

Lev Mazalov

List of Publications by Year in Descending Order

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The third column is the impact factor (IF) of the journal, and the fourth column is the number of citations of the article.

23
papers

84
citations

6
h-index

8
g-index

23
ext. papers

88
ext. citations

2.2
avg, IF

0.98
L-index

#	Paper	IF	Citations
23	XPS experimental and DFT investigations on solid solutions of MoReS (0 Nanoscale, 2018 , 10, 10232-10240)	4.9	17
22	Influence of conformation on the electronic structure of thiacalixarenes according to DFT calculations and X-ray emission spectroscopy. <i>Journal of Molecular Structure</i> , 2011 , 1006, 502-507	3.4	1
21	Ge quantum dots structural peculiarities depending on the preparation conditions. <i>Journal of Synchrotron Radiation</i> , 2003 , 10, 380-3	2.4	7
20	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	3.9	10
19	Investigation of electronic interactions in solid hydrogen fluoride. <i>Journal of Structural Chemistry</i> , 1997 , 38, 570-577	0.9	
18	X-ray emission and X-ray photoelectron study of the electronic structure of polymeric cubanocluster compounds $\text{Re}_4\text{Mo}_x\text{S}_4\text{Te}_4$. <i>Journal of Structural Chemistry</i> , 1996 , 37, 767-772	0.9	3
17	X-ray spectra and electronic structure of solid ammonia. <i>Journal of Structural Chemistry</i> , 1996 , 37, 621-625	0.9	2
16	Synthesis and investigation of the electronic structure of molybdenum tetrasulfide and its lithium intercalates. <i>Journal of Structural Chemistry</i> , 1996 , 37, 626-632	0.9	
15	Quantum chemical and x-ray spectral studies of the structures of superstoichiometric fluorocarbons. <i>Journal of Structural Chemistry</i> , 1996 , 37, 906-912	0.9	1
14	Quantum chemical calculation of $\text{Mo}_4\text{S}_4\text{Cl}_4$ in the cluster approximation. <i>Journal of Structural Chemistry</i> , 1996 , 37, 335-339	0.9	
13	X-ray spectral and quantum chemical analyses of the electronic structure of poly(monofluorocarbon). <i>Journal of Structural Chemistry</i> , 1995 , 36, 572-577	0.9	1
12	The study of interaction of HIV-1 surface GP120 protein with CD4 cell receptor by EXAFS spectroscopy. <i>Nuclear Instruments and Methods in Physics Research, Section A: Accelerators, Spectrometers, Detectors and Associated Equipment</i> , 1995 , 359, 259-262	1.2	
11	MNDO calculations of the electronic structure of lead β -diketonates. <i>Journal of Structural Chemistry</i> , 1994 , 35, 916-919	0.9	2
10	X-ray spectral microinvestigation of the chemical states of atoms in $\text{Cd}_x\text{Hg}_{1-x}\text{Te}$ thin films. <i>Analyst, The</i> , 1992 , 117, 803-805	5	1
9	X-ray spectral microanalysis of the phase composition of high temperature superconductor bismuth lead strontium calcium copper oxygen ceramics using chemometric approaches. <i>Analyst, The</i> , 1992 , 117, 795-802	5	3
8	X-ray spectra and electronic structure of hexafluorobenzene and pentafluoronitrobenzene. <i>Journal of Structural Chemistry</i> , 1989 , 29, 720-727	0.9	1
7	Fragment-by-fragment analysis of molecular orbitals in some substituted benzenes. <i>Journal of Structural Chemistry</i> , 1989 , 30, 194-202	0.9	1

- 6 An SCF X̄W study of the electronic structure and X-ray and photoelectron spectra of Fe(II) and Fe(III) hexacyano complexes in a cluster approach. *Theoretica Chimica Acta*, **1986**, 70, 429-441 5
- 5 Many-electron effects in the x-ray emission spectra and the one-particle Green's function. Correspondence theorem. *Theoretica Chimica Acta*, **1983**, 64, 31-39 6
- 4 Many-electron effects in the X-ray absorption spectra and the one-particle Green's function. *Theoretica Chimica Acta*, **1983**, 62, 537-547 6
- 3 An ab initio investigation of the collective phenomena accompanying the core ionization in the radical NO. *Theoretica Chimica Acta*, **1979**, 52, 311-318 3
- 2 An ab initio investigation of the inner shell excited states of the molecule Cl₂. *Theoretica Chimica Acta*, **1979**, 54, 179-185 7
- 1 An ab initio calculation of K-spectra in molecules HCl and HF. *Theoretica Chimica Acta*, **1977**, 44, 257-263 7