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# **Aerosol modeling with WRF/Chem**

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**WRF/Chem Tutorial, 22 July 2013**

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## **Part I - Introduction**

- **Overview of ...**
  - Aerosol
  - Aerosol processes and life cycle
  - Model treatment of aerosol
  - WRF/Chem aerosol schemes

## **Part II – The details**

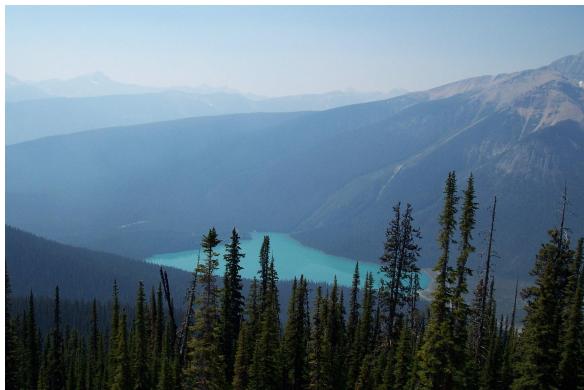
- **Representing the aerosol size distribution**
- **Walk through the WRF/Chem aerosol schemes**
  - How they work and what they do
  - Coupling to other processes
    - ◆ Gas phase chemistry
    - ◆ Aqueous chemistry
    - ◆ ...
- **Hint on how to tell WRF/Chem what to do**
- **Resources**

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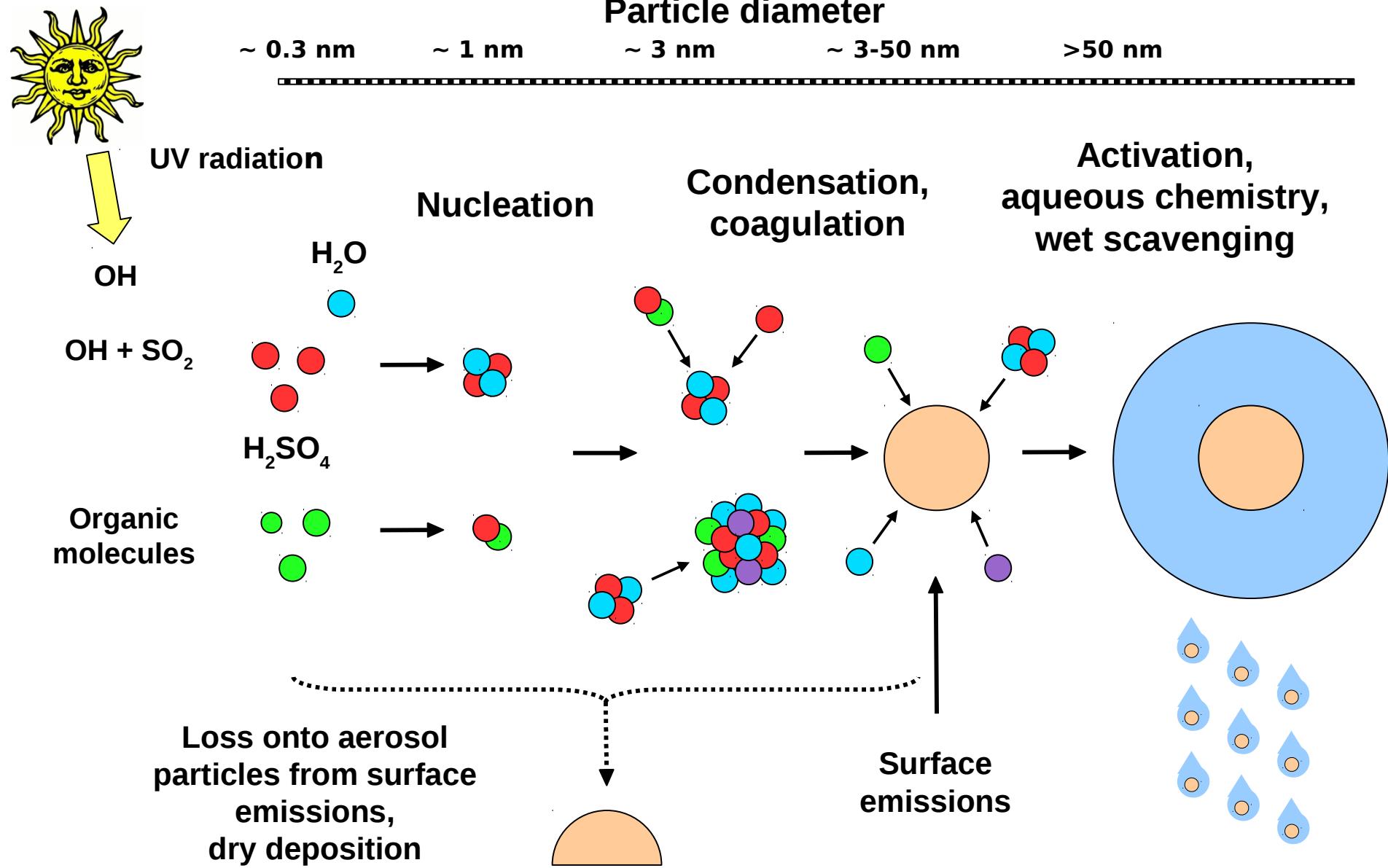
# **Part I – Introduction**

# Aerosol

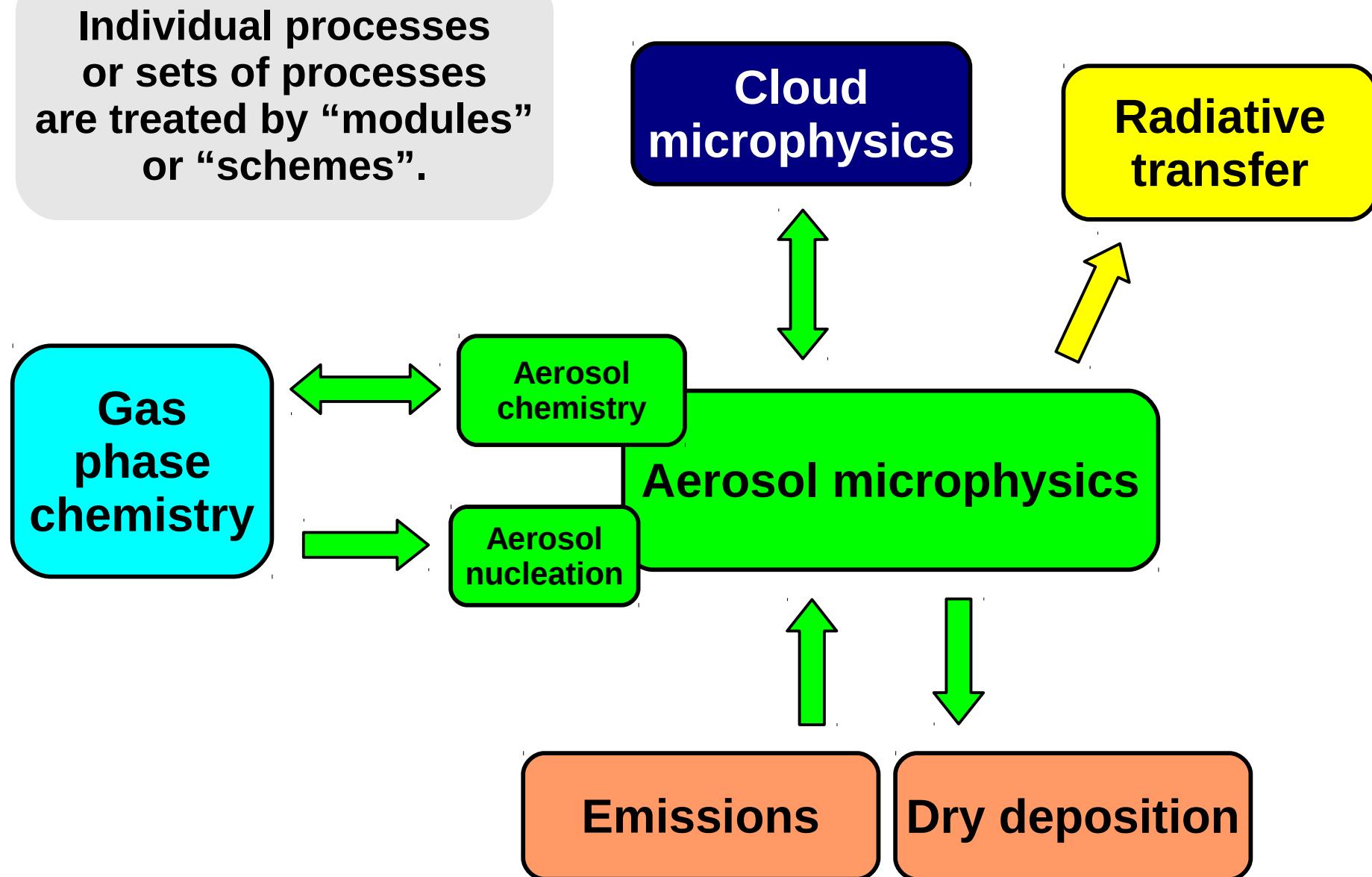
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# Aerosol life cycle and processes



