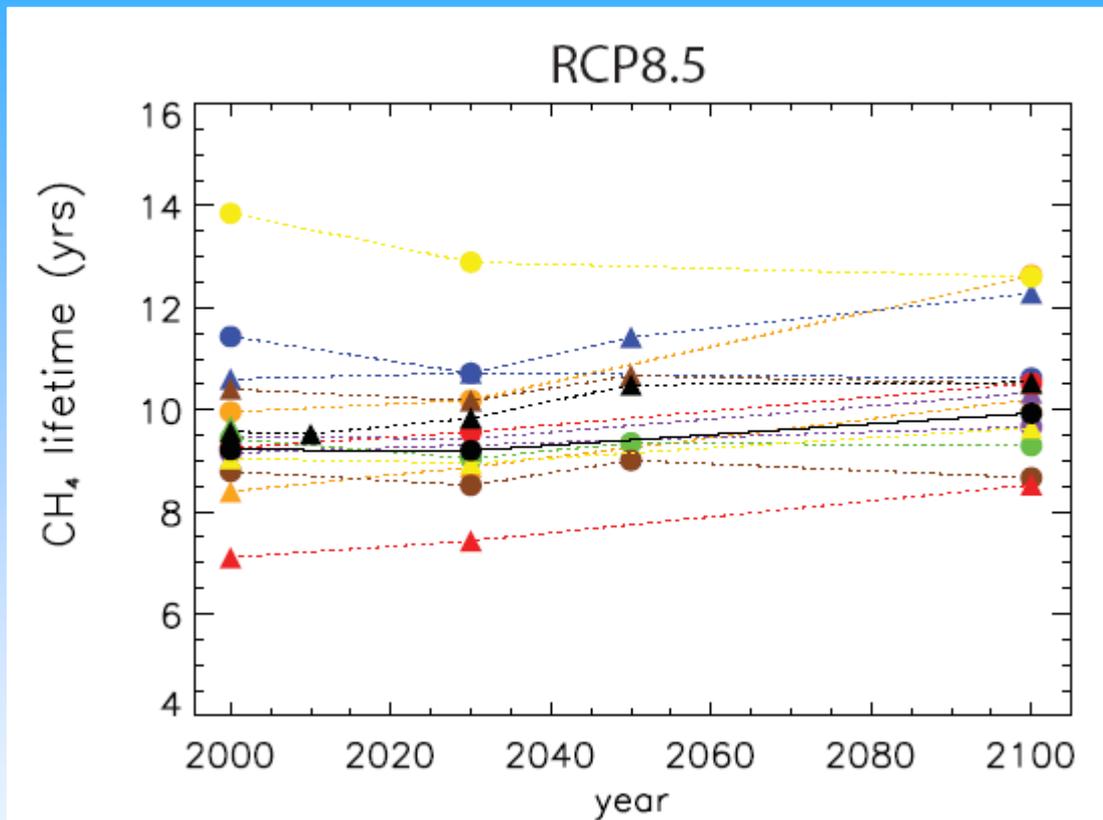


Evaluating Tropospheric OH in Global Models: The Roles of CONTRAST and CCMI

Julie Nicely
University of Maryland
CONTRAST Science Team Meeting
Boulder, CO

21 October 2013

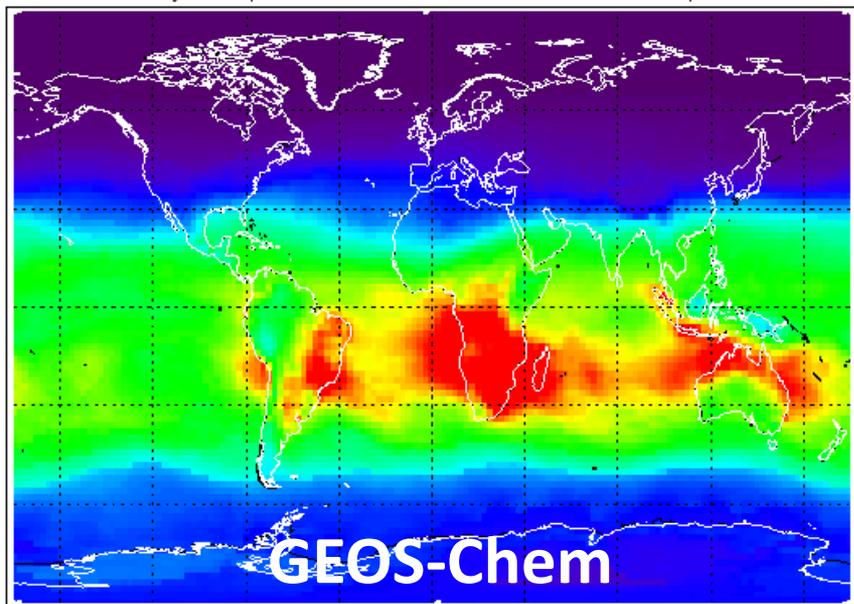
CH₄ lifetimes differ by ±20%



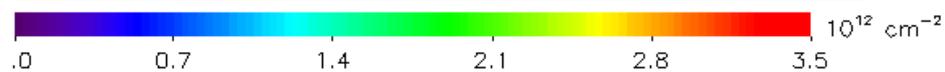
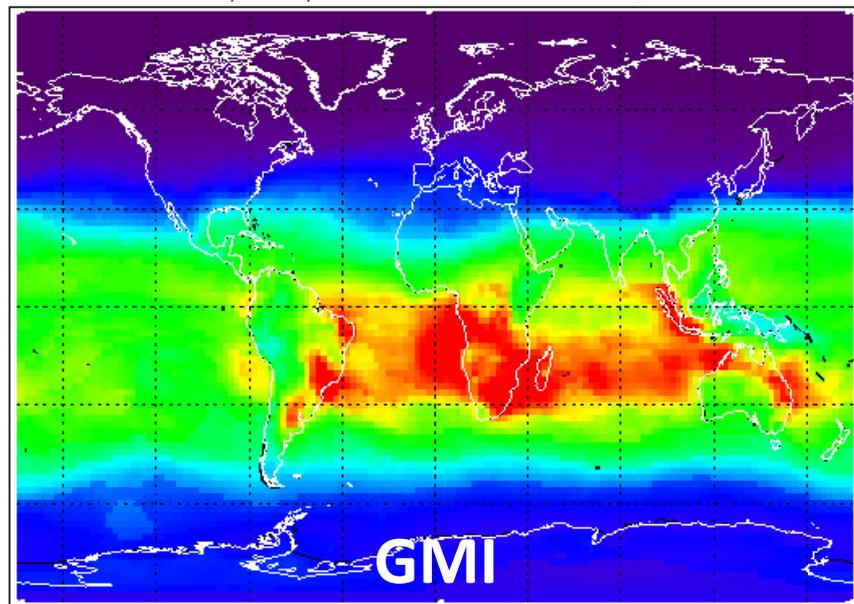
- | | | |
|--------------------|-------------|-------------|
| CESM-CAM-superfast | MIROC-CHEM | LMDzORINCA |
| CICERO-OsloCTM2 | MOCAGE | MIROC-CHEM |
| CMAM | NCAR-CAM3.5 | MOCAGE |
| EMAC | STOC-HadAM3 | NCAR-CAM3.5 |
| GEOSCCM | UM-CAM | STOC-HadAM3 |
| GFDL-AM3 | ACCMIP_mean | UM-CAM |
| GISS-E2-R | MAGICC | ACCMIP_mean |
| HadGEM2 | | MAGICC |

