Atmos. Chem. Phys. Discuss., 11, C12949–C12949, 2011 www.atmos-chem-phys-discuss.net/11/C12949/2011/

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Interactive comment on "Structures and reaction rates of the gaseous oxidation of SO_2 by an $O_3^-(H_2O)_{0-5}$ cluster – a density functional theory investigation" by N. Bork et al.

Anonymous Referee #2

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The comment was uploaded in the form of a supplement: http://www.atmos-chem-phys-discuss.net/11/C12949/2011/acpd-11-C12949-2011-supplement.pdf

Interactive comment on Atmos. Chem. Phys. Discuss., 11, 29647, 2011.

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