

Joshua J Melko

List of Publications by Year in descending order

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papers

649
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567281

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#	ARTICLE	IF	CITATIONS
1	Mechanisms of sequential ion-molecule reactions in protonated methanol using mass spectrometry, ab initio methods, and statistical modeling. <i>Chemical Physics</i> , 2019, 525, 110420.	1.9	5
2	Kinetics of CO ⁺ and CO ₂ ⁺ with N and O atoms. <i>Journal of Chemical Physics</i> , 2018, 148, 084305.	3.0	13
3	Temperature and Isotope Dependent Kinetics of Nickel-Catalyzed Oxidation of Methane by Ozone. <i>Journal of Physical Chemistry A</i> , 2018, 122, 6655-6662.	2.5	12
4	Determining Rate Constants and Mechanisms for Sequential Reactions of Fe ⁺⁺ with Ozone at 500 K. <i>Journal of Physical Chemistry A</i> , 2017, 121, 24-30.	2.5	10
5	Spin-inversion and spin-selection in the reactions FeO ⁺ + H ₂ and Fe ⁺ + N ₂ O. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 19709-19717.	2.8	28
6	Statistical modeling of the reactions Fe ⁺ + N ₂ O → FeO ⁺ + N ₂ and FeO ⁺ + CO → Fe ⁺ + CO ₂ . <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 19700-19708.	2.8	24
7	Coupling an electrospray source and a solids probe/chemical ionization source to a selected ion flow tube apparatus. <i>Review of Scientific Instruments</i> , 2015, 86, 084101.	1.3	6
8	Evaluation of the exothermicity of the chemi-ionization reaction Sm + O → SmO ⁺ + e ⁻ . <i>Journal of Chemical Physics</i> , 2015, 142, 134307.	3.0	44
9	Selected-ion flow tube temperature-dependent measurements for the reactions of O ₂ ⁺ with N atoms and N ₂ ⁺ with O atoms. <i>Journal of Chemical Physics</i> , 2015, 142, 154305.	3.0	9
10	Effect of higher order solvation and temperature on SN ₂ and E ₂ reactivity. <i>International Journal of Mass Spectrometry</i> , 2015, 378, 54-58.	1.5	16
11	Incorporating time-of-flight detection on a selected ion flow tube apparatus. <i>International Journal of Mass Spectrometry</i> , 2015, 377, 479-483.	1.5	11
12	Further Insight into the Reaction FeO ⁺ + H ₂ → Fe ⁺ + H ₂ O: Temperature Dependent Kinetics, Isotope Effects, and Statistical Modeling. <i>Journal of Physical Chemistry A</i> , 2014, 118, 6789-6797.	2.5	38
13	S-P Coupling Induced Unusual Open-Shell Metal Clusters. <i>Journal of the American Chemical Society</i> , 2014, 136, 4821-4824.	13.7	22
14	Temperature-Dependent Kinetics of Charge Transfer, Hydrogen-Atom Transfer, and Hydrogen-Atom Expulsion in the Reaction of CO ⁺ with CH ₄ and CD ₄ . <i>Journal of Physical Chemistry A</i> , 2014, 118, 8141-8146.	2.5	3
15	Activation of Methane by FeO ⁺ : Determining Reaction Pathways through Temperature-Dependent Kinetics and Statistical Modeling. <i>Journal of Physical Chemistry A</i> , 2014, 118, 2029-2039.	2.5	46
16	Photoelectron imaging of small aluminum clusters: quantifying s-p hybridization. <i>Physical Chemistry Chemical Physics</i> , 2013, 15, 3173.	2.8	23
17	Comment on "Role of (NO) ₂ Dimer in Reactions of Fe ⁺ with NO and NO ₂ Studied by ICP-SIFT Mass Spectrometry". <i>Journal of Physical Chemistry A</i> , 2013, 117, 9108-9110.	2.5	2
18	Reactions of Fe ⁺ and FeO ⁺ with C ₂ H ₂ , C ₂ H ₄ , and C ₂ H ₆ : Temperature-Dependent Kinetics. <i>Journal of Physical Chemistry A</i> , 2013, 117, 10178-10185.	2.5	8