Voulgarakis et al., ACP, 2013

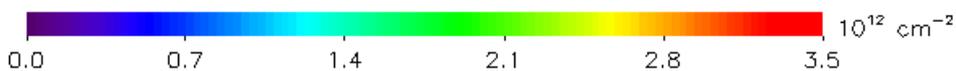
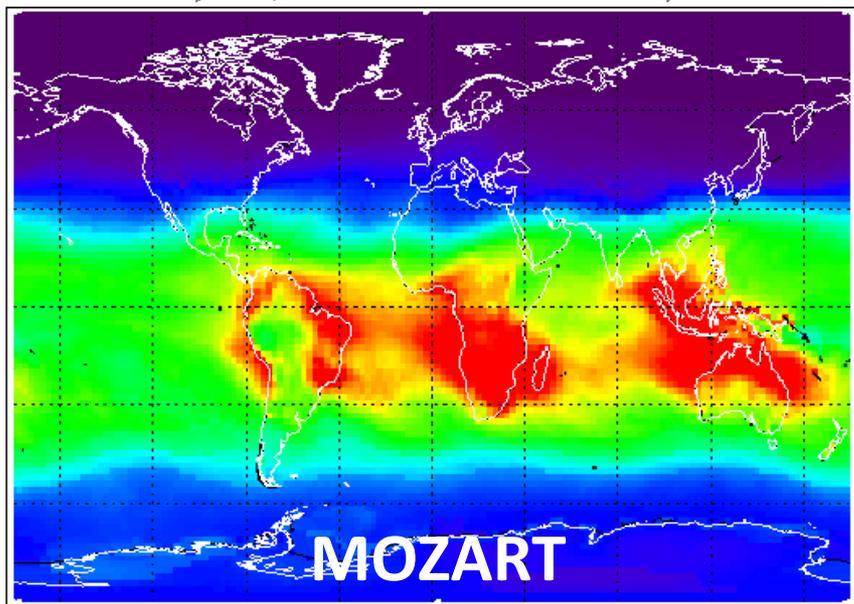
Monthly Trop OH Column, GEOSCHEM, 01/2008



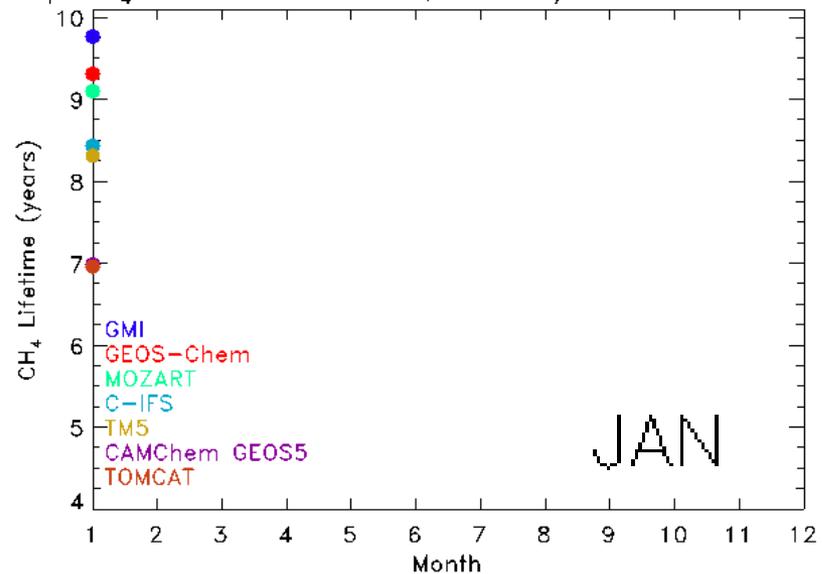
Monthly Trop OH Column, GMI, 01/2008

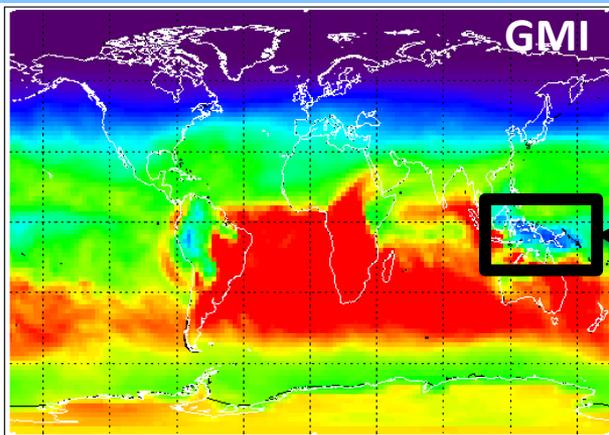
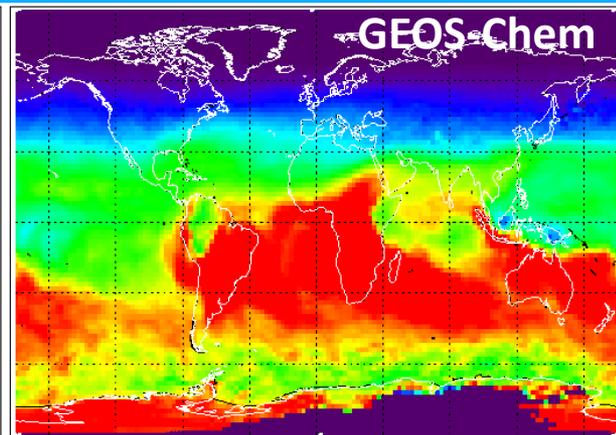
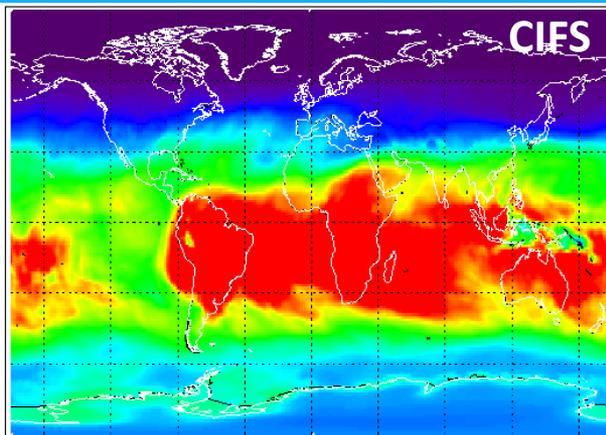
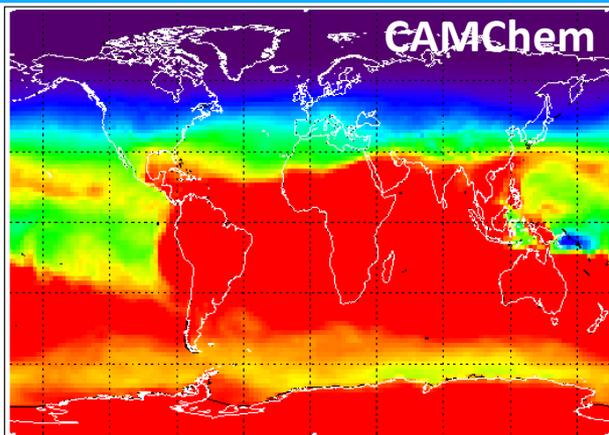


