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Interactive Comment

Interactive comment on "A numerical evaluation of global oceanic emissions of α -pinene and isoprene" by G. Luo and F. Yu

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The authors thank the editor for the helpful comments. Our responses to the comments are given below.

As pointed out during this online review process, the lifetime of α -pinene is longer than that of isoprene. The boundary layer budget equation imposes an additional physical constraint on the relative ratio between isoprene and α -pinene fluxes. [E.g. (Flux(α pinene)/Flux(isoprene) ~ (tau(isoprene)*C(α -pinene))/((tau(α -pinene)*C(isoprene))]. Assuming comparable lifetimes (tau) and concentrations (C), the α -pinene flux would be expected to be in the same range as the isoprene flux. It needs to be explained why the modeled α -pinene flux is more than a factor of 10 higher than the modeled



isoprene flux, while their observed concentrations are very similar.

The explanation for the difference between the global oceanic emissions of α -pinene and isoprene has been discussed in the reply to referee 1. Detailed budget analysis for selected locations has been provided in the supplement.

Also. unable find reference material to how was on monoterpene oxidation is incorporated in GEOS-CHEM (e.g. http://acmg.seas.harvard.edu/geos/wiki docs/chemistry/chemistry_updates_v6.pdf). It will be helpful to the reader to include either a reference or an appendix that documents the chemical reactions which are considered for a-pinene/monoterpene oxidation.

In GESO-Chem, the monoterpene oxidation is treated in the code carbon_mod.f, while the isoprene oxidation is treated in SMVGEAR solver. The reaction rates of α -pinene with OH, O₃ and NO₃ are calculated as following:

```
KO3 = 56.15e-18*EXP( ACT_O3 * ( 1.0/298.0 - 1.0/T ) )
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KOH = 84.4e-12* EXP( ACT_OH * (1.0/298.0 - 1.0/T ) )
```

```
KNO3 = 6.95e-12* EXP( ACT_NO3 * (1.0/298.0 - 1.0/T ) )
```

where ACT_O3= 732.0, ACT_OH= -400.0 and ACT_NO3=-490.0, and T represent temperature in air.

The reaction rates of isoprene in GEOS-Chem with OH, O_3 and NO_3 are calculated as following:

KO3 = 1.05d-14 * EXP(-2000.d0 / T) KOH = 2.70d-11 * EXP(390.d0 / T) KNO3 = 3.03d-12 * EXP(-446.d0 / T)

The above reaction rates have been described in the supplement.

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