

Andrzej KÄdziorski

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/1389828/publications.pdf>

Version: 2024-02-01

24

papers

288

citations

933447

10

h-index

888059

17

g-index

25

all docs

25

docs citations

25

times ranked

408

citing authors

#	ARTICLE	IF	CITATIONS
1	Broadband anti-Stokes white emission of Sr ₂ CeO ₄ nanocrystals induced by laser irradiation. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 27921-27927.	2.8	53
2	Role of the Antenna in Tissue Selective Probes Built of Lanthanide-Organic Chelates. <i>Journal of Physical Chemistry A</i> , 2008, 112, 2397-2407.	2.5	32
3	Ab initio Theoretical Study on the 4f ² and 4f5d Electronic Manifolds of Cubic Defects in CaF ₂ :Pr ³⁺ . <i>Journal of Physical Chemistry A</i> , 2014, 118, 358-368.	2.5	28
4	Efficiency of the energy transfer in lanthanide-organic chelates; spectral overlap integral. <i>Journal of Luminescence</i> , 2010, 130, 1154-1159.	3.1	25
5	The first example of ab initio calculations of f-f transitions for the case of [Eu(DOTP)] ₅ experiment versus theory. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 27808-27817.	2.8	19
6	Enhancement of and interference among higher order multipole transitions in molecules near a plasmonic nanoantenna. <i>Nature Communications</i> , 2019, 10, 5775.	12.8	19
7	New parametrization of spectra of Nd ³⁺ and Sm ³⁺ in glasses. <i>Journal of Alloys and Compounds</i> , 2008, 451, 686-690.	5.5	16
8	electric dipole transitions; old problems in a new light. <i>Journal of Alloys and Compounds</i> , 2009, 488, 586-590.	5.5	16
9	Extended parametrization scheme of f-spectra. <i>Journal of Luminescence</i> , 2007, 127, 552-560.	3.1	12
10	Interatomic potentials of metal dimers: probing agreement between experiment and advanced ab initio calculations for van der Waals dimer Cd ₂ . <i>International Reviews in Physical Chemistry</i> , 2017, 36, 541-620.	2.3	10
11	Magnetic dipole transitions in crystals. <i>Molecular Physics</i> , 2004, 102, 1105-1111.	1.7	9
12	The E 3 Δ 1 + (6 3 S 1) → A 3 Δ 0 + (5 3 P 1) transition in CdAr revisited: The spectrum and new analysis of the E 3 Δ 1 + Rydberg state interatomic potential. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2018, 196, 58-66.	3.9	9
13	Borrowing Intensity in Rare Earth Doped Materials; Magnetic Dipole Transitions. <i>Collection of Czechoslovak Chemical Communications</i> , 2005, 70, 905-922.	1.0	6
14	Rydberg states of the CdAr van der Waals complex. <i>Physical Review A</i> , 2019, 99, .	2.5	6
15	Atomic fine-structure calculations performed with a finite-nuclear-mass approach and with all-electron explicitly correlated Gaussian functions. <i>Chemical Physics Letters</i> , 2020, 751, 137476. xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:mi>l</mml:mi><mml:math>Mn</mml:math>	2.6	6
16	xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:msub><mml:mrow><mml:mn>3</mml:mn></mml:mrow></mml:msub><mml:math>N</mml:math>	3.2	5
17	xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:msub><mml:mrow><mml:math>Hyperfine-induced f Transitions: Effective Operator Formulation. Spectroscopy Letters, 2007, 40, 293-315.	1.0	4
18	Net-value of the relativistic crystal field effect. <i>Journal of Alloys and Compounds</i> , 2004, 380, 151-155.	5.5	3

#	ARTICLE	IF	CITATIONS
19	Experimental and <i>Ab Initio</i> Study on the Intensities of f-f Transitions for the Molecular Eu(III)-DOTP System. <i>ChemistrySelect</i> , 2019, 4, 1394-1402.	1.5	3
20	Magnetic dipole transitions in crystals. <i>Journal of Alloys and Compounds</i> , 2008, 451, 18-34.	5.5	2
21	Influence of dopant concentration on spectroscopic properties of Sr ₂ CeO ₄ :Yb nanocrystals. <i>Optical Materials</i> , 2017, 74, 34-40.	3.6	2
22	Rydberg states of ZnAr complex. <i>Molecular Physics</i> , 2022, 120, .	1.7	2
23	Fine structure of the beryllium Be^{2+} states calculated with all-electron explicitly correlated Gaussian functions. <i>Physical Review A</i> , 2022, 105, 022501.	2.5	1
24	Interatomic potentials of van der Waals dimers Hg ₂ and Cd ₂ : Probing discrepancies between theory and experiment. <i>Journal of Physics: Conference Series</i> , 2017, 810, 012018.	0.4	0