

Damien Thompson

List of Publications by Year in Descending Order

Source: <https://exaly.com/author-pdf/8301562/damien-thompson-publications-by-year.pdf>

Version: 2024-04-23

This document has been generated based on the publications and citations recorded by exaly.com. For the latest version of this publication list, visit the link given above.

The third column is the impact factor (IF) of the journal, and the fourth column is the number of citations of the article.

127
papers

2,738
citations

28
h-index

46
g-index

145
ext. papers

3,501
ext. citations

9.5
avg, IF

5.45
L-index

#	Paper	IF	Citations
127	Balanced lipase interactions for degradation-controlled paclitaxel release from lipid cubic phase formulations. <i>Journal of Colloid and Interface Science</i> , 2022 , 607, 978-991	9.3	0
126	Predictive Modeling of Neurotoxic β Synuclein Polymorphs.. <i>Methods in Molecular Biology</i> , 2022 , 2340, 379-399	1.4	0
125	Characterization of Amyloidogenic Peptide Aggregability in Helical Subspace.. <i>Methods in Molecular Biology</i> , 2022 , 2340, 401-448	1.4	0
124	Stable Universal 1- and 2-Input Single-Molecule Logic Gates.. <i>Advanced Materials</i> , 2022 , e2202135	2.4	0
123	Large cooperative effects in tunneling rates across van der Waals coupled binary self-assembled monolayers. <i>Nano Today</i> , 2022 , 44, 101497	17.9	2
122	Modulating the pro-apoptotic activity of cytochrome c at a biomimetic electrified interface. <i>Science Advances</i> , 2021 , 7, eabg4119	14.3	0
121	Extended Lifetime of Molecules Adsorbed onto Excipients Drives Nucleation in Heterogeneous Crystallization.. <i>Crystal Growth and Design</i> , 2021 , 21, 2101-2112	3.5	1
120	Understanding solid-state processing of pharmaceutical cocrystals via milling: Role of tablet excipients. <i>International Journal of Pharmaceutics</i> , 2021 , 601, 120514	6.5	8
119	Piezoelectricity of the Transmembrane Protein β 3 Cytochrome c Oxidase. <i>Advanced Functional Materials</i> , 2021 , 31, 2100884	15.6	1
118	A practical approach for standardization of converse piezoelectric constants obtained from piezoresponse force microscopy. <i>Journal of Applied Physics</i> , 2021 , 129, 185104	2.5	1
117	Molecular engineering of piezoelectricity in collagen-mimicking peptide assemblies. <i>Nature Communications</i> , 2021 , 12, 2634	17.4	11
116	A single atom change turns insulating saturated wires into molecular conductors. <i>Nature Communications</i> , 2021 , 12, 3432	17.4	5
115	Modulation of physical properties of organic cocrystals by amino acid chirality. <i>Materials Today</i> , 2021 , 42, 29-40	21.8	9
114	Nanoconfined water governs polarization-related properties of self-assembled peptide nanotubes. <i>Nano Select</i> , 2021 , 2, 817-829	3.1	6
113	Molecular Modelling Guided Modulation of Molecular Shape and Charge for Design of Smart Self-Assembled Polymeric Drug Transporters. <i>Pharmaceutics</i> , 2021 , 13,	6.4	2
112	Self-Assembled Pyrene Stacks and Peptide Monolayers Tune the Electronic Properties of Functionalized Electrolyte-Gated Graphene Field-Effect Transistors. <i>ACS Applied Materials & Interfaces</i> , 2021 , 13, 9134-9142	9.5	3
111	Energy-Level Alignment and Orbital-Selective Femtosecond Charge Transfer Dynamics of Redox-Active Molecules on Au, Ag, and Pt Metal Surfaces. <i>Journal of Physical Chemistry C</i> , 2021 , 125, 18474-18482	3.8	0

