

Ester Livshits

List of Publications by Year in descending order

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Version: 2024-02-01

15
papers

4,140
citations

759233

12
h-index

996975

15
g-index

15
all docs

15
docs citations

15
times ranked

5231
citing authors

#	ARTICLE	IF	CITATIONS
1	Two pathways and an isotope effect in H ₃ ⁺ formation following double ionization of methanol. <i>Natural Sciences</i> , 2021, 1, e10022.	2.1	6
2	Absence of Triplets in Single-Photon Double Ionization of Methanol. <i>Journal of Physical Chemistry Letters</i> , 2020, 11, 8108-8113.	4.6	10
3	Time-resolving the ultrafast H ₂ roaming chemistry and H ₃ ⁺ formation using extreme-ultraviolet pulses. <i>Communications Chemistry</i> , 2020, 3, .	4.5	31
4	Making Sense of Coulomb Explosion Imaging. <i>Journal of Physical Chemistry Letters</i> , 2019, 10, 1361-1367.	4.6	36
5	Single-photon Coulomb explosion of methanol using broad bandwidth ultrafast EUV pulses. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 13488-13495.	2.8	18
6	Advances in molecular quantum chemistry contained in the Q-Chem 4 program package. <i>Molecular Physics</i> , 2015, 113, 184-215.	1.7	2,561
7	A Density Functional Theory for Studying Ionization Processes in Water Clusters. <i>Journal of Physical Chemistry A</i> , 2011, 115, 5735-5744.	2.5	51
8	Tuned Range-Separated Hybrids in Density Functional Theory. <i>Annual Review of Physical Chemistry</i> , 2010, 61, 85-109.	10.8	661
9	Deleterious Effects of Long-Range Self-Repulsion on the Density Functional Description of O ₂ Sticking on Aluminum. <i>Journal of Physical Chemistry A</i> , 2009, 113, 7521-7527.	2.5	27
10	A Density Functional Theory for Symmetric Radical Cations from Bonding to Dissociation. <i>Journal of Physical Chemistry A</i> , 2008, 112, 12789-12791.	2.5	45
11	A well-tempered density functional theory of electrons in molecules. <i>Physical Chemistry Chemical Physics</i> , 2007, 9, 2932.	2.8	344
12	The Role of Charge Localization in Current-Driven Dynamics. <i>Israel Journal of Chemistry</i> , 2007, 47, 99-104.	2.3	5
13	Time-Dependent Density-Functional Studies of the D ₂ Coulomb Explosion. <i>Journal of Physical Chemistry A</i> , 2006, 110, 8443-8450.	2.5	28
14	Avoiding self-repulsion in density functional description of biased molecular junctions. <i>Chemical Physics</i> , 2006, 329, 266-275.	1.9	31
15	Electrical or Photocontrol of the Rotary Motion of a Metallocarborane. <i>Science</i> , 2004, 303, 1849-1851.	12.6	286