# Model treatment of aerosol



# WRF/Chem aerosol schemes

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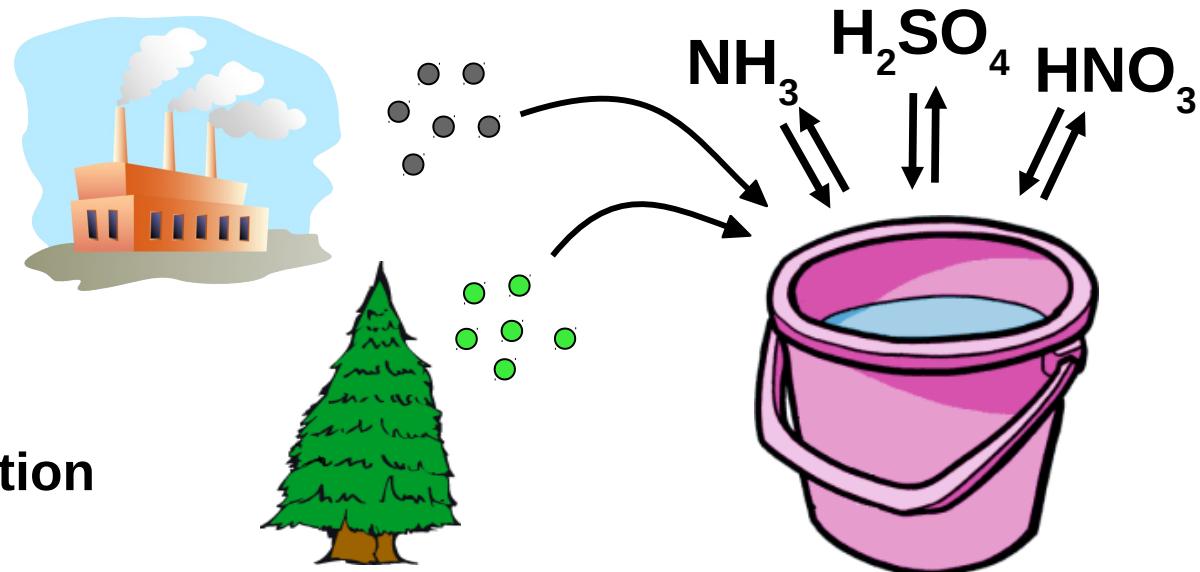
- An efficient aerosol scheme from the GOCART model
  - No size information for sulfate, BC, OC
  - Size information for dust and sea salt
  - No secondary organic aerosol (SOA)
- Modal Aerosol Dynamics Model for Europe – MADE
  - 3 log-normal modes
  - Inorganic, organic aerosol, SOA
- Model for Simulating Aerosol Interactions and Chemistry (MOSAIC)
  - Sectional model, 4 or 8 bins
  - Inorganic, organic aerosol, SOA
- MAM – Modal Aerosol Model (new in WRF/Chem 3.5) from CAM5
  - 3 or 7 log-normal modes
  - Inorganic, organic aerosol, SOA, sea salt, BC, mineral dust
- Simple sectional (bin) scheme for volcanic ash aerosol

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## **Part II – The details**

# Bulk aerosol schemes

- Only total mass of aerosol compounds is known



- No information on
  - Particle number
  - Aerosol size distribution

Aerosol size distribution needs to be assumed for:

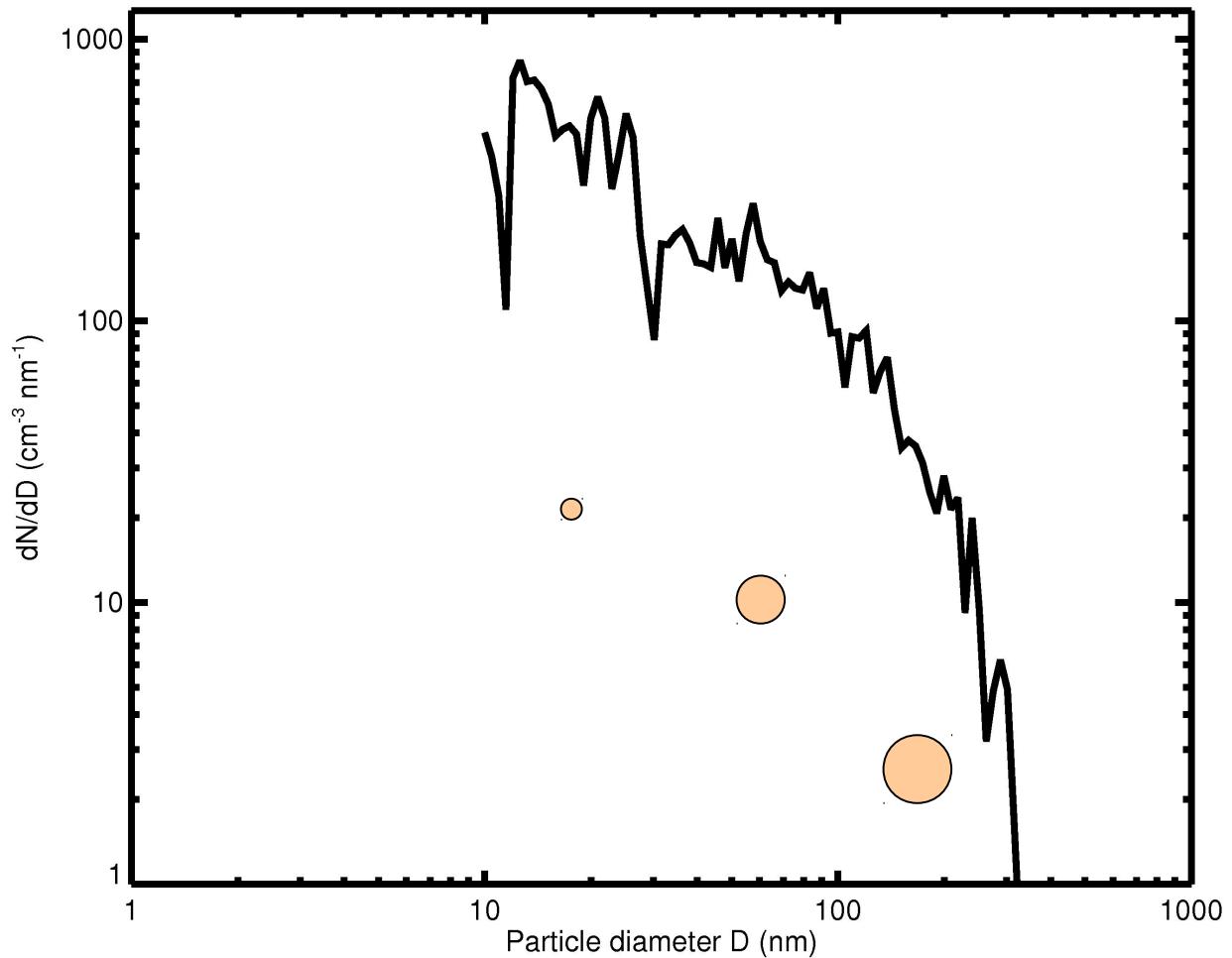
- radiative transfer
- response of cloud properties to aerosol number



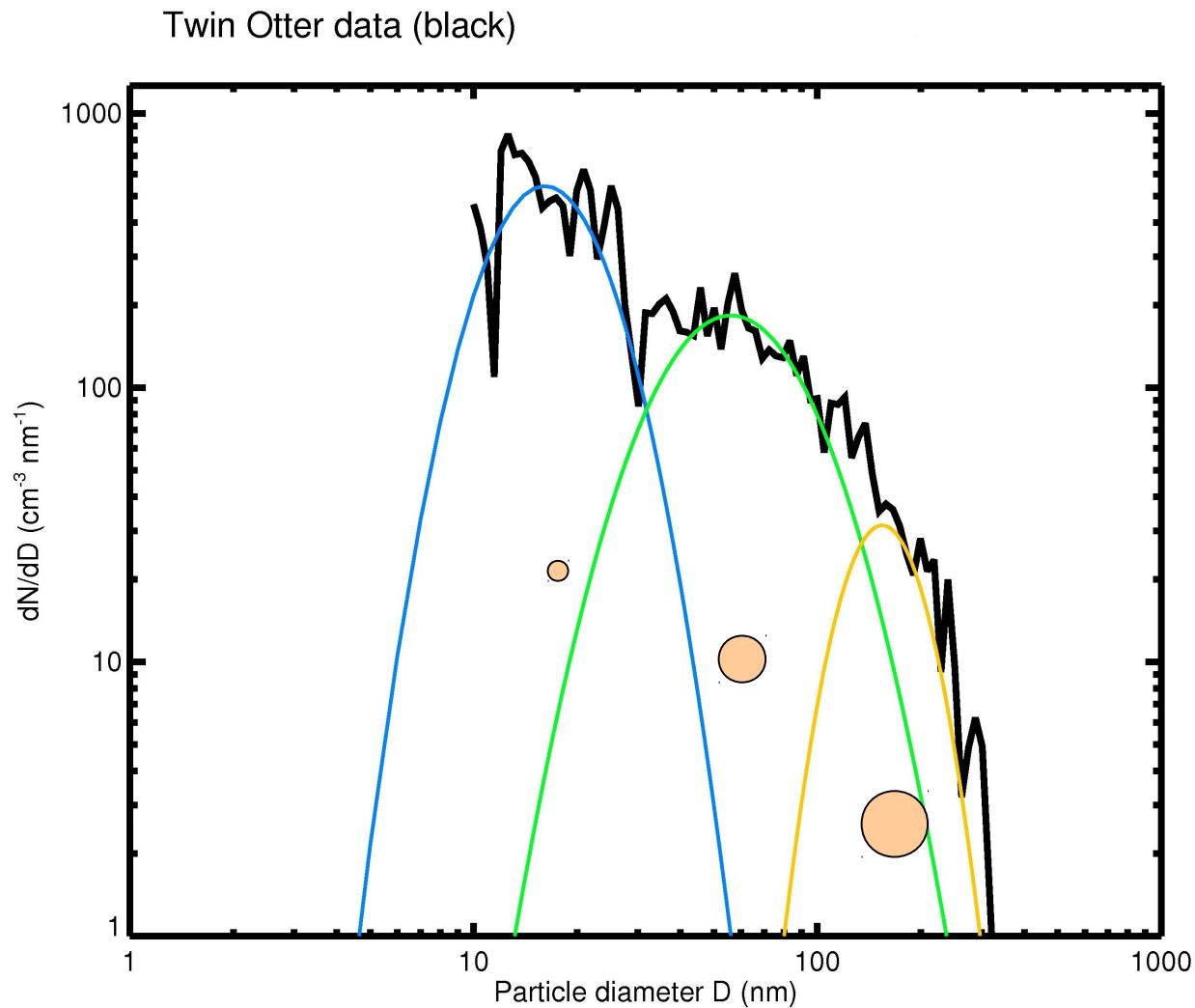
- Numerically efficient
  - Useful when focus is on complex gas phase chemistry
- **GOCART (+ size resolved dust and sea salt)**

