## Alvaro Lobato

## List of Publications by Year in descending order

Source: https://exaly.com/author-pdf/3735532/publications.pdf

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		1307594	1281871	
20	125	7	11	
papers	citations	h-index	g-index	
20	20	20	162	
20	20	20	162	
all docs	docs citations	times ranked	citing authors	

#	Article	IF	Citations
1	The structures of inorganic crystals: A rational explanation from the chemical pressure approach and the anions in metallic matrices model., 2023,, 238-261.		1
2	Structural and Electronic Effect Driven Distortions in Visible Light Absorbing Polar Materials $\langle i\rangle A\langle  i\rangle Ta\langle sub\rangle 2\langle  sub\rangle V\langle sub\rangle 2\langle  sub\rangle O\langle sub\rangle 11\langle  sub\rangle (\langle i\rangle A\langle  i\rangle = Sr, Pb)$ . Journal of Physical Chemistry C, 2022, 126, 8047-8055.	3.1	0
3	Compressing the Channels in the Crystal Structure of Copper Squarate Metal-Organic Framework. Solids, 2022, 3, 374-384.	2.4	2
4	Temperature and pressure-induced strains in anhydrous iron trifluoride polymorphs. Physical Chemistry Chemical Physics, 2021, 23, 2825-2835.	2.8	1
5	Controlling the off-center positions of anions through thermodynamics and kinetics in flexible perovskite-like materials. Physical Chemistry Chemical Physics, 2021, 23, 4491-4499.	2.8	2
6	Mechanical Response of Polymer Epoxy/BMI Composites with Graphene and a Boron Nitride Monolayer from First Principles. ACS Applied Polymer Materials, 2021, 3, 1052-1059.	4.4	10
7	Highs and Lows of Bond Lengths: Is There Any Limit?. Angewandte Chemie, 2021, 133, 17165-17173.	2.0	5
8	Highs and Lows of Bond Lengths: Is There Any Limit?. Angewandte Chemie - International Edition, 2021, 60, 17028-17036.	13.8	13
9	Understanding the Pressure Effect on the Elastic, Electronic, Vibrational, and Bonding Properties of the CeScO <sub>3</sub> Perovskite. Journal of Physical Chemistry C, 2021, 125, 107-119.	3.1	17
10	Comment on "Uncommon structural and bonding properties in Ag <sub>16</sub> 8 <sub>4</sub> 0 <sub>10</sub> ―by A. Kovalevskiy, C. Yin, J. Nuss, U. Wedig, and M. Jansen, <i>Chem. Sci.</i> ), 2020, <b>11</b> , 962. Chemical Science, 2021, 12, 13588-13592.	7.4	6
11	Theoretical (DFT) and experimental (Raman and FTIR) spectroscopic study on communic acids, main components of fossil resins. Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy, 2020, 224, 117405.	3.9	8
12	Generalized Stress-Redox Equivalence: A Chemical Link between Pressure and Electronegativity in Inorganic Crystals. Inorganic Chemistry, 2020, 59, 5281-5291.	4.0	21
13	Computational Modeling of Tensile Stress Effects on the Structure and Stability of Prototypical Covalent and Layered Materials. Nanomaterials, 2019, 9, 1483.	4.1	3
14	Chemical pressure–chemical knowledge: squeezing bonds and lone pairs within the valence shell electron pair repulsion model. Physical Chemistry Chemical Physics, 2019, 21, 12585-12596.	2.8	12
15	The self-absorption phenomenon in quantitative Raman spectroscopy and how to correct its effects. Microchemical Journal, 2018, 139, 134-138.	4.5	9
16	Infrared spectroscopic study of the formation of fossil resin analogs with temperature using trans-communic acid as precursor. Microchemical Journal, 2018, 141, 294-300.	4.5	4
17	Temperature effects on the friction-like mode of graphite. Theoretical Chemistry Accounts, 2017, 136, 1.	1.4	O
18	Local, solvation pressures and conformational changes in ethylenediamine aqueous solutions probed using Raman spectroscopy. Physical Chemistry Chemical Physics, 2016, 18, 26192-26198.	2.8	8

#	Article	IF	CITATIONS
19	Anharmonicity effects in the frictionlike mode of graphite. Physical Review B, 2016, 93, .	3.2	1
20	Raman Spectroscopy, a Useful Tool to Study Nuclear Materials. , 0, , .		2