



# Update on carbon monoxide inversions

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**Johann Rasmus Nüß<sup>1</sup>**

Nikos Daskalakis<sup>1</sup>, Oliver Schneising<sup>2</sup>, Michael Buchwitz<sup>2</sup>,  
Maarten C. Krol<sup>3,4</sup> and Mihalis Vrekoussis<sup>1</sup>

<sup>1</sup> LAMOS group, IUP, University Bremen

<sup>2</sup> Carbon and Greenhouse Gas Group, IUP, University Bremen

<sup>3</sup> MAQ, WIMEK, Wageningen University and Research, the Netherlands

<sup>4</sup> IMAU, Utrecht University, the Netherlands

- 1 Californian Fires
  - Objective and Motivation
  - Intermediate Results
- 2 Novel chemical-scheme development for carbon monoxide inversions (CHEMFORCER)
  - Background
  - Methods
- 3 Summary & Outlook

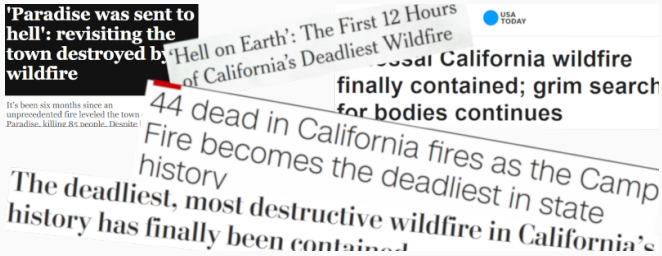
# Californian Fires

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# The bigger plan

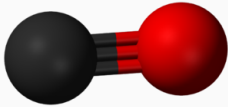
- **Use high-spatial-resolution satellite data**
- Extend the chemical scheme of the model for using satellite data products for multiple species
- Perform inversions using data from multiple satellites instruments

# Californian wild fires

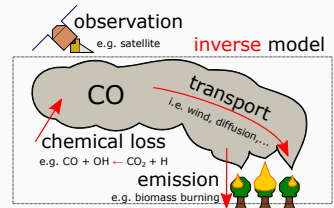
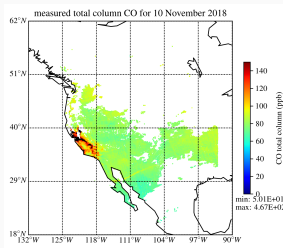


- Warmer and dryer than usual → wildfires more likely
- November and December 2018: major burning events
- Focus on Camp and Woolsey fires, raging in the weeks after November 8th
  - Devastated area about 1000 km<sup>2</sup>
  - Direct damage: 88 dead, burned land and structures, forced evacuation of multiple towns
  - Indirect damage due to pollution

# Objective and Motivation



Retrieve CO emissions from biomass burning events in California using TROPOMI observation in the TM5 4DVAR model.



Images: Fire: Mark McKenna / Zuma Press, LA-Times; S5P: ESA

# TROPOMI observations

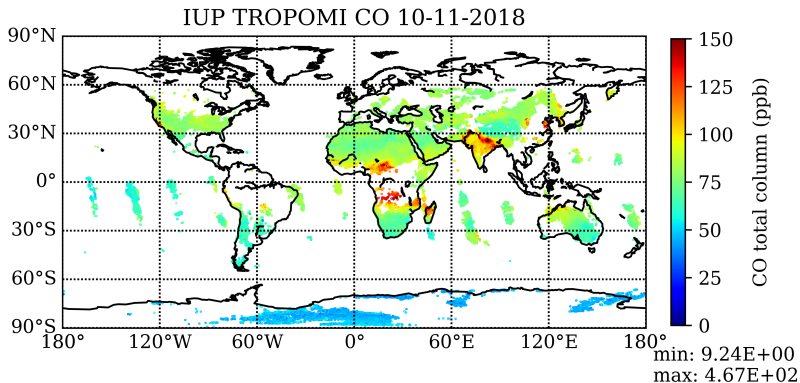
- **TROPO**spheric **M**onitoring **I**nstrument onboard of **Sentinel-5 P**recursor
- Daily global coverage
- Local overpass time 13:30
- High resolution (up to  $7 \times 7 \text{ km}^2$ )
  - Still useful for  $1^\circ \times 1^\circ$  model boxes: lower error, chance to have at least some cloud free pixels
- Especially sensitive to troposphere/boundary layer