# Modal aerosol schemes

Twin Otter data (black)



# Modal aerosol schemes



$$\frac{dN}{dD} = \frac{N}{\sqrt{2\pi} \ln(\sigma) D} e^{-\frac{1}{2} \left[ \frac{\ln(D/\mu)}{\ln(\sigma)} \right]^2}$$

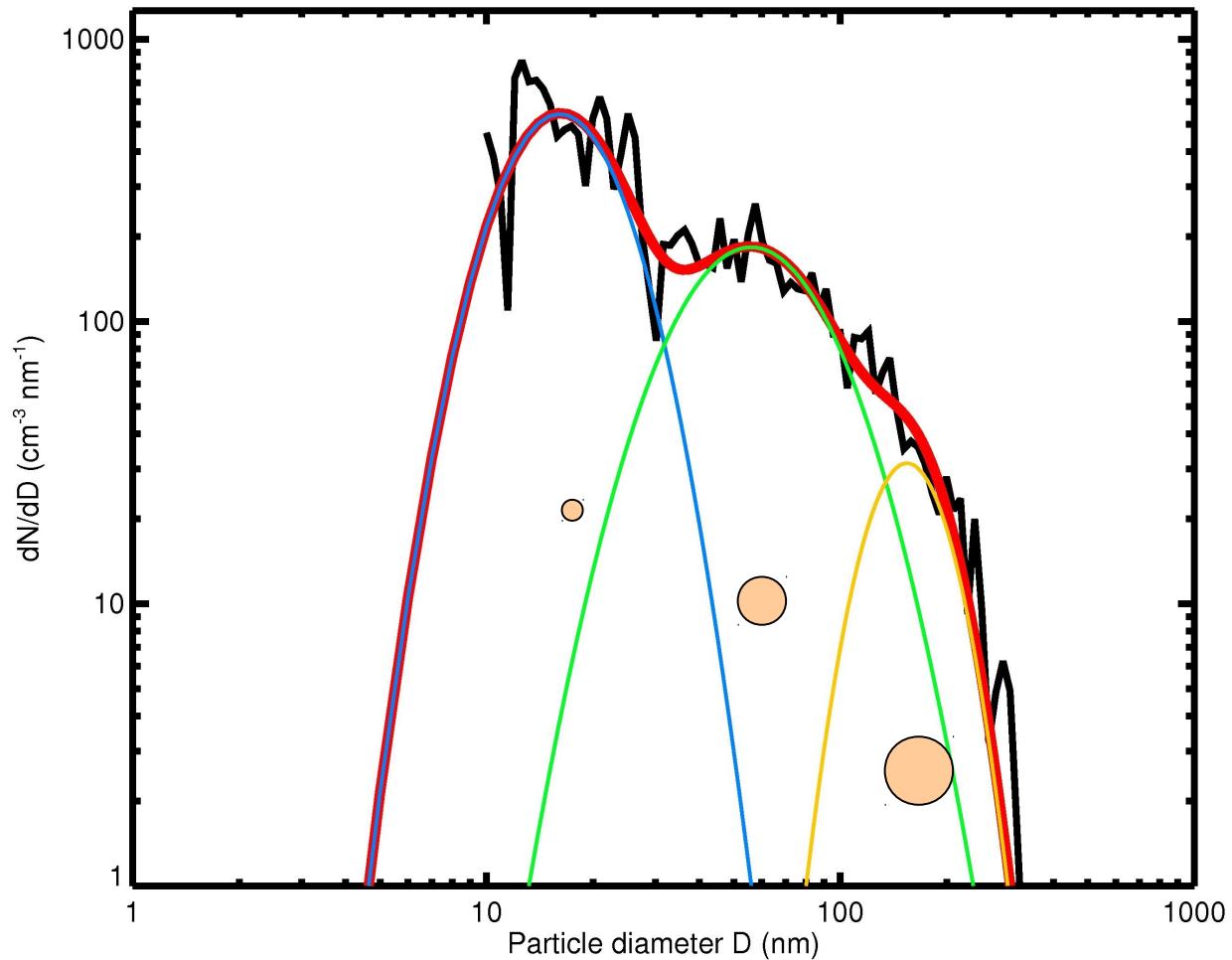
$$\frac{dN}{dD} \rightarrow N = 8195 \text{ cm}^{-3}$$
$$\mu = 18.22 \text{ nm}$$
$$\sigma = 1.42$$

$$\frac{dN}{dD} \rightarrow N = 12732 \text{ cm}^{-3}$$
$$\mu = 68.44 \text{ nm}$$
$$\sigma = 1.57$$

$$\frac{dN}{dD} \rightarrow N = 3140 \text{ cm}^{-3}$$
$$\mu = 164.41 \text{ nm}$$
$$\sigma = 1.28$$

# Modal aerosol schemes

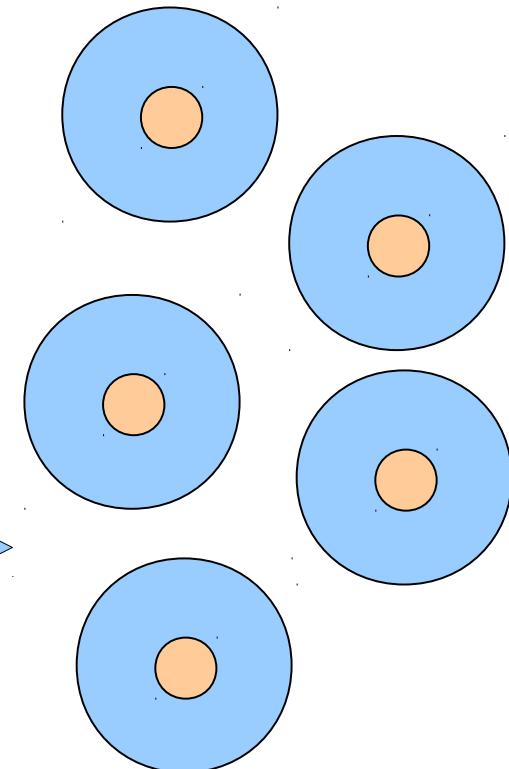
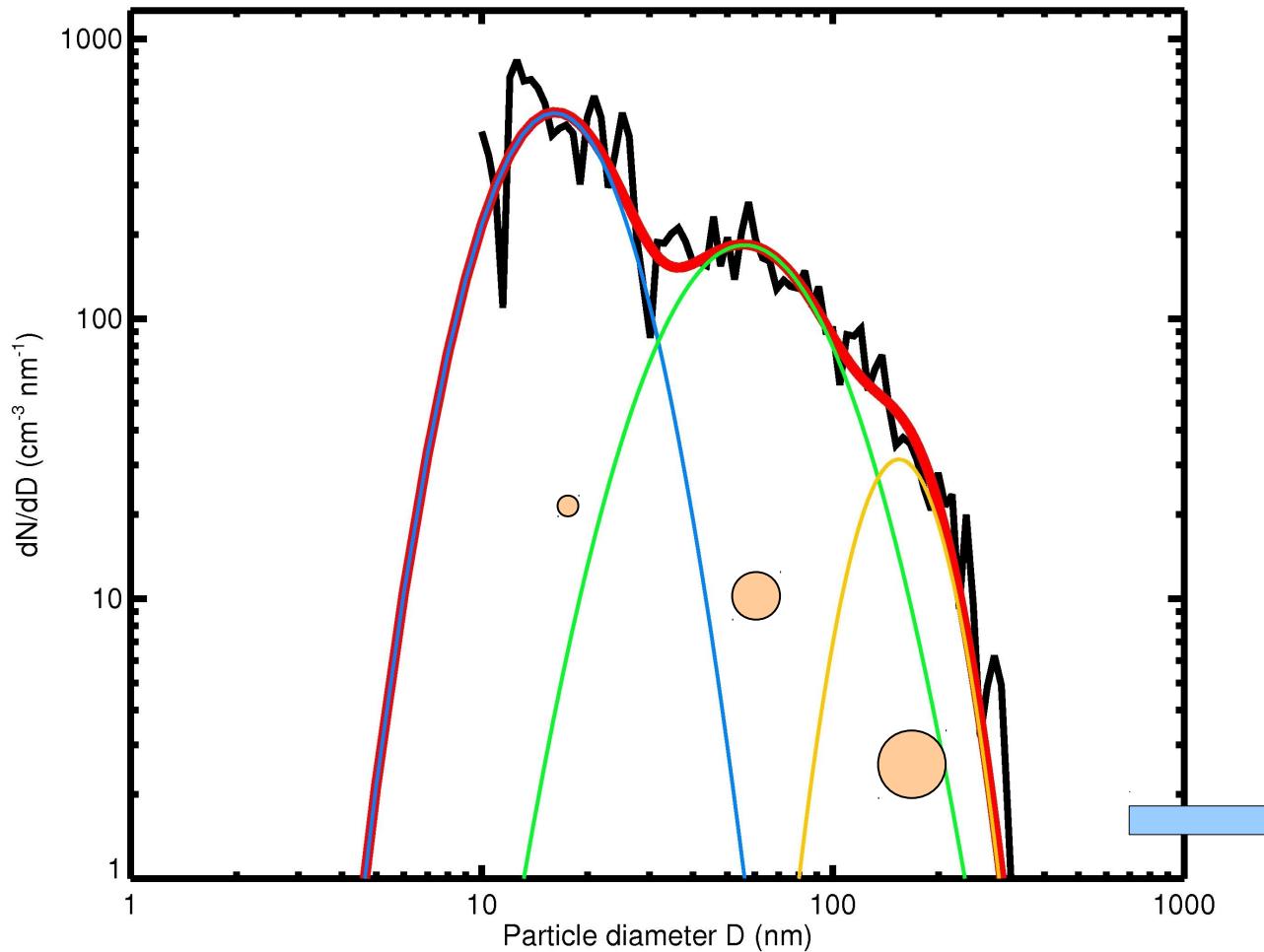
Twin Otter data (black)



