

Jennifer Meyer

List of Publications by Year in descending order

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26
papers

730
citations

687335

13
h-index

552766

26
g-index

26
all docs

26
docs citations

26
times ranked

644
citing authors

#	ARTICLE	IF	CITATIONS
1	Influence of the leaving group on the dynamics of a gas-phase SN2 reaction. <i>Nature Chemistry</i> , 2016, 8, 151-156.	13.6	116
2	Spin and Orbital Magnetic Moments of Free Nanoparticles. <i>Physical Review Letters</i> , 2011, 107, 233401.	7.8	86
3	Ion-Induced Molecule Reaction Dynamics. <i>Annual Review of Physical Chemistry</i> , 2017, 68, 333-353.	10.8	61
4	Imaging dynamic fingerprints of competing E2 and SN2 reactions. <i>Nature Communications</i> , 2017, 8, 25.	12.8	59
5	The spin and orbital contributions to the total magnetic moments of free Fe, Co, and Ni clusters. <i>Journal of Chemical Physics</i> , 2015, 143, 104302.	3.0	54
6	Imaging the dynamics of ion-induced molecule reactions. <i>Chemical Society Reviews</i> , 2017, 46, 7498-7516.	38.1	51
7	Atomistic dynamics of elimination and nucleophilic substitution disentangled for the $F\dot{\sim} + CH_3CH_2Cl$ reaction. <i>Nature Chemistry</i> , 2021, 13, 977-981.	13.6	43
8	Stretching vibration is a spectator in nucleophilic substitution. <i>Science Advances</i> , 2018, 4, eaas9544.	10.3	37
9	Conservation of direct dynamics in sterically hindered $S_N2/E2$ reactions. <i>Chemical Science</i> , 2018, 9, 693-701.	7.4	31
10	Imaging Proton Transfer and Dihalide Formation Pathways in Reactions of $F\dot{\sim} + CH_3I$. <i>Journal of Physical Chemistry A</i> , 2016, 120, 4711-4719.	2.5	26
11	Nucleophilic substitution with two reactive centers: The $CN\dot{\sim} + CH_3I$ case. <i>Journal of Chemical Physics</i> , 2015, 143, 184309.	3.0	20
12	Unexpected Indirect Dynamics in Base-Induced Elimination. <i>Journal of the American Chemical Society</i> , 2019, 141, 20300-20308.	13.7	19
13	X-ray absorption spectroscopy of mass-selected transition metal clusters using a cyclotron ion trap: An experimental setup for measuring XMCD spectra of free clusters. <i>Journal of Electron Spectroscopy and Related Phenomena</i> , 2011, 184, 113-118.	1.7	18
14	Influence of Vibrational Excitation on the Reaction of $F\dot{\sim}$ with CH_3I : Spectator Mode Behavior, Enhancement, and Suppression. <i>Journal of Physical Chemistry Letters</i> , 2020, 11, 4331-4336.	4.6	15
15	Orbit and spin resolved magnetic properties of size selected $[Co_nRh]^{+}$ and $[Co_nAu]^{+}$ nanoalloy clusters. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 28372-28378.	2.8	13
16	Imaging Reaction Dynamics of $F\dot{\sim}(H_2O)$ and $Cl\dot{\sim}(H_2O)$ with CH_3I . <i>Journal of Physical Chemistry A</i> , 2020, 124, 1929-1939.	2.5	13
17	Imaging state-to-state reactive scattering in the $Ar^+ + H_2$ charge transfer reaction. <i>Journal of Chemical Physics</i> , 2017, 147, 013940.	3.0	11
18	Threshold photodetachment spectroscopy of the astrochemical anion $CN\dot{\sim}$. <i>Journal of Chemical Physics</i> , 2020, 153, 184309.	3.0	11

#	ARTICLE	IF	CITATIONS
19	Anionic Carbon Chain Growth in Reactions of C_2H^+ , C_4H^+ , and C_6H^+ with C_2H_2 . <i>Astrophysical Journal</i> , 2019, 878, 162.	4.5	8
20	Proton transfer dynamics modified by CH-stretching excitation. <i>Physical Chemistry Chemical Physics</i> , 2020, 22, 12382-12388.	2.8	8
21	Atomic and electronic structure of free niobium nanoclusters: Simulation of the M 4,5 -XANES spectrum of Nb 13+. <i>Journal of Electron Spectroscopy and Related Phenomena</i> , 2014, 195, 189-194.	1.7	7
22	$Fa^+(H_2O)+CH_3I$ ligand exchange reaction dynamics. <i>Chinese Journal of Chemical Physics</i> , 2020, 33, 210-216.	1.3	7
23	Dynamics of proton transfer from ArH^+ to CO. <i>International Journal of Mass Spectrometry</i> , 2019, 438, 175-185.	1.5	5
24	Methotrexate-gelonin conjugate as an inhibitor of MCF-7 cells expressing the dihydrofolate receptor. <i>Biological Chemistry</i> , 2014, 395, 1461-1466.	2.5	4
25	Suppression of low product kinetic energies in reactions of FHO^+ and $Cl^+(H_2O)$ with CH_3I . <i>International Journal of Mass Spectrometry</i> , 2021, 462, 116526.	1.5	4
26	Mn^{12} Acetate Complexes Studied as Single Molecules. <i>Chemistry - A European Journal</i> , 2022, 28, .	3.3	3