## Said Bouzakraoui

List of Publications by Year in descending order

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933447 996975 17 268 10 15 citations h-index g-index papers 17 17 17 387 docs citations times ranked citing authors all docs

#	Article	IF	CITATIONS
1	Energy absorption and damage characterization of GFRP laminated and PVC-foam sandwich composites under repeated impacts with reduced energies and quasi-static indentation. Case Studies in Construction Materials, 2022, 16, e00844.	1.7	3
2	Adsorption of a cationic dye (Methylene bleu) by Typha Latifolia: Equilibrium, kinetic, thermodynamic and DFT calculations. Chemical Data Collections, 2022, 38, 100834.	2.3	30
3	Carbazole-based hole-transport materials for efficient Perovskite solar cells. A computational study. Optik, 2022, 257, 168793.	2.9	4
4	Structural and thermodynamical properties of early human amylin oligomers using replica exchange molecular dynamics: mutation effect of three key residues F15, H18 and F23. Physical Chemistry Chemical Physics, 2017, 19, 31290-31299.	2.8	10
5	Bianthrone at a metal surface: Conductance switching with a bistable molecule made feasible by image charge effects. , 2015, , .		O
6	Multiscale Simulation of Polyglutamine and the Effect of Neighboring Amino Acids on Oligomerization. Biophysical Journal, 2012, 102, 733a.	0.5	0
7	Dimers of Anthrathiophene and Anthradithiophene Derivatives: Synthesis and Characterization. Organic Letters, 2011, 13, 548-551.	4.6	17
8	Donor/acceptor-substituted anthradithiophene materials: synthesis, optical and electrochemical properties. Tetrahedron, 2011, 67, 7156-7161.	1.9	7
9	Synthesis of soluble oligothiophenes bearing cyano groups, their optical and electrochemical properties. Tetrahedron, 2010, 66, 9560-9572.	1.9	16
10	Bianthrone in a Single-Molecule Junction: Conductance Switching with a Bistable Molecule Facilitated by Image Charge Effects. Journal of Physical Chemistry C, 2010, 114, 20686-20695.	3.1	19
11	Structural and Charge-Transport Properties of a Liquid-Crystalline $\hat{l}\pm, \hat{l}\%$ -Disubstituted Thiophene Derivative: A Joint Experimental and Theoretical Study. Journal of Physical Chemistry C, 2010, 114, 4617-4627.	3.1	18
12	Theoretical study of structural and electronic properties of oligo(thiophene-phenylene)s in comparison with oligothiophenes and oligophenylenes. Chinese Chemical Letters, 2008, 19, 123-126.	9.0	21
13	Theoretical investigation of electroluminescent alkoxy substituted 4,4′-bis(2-phenylethenyl)biphenyls as guest in blue OLEDs. Solar Energy Materials and Solar Cells, 2006, 90, 1393-1402.	6.2	12
14	Prediction of electropolymerization mechanisms of two substituted phenylene: Poly-3-methoxy-toluenes (P3mt1 and P3mt2). Journal of Applied Polymer Science, 2006, 100, 57-64.	2.6	10
15	Density functional theory study of conformational and opto-electronic properties of oligo-para-phenylenes. Computational and Theoretical Chemistry, 2005, 725, 39-44.	1.5	29
16	Density functional theory (B3LYP/6-31G*) study of oligothiophenes in their aromatic and polaronic states. Computational and Theoretical Chemistry, 2005, 726, 271-276.	1.5	61
17	Synthesis and characterization of co-polymers involving various thiophene and phenylene monomers. Synthetic Metals, 2004, 145, 237-243.	3.9	11