$$\frac{dN}{dD} = \frac{dN}{dD} + \frac{dN}{dD} + \frac{dN}{dD}$$

# Modal aerosol schemes

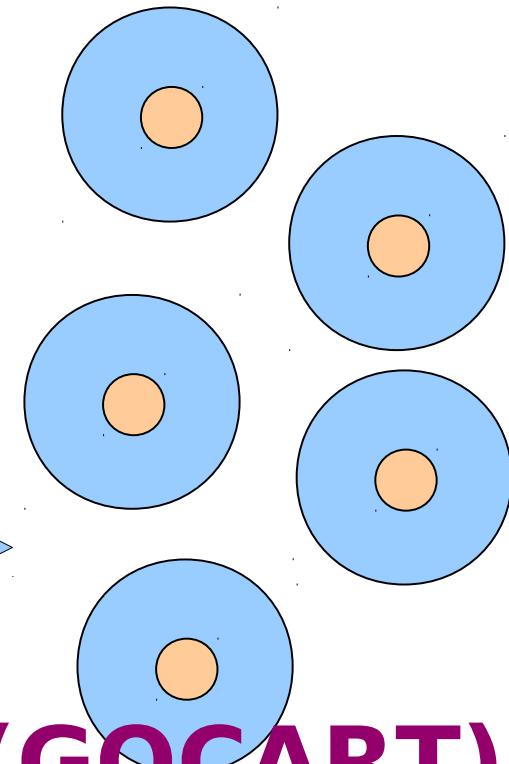
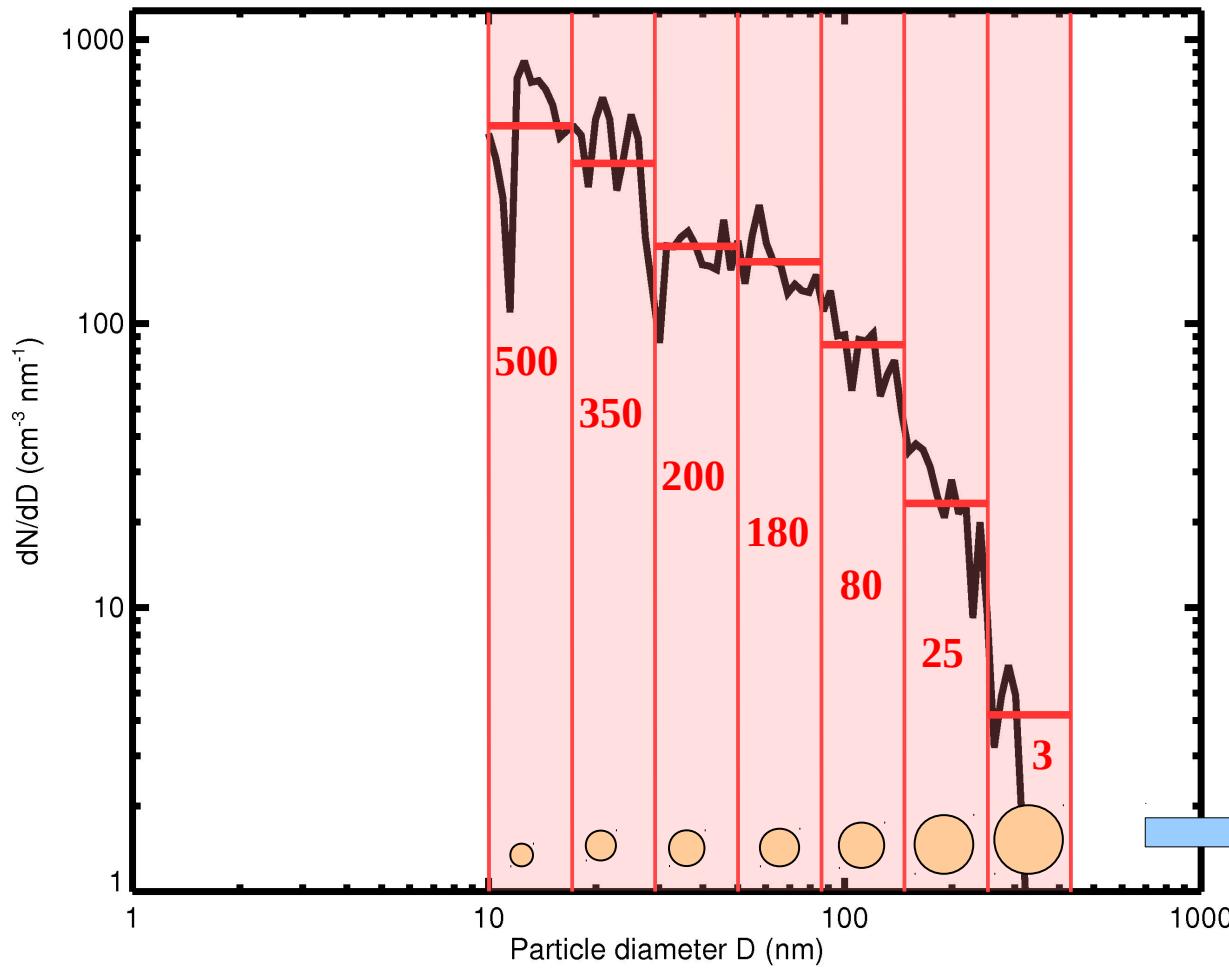
Twin Otter data (black)



→ MADE and MAM

# Sectional aerosol schemes

Twin Otter data (black)



→ MOSAIC, volcanic ash, (GOCART)

# GOCART aerosol module

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- Georgia Tech/Goddard **Global Ozone Chemistry Aerosol Radiation and Transport model** (Chin et al., JGR, 2000)
  - Bulk aerosol:
    - ◆ Hydrophobic black carbon (fresh soot)
    - ◆ Hydrophilic black carbon (aged/coated soot)
    - ◆ Hydrophobic organic carbon (fresh burnt biomass)
    - ◆ Hydrophilic organic carbon (aged/coated burnt biomass)
      - Fresh → aged conversion time 2.5 days
    - ◆ Other GOCART primary PM2.5
    - ◆ Other GOCART primary PM10
    - ◆ Sulfate (only secondary aerosol species)
  - Sectional scheme for dust and sea salt:
    - ◆ Dust: 0.5, 1.4, 2.4, 4.5, 8.0  $\mu\text{m}$  effective radius
    - ◆ Sea salt: 0.3, 1.0, 3.2, 7.5  $\mu\text{m}$  effective radius

# GOCART aerosol module

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**GOCART comes with sulfur gas phase chemistry:**

- DMS + OH → SO<sub>2</sub> + ...
- DMS + OH → MSA + ...
- DMS + NO<sub>3</sub> → SO<sub>2</sub> + ...
- SO<sub>2</sub> + OH → SO<sub>4</sub><sup>=</sup> + ...