Monthly Trop OH Column, MOZART, 01/2008

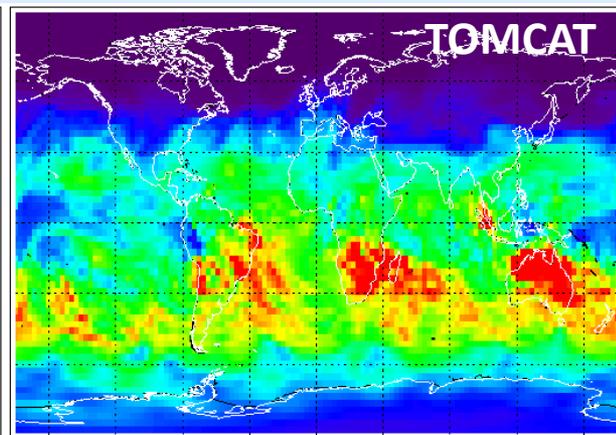
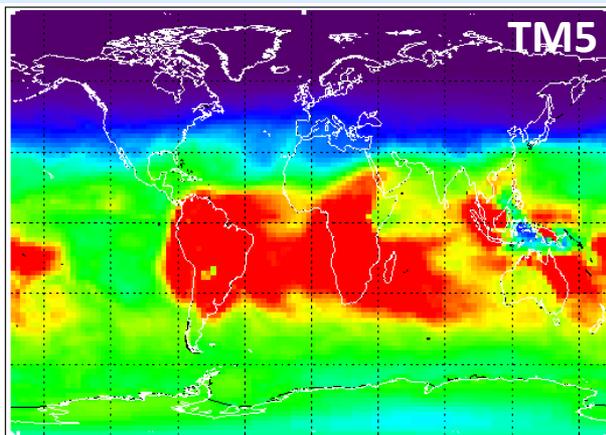
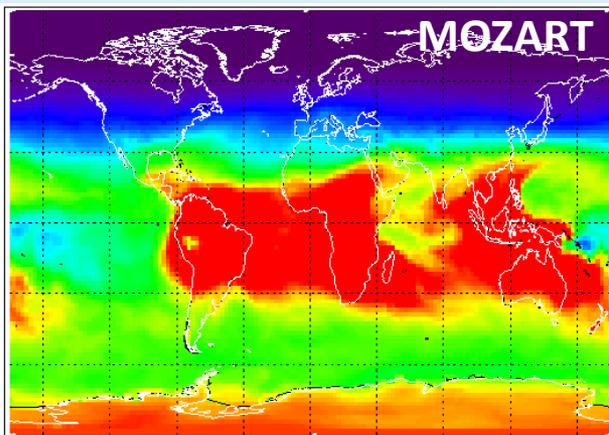


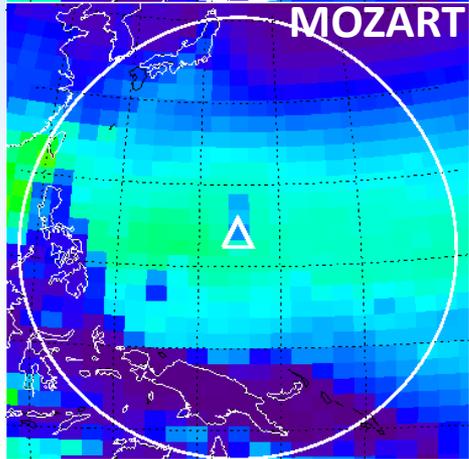
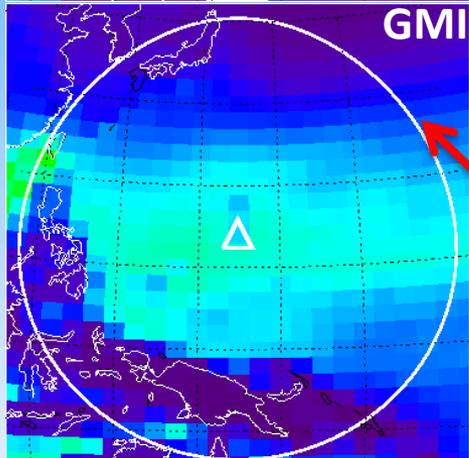
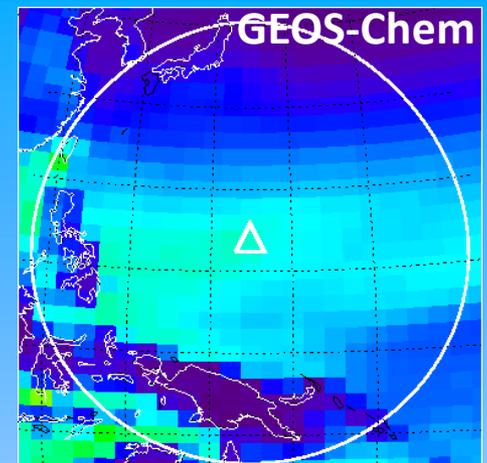
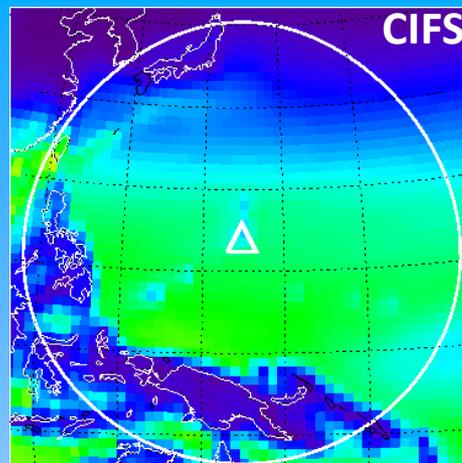
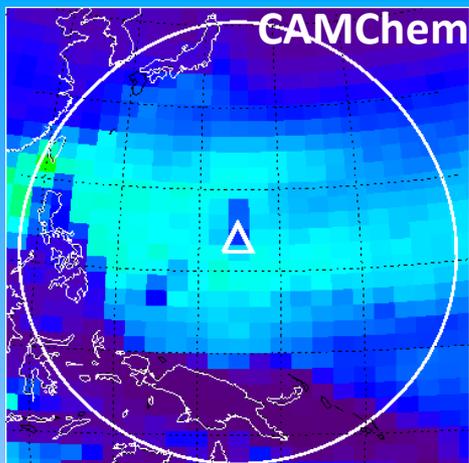
Trop CH₄ Lifetime vs. Month, monthly mean POLMIP output





