## James G M Hooper

List of Publications by Year in descending order

Source: https://exaly.com/author-pdf/9028965/publications.pdf

Version: 2024-02-01

46 1,372 19
papers citations h-index

52 52 52 1542 all docs docs citations times ranked citing authors

36

g-index

#	Article	IF	CITATIONS
1	Graphene-like Boron–Carbon–Nitrogen Monolayers. ACS Nano, 2017, 11, 2486-2493.	14.6	154
2	Metallization of magnesium polyhydrides under pressure. Physical Review B, 2013, 87, .	3.2	102
3	Polyhydrides of the Alkaline Earth Metals: A Look at the Extremes under Pressure. Journal of Physical Chemistry C, 2013, 117, 2982-2992.	3.1	84
4	Layered single-bonded nonmolecular phase of nitrogen from first-principles simulation. Physical Review B, 2005, 72, .	3.2	82
5	Rubidium Polyhydrides Under Pressure: Emergence of the Linear H <sub>3</sub> <sup>â^'</sup> Species. Chemistry - A European Journal, 2012, 18, 5013-5021.	3.3	68
6	High Pressure Potassium Polyhydrides: A Chemical Perspective. Journal of Physical Chemistry C, 2012, 116, 13322-13328.	3.1	63
7	Low-pressure metastable phase of single-bonded polymeric nitrogen from a helical structure motif and first-principles calculations. Physical Review B, 2007, 75, .	3.2	60
8	Systematic Method to New Phases of Polymeric Nitrogen under High Pressure. Physical Review Letters, 2006, 97, 155503.	7.8	59
9	Composition and Constitution of Compressed Strontium Polyhydrides. Journal of Physical Chemistry C, 2014, 118, 6433-6447.	3.1	59
10	Locking and Unlocking the Molecular Spin Crossover Transition. Advanced Materials, 2017, 29, 1702257.	21.0	55
11	Compressed Cesium Polyhydrides: Cs <sup>+</sup> Sublattices and H <sub>3</sub> <sup>–</sup> Three-Connected Nets. Inorganic Chemistry, 2012, 51, 9333-9342.	4.0	54
12	A DFT+U study of defect association and oxygen migration in samarium-doped ceria. Physical Chemistry Chemical Physics, 2011, 13, 6116.	2.8	51
13	Self-assembly of strongly dipolar molecules on metal surfaces. Journal of Chemical Physics, 2015, 142, 101921.	3.0	38
14	Lithium Subhydrides under Pressure and Their Superatomâ€like Building Blocks. ChemPlusChem, 2012, 77, 969-972.	2.8	30
15	Polymorphism driven optical properties of an anil dye. CrystEngComm, 2016, 18, 7249-7259.	2.6	29
16	Origin of Diastereoselectivity in the Tandem Oxy-Cope/Claisen/Ene Reaction:Â Experimental and Theoretical Studies of the Ring Inversion Mechanism. Journal of the American Chemical Society, 2007, 129, 2112-2119.	13.7	25
17	Pressure induced structural transitions in KH, RbH, and CsH. Journal of Applied Physics, 2012, 111, 112611.	2.5	23
18	Proton transfer in surface-stabilized chiral motifs of croconic acid. Physical Review B, 2013, 87, .	3.2	22

