

KartEEK K Bejagam

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/6540081/publications.pdf>

Version: 2024-02-01

30
papers

1,050
citations

394421

19
h-index

434195

31
g-index

31
all docs

31
docs citations

31
times ranked

1314
citing authors

#	ARTICLE	IF	CITATIONS
1	Biomimetic temporal self-assembly via fuel-driven controlled supramolecular polymerization. <i>Nature Communications</i> , 2018, 9, 1295.	12.8	148
2	Dipole-Moment-Driven Cooperative Supramolecular Polymerization. <i>Journal of the American Chemical Society</i> , 2015, 137, 3924-3932.	13.7	115
3	Solvent Clathrate Driven Dynamic Stereomutation of a Supramolecular Polymer with Molecular Pockets. <i>Journal of the American Chemical Society</i> , 2017, 139, 13867-13875.	13.7	86
4	Supramolecular Polymerization of Benzene-1,3,5-tricarboxamide: A Molecular Dynamics Simulation Study. <i>Journal of Physical Chemistry B</i> , 2014, 118, 5218-5228.	2.6	61
5	Autoresolution of Segregated and Mixed π Stacks by Stereoselective Supramolecular Polymerization in Solution. <i>Angewandte Chemie - International Edition</i> , 2015, 54, 13053-13057.	13.8	61
6	Machine-Learned Coarse-Grained Models. <i>Journal of Physical Chemistry Letters</i> , 2018, 9, 4667-4672.	4.6	48
7	Dissolution of Cellulose in Room Temperature Ionic Liquids: Anion Dependence. <i>Journal of Physical Chemistry B</i> , 2015, 119, 1654-1659.	2.6	44
8	PSO-Assisted Development of New Transferable Coarse-Grained Water Models. <i>Journal of Physical Chemistry B</i> , 2018, 122, 1958-1971.	2.6	39
9	Supramolecular Polymerization: A Coarse Grained Molecular Dynamics Study. <i>Journal of Physical Chemistry B</i> , 2015, 119, 5738-5746.	2.6	38
10	Machine-Learning Enabled New Insights into the Coil-to-Globule Transition of Thermosensitive Polymers Using a Coarse-Grained Model. <i>Journal of Physical Chemistry Letters</i> , 2018, 9, 6480-6488.	4.6	34
11	Host-Guest [2+2] Cycloaddition Reaction: Postsynthetic Modulation of CO ₂ Selectivity and Magnetic Properties in a Bimodal Metal-Organic Framework. <i>Chemistry - A European Journal</i> , 2016, 22, 7792-7799.	3.3	30
12	Nanoparticle activated and directed assembly of graphene into a nanoscroll. <i>Carbon</i> , 2018, 134, 43-52.	10.3	29
13	Development of New Transferable Coarse-Grained Models of Hydrocarbons. <i>Journal of Physical Chemistry B</i> , 2018, 122, 7143-7153.	2.6	28
14	Unraveling the Conformations of Backbone and Side Chains in Thermosensitive Bottlebrush Polymers. <i>Macromolecules</i> , 2019, 52, 9398-9408.	4.8	28
15	Machine-Learning Based Stacked Ensemble Model for Accurate Analysis of Molecular Dynamics Simulations. <i>Journal of Physical Chemistry A</i> , 2019, 123, 5190-5198.	2.5	26
16	External electric field reverses helical handedness of a supramolecular columnar stack. <i>Chemical Communications</i> , 2015, 51, 16049-16052.	4.1	22
17	Development of an Accurate Coarse-Grained Model of Poly(acrylic acid) in Explicit Solvents. <i>Macromolecules</i> , 2019, 52, 4875-4887.	4.8	21
18	Solvation dynamics of <i>N</i> -substituted acrylamide polymers and the importance for phase transition behavior. <i>Soft Matter</i> , 2020, 16, 1582-1593.	2.7	20

#	ARTICLE	IF	CITATIONS
19	Molecular dynamics simulations for glass transition temperature predictions of polyhydroxyalkanoate biopolymers. <i>Physical Chemistry Chemical Physics</i> , 2020, 22, 17880-17889.	2.8	19
20	Development of non-bonded interaction parameters between graphene and water using particle swarm optimization. <i>Journal of Computational Chemistry</i> , 2018, 39, 721-734.	3.3	18
21	Development of transferable coarse-grained models of amino acids. <i>Molecular Systems Design and Engineering</i> , 2020, 5, 675-685.	3.4	16
22	Machine Learning for Melting Temperature Predictions and Design in Polyhydroxyalkanoate-Based Biopolymers. <i>Journal of Physical Chemistry B</i> , 2022, 126, 934-945.	2.6	15
23	Development of Transferable Nonbonded Interactions between Coarse-Grained Hydrocarbon and Water Models. <i>Journal of Physical Chemistry B</i> , 2019, 123, 909-921.	2.6	12
24	Understanding the self-assembly of amino ester-based benzene-1,3,5-tricarboxamides using molecular dynamics simulations. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 258-266.	2.8	11
25	Composition and Configuration Dependence of Glass-Transition Temperature in Binary Copolymers and Blends of Polyhydroxyalkanoate Biopolymers. <i>Macromolecules</i> , 2021, 54, 5618-5628.	4.8	11
26	Supramolecular Polymerization of <i>N,N,N',N'-tetra-(Tetradecyl)-1,3,6,8-pyrenetetra-carboxamide: A Computational Study</i> . <i>Journal of Physical Chemistry B</i> , 2017, 121, 11492-11503.	2.6	10
27	Development of non-bonded interaction parameters between hexagonal boron-nitride and water. <i>Computational Materials Science</i> , 2019, 161, 339-345.	3.0	10
28	Durable and highly selective ion transport of a sulfonated Diels Alder Poly(phenylene) for vanadium redox flow batteries. <i>Journal of Power Sources</i> , 2022, 520, 230805.	7.8	9
29	Predicting the Mechanical Response of Polyhydroxyalkanoate Biopolymers Using Molecular Dynamics Simulations. <i>Polymers</i> , 2022, 14, 345.	4.5	7
30	Dehydration of polymer chains initiates graphene folding in water. <i>Carbon</i> , 2021, 180, 244-253.	10.3	5