

MICM chemistry solvers

Advance the chemical concentrations from TimeStart to TimeEnd, based on the ODE derived from the chemical reactions in the mechanism.

User may choose between either of two solver schemes

- chemistry_driver_ros Rosenbrock (Similar to what is used in some WRF configurations)
- chemistry_driver_moz Mozart (Similar to what is used in CESM with an error control on timestep)

The inputs/outputs are identical for both of them, however runtime options will allow choices of tolerance and configurations of those solvers.

Source code locations

- MICM_chemistry/src/chemistry_driver_ros.F90
- MICM_chemistry/src/chemistry_driver_moz.F90

Metadata for chemistry_driver routines

chemistry_driver_xxx_init routine (will show mozart, but Rosenbrock values are identical)

!> \section arg_table_chemistry_driver_moz_init Argument Table

!! local_name standard_name					long_name
units rank type	kind	intent	optional		
!! ----- ----- ----- ----- ----- -----					
!! TimeStart chem_step_start_time					Chem step start time
s 0 real	kind_phys	in	F		
!! TimeEnd chem_step_end_time					Chem step end time
s 0 real	kind_phys	in	F		
!! dt time_step_for_physics					time_step_for_physics
s 0 real	kind_phys	in	F		
!! errmsg ccpp_error_message					CCPP error message
none 0 character	len=512	out	F		
!! errflg ccpp_error_flag					CCPP error flag
flag 0 integer		out	F		
!!					

chemistry_driver_xxx_run routine (will show mozart, but Rosenbrock values are identical)

!> \section arg_table_chemistry_driver_moz_run Argument Table

!! local_name standard_name					long_name
units rank type	kind	intent	optional		
!! ----- ----- ----- ----- ----- -----					
!! vmr concentration					species concentration
mole/mole 1 real	kind_phys	inout	F		
!! TimeStart chem_step_start_time					Chem step start time
s 0 real	kind_phys	in	F		
!! TimeEnd chem_step_end_time					Chem step end time
s 0 real	kind_phys	in	F		
!! j_rateConst photo_rate_constants					photochemical rates constants
1 1 real	kind_phys	in	F		s-
!! k_rateConst gasphase_rate_constants					k rate constants
1 1 real	kind_phys	in	F		s-
!! errmsg ccpp_error_message					CCPP error message
none 0 character	len=512	out	F		
!! errflg ccpp_error_flag					CCPP error flag
flag 0 integer		out	F		
!!					

chemistry_driver_xxx_finalize routine (will show mozart, but Rosenbrock values are identical)

