

***Interactive comment on “HULIS in nanoaerosol clusters; investigations of surface tension and aggregate formation using molecular dynamics simulations” by T. Hede et al.***

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The comment was uploaded in the form of a supplement:  
<http://www.atmos-chem-phys-discuss.net/11/C4552/2011/acpd-11-C4552-2011-supplement.pdf>

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