

Mychel E Varner

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/1262386/publications.pdf>

Version: 2024-02-01

19
papers

902
citations

623734

14
h-index

794594

19
g-index

20
all docs

20
docs citations

20
times ranked

1233
citing authors

#	ARTICLE	IF	CITATIONS
1	Concerted transfer of multiple protons in acid-water clusters: $[(\text{HCl})(\text{H}_2\text{O})_2]$ and $[(\text{HF})(\text{H}_2\text{O})_2]$. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 20641-20646.	2.8	3
2	Reactions of Methanesulfonic Acid with Amines and Ammonia as a Source of New Particles in Air. <i>Journal of Physical Chemistry B</i> , 2016, 120, 1526-1536.	2.6	115
3	Computational Studies of Atmospherically-Relevant Chemical Reactions in Water Clusters and on Liquid Water and Ice Surfaces. <i>Accounts of Chemical Research</i> , 2015, 48, 399-406.	15.6	89
4	New particle formation and growth from methanesulfonic acid, trimethylamine and water. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 13699-13709.	2.8	88
5	Amine-Amine Exchange in Ammonium-Methanesulfonate Aerosols. <i>Journal of Physical Chemistry C</i> , 2014, 118, 29431-29440.	3.1	31
6	Raman spectroscopy of solutions and interfaces containing nitrogen dioxide, water, and 1,4 dioxane: Evidence for repulsion of surface water by NO ₂ gas. <i>Journal of Chemical Physics</i> , 2014, 140, 184702.	3.0	3
7	Reaction of a charge-separated ONONO ₂ species with water in the formation of HONO: an MP2 Molecular Dynamics study. <i>Physical Chemistry Chemical Physics</i> , 2014, 16, 4483.	2.8	31
8	Ab initio and semi-empirical Molecular Dynamics simulations of chemical reactions in isolated molecules and in clusters. <i>Physical Chemistry Chemical Physics</i> , 2014, 16, 9760-9775.	2.8	35
9	Nitrogen dioxide at the air-water interface: trapping, absorption, and solvation in the bulk and at the surface. <i>Physical Chemistry Chemical Physics</i> , 2013, 15, 204-212.	2.8	33
10	Photooxidation of Ammonia on TiO ₂ as a Source of NO and NO ₂ under Atmospheric Conditions. <i>Journal of the American Chemical Society</i> , 2013, 135, 8606-8615.	13.7	72
11	Simplified mechanism for new particle formation from methanesulfonic acid, amines, and water via experiments and ab initio calculations. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2012, 109, 18719-18724.	7.1	173
12	Isomerization and ionization of N ₂ O ₄ on model ice and silica surfaces. <i>Chemical Physics</i> , 2012, 405, 52-59.	1.9	13
13	Dissociation Energy of the HOOO Radical. <i>Journal of Physical Chemistry A</i> , 2009, 113, 11238-11241.	2.5	46
14	On the geometry of the HO ₃ radical. <i>Chemical Physics</i> , 2008, 346, 53-55.	1.9	32
15	Vibrational overtone spectrum of matrix isolated cis, cis-HOONO. <i>Journal of Chemical Physics</i> , 2007, 126, 174308.	3.0	10
16	Molecular dynamics simulations of human Formula: the role of modified bases in mRNA recognition. <i>Nucleic Acids Research</i> , 2006, 34, 5361-5368.	14.5	37
17	A comparison of experimental and calculated spectra of HNO ₃ in the near-infrared using Fourier transform infrared spectroscopy and vibrational perturbation theory. <i>Journal of Chemical Physics</i> , 2006, 124, 124323.	3.0	33
18	Propargyl Radical: Ab Initio Anharmonic Modes and the Polarized Infrared Absorption Spectra of Matrix-Isolated HCCCH ₂ . <i>Journal of Physical Chemistry A</i> , 2005, 109, 3812-3821.	2.5	55

#	ARTICLE	IF	CITATIONS
19	Oxidation mechanism of ammonia in water clusters. <i>Molecular Physics</i> , 0, , .	1.7	0