110	Decision trees within a molecular memristor. <i>Nature</i> , 2021 , 597, 51-56	50.4	19
109	A Piezoelectric Ionic Cocystal of Glycine and Sulfamic Acid. <i>Crystal Growth and Design</i> , 2021 , 21, 5818-5827	3.7	3
108	Predictive Modeling of Ceramic Materials 2021 , 475-480		
107	Restriction boosts piezoelectricity. <i>Nature Materials</i> , 2021 , 20, 574-575	27	2
106	Electric-field-driven dual-functional molecular switches in tunnel junctions. <i>Nature Materials</i> , 2020 , 19, 843-848	27	54
105	Long-range Regulation of Partially Folded Amyloidogenic Peptides. <i>Scientific Reports</i> , 2020 , 10, 7597	4.9	5
104	Charge disproportionate molecular redox for discrete memristive and memcapacitive switching. <i>Nature Nanotechnology</i> , 2020 , 15, 380-389	28.7	37
103	Colossal current and voltage tunability in an organic memristor via electrode engineering. <i>Applied Materials Today</i> , 2020 , 19, 100626	6.6	11
102	Complete aggregation pathway of amyloid α (1-40) and (1-42) resolved on an atomically clean interface. <i>Science Advances</i> , 2020 , 6, eaaz6014	14.3	28
101	Diphenylalanine-Derivative Peptide Assemblies with Increased Aromaticity Exhibit Metal-like Rigidity and High Piezoelectricity. <i>ACS Nano</i> , 2020 , 14, 7025-7037	16.7	18
100	Anchoring and packing of self-assembled monolayers of semithio-bambusurils on Au(111). <i>Molecular Systems Design and Engineering</i> , 2020 , 5, 511-520	4.6	1
99	Accelerated charge transfer in water-layered peptide assemblies. <i>Energy and Environmental Science</i> , 2020 , 13, 96-101	35.4	21
98	Quantitative Polarization-Resolved Second-Harmonic-Generation Microscopy of Glycine Microneedles. <i>Advanced Materials</i> , 2020 , 32, e2002873	24	3
97	Design principles of dual-functional molecular switches in solid-state tunnel junctions. <i>Applied Physics Letters</i> , 2020 , 117, 030502	3.4	8
96	Piezoelectricity in the proteinogenic amino acid L-leucine: A novel piezoactive bioelectret. <i>IEEE Transactions on Dielectrics and Electrical Insulation</i> , 2020 , 27, 1465-1468	2.3	1
95	Tunable Mechanical and Optoelectronic Properties of Organic Cocystals by Unexpected Stacking Transformation from H- to J- and X-Aggregation. <i>ACS Nano</i> , 2020 , 14, 10704-10715	16.7	18
94	Atomistic-Benchmarking towards a protocol development for rapid quantitative metrology of piezoelectric biomolecular materials. <i>Applied Materials Today</i> , 2020 , 21, 100818	6.6	8
93	Racemic Amino Acid Piezoelectric Transducer. <i>Physical Review Letters</i> , 2019 , 122, 047701	7.4	27