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Image: ESA

# Satellite observations



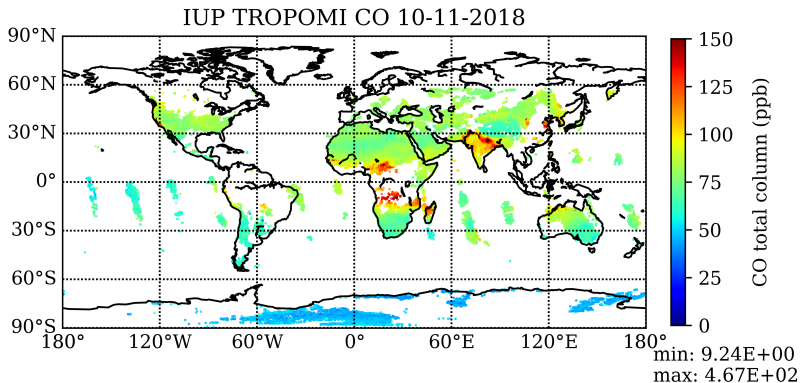
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Satellite data courtesy of Oliver Schneising and Michael Buchwitz of IUP  
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# Satellite observations



- Given: TROPOMI CO total column observation
- Wanted: Location and temporal development of emissions

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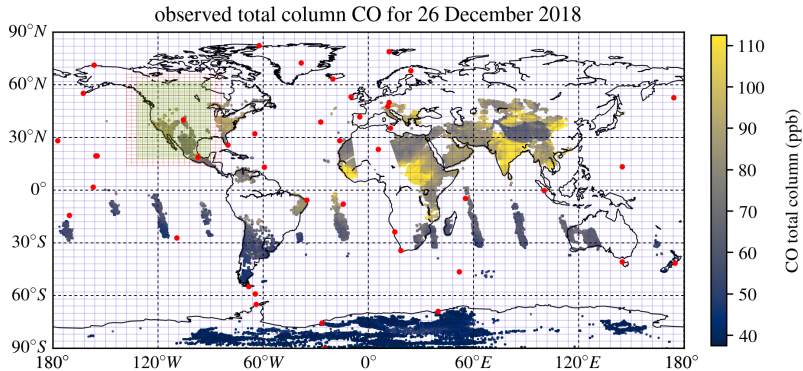
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- M1qN3 optimizer
- Zoom over California

# Zooming, flask measurements and satellite observations



- Start from TM5MP full chemistry fields
- Spin-up inversion for 5 months to flasks and global IASI TROPOMI observations on  $6^\circ \times 4^\circ$
- 3 month main inversion period around 2 week time of interest



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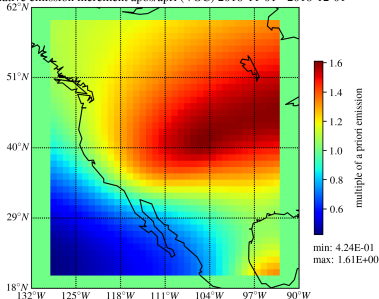


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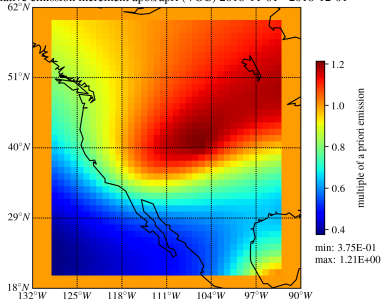
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- Spin up works as intended

# Aliasing between biomass burning and VOC emissions

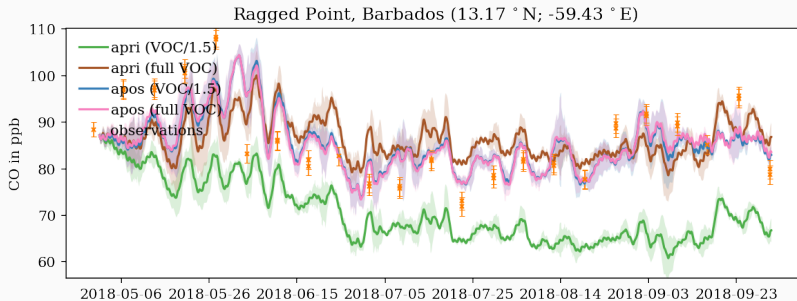
Relative emission increment apos/apri (VOC) 2018-11-01 - 2018-12-01