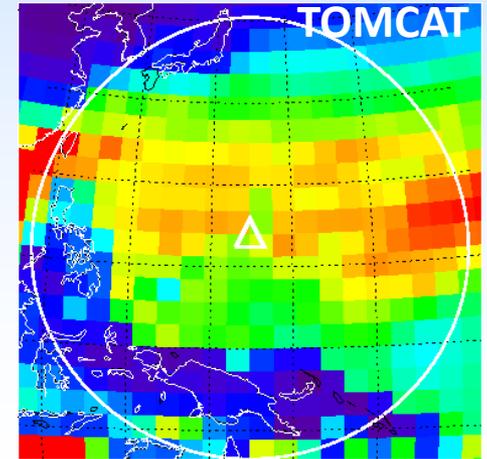
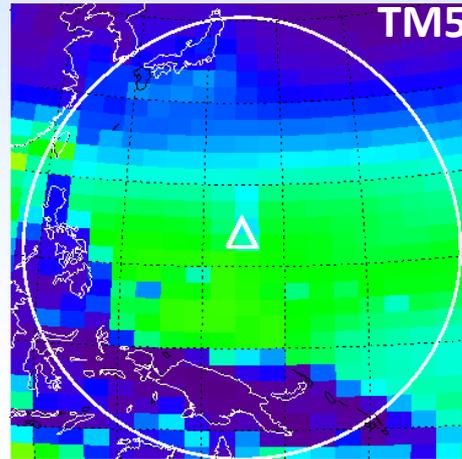
Monthly Mean OH M.R. at ~250 hPa
January, 2008
POLMIP

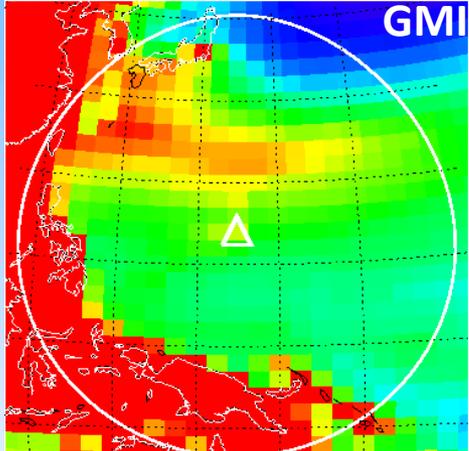
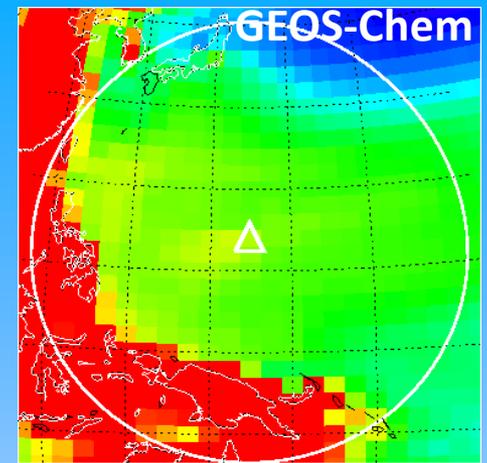
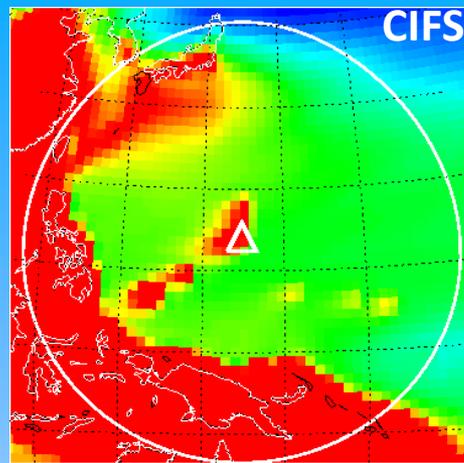
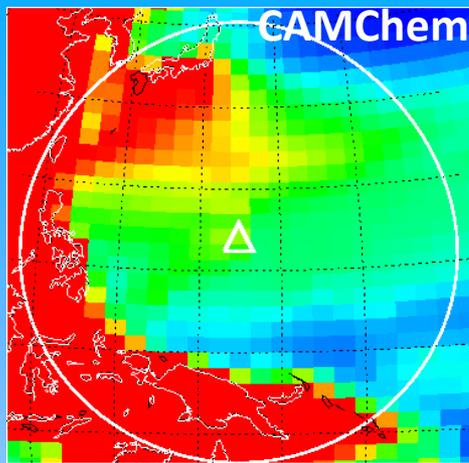




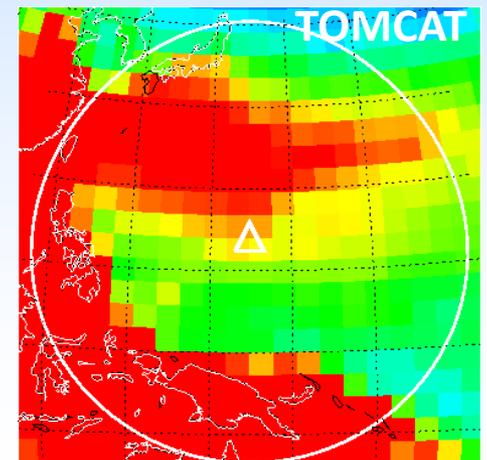
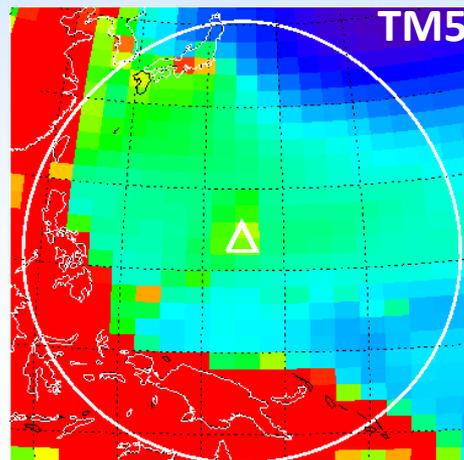
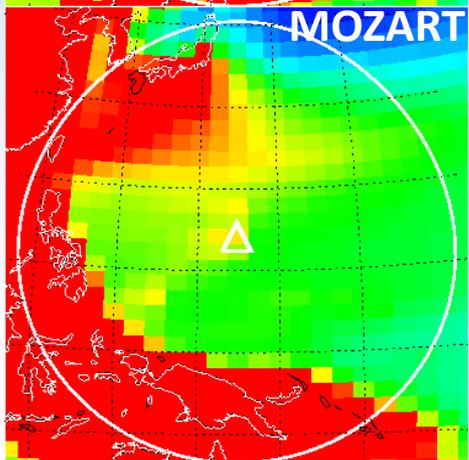
Monthly Mean OH Mixing Ratios at 957.7 hPa
January, 2008
POLMIP