92	Bioinspired Stable and Photoluminescent Assemblies for Power Generation. <i>Advanced Materials</i> , 2019 , 31, e1807481	24	41
91	Phase-change memories (PCM) - Experiments and modelling: general discussion. <i>Faraday Discussions</i> , 2019 , 213, 393-420	3.6	3
90	On the ubiquity of helical β -synuclein tetramers. <i>Physical Chemistry Chemical Physics</i> , 2019 , 21, 12036-12043	3.6	14
89	Probing Crystal Nucleation of Fenoxycarb from Solution through the Effect of Solvent. <i>Crystal Growth and Design</i> , 2019 , 19, 2037-2049	3.5	22
88	Organic piezoelectric materials: milestones and potential. <i>NPG Asia Materials</i> , 2019 , 11,	10.3	57
87	Direct measurement of the local field within alkyl-ferrocenyl-alkanethiolate monolayers: Importance of the supramolecular and electronic structure on the voltammetric response and potential profile. <i>Electrochimica Acta</i> , 2019 , 311, 92-102	6.7	9
86	The supramolecular structure and van der Waals interactions affect the electronic structure of ferrocenyl-alkanethiolate SAMs on gold and silver electrodes. <i>Nanoscale Advances</i> , 2019 , 1, 1991-2002	5.1	7
85	Molecular Simulations Reveal Terminal Group Mediated Stabilization of Helical Conformers in Both Amyloid- β 2 and β -synuclein. <i>ACS Chemical Neuroscience</i> , 2019 , 10, 2830-2842	5.7	19
84	Self-Assembled Cationic β -Cyclodextrin Nanostructures for siRNA Delivery. <i>Molecular Pharmaceutics</i> , 2019 , 16, 1358-1366	5.6	28
83	Monitoring transient changes in the structure of water at a polarised liquid-liquid interface using electrocapillary curves. <i>Electrochemistry Communications</i> , 2019 , 109, 106564	5.1	3
82	Molecular Diodes: Stable Molecular Diodes Based on π -Interactions of the Molecular Frontier Orbitals with Graphene Electrodes (Adv. Mater. 10/2018). <i>Advanced Materials</i> , 2018 , 30, 1870069	24	
81	On the distinct binding modes of expansin and carbohydrate-binding module proteins on crystalline and nanofibrous cellulose: implications for cellulose degradation by designer cellulosomes. <i>Physical Chemistry Chemical Physics</i> , 2018 , 20, 8278-8293	3.6	6
80	Graphene wrinkle effects on molecular resonance states. <i>Npj 2D Materials and Applications</i> , 2018 , 2,	8.8	7
79	Revisiting the earliest signatures of amyloidogenesis: Roadmaps emerging from computational modeling and experiment. <i>Wiley Interdisciplinary Reviews: Computational Molecular Science</i> , 2018 , 8, e1359	7.9	9
78	The fold preference and thermodynamic stability of β -synuclein fibrils is encoded in the non-amyloid- β component region. <i>Physical Chemistry Chemical Physics</i> , 2018 , 20, 4502-4512	3.6	11
77	Stable Molecular Diodes Based on π -Interactions of the Molecular Frontier Orbitals with Graphene Electrodes. <i>Advanced Materials</i> , 2018 , 30, 1706322	24	23
76	Ordering of Air-Oxidized Decanethiols on Au(111). <i>Journal of Physical Chemistry C</i> , 2018 , 122, 8430-8436	3.8	9
75	Longitudinal Piezoelectricity in Orthorhombic Amino Acid Crystal Films. <i>Crystal Growth and Design</i> , 2018 , 18, 4844-4848	3.5	16

74	. <i>IEEE Transactions on Dielectrics and Electrical Insulation</i> , 2018 , 25, 803-807	2.3	8
73	Subcellular Imaging of Liquid Silicone Coated-Intestinal Epithelial Cells. <i>Scientific Reports</i> , 2018 , 8, 107634.9	4.9	3
72	Deconstructing collagen piezoelectricity using alanine-hydroxyproline-glycine building blocks. <i>Nanoscale</i> , 2018 , 10, 9653-9663	7.7	20
71	Steered molecular dynamics simulations reveal the role of Ca in regulating mechanostability of cellulose-binding proteins. <i>Physical Chemistry Chemical Physics</i> , 2018 , 20, 22674-22680	3.6	4
70	Control of piezoelectricity in amino acids by supramolecular packing. <i>Nature Materials</i> , 2018 , 17, 180-186.7	6.7	118
69	Molecular simulations reveal that a short helical loop regulates thermal stability of type I cohesin-dockerin complexes. <i>Physical Chemistry Chemical Physics</i> , 2018 , 20, 28445-28451	3.6	2
68	The laterally acquired GH5 EngA from the marine bacterium is dedicated to hemicellulose hydrolysis. <i>Biochemical Journal</i> , 2018 , 475, 3609-3628	3.8	4
67	Re-designing the β -synuclein tetramer. <i>Chemical Communications</i> , 2018 , 54, 8080-8083	5.8	16
66	Familial Mutations May Switch Conformational Preferences in β -Synuclein Fibrils. <i>ACS Chemical Neuroscience</i> , 2017 , 8, 837-849	5.7	25
65	Motion of Fullerenes around Topological Defects on Metals: Implications for the Progress of Molecular Scale Devices. <i>ACS Applied Materials & Interfaces</i> , 2017 , 9, 7897-7902	9.5	3
64	Supramolecular Structure of the Monolayer Triggers Odd-Even Effects in the Tunneling Rates across Noncovalent Junctions on Graphene. <i>Journal of Physical Chemistry C</i> , 2017 , 121, 4172-4180	3.8	13
63	Controlling Protein Surface Orientation by Strategic Placement of Oligo-Histidine Tags. <i>ACS Nano</i> , 2017 , 11, 9068-9083	16.7	31
62	Non-local effects of point mutations on the stability of a protein module. <i>Journal of Chemical Physics</i> , 2017 , 147, 105101	3.9	4
61	The length but not the sequence of peptide linker modules exerts the primary influence on the conformations of protein domains in cellulosome multi-enzyme complexes. <i>Physical Chemistry Chemical Physics</i> , 2017 , 19, 21414-21425	3.6	13
60	Restoring the Electrical Properties of CVD Graphene via Physisorption of Molecular Adsorbates. <i>ACS Applied Materials & Interfaces</i> , 2017 , 9, 25014-25022	9.5	19
59	Molecular diodes with rectification ratios exceeding 10 driven by electrostatic interactions. <i>Nature Nanotechnology</i> , 2017 , 12, 797-803	28.7	148
58	Harnessing Nanoscale Physics for Next-Generation Electronic Medical Devices 2016 , 491-509		
57	A robust molecular probe for β gstrom-scale analytics in liquids. <i>Nature Communications</i> , 2016 , 7, 12403	17.4	3