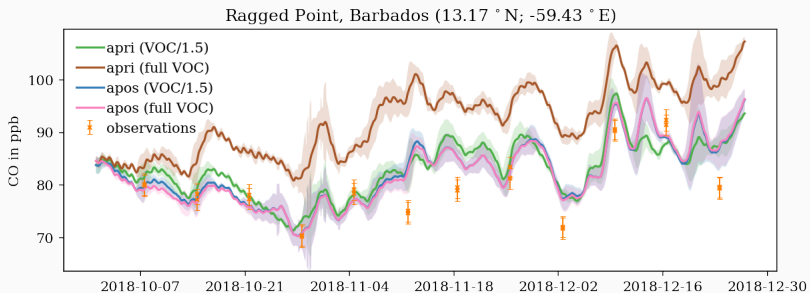


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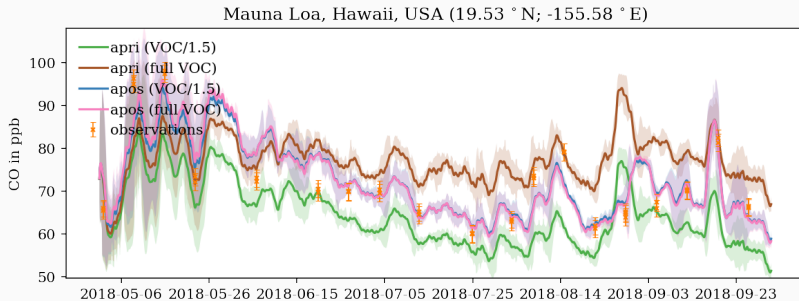


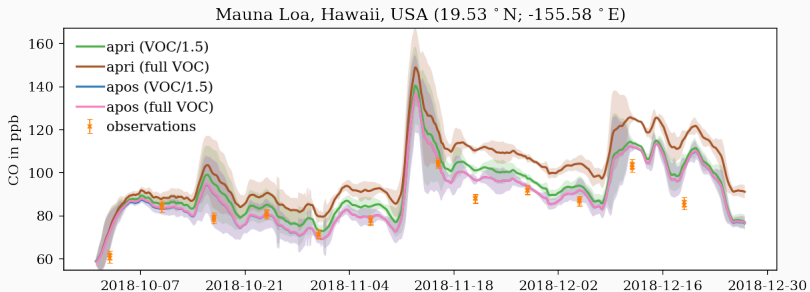
- Identical setups except for BB prior (left: GFED, right: FINN)
- Simultaneous inversion of BB CO and CO from VOC/CH<sub>4</sub>
- A posterior production from VOC/CH<sub>4</sub> strongly affected by a priori biomass emissions → simultaneous inversion not feasible

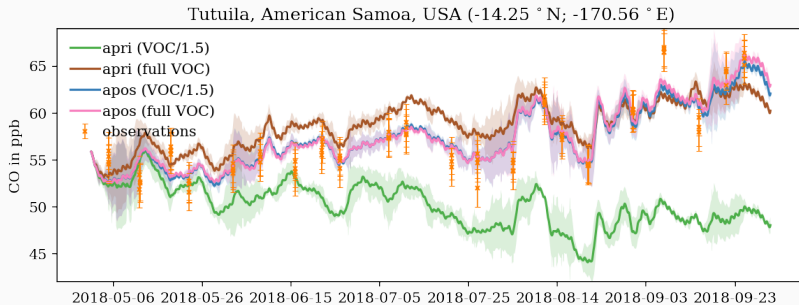




# MLO - NH-summer

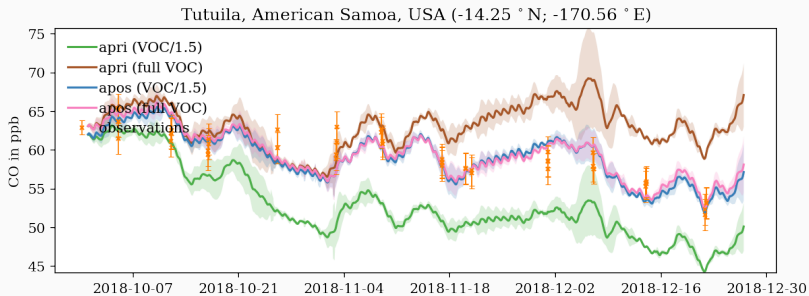








# SMO - NH-winter



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- Consider more complex chemistry, like HCHO

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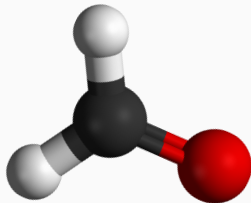
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- Collaboration with Prof. Dr. Maarten C. Krol, one of the model's original creators, and other students working on related topics

# The bigger plan

- Use high-spatial-resolution satellite data
- **Extend the chemical scheme of the model for using satellite data products for multiple species**
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# Formaldehyde (HCHO)



