Luc M Leblanc

List of Publications by Year
in descending order

21 papers	443 citations	759233 12 h-index	713466 21 g-index
21 all docs	21 docs citations	21 times ranked	661 citing authors

#	Article	IF	Citations
1	Caffeic Acid Phenethyl Ester and Its Amide Analogue Are Potent Inhibitors of Leukotriene Biosynthesis in Human Polymorphonuclear Leukocytes. PLoS ONE, 2012, 7, e31833.	2.5	58
2	Synthesis and Antiradical/Antioxidant Activities of Caffeic Acid Phenethyl Ester and Its Related Propionic, Acetic, and Benzoic Acid Analoguesc. Molecules, 2012, 17, 14637-14650.	3.8	50
3	A computational exploration of the crystal energy and charge-carrier mobility landscapes of the chiral [6]helicene molecule. Nanoscale, 2018, 10, 1865-1876.	5.6	48
4	Pervasive Delocalisation Error Causes Spurious Proton Transfer in Organic Acidâ€“Base Coâ€Crystals. Angewandte Chemie - International Edition, 2018, 57, 14906-14910.	13.8	45
5	What is â€œmany-bodyâ€•dispersion and should I worry about it?. Physical Chemistry Chemical Physics, 2020, 22, 8266-8276.	2.8	30
6	Torquoselectivity in the Nazarov Reactions of Allenyl Vinyl Ketones. Journal of Organic Chemistry, 2015, 80, 1042-1051.	3.2	29
7	Structureâ€“activity relationship of caffeic acid phenethyl ester analogs as new 5â€lipoxygenase inhibitors. Chemical Biology and Drug Design, 2017, 89, 514-528.	3.2	24
8	Dispersion XDM with Hybrid Functionals: Delocalization Error and Halogen Bonding in Molecular Crystals. Journal of Chemical Theory and Computation, 2019, 15, 4933-4944.	5.3	22
9	NMR Metabolomics Analysis of the Effects of 5-Lipoxygenase Inhibitors on Metabolism in Glioblastomas. Journal of Proteome Research, 2013, 12, 2165-2176.	3.7	19
10	Composite and Low-Cost Approaches for Molecular Crystal Structure Prediction. Journal of Chemical Theory and Computation, 2018, 14, 2265-2276.	5.3	18
11	Evaluation of Shear-Slip Transitions in Crystalline Aspirin by Density-Functional Theory. Crystal Growth and Design, 2016, 16, 6867-6873.	3.0	17
12	Asymptotic Pairwise Dispersion Corrections Can Describe Layered Materials Accurately. Journal of Physical Chemistry Letters, 2020, 11, 2298-2302.	4.6	17
13	Crystal-energy landscapes of active pharmaceutical ingredients using composite approaches. CrystEngComm, 2019, 21, 5995-6009.	2.6	14
14	Pervasive Delocalisation Error Causes Spurious Proton Transfer in Organic Acidâ€“Base Coâ€Crystals. Angewandte Chemie, 2018, 130, 15122-15126.	2.0	10
15	A Universal Force Field for Materials, Periodic GFN-FF: Implementation and Examination. Journal of Chemical Theory and Computation, 2021, 17, 7827-7849.	5.3	10
16	Non-Covalent Interactions in Molecular Crystals: Exploring the Accuracy of the Exchange-Hole Dipole Moment Model with Local Orbitals. Journal of Chemical Theory and Computation, 2018, 14, 5715-5724.	5.3	9
17	Clusters in Liquid Fatty Acids: Structure and Role in Nucleation. Journal of Physical Chemistry B, 2019, 123, 7043-7054.	2.6	8
18	Density Functional Theory Study of BF₃-Mediated Additions of Enols and [(Trimethylsilyl)oxy]alkenes to an Oxyallyl Cation: Homologous Mukaiyama Reactions. Journal of Physical Chemistry A, 2015, 119, 6714-6722.	2.5	6

#	Article	IF	Citations
19	Competing fragmentation processes of Î²â€substituted propanoate ions upon collisionâ€induced dissociation. Rapid Communications in Mass Spectrometry, 2016, 30, 2133-2144.	1.5	4
20	Phenyl group participation in rearrangements during collision-induced dissociation of deprotonated phenoxyacetic acid. Rapid Communications in Mass Spectrometry, 2015, 29, 2293-2301.	1.5	3
21	Ketone Analog of Caffeic Acid Phenethyl Ester Exhibits Antioxidant Activity via Activation of ERK-Dependent Nrf2 Pathway. Applied Sciences (Switzerland), 2022, 12, 3062.	2.5	2