56	Noncovalent Self-Assembled Monolayers on Graphene as a Highly Stable Platform for Molecular Tunnel Junctions. <i>Advanced Materials</i> , 2016 , 28, 631-9	24	35
55	Nanoscale Engineering of Designer Cellulosomes. <i>Advanced Materials</i> , 2016 , 28, 5619-47	24	35
54	Fingerprinting Electronic Molecular Complexes in Liquid. <i>Scientific Reports</i> , 2016 , 6, 19009	4.9	8
53	Molecular Electronics: Noncovalent Self-Assembled Monolayers on Graphene as a Highly Stable Platform for Molecular Tunnel Junctions (Adv. Mater. 4/2016). <i>Advanced Materials</i> , 2016 , 28, 784-784	24	3
52	Even the Odd Numbers Help: Failure Modes of SAM-Based Tunnel Junctions Probed via Odd-Even Effects Revealed in Synchrotrons and Supercomputers. <i>Accounts of Chemical Research</i> , 2016 , 49, 2061-2069	24.3	56
51	Controlling the direction of rectification in a molecular diode. <i>Nature Communications</i> , 2015 , 6, 6324	17.4	153
50	Formation Mechanism of Metal-Molecule-Metal Junctions: Molecule-Assisted Migration on Metal Defects. <i>Journal of Physical Chemistry C</i> , 2015 , 119, 19438-19451	3.8	10
49	One Carbon Matters: The Origin and Reversal of Odd-Even Effects in Molecular Diodes with Self-Assembled Monolayers of Ferrocenyl-Alkanethiolates. <i>Journal of Physical Chemistry C</i> , 2015 , 119, 17910-17919	3.8	50
48	One-Nanometer Thin Monolayers Remove the Deleterious Effect of Substrate Defects in Molecular Tunnel Junctions. <i>Nano Letters</i> , 2015 , 15, 6643-9	11.5	38
47	Nonideal Electrochemical Behavior of Ferrocenyl-Alkanethiolate SAMs Maps the Microenvironment of the Redox Unit. <i>Journal of Physical Chemistry C</i> , 2015 , 119, 21978-21991	3.8	41
46	On the hydration of subnanometric antifouling organosilane adlayers: a molecular dynamics simulation. <i>Journal of Colloid and Interface Science</i> , 2015 , 437, 197-204	9.3	28
45	Peptide Recognition Capabilities of Cellulose in Molecular Dynamics Simulations. <i>Journal of Physical Chemistry C</i> , 2015 , 119, 24404-24416	3.8	13
44	Capturing the embryonic stages of self-assembly - design rules for molecular computation. <i>Scientific Reports</i> , 2015 , 5, 10116	4.9	14
43	Physisorption Controls the Conformation and Density of States of an Adsorbed Porphyrin. <i>Journal of Physical Chemistry C</i> , 2015 , 119, 27982-27994	3.8	24
42	On the remarkable role of surface topography of the bottom electrodes in blocking leakage currents in molecular diodes. <i>Journal of the American Chemical Society</i> , 2014 , 136, 6554-7	16.4	77
41	Nanoelectrical analysis of single molecules and atomic-scale materials at the solid/liquid interface. <i>Nature Materials</i> , 2014 , 13, 947-53	27	27
40	Reassigning the most stable surface of hydroxyapatite to the water resistant hydroxyl terminated (010) surface. <i>Surface Science</i> , 2014 , 623, 55-63	1.8	17
39	Mechanostability of cohesin-dockerin complexes in a structure-based model: anisotropy and lack of universality in the force profiles. <i>Journal of Chemical Physics</i> , 2014 , 141, 245103	3.9	7

