

# ***Interactive comment on “Molecular Dynamics Simulation of the Surface Tension of Aqueous Sodium Chloride: from Dilute to Highly Supersaturated Solutions and Molten Salt” by Xiaoxiang Wang et al.***

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Received and published: 29 October 2018

Dear Editor and Reviewers,

Many thanks for the kind effort and help in the peer review process. Our responses to the comments are provided in the attached file. The revised manuscript is also provided.

Please also note the supplement to this comment:

<https://www.atmos-chem-phys-discuss.net/acp-2017-1013/acp-2017-1013-AC1-supplement.pdf>

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Interactive comment on Atmos. Chem. Phys. Discuss., <https://doi.org/10.5194/acp-2017-1013>, 2017.

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