

Mikhail V Ryzhkov

List of Publications by Year in descending order

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13
papers

163
citations

1478505

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1281871

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g-index

13
all docs

13
docs citations

13
times ranked

190
citing authors

#	ARTICLE	IF	CITATIONS
1	The electronic structure and the nature of the chemical bond in CeO ₂ . Physical Chemistry Chemical Physics, 2018, 20, 16167-16175.	2.8	45
2	Electronic structure of endohedral fullerenes An@C28 (An=Th–Md). Computational and Theoretical Chemistry, 2012, 985, 46-52.	2.5	36
3	Electronic structure and geometry optimization of nanoparticles Fe2C, FeC2, Fe3C, FeC3 and Fe2C2. Chemical Physics Letters, 2005, 404, 400-408.	2.6	34
4	Electronic structure of predicted endohedral fullerenes An@C40 (An=Th–Md). Computational and Theoretical Chemistry, 2013, 1013, 70-77.	2.5	19
5	Valence electronic state density in thorium dioxide. Nuclear Technology and Radiation Protection, 2008, 23, 34-42.	0.8	10
6	X-ray photoelectron spectra structure and chemical bonding in AmO ₂ . Nuclear Technology and Radiation Protection, 2015, 30, 83-98.	0.8	10
7	The nature of the chemical bond in UO ₂ . International Journal of Quantum Chemistry, 2019, 119, e26040.	2.0	4
8	Plutonium complexes in water: new approach to ab initio modeling. Radiochimica Acta, 2021, 109, 327-342.	1.2	2
9	Quantum-chemical simulation of the electronic structure and chemical bonding in the new δ -superstoichiometric titanium carbonitride Ti ₂ CN ₄ . Mendeleev Communications, 2001, 11, 184-185.	1.6	1
10	Electronic Structure and Chemical Bonding in β -Sialons. Journal of Structural Chemistry, 2002, 43, 18-25.	1.0	1
11	Electronic structure and effective charges on atoms near anion point defects in uranium dioxide. Computational Condensed Matter, 2019, 18, e00353.	2.1	1
12	Transformation of electron density distribution induced by the cation point defects in uranium dioxide. Journal of Radioanalytical and Nuclear Chemistry, 2020, 325, 253-262.	1.5	0
13	First-principles study on the plutonium ions interaction with diamide molecules in acid solutions. International Journal of Quantum Chemistry, 2021, 121, e26681.	2.0	0