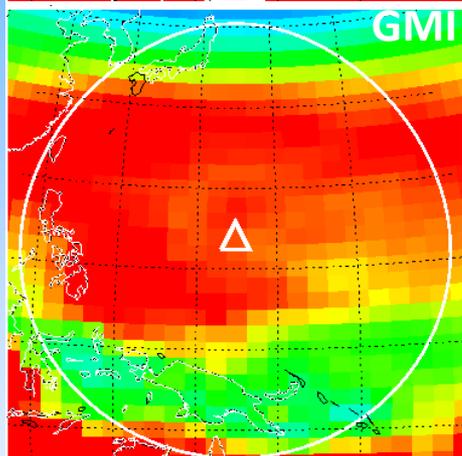
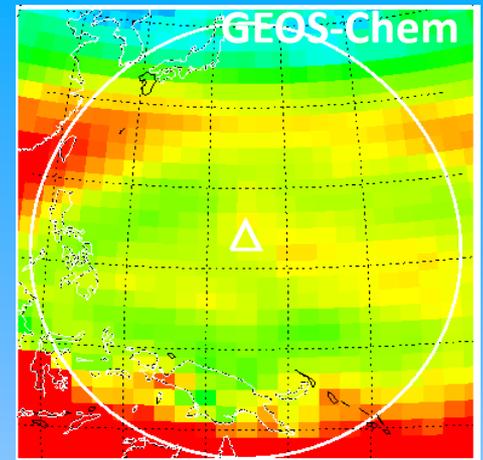
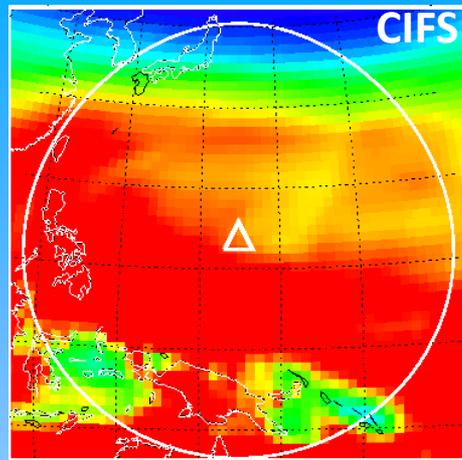
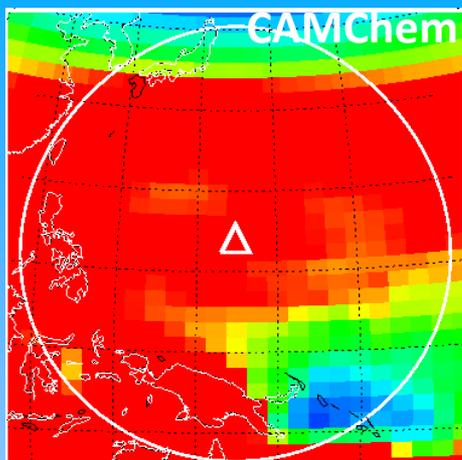
Circles designate
GV flight range



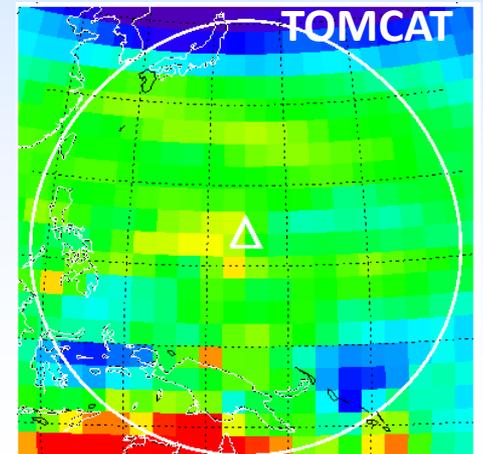
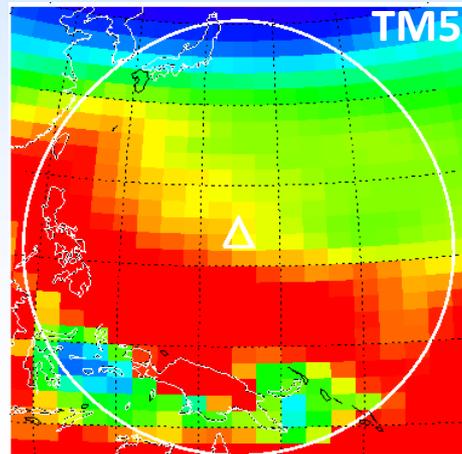
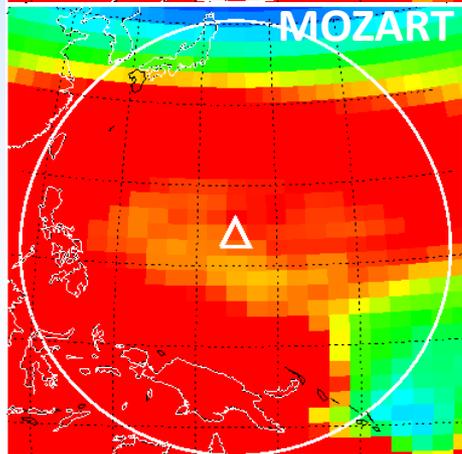
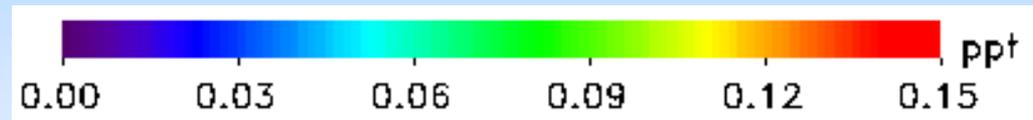


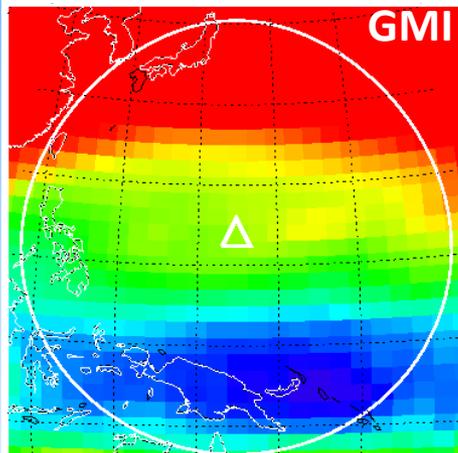
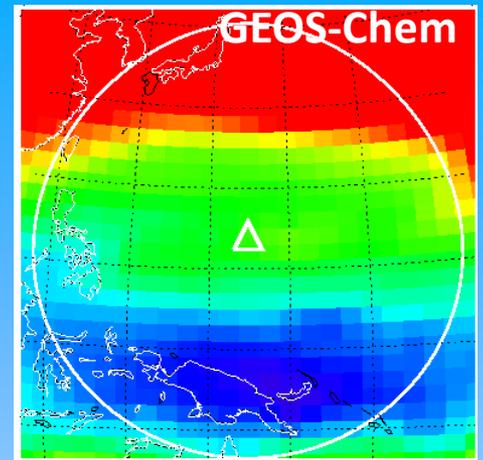
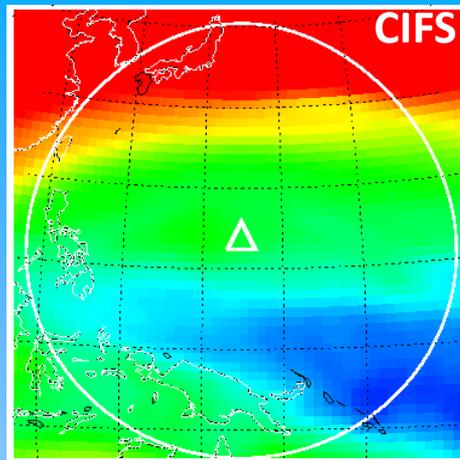
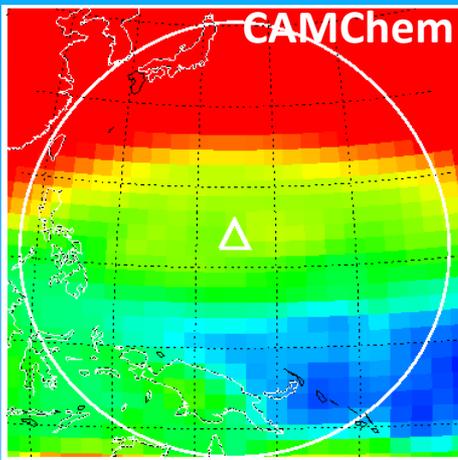
Monthly Mean HCHO Mixing Ratios at 957.7 hPa
January, 2008
POLMIP



