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DOI: 10.1021/ct800229h

Journal of Chemical Theory and Computation, 2008, 4, 1829-3

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227	Benzene Dimer: Dynamic Structure and Thermodynamics Derived from On-the-Fly ab initio DFT-D Molecular Dynamic Simulations. <i>Journal of Chemical Theory and Computation</i> , 2008 , 4, 1835-40	6.4	36
226	Stacking of polycyclic aromatic hydrocarbons as prototype for graphene multilayers, studied using density functional theory augmented with a dispersion term. <i>Journal of Chemical Physics</i> , 2009 , 131, 1947-52	3.9	41
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