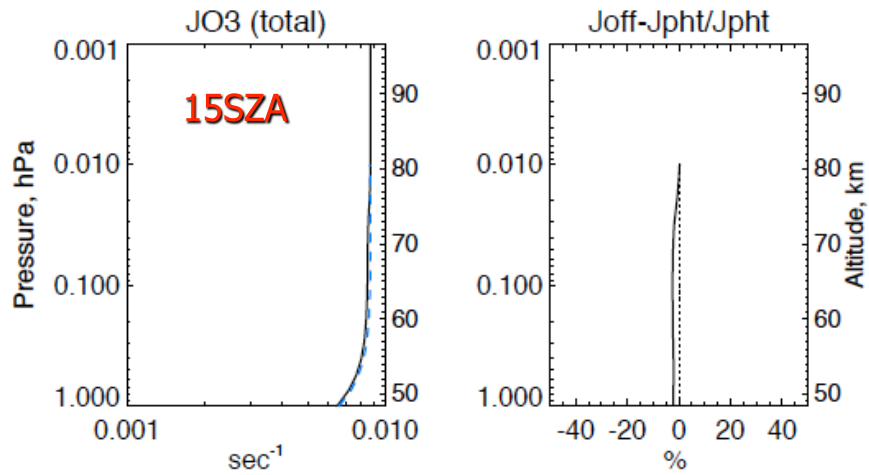
Courtesy of Stan Solomon



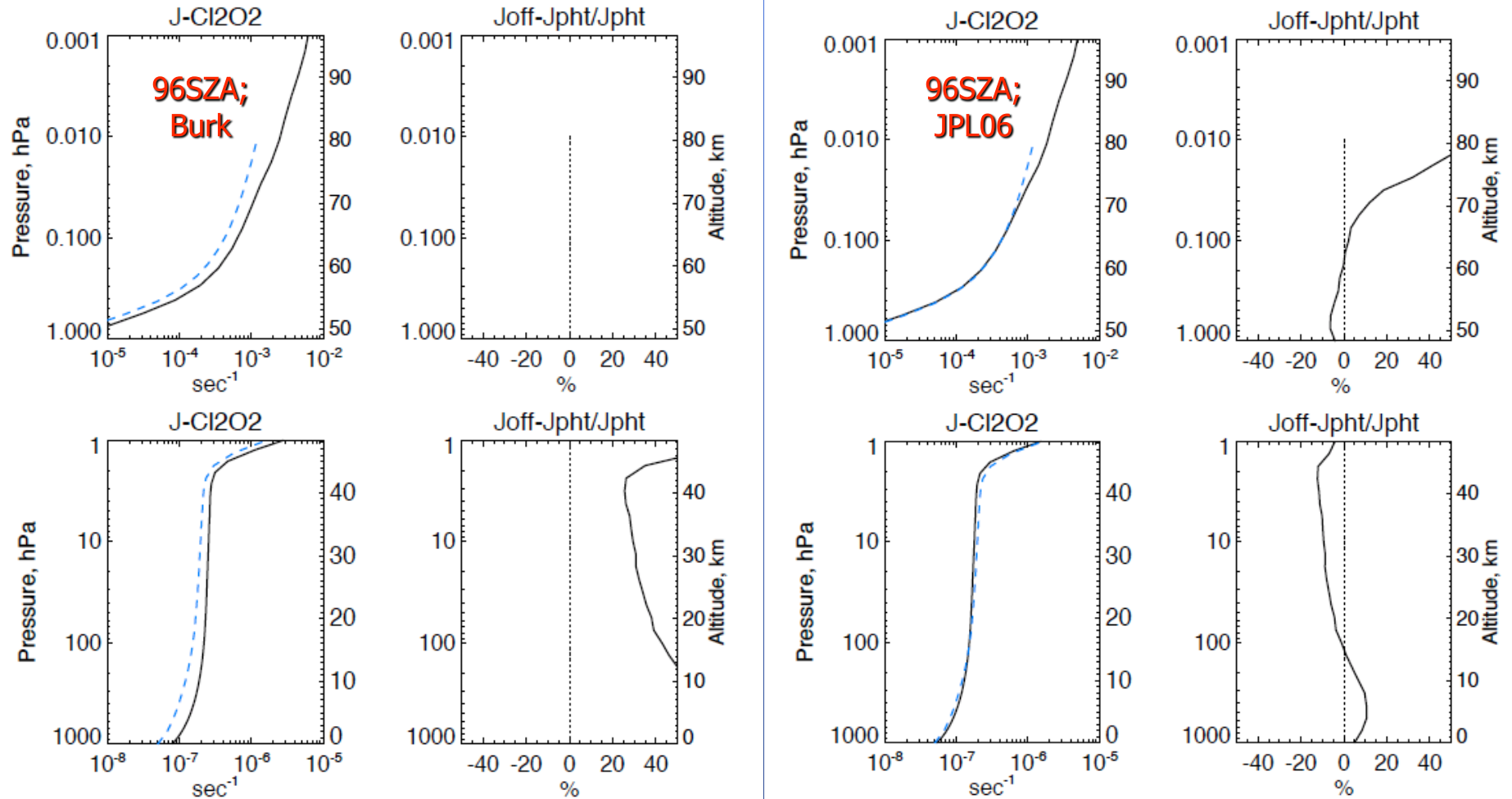
# Photolysis

- **Updated the preprocessor – tagging each photolysis reaction**
  - [jno2]  $\text{NO}_2 + h\nu \rightarrow \text{NO} + \text{O}$ .
  - The current LUT has >95 J's in the master list [code only reads what it needs].
- **Created an “offline” photolysis module that can calculate all J's given ~ 12 constituents [e.g., O<sub>3</sub>, T, CLOUD, ASDIR, etc...].**
- **Ran the “offline” photolysis model for a CCMVal Photocomp.**
  - Clear Sky J's at 15°, 84°, 96° SZA
  - 24-hr avg J's at Equinox, 84N.
  - 15° SZA without Rayleigh Scattering

# Photolysis



# Photolysis



## **SD-WACCM3**

- **Used extensively in testing updated chemistry module changes.**
- **Support ACD lead Aircraft Campaigns [e.g., START08] and Satellite program [e.g., HIRDLS].**
- **Supports ACD UTLS Science Studies [Randel's group].**
- **Will be made available to the University Community in the next WACCM release.**

