

M Romero

List of Publications by Year in descending order

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18
papers

400
citations

1040056

9
h-index

839539

18
g-index

19
all docs

19
docs citations

19
times ranked

507
citing authors

#	ARTICLE	IF	CITATIONS
1	Immunophenotypic heterogeneity of multiple myeloma: influence on the biology and clinical course of the disease. <i>British Journal of Haematology</i> , 1991, 77, 185-190.	2.5	94
2	Lymphoid subsets and prognostic factors in multiple myeloma. <i>British Journal of Haematology</i> , 1992, 80, 305-309.	2.5	64
3	Synthesis by molten salt method of the AFeO ₃ system (A=La, Gd) and its structural, vibrational and internal hyperfine magnetic field characterization. <i>Physica B: Condensed Matter</i> , 2014, 443, 90-94.	2.7	61
4	First-principles calculations of structural, elastic and electronic properties of Nb ₂ SnC under pressure. <i>Computational Materials Science</i> , 2012, 55, 142-146.	3.0	37
5	N-Arylpiperazinyl-N ⁺ -propylamino Derivatives of Heteroaryl Amides as Functional Uroselective \pm 1-Adrenoceptor Antagonists. <i>Journal of Medicinal Chemistry</i> , 1997, 40, 2674-2687.	6.4	33
6	X-ray photoelectron spectroscopy studies of the electronic structure of superconducting Nb ₂ SnC and Nb ₂ SC. <i>Journal of Alloys and Compounds</i> , 2013, 579, 516-520.	5.5	23
7	Direct N3 Alkylation of Uracil and Derivatives via n1-[2-(trimethylsilyl)ethoxymethyl] Protection. <i>Synlett</i> , 1997, 1997, 1233-1234.	1.8	16
8	Synthesis, pharmacology and pharmacokinetics of 3-(4-Aryl-piperazin-1-ylalkyl)-uracils as uroselective \pm 1A-antagonists. <i>Bioorganic and Medicinal Chemistry Letters</i> , 2003, 13, 1873-1878.	2.2	15
9	Structural and mechanic properties of RFeO ₃ with R = Y, Eu and La perovskites: a first-principles calculation. <i>European Physical Journal D</i> , 2015, 69, 1.	1.3	10
10	Using Wikipedia concepts and frequency in language to extract key terms from support documents. <i>Expert Systems With Applications</i> , 2012, 39, 13480-13491.	7.6	8
11	First-principles study of the structural, elastic, vibrational, thermodynamic and electronic properties of the Mo ₂ B intermetallic under pressure. <i>Journal of Molecular Structure</i> , 2016, 1125, 350-357.	3.6	7
12	A cloud of FAQ: A highly-precise FAQ retrieval system for the Web 2.0. <i>Knowledge-Based Systems</i> , 2013, 49, 81-96.	7.1	6
13	Ab initio study of structural, elastic, and electronic properties of Mo ₃ 46B ₁₂ under high pressure. <i>European Physical Journal B</i> , 2019, 92, 1.	1.5	6
14	Exposed Surface and Confinement Effects on the Electronic, Magnetic, and Mechanical Properties of LaTiO ₃ Slabs. <i>IEEE Transactions on Magnetics</i> , 2021, 57, 1-4.	2.1	5
15	LDA+U study of hydrostatic pressure effect on double perovskite Sr ₂ FeNbO ₆ : crystal structure, mechanical and electronic properties. <i>Physica Scripta</i> , 2020, 95, 115704.	2.5	5
16	Effects of the phase transition on the structural, mechanical, electronic and vibrational properties of the CaSnO ₃ perovskite: Study under hydrostatic pressure. <i>Journal of Physics and Chemistry of Solids</i> , 2022, 163, 110594.	4.0	5
17	FAQtory: A framework to provide high-quality FAQ retrieval systems. <i>Expert Systems With Applications</i> , 2012, 39, 11525-11534.	7.6	4
18	Ab Initio study of the crystal structure and the elastic properties of the Mo _{0.85} Nb _{0.15} B ₃ compound under pressure. <i>MRS Advances</i> , 2019, 4, 3453-3461.	0.9	1