

Kono H Lemke

List of Publications by Year in descending order

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14
papers

500
citations

1040056

9
h-index

1058476

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g-index

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all docs

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docs citations

14
times ranked

668
citing authors

#	ARTICLE	IF	CITATIONS
1	Structure and solvation dynamics of the hydroxide ion in ice-like water clusters: a CCSD(T) and carâ€“parrinello molecular dynamics study. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 18990-18998.	2.8	1
2	The ozoneâ€“water complex: CCSD(T)/CBS structures and anharmonic vibrational spectroscopy of O ₃ (H ₂ O) (ν ₁ = 1 \hat{a} 2). <i>Journal of Chemical Physics</i> , 2020, 153, 084302.	3.0	2
3	Molecular Clusters and Solvation in Volcanic and Hydrothermal Vapors. <i>Reviews in Mineralogy and Geochemistry</i> , 2018, 84, 57-83.	4.8	5
4	Structure and binding energy of the H ₂ S dimer at the CCSD(T) complete basis set limit. <i>Journal of Chemical Physics</i> , 2017, 146, 234301.	3.0	23
5	ESI FT-MS and DFT studies of aqueous Au(I) complexes with cysteine and its disulfide form cystine. <i>International Journal of Mass Spectrometry</i> , 2016, 406, 12-19.	1.5	4
6	Calculations of the water-dimer encapsulations into C ₈₄ . <i>Fullerenes Nanotubes and Carbon Nanostructures</i> , 2016, 24, 1-7.	2.1	15
7	Systematic testing of Gaussian and complete basis set methods with dispersion corrections for environmentally relevant clusters. <i>Chemical Physics Letters</i> , 2014, 615, 50-55.	2.6	1
8	Gold chloride clusters with Au(III) and Au(I) probed by FT-ICR mass spectrometry and MP2 theory. <i>Physical Chemistry Chemical Physics</i> , 2014, 16, 7813.	2.8	10
9	Thermodynamic properties of carbon dioxide clusters by M06-2X and dispersion-corrected B2PLYP-D theory. <i>Chemical Physics Letters</i> , 2013, 573, 19-23.	2.6	30
10	Peptide Synthesis in Early Earth Hydrothermal Systems. <i>Astrobiology</i> , 2009, 9, 141-146.	3.0	50
11	Ab initio investigation of the structure, stability, and atmospheric distribution of molecular clusters containing H ₂ O, CO ₂ , and N ₂ O. <i>Journal of Geophysical Research</i> , 2008, 113, .	3.3	15
12	Solvation processes in steam: Ab initio calculations of ionâ€“solvent structures and clustering equilibria. <i>Geochimica Et Cosmochimica Acta</i> , 2008, 72, 3293-3310.	3.9	16
13	Mass spectrometric and quantum chemical determination of proton water clustering equilibria. <i>Geochimica Et Cosmochimica Acta</i> , 2007, 71, 2436-2447.	3.9	21
14	Extreme accumulation of nucleotides in simulated hydrothermal pore systems. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2007, 104, 9346-9351.	7.1	307