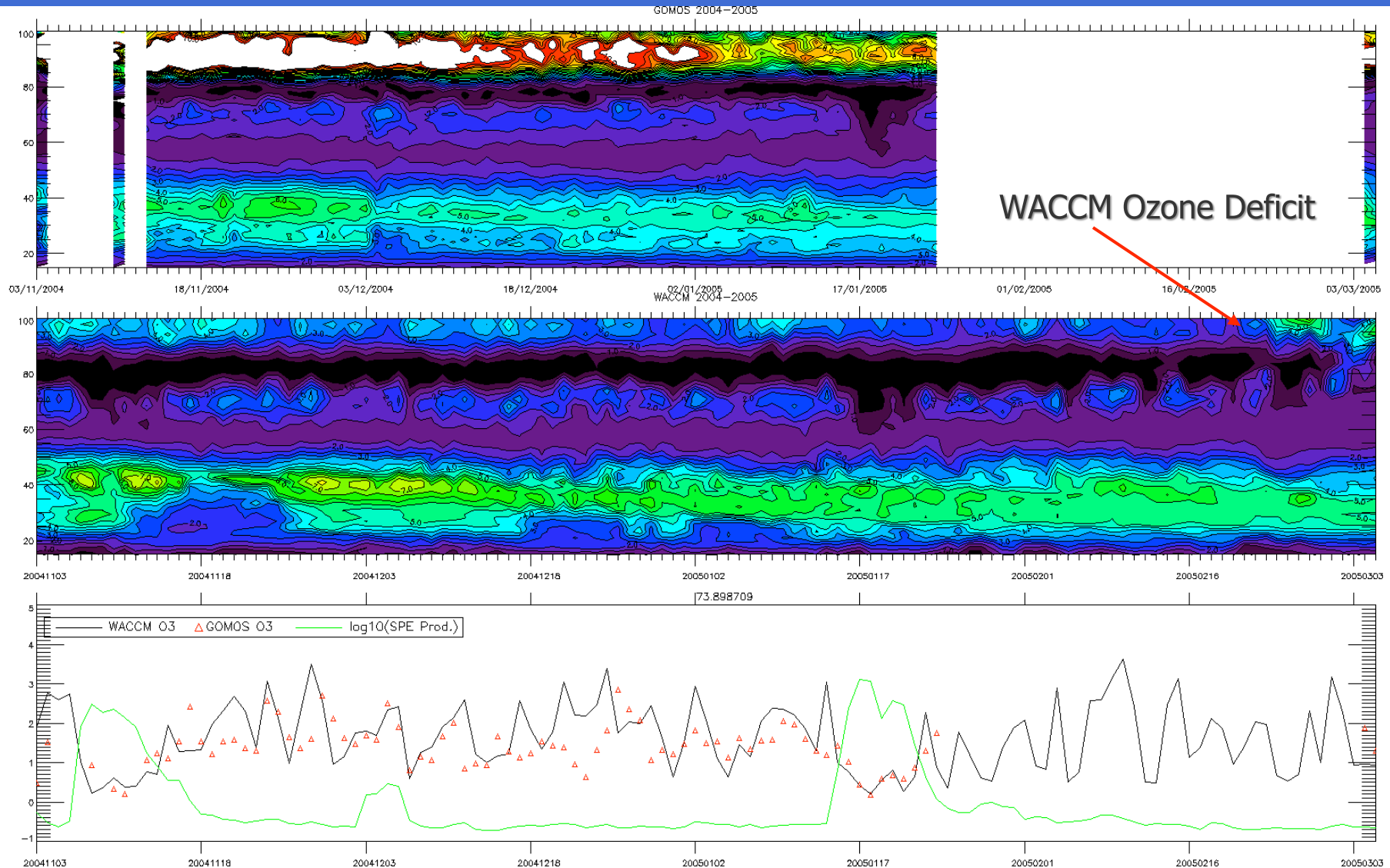
# SD-WACCM3

- **Logic existed to nudge CAM3 [WACCM3] met fields with offline met fields every time step [Hess, Rasch, Vitt].**
- **Created a procedure to convert GEOS5.1 met fields from 0.5x0.66 horiz. resolution to 1.9x2.5 (or 4x5). [Yudin, Kinnison, Tilmes]**
- **Transition from nudged met fields to fully interactive dynamics between a specified altitude region was added for WACCM.**
- **Current Studies...**
  - CU/LASP Randall et al. [O<sub>3</sub> loss]
  - Yudin et al. , [Lamina Statistics with HIRDLS]
  - Tilmes et al., [START08 / ARCTAS].
  - Marsh et al., [MLT Ozone / SPE / GOMOS].

# SD-WACCM3 \*\*\* with SPE's

## Comparison with GOMOS data (Viktoria Sofieva)

- GOMOS is stellar occultation and so these are nighttime ozone values.
- Average of data from 70 to 76 N



# CAM-CHEM / WACCM3 Merge (>CAM3.6)

Expertise lies in both group for common modules:

- Wet and dry deposition is more advanced in CAM-CHEM
- Emission databases are evaluated in the CAM-CHEM community.
- Lower boundary condition modules are used by both teams.
- Tropospheric mechanisms have been developed and used by the CAM-CHEM team. Stratospheric and MLT mechanism have been developed and used by the WACCM team.
- LUT/parameterization approaches have been developed by the WACCM team and used by CAM-CHEM team.
- Heterogeneous processes on stratospheric aerosols have been developed by the WACCM team and used by CAM-CHEM team.

Maintaining one chemistry pre-processor and common code would increase efficiency for both groups .



# Just Remember: Chemistry is Really Important!

JCl<sub>2</sub>O<sub>2</sub> (Burkholder et al., 1990)

82S \*\*\* 43hPa

