

# WACCM Versions

- 3.1.9
  - Released version
- 3.5.48
  - CCMVal simulations
  - Gravity wave improvements
  - Chemistry updated to JPL-2006
  - <https://wiki.ucar.edu/display/~fvitt/WACCM+Versions>
- 3.6.16
  - WACCM/CAM-Chem codes combined (mostly)
  - More flexible

# WACCM3.6 Developments

- Combined WACCM and CAM-Chem chemistry codes
  - WACCM adopted CAM-Chem flexibilities
  - Common namelist specifications
- Configure changes
  - Invokes the chemistry preprocessor
  - Easier to change chemical mechanisms
- Flexible photolysis scheme
- Ability to specify non-transported short lived species

# WACCM/CAM-Chem Code Merge

- Share common routines
  - Easier to maintain
  - Examples:
    - Surface emissions
    - Lightning NOx production
    - Dry Deposition
  - WACCM adopted CAM-Chem flexibilities
    - Use helper functions
    - Flexible namelist specifications
    - Easier to change mechanisms

# Surface Emissions

- Easy to add/remove tracer emissions without code changes
- Can be 'CYCLICAL', 'SERIAL', or 'FIXED'

```
srf_emis_specifier =  
    'CH2O -> /inputdata/atm/wacm/emis/emis.ch2o.c080919.nc',  
    'CO    -> /inputdata/atm/wacm/emis/emis.co.c080919.nc',  
    'NO    -> /inputdata/atm/wacm/emis/emis.nox.c080919.nc'  
srf_emis_type      = 'CYCLICAL'  
srf_emis_date      = 19970101
```

# Elevated Emissions

- Ability to specify 3D time dependent emissions
- Example aircraft emissions
- Can be 'CYCLICAL', 'SERIAL', or 'FIXED'

```
ext_frc_specifier =  
    'CO    -> /inputdata/atm/waccm/emis/vrt_emis.co.nc',  
    'NO    -> /inputdata/atm/waccm/emis/vrt_emis.nox.nc'  
ext_frc_type      = 'CYCLICAL'  
ext_frc_date      = 19970101
```

# Addition Namelist Changes

## Flexible specifications:

```
drydep_list = 'NO', 'NO2', 'HNO3', 'HO2NO2', 'O3', 'H2O2',  
              'CH2O', 'CH3OOH', 'CO'
```

```
flbc_file = '../lb/LBC_1850-2100_1.9x2.5_REF2_za_c080114.nc'
```

```
flbc_list = 'N2O', 'HCFC22', 'H2', 'CO2', 'CH4', 'CH3CL',  
            'CH3CCL3', 'CH3BR', 'CFC11', 'CFC113', 'CF3BR',  
            'CF2CLBR', 'CFC12', 'CCL4'
```

```
flbc_type = 'CYCLICAL' | 'SERIAL' | 'FIXED'
```

```
flbc_date = 20000101
```

```
sad_file = '../sulf/SAD_SULF_1850-2100_1.9x2.5_c080220.nc'
```

```
sad_type = 'CYCLICAL' | 'SERIAL' | 'FIXED'
```

```
sad_date = 20000101
```

# Changes to Configure Utility

- Easier to customize chemical mechanism
- Invokes the chemistry preprocessor
- Option added:
  - **-usr\_mech\_infile \$mechanism\_file**
    - Allows user to specify a customized preprocessor input file
    - Determines the number of advected tracers
- Ability to configure with non-transported chemical tracers via preprocessor

## Preprocessor input file

```

List solution species      Solution
                           End Solution

List prescribed species   Fixed
                           End Fixed

List species not transported Not-transported
                           End Not-transported

                           END SPECIES

                           SOLUTION CLASSES

Longer lived species      Explicit
                           End Explicit

Shorter lived species     Implicit
                           End Implicit

                           END SOLUTION CLASSES

                           CHEMISTRY

Photolysis reactions      Photolysis
                           End Photolysis

Gas phase reactions       Reactions
                           End Reactions

Gas phase wet removal     Heterogeneous
                           End Heterogeneous

Prescribed prod/loss rates Ext Forcing
                           End Ext Forcing

                           END CHEMISTRY
```



# Non-transported chemical tracers

mechanism file:

```
SPECIES  
  
  . . .  
  
  NOT-TRANSPORTED  
    OH, e, ..  
  END NOT-TRANSPORTED  
  
END SPECIES
```

build script:

```
$cfgdir/configure \  
  -chem waccm_mozart \  
  -usr_mech_infile $mechanism_file \  
  -usr_mech_outfile $mechanism_file \  
  -usr_mech_infile $mechanism_file \  
  -usr_mech_outfile $mechanism_file \  
  -usr_mech_infile $mechanism_file \  
  -usr_mech_outfile $mechanism_file \  
  -usr_mech_infile $mechanism_file \  
  -usr_mech_outfile $mechanism_file \  
  -usr_mech_infile $mechanism_file \  
  -usr_mech_outfile $mechanism_file
```

# Flexible Photolysis

- Easier to change photolysis
- Only the cross sections needed are read in
- Re-assignment of rates are straight forward

```
      Photolysis
[ jo2_a=userdefined, ] O2 + hv -> O + O1D
[ jo2_b=userdefined, ] O2 + hv -> 2*O
[ jo3_a ]              O3 + hv -> O1D + O2_1D
[ jo3_b ]              O3 + hv -> O + O2
[ jn2o ]              N2O + hv -> O1D + N2
...
[ jh2o2 ]             H2O2 + hv -> 2*OH
[ jch3ooh ]           CH3OOH + hv -> CH2O + H + OH
[ jpooh->,jch3ooh ]   POOH + hv -> CH3CHO + CH2O + HO2 + OH
[ jch3co3h->,0.28*jh2o2 ] CH3COOOH + hv -> CH3O2 + OH + CO2
...
[ jeuv_1=userdefined,userdefined ] O + hv -> Op + e
[ jeuv_2=userdefined,userdefined ] O + hv -> Op + e
...
      End Photolysis
```