

Interactive comment on “Structures and reaction rates of the gaseous oxidation of SO₂ by an O₃[−](H₂O)_{0–5} cluster – a density functional theory investigation” by N. Bork et al.

Anonymous Referee #2

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<http://www.atmos-chem-phys-discuss.net/11/C12949/2011/acpd-11-C12949-2011-supplement.pdf>

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