

Vesa Hänninen

List of Publications by Year in descending order

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Version: 2024-02-01

10
papers

122
citations

1937685

4
h-index

1588992

8
g-index

10
all docs

10
docs citations

10
times ranked

132
citing authors

#	ARTICLE	IF	CITATIONS
1	An Ab Initio Molecular Dynamics Study of the Hydrolysis Reaction of Sulfur Trioxide Catalyzed by a Formic Acid or Water Molecule. <i>Journal of Physical Chemistry A</i> , 2020, 124, 1922-1928.	2.5	1
2	Photoacoustic Optical Frequency Comb Spectroscopy of Radioactive Methane in the Mid-Infrared Region. , 2019, , .		0
3	Ab Initio Molecular Dynamics Simulations of the Influence of Lithium Bromide on the Structure of the Aqueous Solution's Air Interface. <i>Journal of Physical Chemistry B</i> , 2019, 123, 729-737.	2.6	4
4	Ab initio molecular dynamics studies of formic acid dimer colliding with liquid water. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 23717-23725.	2.8	15
5	Effects of Global and Local Anharmonicities on the Thermodynamic Properties of Sulfuric Acid Monohydrate. <i>Journal of Chemical Theory and Computation</i> , 2016, 12, 5511-5524.	5.3	3
6	Generalized Intermolecular Interaction Tensor Applied to Long-Range Interactions in Hydrogen and Coinage Metal (Cu, Ag, and Au) Clusters. <i>Journal of Physical Chemistry A</i> , 2015, 119, 11729-11736.	2.5	2
7	The effect of large amplitude motions on the vibrational intensities in hydrogen bonded complexes. <i>Journal of Chemical Physics</i> , 2015, 142, 094304.	3.0	39
8	The effect of large amplitude motions on the transition frequency redshift in hydrogen bonded complexes: A physical picture. <i>Journal of Chemical Physics</i> , 2014, 140, 184309.	3.0	47
9	An ab Initio Study of van der Waals Potential Energy Parameters for Silver Clusters. <i>Journal of Physical Chemistry A</i> , 2011, 115, 2332-2339.	2.5	6
10	Path sampling for atmospheric reactions: formic acid catalysed conversion of $\text{SO}_3 + \text{H}_2\text{O}$ to H_2SO_4 . <i>PeerJ Physical Chemistry</i> , 0, 2, e7.	0.0	5