

Lev Mazalov

List of Publications by Year in descending order

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23
papers

97
citations

1477746

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1473754

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23
all docs

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docs citations

23
times ranked

155
citing authors

#	ARTICLE	IF	CITATIONS
1	XPS experimental and DFT investigations on solid solutions of $\text{Mo}_{1-x}\text{Re}_x\text{S}_2$ ($0 \leq x \leq 0.20$). <i>Nanoscale</i> , 2018, 10, 10232-10240.	2.8	23
2	Synthesis and study of the electronic structure of molybdenum tetrasulfide and its lithium intercalates. <i>Journal of Physics and Chemistry of Solids</i> , 1998, 59, 283-288.	1.9	12
3	An ab initio calculation of K-spectra in molecules HCl and HF. <i>Theoretica Chimica Acta</i> , 1977, 44, 257-263.	0.9	7
4	An ab initio investigation of the inner shell excited states of the molecule Cl_2 . <i>Theoretica Chimica Acta</i> , 1979, 54, 179-185.	0.9	7
5	Ge quantum dots structural peculiarities depending on the preparation conditions. <i>Journal of Synchrotron Radiation</i> , 2003, 10, 380-383.	1.0	7
6	Many-electron effects in the x-ray emission spectra and the one-particle Green's function. Correspondence theorem. <i>Theoretica Chimica Acta</i> , 1983, 64, 31-39.	0.9	6
7	Many-electron effects in the X-ray absorption spectra and the one-particle Green's function. <i>Theoretica Chimica Acta</i> , 1983, 62, 537-547.	0.9	6
8	An SCF XPS study of the electronic structure and X-ray and photoelectron spectra of Fe(II) and Fe(III) hexacyano complexes in a cluster approach. <i>Theoretica Chimica Acta</i> , 1986, 70, 429-441.	0.9	5
9	X-ray emission and X-ray photoelectron study of the electronic structure of polymeric cubanocluster compounds $\text{Re}_4 \times \text{Mo}_x \text{S}_4 \text{Te}_4$. <i>Journal of Structural Chemistry</i> , 1996, 37, 767-772.	0.3	5
10	X-ray spectral microanalysis of the phase composition of high temperature superconductor bismuth-lead-strontium-calcium-copper-oxygen ceramics using chemometric approaches. <i>Analyst</i> , 1992, 117, 795-802.	1.7	4
11	An ab initio investigation of the collective phenomena accompanying the core ionization in the radical NO. <i>Theoretica Chimica Acta</i> , 1979, 52, 311-318.	0.9	3
12	X-ray spectral microinvestigation of the chemical states of atoms in $\text{Cd}_x\text{Hg}_{1-x}\text{Te}$ GaAs thin films. <i>Analyst</i> , 1992, 117, 803-805.	1.7	2
13	MNDO calculations of the electronic structure of lead P^2 -diketonates. <i>Journal of Structural Chemistry</i> , 1994, 35, 916-919.	0.3	2
14	X-ray spectra and electronic structure of solid ammonia. <i>Journal of Structural Chemistry</i> , 1996, 37, 621-625.	0.3	2
15	X-ray spectra and electronic structure of hexafluorobenzene and pentafluoronitrobenzene. <i>Journal of Structural Chemistry</i> , 1989, 29, 720-727.	0.3	1
16	Fragment-by-fragment analysis of molecular orbitals in some substituted benzenes. <i>Journal of Structural Chemistry</i> , 1989, 30, 194-202.	0.3	1
17	X-ray spectral and quantum chemical analyses of the electronic structure of poly(monofluorocarbon). <i>Journal of Structural Chemistry</i> , 1995, 36, 572-577.	0.3	1
18	Synthesis and investigation of the electronic structure of molybdenum tetrasulfide and its lithium intercalates. <i>Journal of Structural Chemistry</i> , 1996, 37, 626-632.	0.3	1

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19	Quantum chemical and x-ray spectral studies of the structures of superstoichiometric fluorocarbons. Journal of Structural Chemistry, 1996, 37, 906-912.	0.3	1
20	Influence of conformation on the electronic structure of thiacalixarenes according to DFT calculations and X-ray emission spectroscopy. Journal of Molecular Structure, 2011, 1006, 502-507.	1.8	1
21	The study of interaction of HIV-1 surface GP120 protein with CD4 cell receptor by EXAFS spectroscopy. Nuclear Instruments and Methods in Physics Research, Section A: Accelerators, Spectrometers, Detectors and Associated Equipment, 1995, 359, 259-262.	0.7	0
22	Quantum chemical calculation of Mo ₄ S ₄ Cl ₄ in the cluster approximation. Journal of Structural Chemistry, 1996, 37, 335-339.	0.3	0
23	Investigation of electronic interactions in solid hydrogen fluoride. Journal of Structural Chemistry, 1997, 38, 570-577.	0.3	0