38	Frontiers of Cu Electrodeposition and Electroless Plating for On-chip Interconnects. <i>Nanostructure Science and Technology</i> , 2014 , 99-113	0.9	2
37	Large directional conductivity change in chemically stable layered thin films of vanadium oxide and a 1D metal complex. <i>Journal of Materials Chemistry C</i> , 2013 , 1, 5675	7.1	19
36	Nanoscale dynamics and protein adhesivity of alkylamine self-assembled monolayers on graphene. <i>Langmuir</i> , 2013 , 29, 7271-82	4	23
35	The role of van der Waals forces in the performance of molecular diodes. <i>Nature Nanotechnology</i> , 2013 , 8, 113-8	28.7	245
34	A multi-scale molecular dynamics study of the assembly of micron-size supraparticles from 30 nm alkyl-coated nanoparticles. <i>Physical Chemistry Chemical Physics</i> , 2013 , 15, 8132-43	3.6	7
33	Non-Covalent Functionalization of Graphene Using Self-Assembly of Alkane-Amines. <i>Advanced Functional Materials</i> , 2012 , 22, 717-725	15.6	57
32	Nanoparticle-based drug delivery: case studies for cancer and cardiovascular applications. <i>Cellular and Molecular Life Sciences</i> , 2012 , 69, 389-404	10.3	64
31	Scanning the potential energy surface for synthesis of dendrimer-wrapped gold clusters: design rules for true single-molecule nanostructures. <i>ACS Nano</i> , 2012 , 6, 3007-17	16.7	25
30	Monofunctionalized gold nanoparticles stabilized by a single dendrimer form dumbbell structures upon homocoupling. <i>Journal of the American Chemical Society</i> , 2012 , 134, 14674-7	16.4	39
29	Accommodating Curvature in a Highly Ordered Functionalized Metal Oxide Nanofiber: Synthesis, Characterization, and Multiscale Modeling of Layered Nanosheets. <i>Chemistry of Materials</i> , 2012 , 24, 3987-3992	9.6	15
28	Gradient-driven motion of multivalent ligand molecules along a surface functionalized with multiple receptors. <i>Nature Chemistry</i> , 2011 , 3, 317-22	17.6	86
27	Molecular dynamics of the "hydrophobic patch" that immobilizes hydrophobin protein HFBII on silicon. <i>Journal of Molecular Modeling</i> , 2011 , 17, 2227-35	2	14
26	Gold nanoparticles stabilized by thioether dendrimers. <i>Chemistry - A European Journal</i> , 2011 , 17, 13473-81	4.8	37
25	Electronic structure calculations and physicochemical experiments quantify the competitive liquid ion association and probe stabilisation effects for nitrobenzospirropyran in phosphonium-based ionic liquids. <i>Physical Chemistry Chemical Physics</i> , 2011 , 13, 6156-68	3.6	16
24	Interdigitating organic bilayers direct the short interlayer spacing in hybrid organic-inorganic layered vanadium oxide nanostructures. <i>Journal of Physical Chemistry B</i> , 2011 , 115, 14518-25	3.4	17
23	Linker-mediated assembly of gold nanoparticles into multimeric motifs. <i>Nanotechnology</i> , 2011 , 22, 44560-4	9.1	16
22	Molecular dynamics study of naturally occurring defects in self-assembled monolayer formation. <i>ACS Nano</i> , 2010 , 4, 921-32	16.7	33
21	Alchemical free energy simulations for biological complexes: powerful but temperamental. <i>Journal of Molecular Recognition</i> , 2010 , 23, 117-27	2.6	46