**Extended gas phase chemistry can be used:**

- MOZART (with KPP)
- RACM (with KPP)
- RADM (with and without KPP)

# GOCART aerosol module

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- **Interaction with radiation:**
  - Direct effect for some model setups
  - Effect on photochemistry
- **Interaction with clouds:**
  - Aqueous chemistry
    - ◆  $\text{SO}_2 + \text{H}_2\text{O}_2 \rightarrow \text{SO}_4^=$
    - ◆  $\text{SO}_2 + \text{O}_3 \rightarrow \text{SO}_4^=$
  - No indirect effect
  - No wet scavenging/deposition
- **No secondary organic aerosol (SOA)**

# MADE aerosol module

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## Modal Aerosol Dynamics Model for Europe

(Ackermann et al., Atm. Env., 1998)

- **3 log-normal aerosol modes: Aitken, accumulation, coarse**
- Mode width  $\sigma$  is fixed
- Aerosol number and mass variable
- (Currently no nucleation mode)
- **Interaction with radiation:**
  - Direct aerosol effect
  - Effect on photolysis
- **Interaction with clouds:**
  - Aerosol number determines cloud drop number and size
  - Radiative response → 1<sup>st</sup> indirect aerosol effect
    - ◆ only for resolved clouds (Sc)
  - Aqueous chemistry
  - Wet removal (scavenging)

# MADE aerosol module

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## Aitken and accumulation modes:

- $\text{SO}_4^{=}$ ,  $\text{NH}_4^+$ ,  $\text{NO}_3^-$ ,  $\text{H}_2\text{O}$
- NaCl (sea salt)
- Anthropogenic SOA from oxidation of ...
  - Alkanes
  - Alkenes
  - Aromatics
- Biogenic SOA from oxidation of ...
  - Alpha-pinene
  - Limonene
  - Isoprene
- Anthropogenic POA
- Elemental carbon (soot)
- Primary PM2.5

# MADE aerosol module

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## Coarse mode:

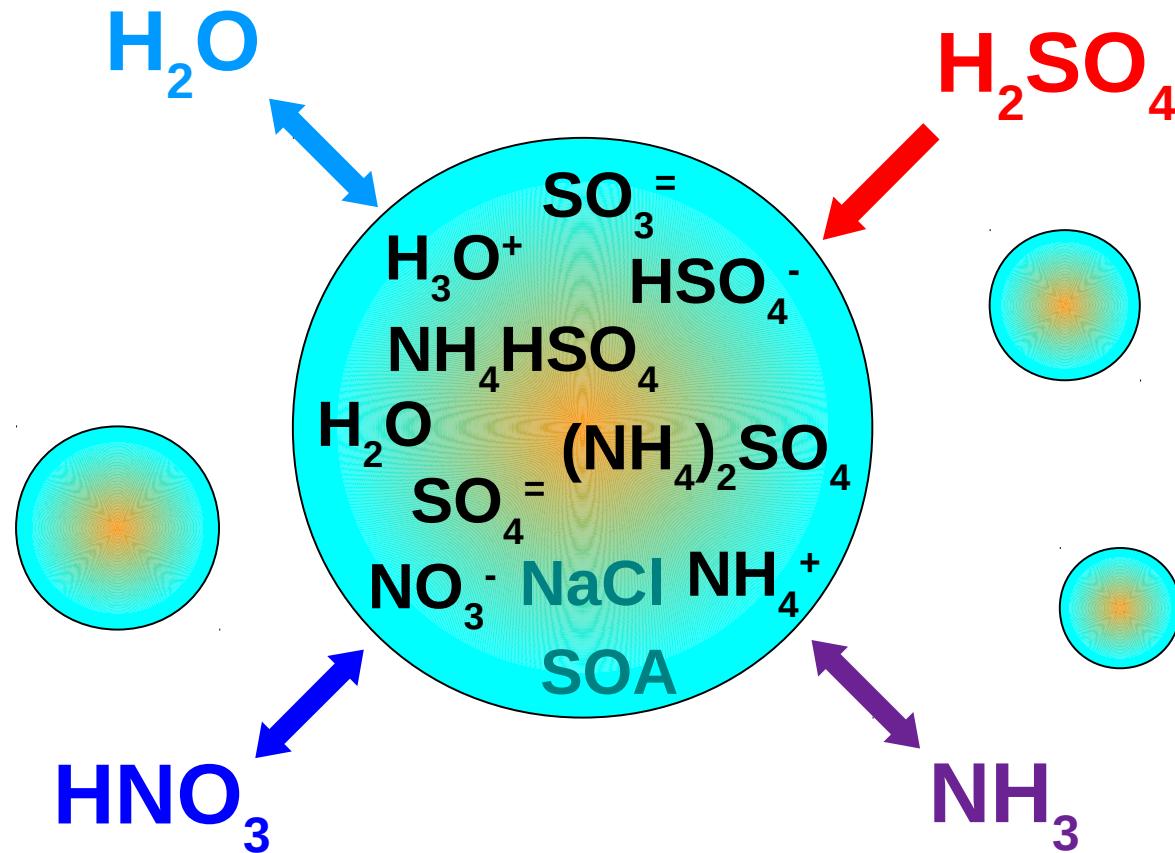
- Anthropogenic primary aerosol – e.g. from
  - Coal combustion
  - Cement manufacturing
  - Metallurgy
  - Waste incineration
- Sea salt
- Soil derived particles (mineral dust)

# MADE aerosol coupling with chemistry

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- **Gas phase chemistry:**
  - **RADM2 (Regional Acid Deposition Model version 2)**
  - **RACM (Regional Atmospheric Chemistry Mechanism)**
  - **RACM NOAA/ESRL version**
  - **CBMZ (Carbon-Bond Mechanism version Z)**
- **Gas phase/particle partitioning (aerosol chemistry):**
  - **MARS (Model for an Aerosol Reacting System)**
  - **SORGAM (Secondary Organic Aerosol Model)**
  - **VBS (Volatility Basis Set)**
- **Aqueous chemistry:**
  - CMU aqueous chemistry
  - CMAQ (EPA) aqueous chemistry
  - Only for Aitken and accumulation mode
  - Only for selected gas phase chemistry options

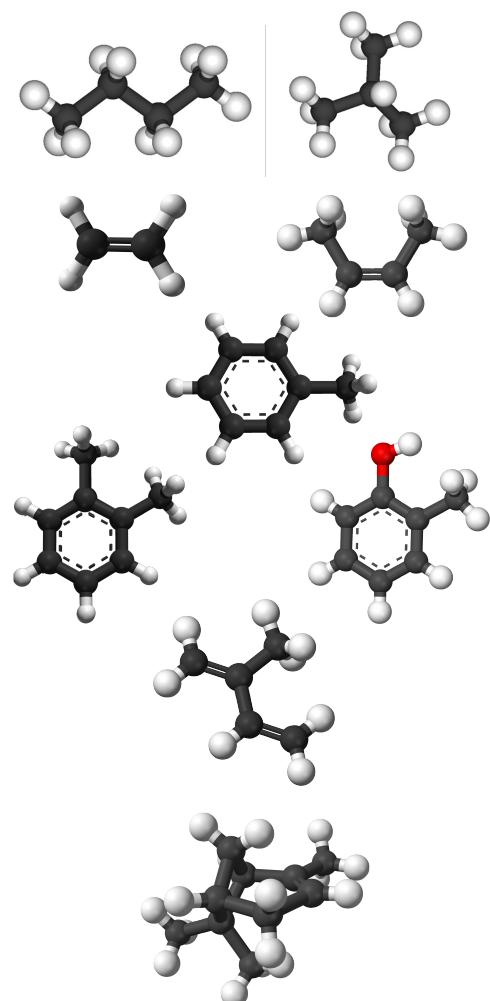
# MADE and MARS: Inorganic aerosol chemistry



MARS (Model for an Aerosol Reacting System),  
Saxena et al., Atm. Env., 1986

# MADE/SORGAM

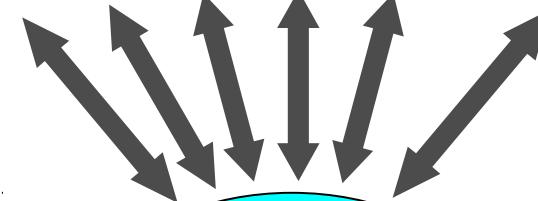
| Gas phase scheme<br>(RADM2, RACM) |  |
|-----------------------------------|--|
| Alkanes                           |  |
| Alkenes                           |  |
| Toluene                           |  |
| Xylene, cresole, ...              |  |
| Isoprene                          |  |
| Sesquiterpene                     |  |
| Alpha-pinene, limonene            |  |



