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Title:	Modelling atmospheric OH-reactivity in a boreal forest ecosystem
Authors:	Sinha, V. (/jspui/browse?type=author&value=Sinha%2C+V.)
Keywords:	Atmospheric chemistry Atmospheric modeling
Issue Date:	2012
Citation:	Atmospheric Chemistry and Physics, 11 (18), pp. 9709-9719.
Abstract:	We have modelled the total atmospheric OH-reactivity in a boreal forest and investigated the individual contributions from gas phase inorganic species, isoprene, monoterpenes, and methane along with other important VOCs. Daily and seasonal variation in OH-reactivity for the year 2008 was examined as well as the vertical OH-reactivity profile. We have used SOSA; a one dimensional vertical chemistry-transport model (Boy et al., 2011a) together with measurements from Hyytiälä, SMEAR II station, Southern Finland, conducted in August 2008. Model simulations only account for ~30-50% of the total measured OH sink, and in our opinion, the reason for missing OH-reactivity is due to unmeasured unknown BVOCs, and limitations in our knowledge of atmospheric chemistry including uncertainties in rate constants. Furthermore, we found that the OH-reactivity correlates with both organic and inorganic compounds and increases during summer. The summertime canopy level OH-reactivity peaks during night and the vertical OH-reactivity decreases with height.
URI:	<a href="http://www.atmos-chem-phys.net/11/9709/2011/acp-11-9709-2011.html">http://www.atmos-chem-phys.net/11/9709/2011/acp-11-9709-2011.html</a> ( <a href="http://www.atmos-chem-phys.net/11/9709/2011/acp-11-9709-2011.html">http://www.atmos-chem-phys.net/11/9709/2011/acp-11-9709-2011.html</a> )
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