- Volatile, colorless and toxic gas
- Intermediate species in oxidation of  $\text{CH}_4$  and VOCs to CO
  - Specifically, proxy species for the isoprene produced by plants
  - As such part of and proxy for the natural/biogenic CO source
- Shortlived, lifetime of a few hours in the sunny atmosphere



## Extended inversion scheme

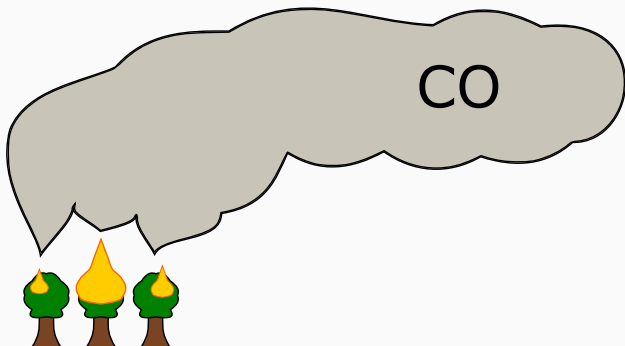
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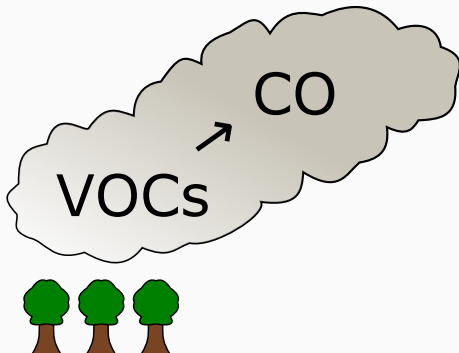
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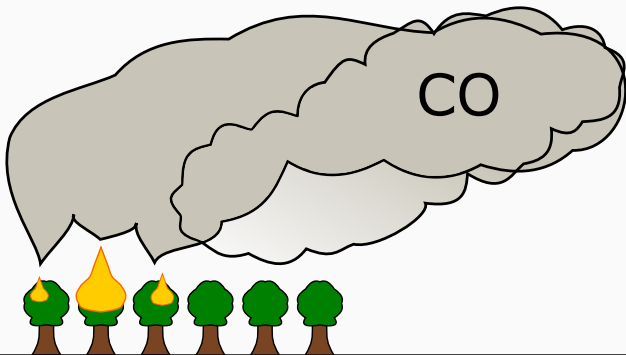
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Solution:

- Use observations for CO **and** HCHO
  - HCHO sufficient proxy for the whole natural source
- If ambient and local HCHO are known, the biogenic source can be estimated and distinguished from the biomass burning source

## Extended inversion scheme (2)

Multiple approaches:

- Extend the currently very simplified CO chemistry in TM5-4DVAR to include HCHO
- Solving more complex chemistry → severely increased computation time
- Create look-up tables for loss and production of CO and HCHO under different ambient conditions
  - Use TM5MP to get loss rates..
  - .. either use a simplified scaling approach
  - .. or do a dedicated HCHO inversion

## Scaling approach

$$P_{\text{CO}} \propto P_{\text{HCHO}} = L_{\text{HCHO}} \cdot [\text{HCHO}]_{\text{TM5}} \cdot \text{scaling\_factor}$$

- $P_{\text{HCHO}}$ : Biogenic source of HCHO to be used in the inversions
- $L_{\text{HCHO}}$ : Chemical loss rate of HCHO, to be taken from TM5MP
- $[\text{HCHO}]_{\text{TM5}}$ : Current concentration of HCHO, to be taken from TM5MP and compared to satellite observations to obtain scaling\_factor

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- $\text{scaling\_factor} = [\text{HCHO}]_{\text{obs}} / [\text{HCHO}]_{\text{TM5}}$
- Varies in time and space
- Each grid box has its own factor, default 1
- Updated whenever there is a measurement or monthly, to be tested



Multiple challenges for HCHO inversions:

- Generally short, but variable lifetime
- Pronounced diurnal cycle
- Little, but non-negligible transport
- Diurnal cycle in vertical distribution of precursors
- Fast chemistry, highly depended on exact ambient conditions (sun, OH, lots of other VOCs)

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- Lots of coding left to be done

# Acknowledgments

- The computations were performed on the HPC cluster Aether at the University of Bremen, financed by DFG in the scope of the Excellence Initiative.
- The PhD position is paid for by the University Bremen.
- The travel expenses were covered by the BremenIDEA program of the University of Bremen.
- Special thanks to the TM5 community, especially Maarten Krol and Sourish Basu for provision of and help with the TM5-4DVAR model.
  
- ... and of course thank You for your attention