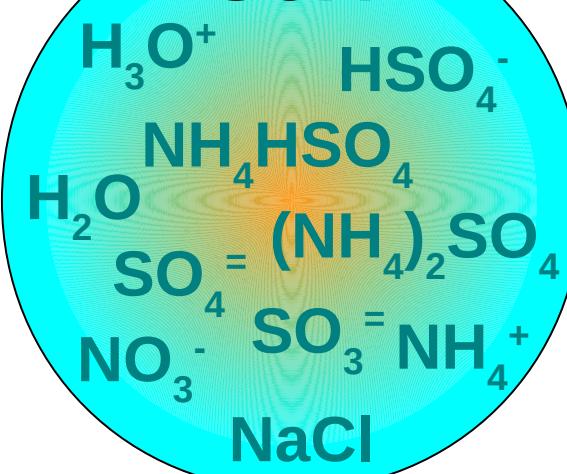
**OH, O<sub>3</sub>, NO<sub>3</sub>**

**Semi-volatile organics**

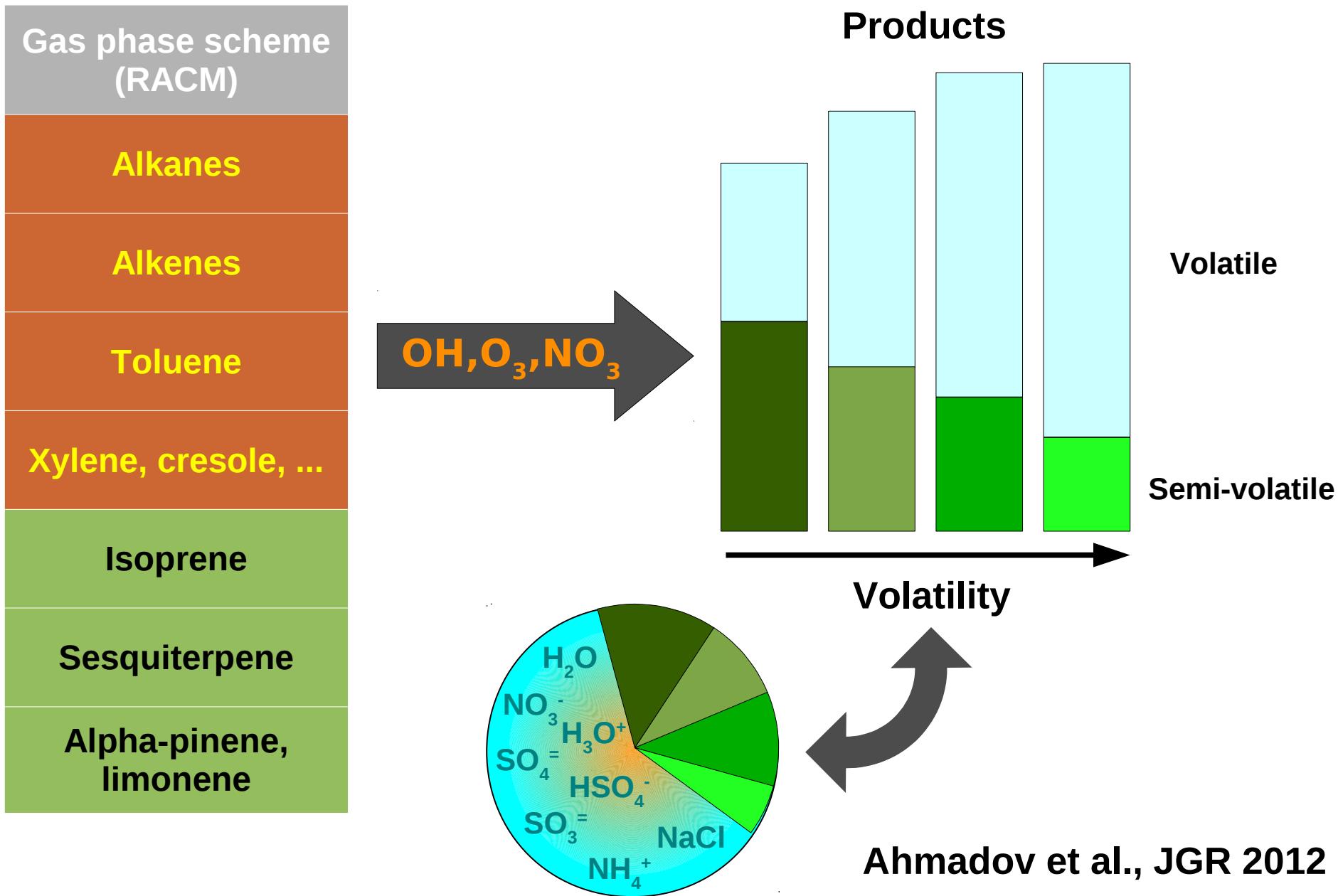
$X_1, X_2, X_3, X_4, X_5, \dots, X_n$



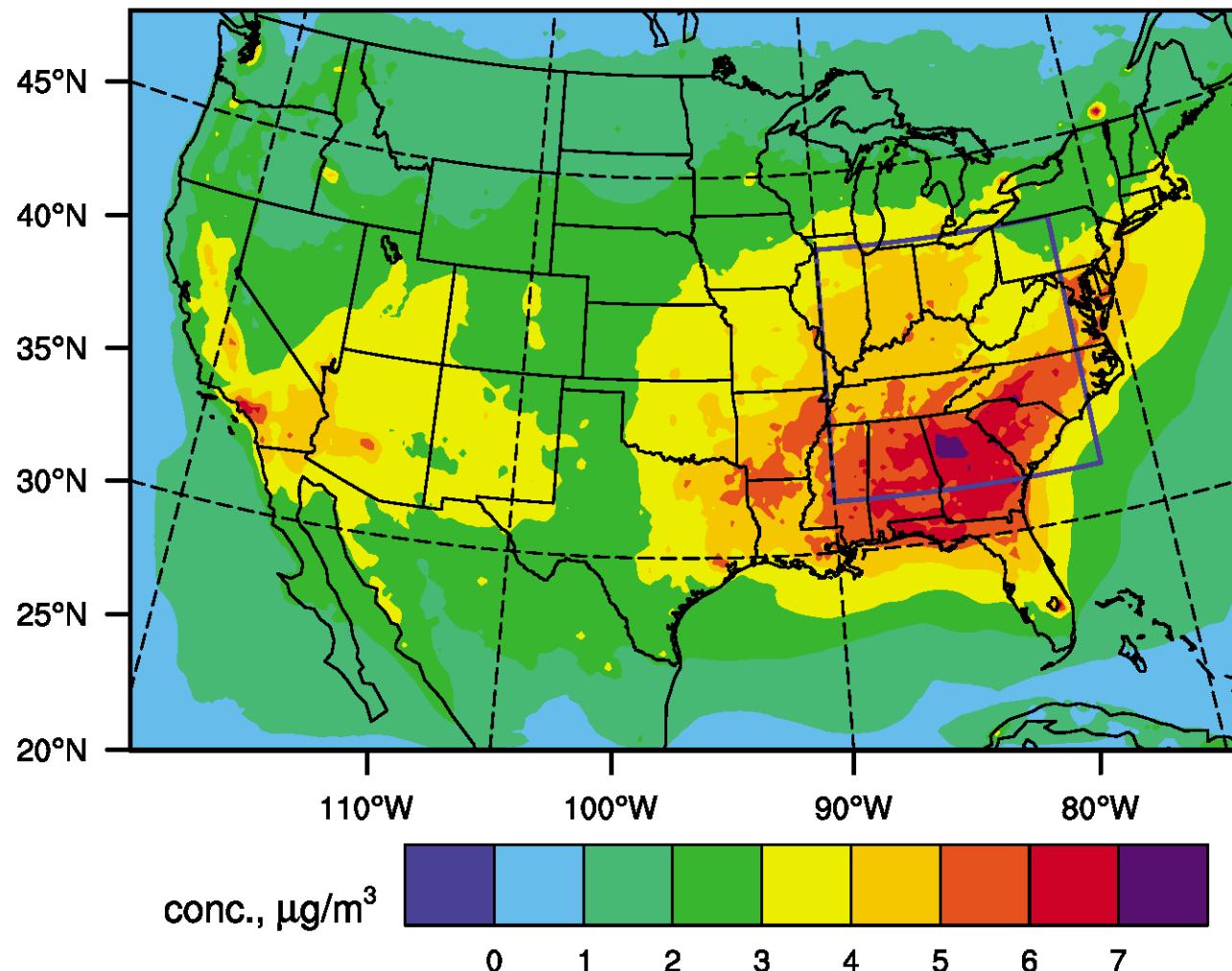
**SOA**



# MADE/VBS (Volatility Basis Set)



# MADE/VBS (Volatility Basis Set)



Organic aerosol mass in the surface layer  
(August - September 2006)

Ahmadov et al., JGR 2012

# MADE and aqueous chemistry

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- **CMU aqueous chemistry** (Fahey & Pandis, Atm. Env., 2001)
  - Slow
  - Only for resolved clouds (Sc)
  - Does not really conserve mass
- **CMAQ aqueous chemistry** (Walcek & Taylor, JAS, 1986)
  - Relatively fast
  - In both resolved (Sc) and parameterized (Cu) clouds
  - Coupled to wet deposition of  $\text{SO}_4^=$  and  $\text{NO}_3^-$

MADE + CMAQ aqueous chemistry

+ wet deposition of  $\text{SO}_4^=$  and  $\text{NO}_3^-$  :

- **chem\_opt = 41 : RADM2 gas phase chemistry**

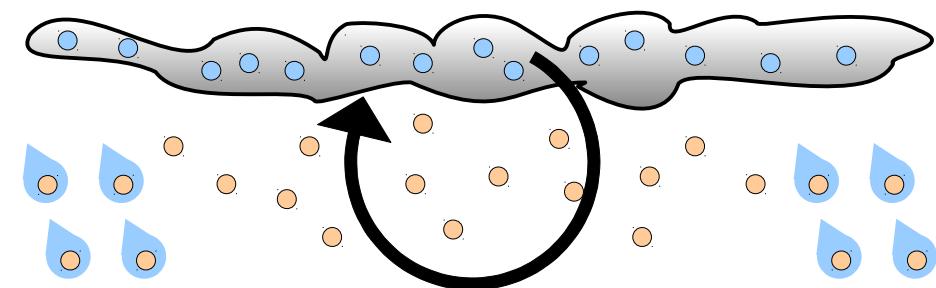
Planned for WRF/Chem 3.5.1:

- **chem\_opt = 42 : RACM gas phase chemistry (KPP)**
- **chem\_opt = 43 : RACM ESRL gas phase chemistry (KPP)**