20	Interaction of acridine-calix[4]arene with DNA at the electrified liquid liquid interface. <i>Electrochimica Acta</i> , 2010 , 55, 3348-3354	6.7	22
19	Quantification of ink diffusion in microcontact printing with self-assembled monolayers. <i>Langmuir</i> , 2009 , 25, 242-7	4	22
18	Monolayer Packing, Dehydration, and Ink-Binding Dynamics at the Molecular Printboard. <i>Journal of Physical Chemistry C</i> , 2009 , 113, 7298-7304	3.8	5
17	Guanidinium chloride molecular diffusion in aqueous and mixed water-ethanol solutions. <i>Journal of Physical Chemistry B</i> , 2008 , 112, 8906-11	3.4	16
16	The effective concentration of unbound ink anchors at the molecular printboard. <i>Journal of Physical Chemistry B</i> , 2008 , 112, 4994-9	3.4	5
15	Computer simulations reveal a novel nucleotide-type binding orientation for ellipticine-based anticancer c-kit kinase inhibitors. <i>Biochemistry</i> , 2008 , 47, 10333-44	3.2	17
14	Coarse-grained molecular dynamics simulations of nanopatterning with multivalent inks. <i>Journal of Chemical Physics</i> , 2008 , 128, 234906	3.9	11
13	Probing electrostatic interactions and ligand binding in aspartyl-tRNA synthetase through site-directed mutagenesis and computer simulations. <i>Proteins: Structure, Function and Bioinformatics</i> , 2008 , 71, 1450-60	4.2	12
12	Hydrocarbon Selective Oxidation on Vanadium Phosphorus Oxide Catalysts: Insights from Electronic Structure Calculations. <i>Topics in Catalysis</i> , 2008 , 50, 116-123	2.3	6
11	Free energy balance predicates dendrimer binding multivalency at molecular printboards. <i>Langmuir</i> , 2007 , 23, 8441-51	4	16
10	In silico engineering of tailored ink-binding ability at molecular printboards. <i>ChemPhysChem</i> , 2007 , 8, 1684-93	3.2	5
9	Ammonium scanning in an enzyme active site. The chiral specificity of aspartyl-tRNA synthetase. <i>Journal of Biological Chemistry</i> , 2007 , 282, 30856-68	5.4	22
8	Free-energy simulations and experiments reveal long-range electrostatic interactions and substrate-assisted specificity in an aminoacyl-tRNA synthetase. <i>ChemBioChem</i> , 2006 , 7, 337-44	3.8	34
7	Molecular dynamics simulations show that bound Mg ²⁺ contributes to amino acid and aminoacyl adenylate binding specificity in aspartyl-tRNA synthetase through long range electrostatic interactions. <i>Journal of Biological Chemistry</i> , 2006 , 281, 23792-803	5.4	33
6	Modeling competitive guest binding to beta-cyclodextrin molecular printboards. <i>Journal of Physical Chemistry B</i> , 2006 , 110, 16640-5	3.4	26
5	Electronic structure of the extended vanadyl pyrophosphate (1 0 0) surface. <i>Catalysis Today</i> , 2004 , 91-92, 177-180	5.3	6
4	A DFT periodic study of the vanadyl pyrophosphate (100) surface. <i>Surface Science</i> , 2003 , 547, 438-451	1.8	9
3	Modelling the active sites in vanadyl pyrophosphate. <i>Journal of Molecular Catalysis A</i> , 2003 , 198, 125-137		22

- 2 The interplay of electrostatic and covalent effects in 1-butene oxidation over vanadyl pyrophosphate. *Journal of Molecular Catalysis A*, **2003**, 206, 435-439 8
- 1 Modulating vectored non-covalent interactions for layered assembly with engineerable properties. *Bio-Design and Manufacturing*, 1 4-7 ○