Torsha Moitra

List of Publications by Year in descending order

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		1163117	1281871
11	233	8	11
papers	citations	h-index	g-index
11	11	11	291
all docs	docs citations	times ranked	citing authors

#	Article	IF	CITATIONS
1	Simulating weak-field attosecond processes with a Lanczos reduced basis approach to time-dependent equation-of-motion coupled-cluster theory. Physical Review A, 2022, 105, .	2.5	14
2	Multi-electron excitation contributions towards primary and satellite states in the photoelectron spectrum. Physical Chemistry Chemical Physics, 2022, 24, 8329-8343.	2.8	7
3	Behind the scenes of spin-forbidden decay pathways in transition metal complexes. Physical Chemistry Chemical Physics, 2021, 23, 59-81.	2.8	14
4	Capturing Correlation Effects on Photoionization Dynamics. Journal of Chemical Theory and Computation, 2021, 17, 5064-5079.	5.3	14
5	Inner-shell photoabsorption and photoionisation cross-sections of valence excited states from asymmetric-Lanczos equation-of-motion coupled cluster singles and doubles theory. Molecular Physics, 2021, 119, .	1.7	5
6	$\langle i\rangle$ e $\langle i\rangle$ T $\langle i\rangle$ 1.0: An open source electronic structure program with emphasis on coupled cluster and multilevel methods. Journal of Chemical Physics, 2020, 152, 184103.	3.0	68
7	Accurate Description of Photoionization Dynamical Parameters. Journal of Physical Chemistry Letters, 2020, 11, 5330-5337.	4.6	26
8	Vibrationally resolved coupled-cluster x-ray absorption spectra from vibrational configuration interaction anharmonic calculations. Journal of Chemical Physics, 2020, 153, 234111.	3.0	8
9	Molecular inner-shell photoabsorption/photoionization cross sections at core-valence-separated coupled cluster level: Theory and examples. Journal of Chemical Physics, 2019, 150, 224104.	3.0	33
10	Strong Duschinsky Mixing Induced Breakdown of Kasha's Rule in an Organic Phosphor. Journal of Physical Chemistry Letters, 2019, 10, 369-374.	4.6	28
11	Intersystem crossing rate dependent dual emission and phosphorescence from cyclometalated platinum complexes: a second order cumulant expansion based approach. Physical Chemistry Chemical Physics, 2018, 20, 23244-23251.	2.8	16