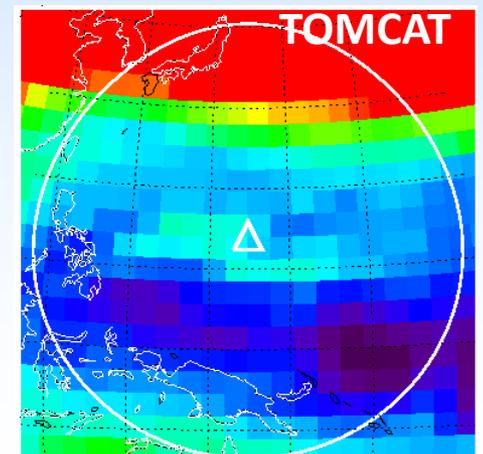
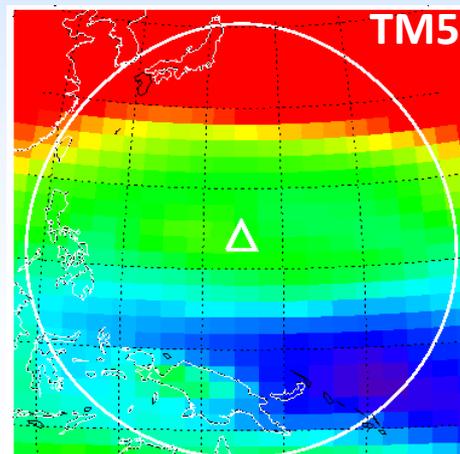
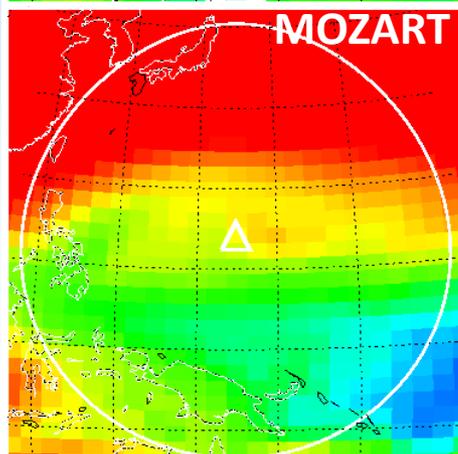


Monthly Mean OH Mixing Ratio
225 hPa
JAN 2008





Monthly Mean O₃ Mixing Ratio
225 hPa
JAN 2008



Box Model

DSMACC: Dynamically Simple Model for Atmospheric Chemical Complexity
tropospheric chemistry box model that can interface to various
chemical mechanisms

Emmerson and Evans, ACP, 2009

Makes use of the:

KPP (Kinetics PreProcessor)

Damian et al., Computers and Chemical Engineering, 2002.

Leeds Master Chemical Mechanism

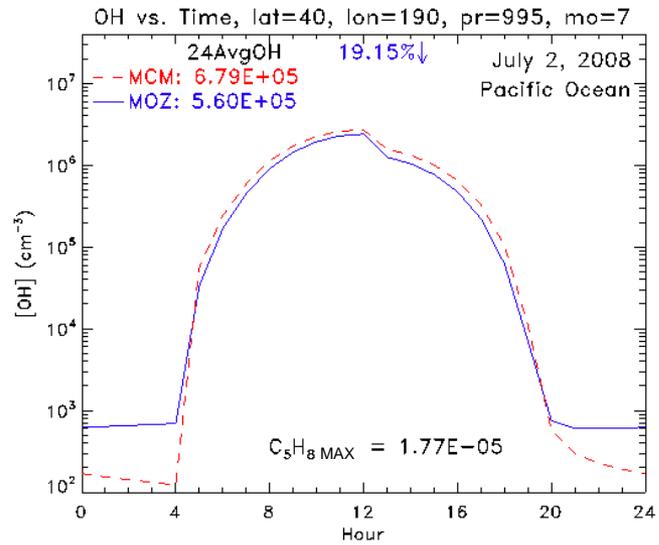
Bloss et al., ACP, 2005

**Here the box model constrained by H_2O , O_3 , CH_4 , CO , C_5H_8 , etc from CTM
as well as NO_2 , $J_{\text{O}1\text{D}}$, & $J_{\text{NO}2}$ from CTM**

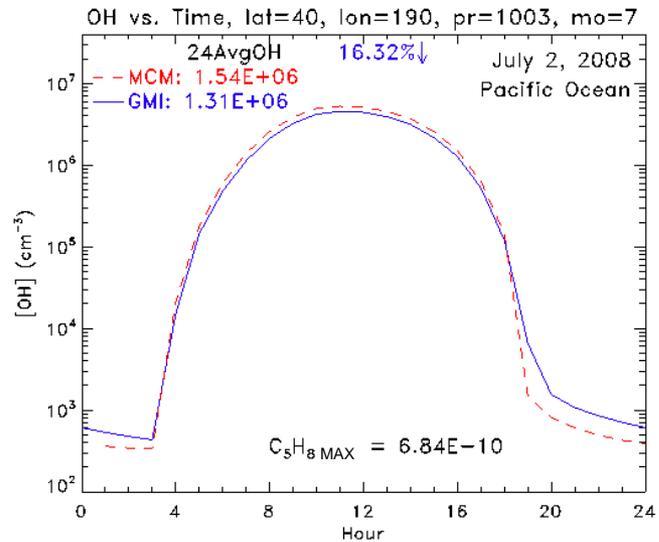
Focus on a few test cases for clear sky conditions

Model framework is amenable to analysis of cloudy conditions !