# MADE and CMAQ aqueous chemistry

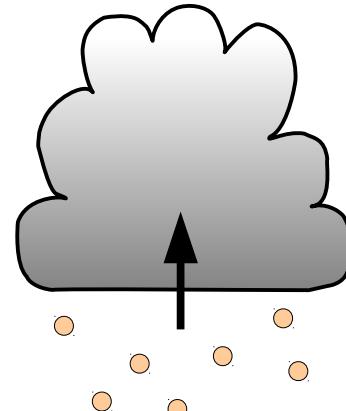
Details of aqueous chemistry depend on cloud type

Stratocumulus



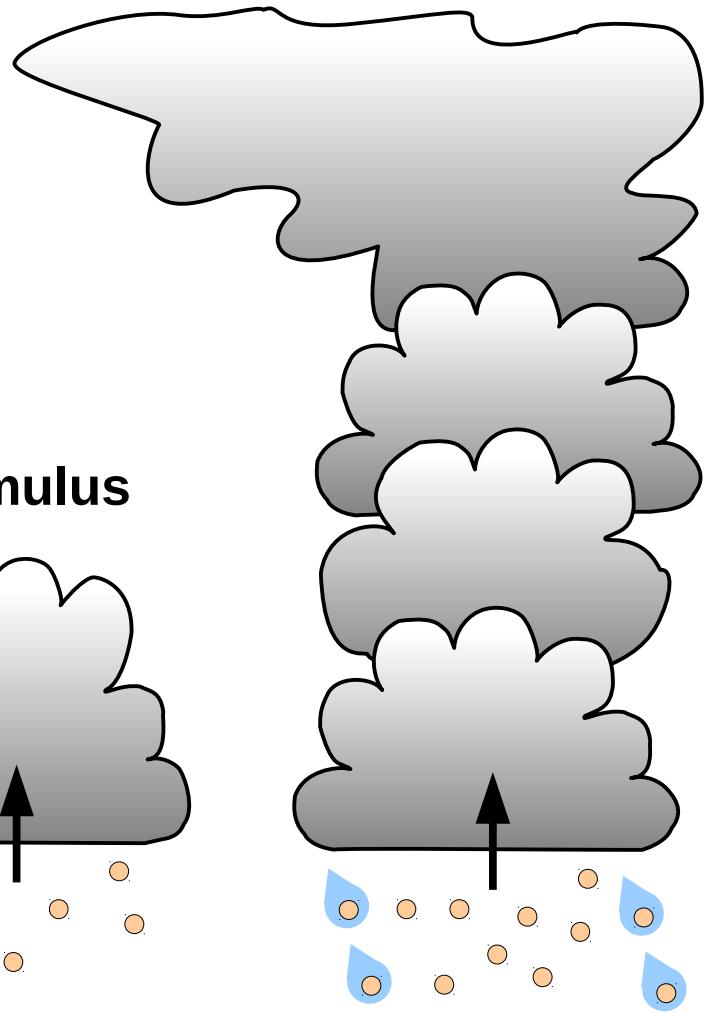
~ 10km

Cumulus



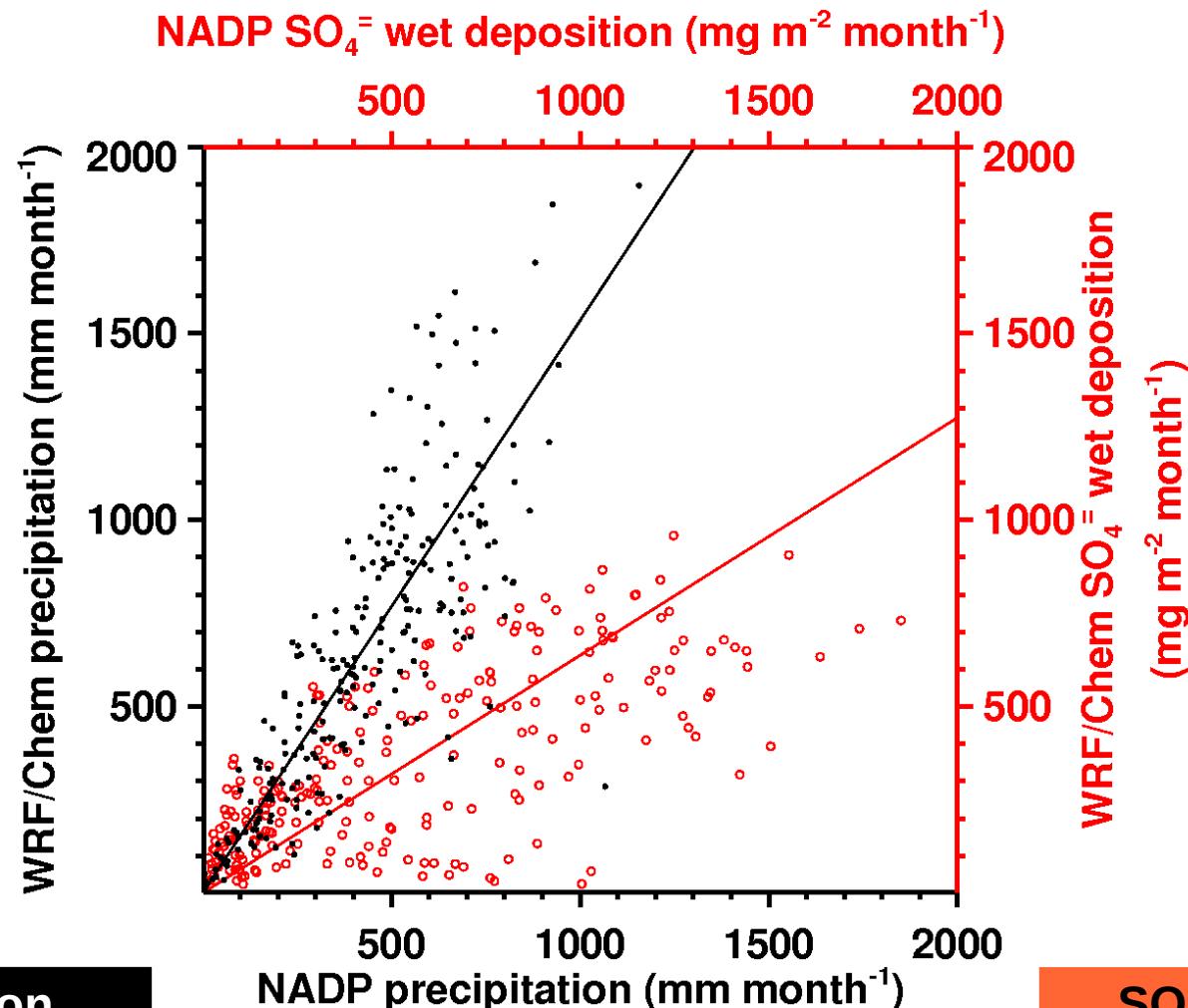
~100m

Cumulonimbus



~1km

# MADE and wet deposition



| Precipitation |            |
|---------------|------------|
| r             | model/obs. |
| 0.81          | 1.25       |

| $\text{SO}_4^{=}$ wet dep. |            |
|----------------------------|------------|
| r                          | model/obs. |
| 0.86                       | 0.53       |

May-September 2006  
(National Atmospheric Deposition Program)

# MOSAIC aerosol module

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**Model for Simulating Aerosol Interactions and Chemistry**  
(Zaveri et al., JGR, 2008)

- Modern aerosol scheme in WRF/Chem
- **4 or 8 aerosol size sections (bins) 39 nm – 10 µm**
- (Lower bin boundary of 39 nm too large for nucleation)
- **Interaction with radiation:**
  - Direct aerosol effect
  - Effect on photolysis
- **Interaction with clouds:**
  - Aerosol number determines cloud drop number and size
  - Radiative response → 1<sup>st</sup> indirect aerosol effect
  - Aqueous chemistry
  - Wet removal (scavenging)
  - **only for resolved clouds (Sc)**

