

Giuseppe Zerbi

List of Publications by Year in descending order

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12
papers

758
citations

1040056

9
h-index

1199594

12
g-index

12
all docs

12
docs citations

12
times ranked

1189
citing authors

#	ARTICLE	IF	CITATIONS
1	Graphite particles induce ROS formation in cell free systems and human cells. <i>Nanoscale</i> , 2017, 9, 13640-13650.	5.6	16
2	Resonant Raman-based cytochrome C biosensor as a tool for evaluating the oxidative properties of the diesel exhaust particulate matter. <i>Journal of Raman Spectroscopy</i> , 2016, 47, 796-800.	2.5	3
3	Fingerprints of polycyclic aromatic hydrocarbons (PAHs) in infrared absorption spectroscopy. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2016, 152, 134-148.	3.9	48
4	A joint Raman and EPR spectroscopic study on ball-milled nanographites. <i>Chemical Physics Letters</i> , 2011, 516, 220-224.	2.6	41
5	Toward carbyne: Synthesis and stability of really long polyynes. <i>Pure and Applied Chemistry</i> , 2010, 82, 891-904.	1.9	59
6	Disposable Electrospun Electrodes Based on Conducting Nanofibers. <i>Electroanalysis</i> , 2008, 20, 1374-1377.	2.9	18
7	Structure of new carbonaceous materials: The role of vibrational spectroscopy. <i>Carbon</i> , 2005, 43, 1593-1609.	10.3	92
8	Resonant Raman spectroscopy of nanostructured carbon-based materials: the molecular approach. <i>AIP Conference Proceedings</i> , 2004, , .	0.4	8
9	Raman spectroscopy of polyconjugated molecules and materials: confinement effect in one and two dimensions. <i>Philosophical Transactions Series A, Mathematical, Physical, and Engineering Sciences</i> , 2004, 362, 2425-2459.	3.4	248
10	Resonance Raman contribution to the D band of carbon materials: Modeling defects with quantum chemistry. <i>Journal of Chemical Physics</i> , 2004, 120, 11889-11900.	3.0	87
11	A Computational Study of the Raman Spectra of Large Polycyclic Aromatic Hydrocarbons: Toward Molecularly Defined Subunits of Graphite. <i>Journal of Physical Chemistry A</i> , 2002, 106, 3306-3317.	2.5	131
12	Excited-State Molecular Dynamics Simulations of Conjugated Oligomers Using the Electronic Density Matrix. <i>Journal of Physical Chemistry A</i> , 2001, 105, 7057-7071.	2.5	7