MOZART

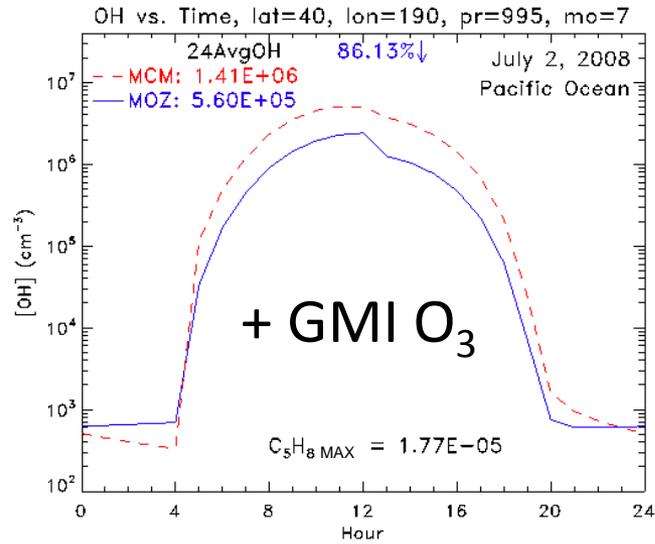


GMI

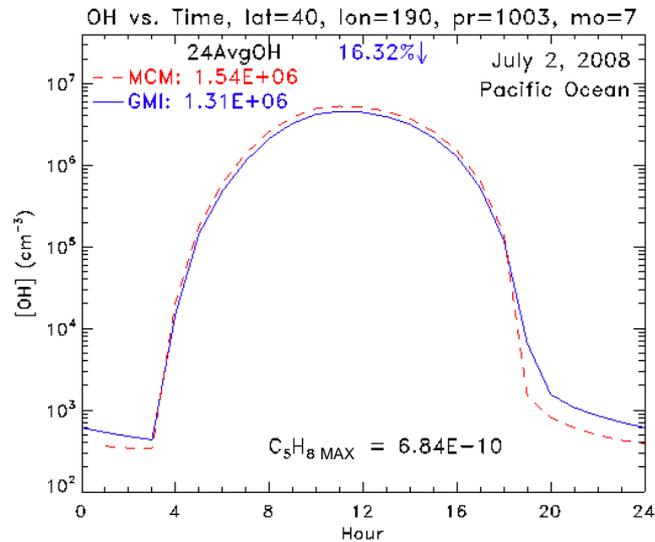


Box Modeling

MOZART

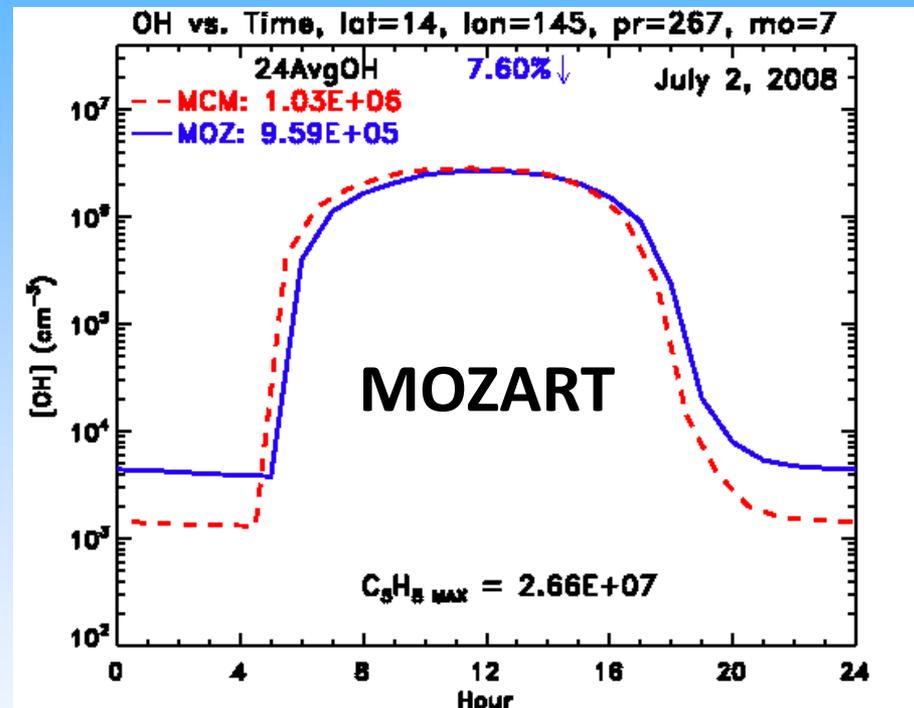
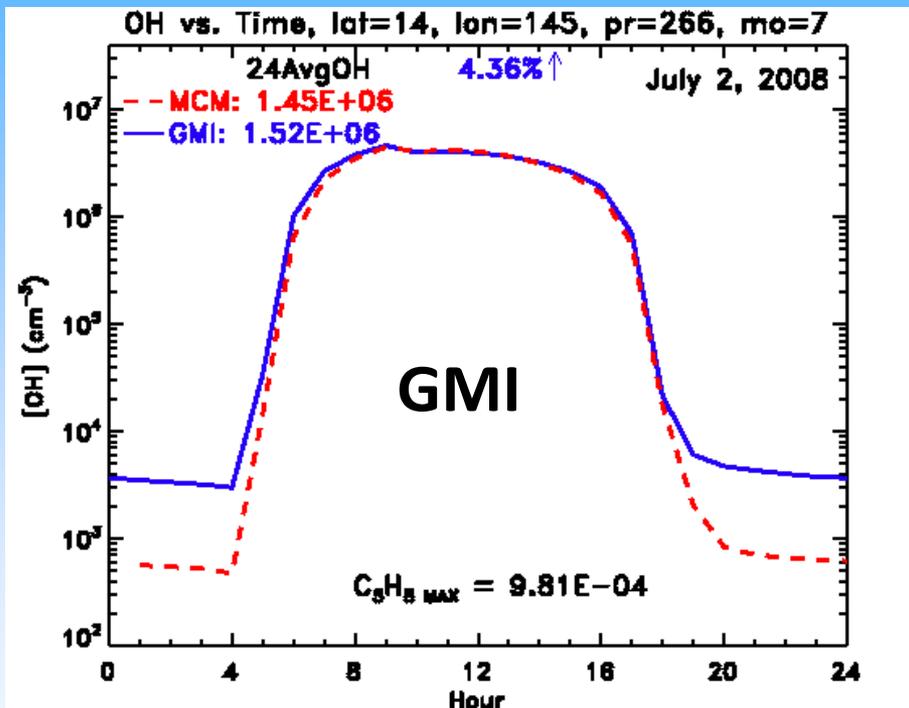


GMI



| Variable | Δ MOZART MCM 24AvgOH | MOZART Range | GMI-Clearsky Range |
|---|----------------------------|--|--|
| O ₃ (ppb) | +107.7% | 4.59 – 7.61 | 13.0 – 16.6 |
| J(NO ₂) & J(O ¹ D) (s ⁻¹) | +15.76% | 0 – 4.69×10 ⁻⁵ 0 – 0.0106 | 0 – 5.05×10 ⁻⁵ 0 – 0.0171 |
| NO _x (ppt) | +5.30% | 0.330 – 0.574 | 0.327 – 1.80 |
| H ₂ O (v.m.r.) | +3.53% | 0.0184 – 0.0193 | 0.0192 – 0.0196 |
| Isoprene (v.m.r.) | +1.18% | 6.78×10 ⁻²⁷ – 7.13×10 ⁻²⁵ | 2.22×10 ⁻²⁹ – 1.91×10 ⁻²⁶ |
| CO (ppb) | -2.50% | 68.1 – 69.2 | 72.3 – 82.2 |