# MOSAIC aerosol module

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## Aerosol composition

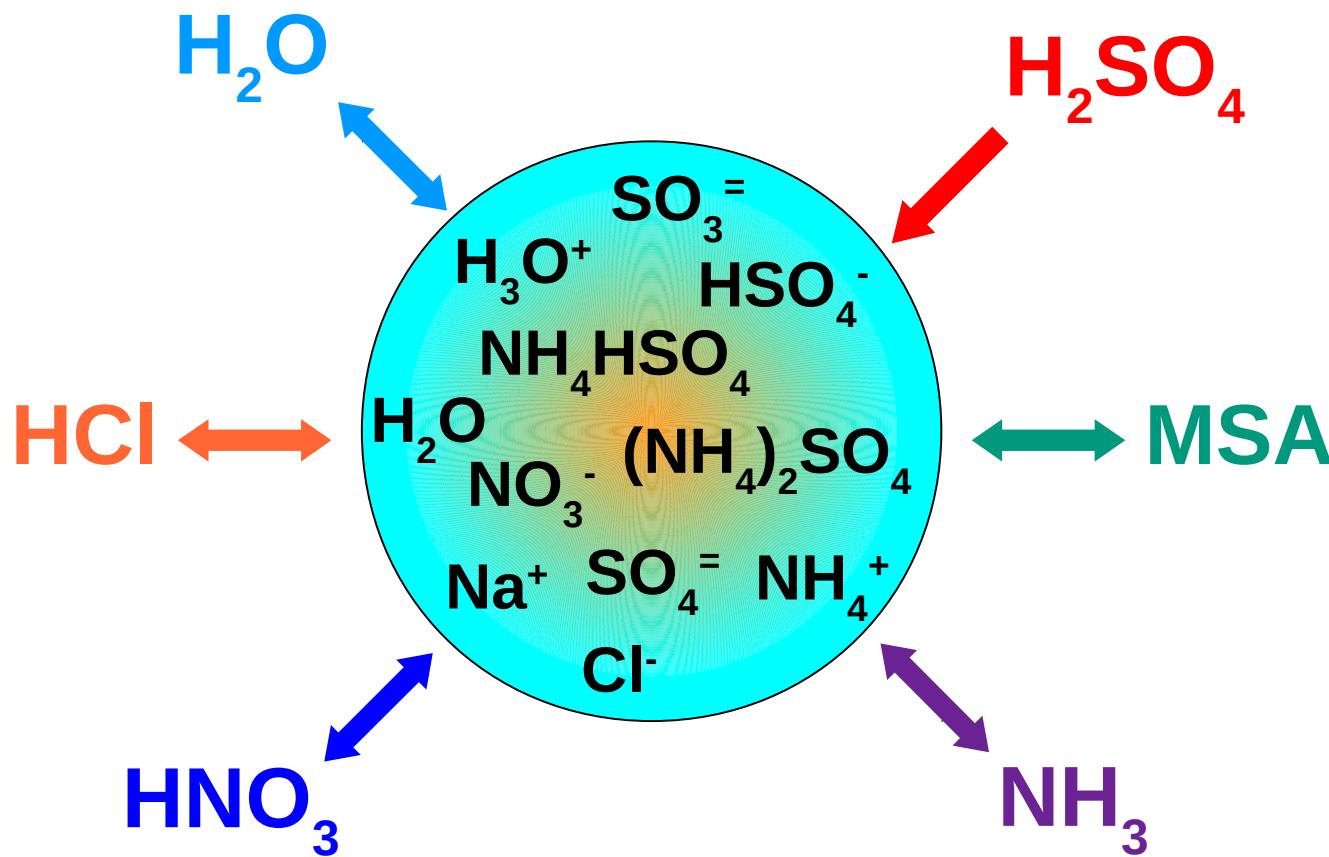
- $\text{SO}_4^{=}$ ,  $\text{NH}_4^+$ ,  $\text{NO}_3^-$ ,  $\text{H}_2\text{O}$
- NaCl (sea salt)
- $\text{CH}_3\text{SO}_3$  (methanesulfonate)
- carbonate ( $\text{CO}_3$ )
- calcium (Ca)
- black carbon (BC)
- primary organic mass (OC)
- other inorganic mass (minerals, trace metals)

# MOSAIC aerosol coupling with chemistry

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- **Gas phase chemistry:**
  - **CBMZ (Carbon-Bond Mechanism version Z)**
    - ◆ “Standard” gas phase chemical scheme for MOSAIC
  - **SAPRC99 (extensive VOC chemistry)**
    - ◆ Works with the VBS SOA scheme
  - **MOZART (Model for Ozone and Related chem. Tracers)**
    - ◆ Works with the VBS SOA scheme
- **Gas phase/particle partitioning (aerosol chemistry):**
  - **MTEM (Multicomponent Taylor Expansion Method)**
  - **MESA (Multicomponent Equilibrium Solver for Aerosols)**
  - **VBS (Volatility Basis Set)**
- **Aqueous chemistry:**
  - CMU aqueous chemistry, only for resolved clouds (Sc)
  - Not with KPP versions of gas phase chemistry schemes

# MOSAIC, MTEM, and MESA



**MTEM** calculates activity coefficients  
**MESA** solves ion-equilibria in the liquid phase  
For SOA: VBS (Volatility Basis Set) scheme

**MTEM** (Multicomponent Taylor Expansion Method), Zaveri et al., JGR 2005a

**MESA** (Multicomponent Equilibrium Solver for Aerosols), Zaveri et al., JGR 2005b

# MAM aerosol module

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## Modal Aerosol Model from CAM5

- New in WRF/Chem 3.5
- 3 or 7 log-normal aerosol modes: MAM3 and MAM7
- Mode width  $\sigma$  is fixed
- Aerosol number and mass variable
- Liu et al., Geosci. Model Dev., 5, 709-739, 2012

# MAM 3

## Aitken mode

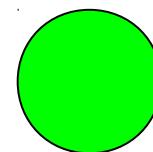


- Sulfate ( $\text{SO}_4^{=}$ )
- SOA
- Sea salt

Coagulation,  
condensation

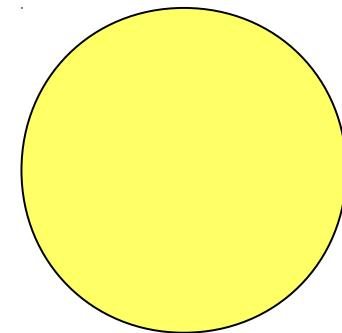


## Accumulation mode



- Sulfate ( $\text{SO}_4^{=}$ )
- SOA
- Primary organic matter
- Sea salt
- Black carbon
- Mineral dust
- Sea salt

## Coarse mode



- Sulfate ( $\text{SO}_4^{=}$ )
- Mineral dust
- Sea salt

15 – 53 nm

58 – 270 nm

0.8 – 3.65  $\mu\text{m}$

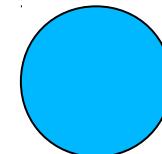
Dry particle diameter

# MAM 7

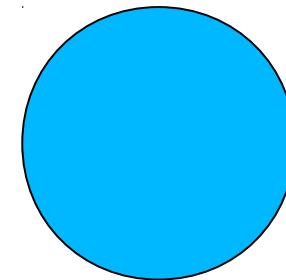
Aitken mode

Accumulation mode

Fine sea salt,  $\text{SO}_4^{=}$



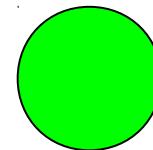
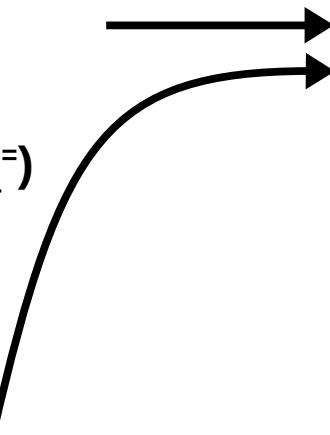
Coarse sea salt,  $\text{SO}_4^{=}$



Coagulation,  
condensation



- Sulfate ( $\text{SO}_4^{=}$ )
- SOA
- Sea salt



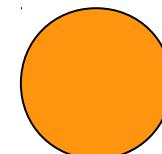
- Sulfate ( $\text{SO}_4^{=}$ )
- SOA
- Primary organic matter
- Sea salt
- Black carbon

Primary carbon

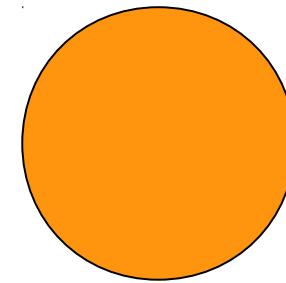


- Primary organic matter
- Black carbon

Fine dust,  $\text{SO}_4^{=}$



Coarse dust,  $\text{SO}_4^{=}$



# MAM aerosol module

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- Currently only one gas phase chemistry scheme
  - CBMZ (Carbon-Bond Mechanism version Z)
- Interaction with radiation:
  - Coupled to RRTMG radiation → Direct aerosol effect

As in CAM5:

- Gas phase/particle partitioning (aerosol chemistry):
  - Condensation of water vapor and of the 4 inorganic trace gase species:  $\text{NH}_3$ ,  $\text{H}_2\text{SO}_4$ ,  $\text{HNO}_3$ ,  $\text{HCl}$
- Interaction with clouds only resolved clouds (Sc):
  - Coupled to Morrison & Gettelman cloud microphysics
  - Radiative response → 1<sup>st</sup> indirect aerosol effect
  - Wet removal (scavenging)
  - Aqueous chemistry
- Dry deposition

# Volcanic ash

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- **10 bins for volcanic ash aerosol**
- **Transport, settling, dry deposition**
- **Currently no other aerosol**
- **Single active volcano**
- **Database of 1535 volcanoes (latitude, longitude, height)**

# How to tell WRF/Chem what to do

..../WRFV3/test/em\_real/real.exe  
| ..../WRFV3/test/em\_real/namelist.input  
  ..../WRFV3/test/em\_real/...  
  ..../WRFV3/test/em\_real/...

```
...
...
&chem
chem_opt      = 42
photdt        = 0.25
chemdt         = 0
...
aerchem_onoff = 1
...
conv_tr_aqchem = 1
```

MADE/SORGAM,  
RACM, CMAQ  
aqueous chemistry

Switches all aerosol  
processes on/off

CMAQ aqueous  
chemistry on in Cu

# Resources

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- **WRF/Chem User's Guide**
  - Model options (namelist parameters)
  - Combinations of physical/chemical schemes
  - ...
- **Papers referenced in the WRF/Chem User's Guide**
- **WRF/Chem source code**
- **WRF/Chem Help ([wrfchemhelp.gsd@noaa.gov](mailto:wrfchemhelp.gsd@noaa.gov))**
- **Yours truly ([jan.kazil@noaa.gov](mailto:jan.kazil@noaa.gov))**