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19	Iron cation catalyzed reduction of N ₂ O by CO: gas-phase temperature dependent kinetics. <i>Physical Chemistry Chemical Physics</i> , 2013, 15, 11257.	2.8	26
20	Temperature dependences for the reactions of Ar ⁺ , O ₂ ⁺ , and C ₇ H ₇ ⁺ with toluene and ethylbenzene. <i>International Journal of Mass Spectrometry</i> , 2013, 353, 60-66.	1.5	3
21	Probing the Electronic Structures and Relative Stabilities of Monomagnesium Oxide Clusters MgO _x (x = 1-4): A Combined Photoelectron Imaging and Theoretical Investigation. <i>Journal of Physical Chemistry A</i> , 2013, 117, 11896-11905.	2.5	11
22	Temperature Dependence of the OH ⁺ + CH ₃ I Reaction Kinetics. Experimental and Simulation Studies and Atomic-Level Dynamics. <i>Journal of Physical Chemistry A</i> , 2013, 117, 14019-14027.	2.5	40
23	A novel technique for measurement of thermal rate constants and temperature dependences of dissociative recombination: CO ₂ ⁺ , CF ₃ ⁺ , N ₂ O ⁺ , C ₇ H ₈ ⁺ , C ₇ H ₇ ⁺ , C ₆ H ₆ ⁺ , C ₆ H ₅ ⁺ , C ₅ H ₆ ⁺ , C ₄ H ₄ ⁺ , and C ₃ H ₃ ⁺ . <i>Journal of Chemical Physics</i> , 2013, 138, 154201.	3.0	22
24	Temperature dependences for the reactions of O ₂ ⁺ and O ⁺ with N and O atoms in a selected-ion flow tube instrument. <i>Journal of Chemical Physics</i> , 2013, 139, 144302.	3.0	17
25	Exploring the Reactions of Fe ⁺ and FeO ⁺ with NO and NO ₂ . <i>Journal of Physical Chemistry A</i> , 2012, 116, 11500-11508.	2.5	20
26	Electron Attachment to C ₇ F ₁₄ , Thermal Detachment from C ₇ F ₁₄ ⁺ , the Electron Affinity of C ₇ F ₁₄ , and Neutralization of C ₇ F ₁₄ ⁺ by Ar ⁺ . <i>Journal of Physical Chemistry A</i> , 2012, 116, 10293-10300.	2.5	8
27	Electronic Structure Similarities in Pb _x Sb _y and Sn _x Bi _y Clusters. <i>Journal of Physical Chemistry A</i> , 2011, 115, 10276-10280.	2.5	12
28	Stability and electronic properties of isoelectronic heteroatomic analogs of. <i>Chemical Physics Letters</i> , 2011, 505, 92-95.	2.6	8
29	Resilient aromaticity in lead-indium clusters. <i>Chemical Physics Letters</i> , 2010, 500, 196-201.	2.6	13
30	Structural Evolution of Triniobium Carbide Clusters: Evidence of Large C _n Chains (n = 3-10) in Nb ₃ C _n ⁺ (n = 5-10) Clusters. <i>Journal of Physical Chemistry A</i> , 2010, 114, 1290-1297.	2.5	6
31	Anion Photoelectron Spectroscopy and First-Principles Study of Pb _x In _y Clusters. <i>Journal of Physical Chemistry C</i> , 2010, 114, 20907-20916.	3.1	15
32	The applicability of three-dimensional aromaticity in BiSnn ⁺ Zintl analogues. <i>Journal of Chemical Physics</i> , 2010, 133, 134302.	3.0	17
33	Origins of Stability in Mixed Bismuth-Indium Clusters. <i>Journal of Physical Chemistry C</i> , 2010, 114, 15963-15972.	3.1	16
34	Combined Experimental and Theoretical Study of Al _n X (n = 1-6; X = As, Sb) Clusters: Evidence of Aromaticity and the Jellium Model. <i>Journal of Physical Chemistry A</i> , 2010, 114, 2045-2052.	2.5	23
35	Structure of Bi ₃ Tj ETQq1 1 0.784314 rgBT /Overlock 10 Tf 50 92 Td (stretchy="false")		
36	Electron delocalization in a non-cyclic all-metal III ⁻ V cluster. <i>Chemical Physics Letters</i> , 2009, 480, 189-192.	2.6	7

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37	Al _n Bi Clusters: Transitions Between Aromatic and Jellium Stability. Journal of Physical Chemistry A, 2008, 112, 13316-13325.	2.5	29
38	Effect of Charge and Composition on the Structural Fluxionality and Stability of Nine Atom Tin-Bismuth Zintl Analogues. Inorganic Chemistry, 2008, 47, 10953-10958.	4.0	22