TTL over Guam



Box Modeling

| Box model constraints | Box model output |
|--------------------------------|-------------------------------|
| p | OH and HO ₂ |
| T | H ₂ O ₂ |
| Overhead O ₃ Column | NO |
| O ₃ | HCHO |
| CO | Acetone |
| H ₂ O | Acetaldehyde |
| NO ₂ | Methanol |
| CH ₄ | Ethanol |
| C ₂ H ₆ | Methyl vinyl ketone |
| C ₃ H ₈ | Methacrolein |
| Isoprene | Methyl butenol |
| J(O ¹ D) | Propanal |
| J(NO ₂) | Butanal |
| Aerosol Surface Area Density | ... |

*Bold, blue text = measured during CONTRAST

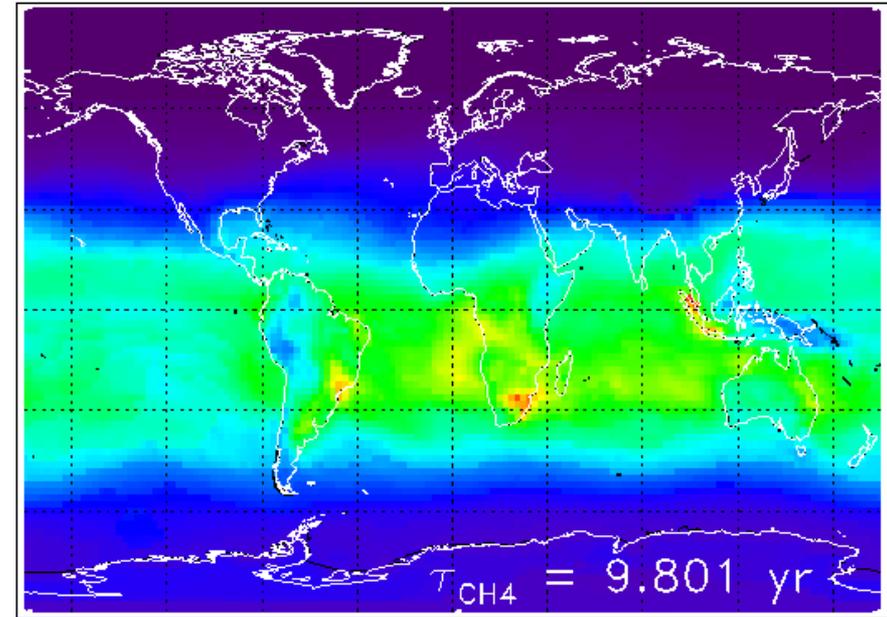
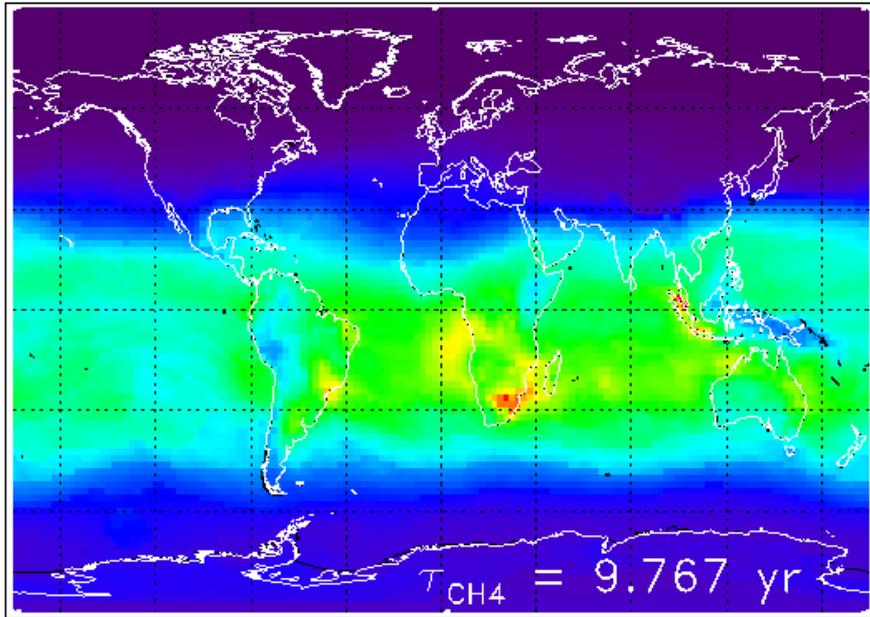
Neural Network Modeling

GMI (Actual Model)

Neural Network

Monthly Trop OH Column, GMI, 01/2008

Monthly Trop OH Column, GMI NN, 01/2008



TROP OH COLUMN

TROP OH COLUMN

Mean neural network $\Delta\tau_{\text{CH}_4}$ values (years) from swapping specified variable between models A/B

| Variable | CAM/GMI | MOZ/CAM | GMI/MOZ | Mean Abs. Value |
|--|-------------------|-------------------|-------------------|-----------------|
| CH ₄ | +0.88±0.24 | -0.51±0.13 | -0.57±0.32 | 0.65 |
| J(O ¹ D) | +0.93±0.21 | -0.13±0.11 | -0.74±0.20 | 0.60 |
| CO | +0.57±0.03 | -0.48±0.04 | -0.17±0.001 | 0.41 |
| O ₃ | -0.30±0.02 | -0.19±0.01 | +0.61±0.02 | 0.37 |
| J(NO ₂) | +0.10±0.05 | -0.32±0.24 | +0.21±0.13 | 0.21 |
| H ₂ O | +0.03±0.02 | -0.23±0.04 | +0.22±0.02 | 0.16 |
| NO _x | +0.14±0.05 | -0.07±0.03 | -0.06±0.01 | 0.09 |
| ISOP | +0.09±0.03 | +0.03±0.0045 | -0.14±0.01 | 0.09 |
| T | -0.002±0.001 | -0.002±0.002 | +0.003±0.002 | 0.002 |
| | | | | |
| Original τ_{CH_4} | 6.99 / 9.80 | 9.12 / 6.99 | 9.80 / 9.12 | - |
| τ_{CH_4} + (total Δ) | 8.70 / 6.80 | 6.18 / 8.63 | 8.45 / 9.06 | - |

Conclusions

- CTMs and CCMs have τ_{CH_4} values that differ by $\pm 20\%$ for reasons that are not well understood
- Box models can identify reasons for differences in OH between 2 CCMs when the CCMs use similar chemical mechanisms
- Neural networks provide method of identifying causes of OH differences between 2 CCMs even when the CCMs use markedly different chemical mechanisms
- CONTRAST will provide data to evaluate the accuracy of models' precursor fields and chemical mechanisms