

# Che R Seabourne

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/10884277/publications.pdf>

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10  
papers

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citations

1040056

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1372567

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docs citations

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times ranked

1140  
citing authors

#	ARTICLE	IF	CITATIONS
1	Probing the Bonding and Electronic Structure of Single Atom Dopants in Graphene with Electron Energy Loss Spectroscopy. <i>Nano Letters</i> , 2013, 13, 4989-4995.	9.1	187
2	Stacking Variants and Superconductivity in the Bi <sup>2</sup> Se <sub>3</sub> System. <i>Journal of the American Chemical Society</i> , 2013, 135, 5372-5374.	13.7	80
3	Electronic Structure Modification of Ion Implanted Graphene: The Spectroscopic Signatures of p- and n-Type Doping. <i>ACS Nano</i> , 2015, 9, 11398-11407.	14.6	75
4	Subangstrom Edge Relaxations Probed by Electron Microscopy in Hexagonal Boron Nitride. <i>Physical Review Letters</i> , 2012, 109, 205502.	7.8	52
5	Incisive Probing of Intermolecular Interactions in Molecular Crystals: Core Level Spectroscopy Combined with Density Functional Theory. <i>Journal of Physical Chemistry B</i> , 2014, 118, 12121-12129.	2.6	28
6	Local Plasmon Engineering in Doped Graphene. <i>ACS Nano</i> , 2018, 12, 1837-1848.	14.6	25
7	A systematic approach to choosing parameters for modelling fine structure in electron energy-loss spectroscopy. <i>Ultramicroscopy</i> , 2009, 109, 1374-1388.	1.9	23
8	Atomic-Scale Surface Roughness of Rutile and Implications for Organic Molecule Adsorption. <i>Langmuir</i> , 2013, 29, 6876-6883.	3.5	16
9	Energy of Step Defects on the TiO <sub>2</sub> Rutile (110) Surface: An ab initio DFT Methodology. <i>Journal of Physical Chemistry C</i> , 2013, 117, 23766-23780.	3.1	13
10	Analysis of computational EELS modelling results for MgO-based systems. <i>Ultramicroscopy</i> , 2010, 110, 1059-1069.	1.9	5