

Jennifer A Ciezak-Jenkins

List of Publications by Year in descending order

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Version: 2024-02-01

10
papers

80
citations

1684188
5
h-index

1474206
9
g-index

10
all docs

10
docs citations

10
times ranked

140
citing authors

#	ARTICLE	IF	CITATIONS
1	The Local Atomic Structures of Liquid CO at 3.6...GPa and Polymerized CO at 0 to 30...GPa from High Pressure Pair Distribution Function Analysis. <i>Chemistry - A European Journal</i> , 2014, 20, 11531-11539.	3.3	17
2	Structural and spectroscopic studies of nitrogen-carbon monoxide mixtures: Photochemical response and observation of a novel phase. <i>Journal of Chemical Physics</i> , 2017, 146, 184309.	3.0	13
3	The High Pressure Characterization of Melt-Castable Energetic Materials: 3,3'-Bis(5,5'-Bis(Methylene Dinitrate). <i>Propellants, Explosives, Pyrotechnics</i> , 2019, 44, 160-165.	1.6	13
4	Mechanochemical induced structural changes in sucrose using the rotational diamond anvil cell. <i>Journal of Applied Physics</i> , 2018, 123, .	2.5	11
5	Characterization of the Isothermal Compression Behavior of LLM-172. <i>Journal of Physical Chemistry A</i> , 2017, 121, 4263-4271.	2.5	7
6	High Pressure Characterization of High-Performance In-sensitive Energetic Materials: Dihydroxylammonium 5,5'-Bis(3-Nitro-1,2,4-Triazolate)-N-Oxide) (MAD-X1). <i>Propellants, Explosives, Pyrotechnics</i> , 2019, 44, 1263-1269.	1.6	1
7	High Pressure Characterization of Melt-Castable Biisoxazole Energetics: 3,3'-Bis(5,5'-Bis(Methylene) Dinitrate and 3,3'-Biisoxazole-4,4',5,5'-Tetrakis(Methylene) Nitrate). <i>Propellants, Explosives, Pyrotechnics</i> , 2019, 44, 1015-1020.	1.6	1
8	Hydrogen-Bonding Modification in Biuret Under Pressure. <i>Journal of Physical Chemistry A</i> , 2017, 121, 762-770.	2.5	4
9	Kinetic effects on the morphology and stability of the pressure-induced extended-solid of carbon monoxide. <i>Journal of Chemical Physics</i> , 2018, 148, 144702.	3.0	3
10	High-pressure polymorphism of the electrochemically active organic molecule tetrahydroxy-p-benzoquinone. <i>Journal of Molecular Structure</i> , 2016, 1119, 71-77.	3.6	1