

Emil J Zak

List of Publications by Year in descending order

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Version: 2024-02-01

17
papers

3,610
citations

1039406

9
h-index

996533

15
g-index

17
all docs

17
docs citations

17
times ranked

4033
citing authors

#	ARTICLE	IF	CITATIONS
1	Controlling rotation in the molecular frame with an optical centrifuge. <i>Physical Review Research</i> , 2021, 3, .	1.3	3
2	Dynamical Semigroups in the Birkhoff Polytope of Order 3 as a Tool for Analysis of Quantum Channels. <i>Open Systems and Information Dynamics</i> , 2020, 27, 2050001.	0.5	1
3	Using collocation and a hierarchical basis to solve the vibrational Schrödinger equation. <i>Journal of Chemical Physics</i> , 2019, 150, 204108.	1.2	8
4	Field-Induced Diastereomers for Chiral Separation. <i>Physical Review Letters</i> , 2019, 123, 243202.	2.9	33
5	Simulating electric field interactions with polar molecules using spectroscopic databases. <i>Scientific Reports</i> , 2017, 7, 45068.	1.6	7
6	From quantum superposition to orbital communication. <i>Computational and Theoretical Chemistry</i> , 2017, 1115, 80-87.	1.1	3
7	Room temperature line lists for CO ₂ asymmetric isotopologues with ab initio computed intensities. <i>Journal of Quantitative Spectroscopy and Radiative Transfer</i> , 2017, 203, 265-281.	1.1	31
8	Room temperature line lists for CO ₂ symmetric isotopologues with ab initio computed intensities. <i>Journal of Quantitative Spectroscopy and Radiative Transfer</i> , 2017, 189, 267-280.	1.1	49
9	Highly accurate intensity factors of pure CO ₂ lines near 2 μ m. <i>Journal of Chemical Physics</i> , 2017, 146, 244309.	1.2	25
10	The HITRAN2016 molecular spectroscopic database. <i>Journal of Quantitative Spectroscopy and Radiative Transfer</i> , 2017, 203, 3-69.	1.1	2,840
11	Ro-vibronic transition intensities for triatomic molecules from the exact kinetic energy operator; electronic spectrum for the Clfâ€™%1B2â€™%â†â€™%Xlfâ€™%1A1 transition in SO ₂ . <i>Journal of Chemical Physics</i> , 2017, 147, 094305.	1.2	5
12	The ExoMol database: Molecular line lists for exoplanet and other hot atmospheres. <i>Journal of Molecular Spectroscopy</i> , 2016, 327, 73-94.	0.4	364
13	A room temperature CO ₂ line list with ab initio computed intensities. <i>Journal of Quantitative Spectroscopy and Radiative Transfer</i> , 2016, 177, 31-42.	1.1	91
14	ROOM TEMPERATURE LINE LISTS FOR CO ₂ ISOTOPOLOGUES WITH AB INITIO COMPUTED INTENSITIES. , 2016, , .		1
15	Electron delocalization index based on bond order orbitals. <i>Chemical Physics Letters</i> , 2014, 593, 154-159.	1.2	37
16	Quasiperiodic Energy Dependence of Exciton Relaxation Kinetics in the Sexithiophene Crystal. <i>Journal of Physical Chemistry A</i> , 2014, 118, 9653-9660.	1.1	0
17	A uniform approach to the description of multicenter bonding. <i>Physical Chemistry Chemical Physics</i> , 2014, 16, 20514-20523.	1.3	112