#	Article	IF	Citations
19	Coverage-Dependent Interactions at the Organics–Metal Interface: Quinonoid Zwitterions on Au(111). Journal of Physical Chemistry C, 2013, 117, 16406-16415.	3.1	21
20	Anion–π recognition between [M(CN) <sub>6</sub> ] <sup>3−</sup> complexes and HAT(CN) <sub>6</sub> : structural matching and electronic charge density modification. Dalton Transactions, 2017, 46, 3482-3491.	3.3	20
21	Mesomorphic phase transitions of 3F7HPhF studied by complementary methods. Phase Transitions, 2018, 91, 186-198.	1.3	20
22	Kagome-like lattice of π–π stacked 3-hydroxyphenalenone on Cu(111). Chemical Communications, 2014, 50, 8659-8662.	4.1	19
23	2D Cocrystallization from H-Bonded Organic Ferroelectrics. Journal of Physical Chemistry Letters, 2016, 7, 435-440.	4.6	19
24	Computational insights into the nature of increased ionic conductivity in concentrated samarium-doped ceria: a genetic algorithm study. Physical Chemistry Chemical Physics, 2010, 12, 12969.	2.8	18
25	Resonance Assisted Hydrogen Bonding Phenomenon Unveiled through Both Experiments and Theory: A New Family of Ethyl Nâ€Salicylideneglycinate Dyes. Chemistry - A European Journal, 2020, 26, 12987-12995.	3.3	18
26	Rhodizonic Acid on Noble Metals: Surface Reactivity and Coordination Chemistry. Journal of Physical Chemistry Letters, 2013, 4, 3413-3419.	4.6	14
27	Genetic algorithm and first-principles DFT study of the high-pressure molecularζphase of nitrogen. Physical Review B, 2009, 80, .	3.2	13
28	Electronic Structure of Iron Porphyrin Adsorbed to the Pt(111) Surface. Journal of Physical Chemistry C, 2016, 120, 29173-29181.	3.1	13
29	On the Origin of Altered Diastereomeric Ratios for Anionic versus Neutral Reaction Conditions in the Oxyâ€Cope/Ene Reaction: An Interplay of Experiment and Computational Modeling. Chemistry - A European Journal, 2010, 16, 14124-14130.	3.3	11
30	Genetic algorithm based approach to investigate doped metal oxide materials: Application to lanthanide-doped ceria. Physical Review B, $2010,81,\ldots$	3.2	11
31	Modulating Bond Lengths via Backdonation: A First-Principles Investigation of a Quinonoid Zwitterion Adsorbed to Coinage Metal Surfaces. Journal of Physical Chemistry C, 2016, 120, 6633-6641.	3.1	11
32	Anion-Ï€ Architectures of HAT(CN) <sub>6</sub> and 5d Polycyanidometalates: [W(CN) <sub>8</sub> ] <sup>3â€"</sup> , [Re(CN) <sub>7</sub> ] <sup>3â€"</sup> , and [Pt(CN) <sub>6</sub> ] <sup>2â€"</sup> . Crystal Growth and Design, 2019, 19, 1215-1225.	3.0	11
33	On the Nature of Ge–Pb Bonding in the Solid State. Synthesis, Structural Characterization, and Electronic Structures of Two Unprecedented Germanide-Plumbides. Journal of the American Chemical Society, 2012, 134, 12708-12716.	13.7	10
34	Molecular Deformation, Charge Flow, and Spongelike Behavior in Anion–π {[M(CN) <sub>4</sub> ] <sup>2â^³</sup> ;[HAT(CN) <sub>6</sub> ]} <sub>â~ž</sub> (M=Ni, Pd, Pt) Supramolecular Stacks. Chemistry - A European Journal, 2018, 24, 16302-16314.	3.3	10
35	Influence of fluorosubstitution on physical properties of the smectogenic chiral ester. Materials Research Bulletin, 2022, 150, 111756.	5.2	10
36	Interplay between Hydrogen Bonding, Epitaxy, and Charge Transfer in the Self-Assembly of Croconic Acid on Au(111) and Ag(111). Journal of Physical Chemistry C, 2015, 119, 26429-26437.	3.1	9

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37	Structure and Proton-Transfer Mechanism in One-Dimensional Chains of Benzimidazoles. Journal of Physical Chemistry C, 2016, 120, 5804-5809.	3.1	8
38	Seeking Out Heterogeneous Hydrogen Bonding in a Self-Assembled 2D Cocrystal of Croconic Acid and Benzimidazole on Au(111). Journal of Physical Chemistry C, 2021, 125, 2403-2410.	3.1	8
39	Chiral surface networks of 3-HPLN — A molecular analog of rounded triangle assembly. Surface Science, 2014, 629, 65-74.	1.9	7
40	On the Border between Low-Nuclearity and One-Dimensional Solids: A Unique Interplay of 1,2,4-Triazolyl-Based {Cu <sup>II</sup> <sub>5</sub> (OH) <sub>2</sub> } Clusters and Mo <sup>VI</sup> -Oxide Matrix. Inorganic Chemistry, 2018, 57, 6076-6083.	4.0	7
41	Urban Particulate Matterâ€Induced Decomposition of <i>S</i> à€Nitrosoglutathione Relevant to Aberrant Nitric Oxide Biological Signaling. ChemSusChem, 2019, 12, 661-671.	6.8	7
42	Binding of anionic Pt( <scp>ii</scp> ) complexes in a dedicated organic matrix: towards new binary crystalline composites. Dalton Transactions, 2021, 50, 170-185.	3.3	7
43	A concerted evolution of supramolecular interactions in a {cation; metal complex; π-acid; solvent} anion-π system. Inorganic Chemistry Frontiers, 2020, 7, 1851-1863.	6.0	6
44	Exploring "Triazole-Thiourea―Based Ligands for the Self-Assembly of Photoluminescent Hg(II) Coordination Compounds. Crystal Growth and Design, 2021, 21, 3562-3581.	3.0	5
45	First principles investigation on how site preference and entropy affect the stability of (Eu <sub><i>x</i></sub> Ge <sub>2</sub> Pb (M = Ca,)	[j <b>ET.Q</b> q1 ]	1 <b>0.₹</b> 84314 rg
46	Molecular Deformation, Charge Flow, and Spongelike Behavior in Anion–π {[M(CN) <sub>4</sub> ] <sup>2â^²</sup> ;[HAT(CN) <sub>6</sub> ]} <sub>â^ž</sub> (M=Ni, Pd, Pt) Supramolecular Stacks. Chemistry - A European Journal, 2018, 24, 16195-16195.	3.3	0