

# Erwan AndrÃ©

## List of Publications by Year in descending order

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22  
papers

509  
citations

759233

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677142

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docs citations

24  
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857  
citing authors

#	ARTICLE	IF	CITATIONS
1	Thin Films of SiP Lamellar Alloys: A First Step toward 2D SiP. <i>Journal of Physical Chemistry C</i> , 2021, 125, 3235-3241.	3.1	6
2	Formation of SiP <sub>2</sub> Nanocrystals Embedded in SiO <sub>2</sub> from Phosphorus-Rich SiO <sub>1.5</sub> Thin Films. <i>Journal of Physical Chemistry C</i> , 2020, 124, 7973-7978.	3.1	4
3	Experimental and Theoretical Infrared Signatures of REMO <sub>3</sub> (RE = La, Pr, Nd, Sm, and M =) <i>Tj ETQq1 1 0,784314 rgBT / Overlock 10 Tf 50 382 T</i>	3.1	12
4	Carbonate-Hydrogenocarbonate Coexistence and Dynamics in Layered Double Hydroxides. <i>Journal of Physical Chemistry C</i> , 2017, 121, 6104-6112.	3.1	23
5	Enhanced photocatalytic ability of Cu, Co doped ZnAl based mixed metal oxides derived from layered double hydroxides. <i>Colloids and Surfaces A: Physicochemical and Engineering Aspects</i> , 2017, 524, 43-52.	4.7	28
6	Probing the Dynamics of Layered Double Hydroxides by Solid-State <sup>27</sup> Al NMR Spectroscopy. <i>Journal of Physical Chemistry C</i> , 2017, 121, 7276-7281.	3.1	8
7	Ternary Layered Double Hydroxides (LDHs) Based on Co-, Cu-Substituted ZnAl for the Design of Efficient Photocatalysts. <i>European Journal of Inorganic Chemistry</i> , 2017, 2017, 669-678.	2.0	43
8	Enhanced catalytic oxidation ability of ternary layered double hydroxides for organic pollutants degradation. <i>Dalton Transactions</i> , 2016, 45, 8224-8235.	3.3	32
9	Molecular Sieving with Vertically Aligned Mesoporous Silica Films and Electronic Wiring through Isolating Nanochannels. <i>Chemistry of Materials</i> , 2016, 28, 2511-2514.	6.7	58
10	Remarkable Structure and Elasticity Relaxation Dynamics of Poly(diallyldimethylammonium) <i>Tj ETQq0 0 0 rgBT / Overlock 10 Tf 50 382 T</i>	3.1	12
11	Properties of rare-earth orthoferrites perovskite driven by steric hindrance. <i>Journal of Alloys and Compounds</i> , 2016, 657, 631-638.	5.5	32
12	Tuning and Investigating the Structure of M <sup>II</sup> -Fe <sup>III</sup> Layered Double Hydroxides (M <sup>II</sup> = Ni <sup>II</sup> , Co <sup>II</sup> ) <i>Tj ETQq0 0 0 rgBT / Overlock 10 Tf 50 382 T</i>	0.2	12
13	Properties. <i>Current Inorganic Chemistry</i> , 2015, 5, 169-183.		
14	Modelling the Structure and Vibrational Properties of Layered Double Hydroxides. , 2015, , 317-323.		1
15	The Raman spectrum of CaCO <sub>3</sub> polymorphs calcite and aragonite: A combined experimental and computational study. <i>Journal of Chemical Physics</i> , 2014, 140, 164509.	3.0	131
16	Versatile Reactivity of Phosphagermaallene Tip( <i>i</i> -Bu)Ge <sup>+</sup> Cl <sup>-</sup> PMes* with $\hat{I}\pm$ -Ethylenic Esters. <i>Organometallics</i> , 2013, 32, 1085-1093.	2.3	5
17	pH influence on the complexation site of Al(III) with protocatechuic acid. A spectroscopic and theoretical approach. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2013, 108, 280-287.	3.9	9
18	Versatile Stereoselective Cycloadditions between Heterocumulenes and Phosphagermaallene Tip( <i>i</i> -Bu)Ge <sup>+</sup> Cl <sup>-</sup> PMes*: Experimental and Theoretical Investigations. <i>Chemistry - A European Journal</i> , 2011, 17, 12763-12772.	3.3	14
19	1,3- $\delta$ -Dipole Behavior of Phosphagermaallene Tip( <i>i</i> -Bu)Ge <sup>+</sup> Cl <sup>-</sup> PMes* Leading to a Phosphagermaheterocyclic Carbene. <i>Angewandte Chemie - International Edition</i> , 2010, 49, 8704-8707.	13.8	20

#	ARTICLE	IF	CITATIONS
19	The Reactivity of Phosphagermaallene $\text{Mes}^*\text{P}=\text{C}=\text{Ge}(\text{i-Bu})_2$ toward Aldehydes and Ketones: an Experimental and Theoretical Study. <i>Organometallics</i> , 2010, 29, 2566-2578.	2.3	26
20	Toward a Better Understanding of the Regioselectivity of the $\text{Al(III)}$ -Protocatechuic Acid Complexation Reaction. <i>Journal of Physical Chemistry A</i> , 2008, 112, 9829-9834.	2.5	10
21	Metal complexation of protocatechuic acid and its derivatives: Determination of the optimal computational conditions for the simulation of electronic spectra. <i>Computational and Theoretical Chemistry</i> , 2007, 806, 131-140.	1.5	12
22	Characterization of the $\text{Al(III)}$ binding site of protocatechuic acid by electronic spectroscopy and quantum chemical calculations. <i>Chemical Physics Letters</i> , 2007, 434, 155-159.	2.6	12