Monica H Lamm

List of Publications by Year in descending order

Source: https://exaly.com/author-pdf/545815/publications.pdf

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40 1,428 21 38 g-index

40 40 40 1813

times ranked

citing authors

docs citations

all docs

#	Article	IF	CITATIONS
1	Sampling Performance of Multiple Independent Molecular Dynamics Simulations of an RNA Aptamer. ACS Omega, 2020, 5, 20187-20201.	3.5	10
2	Common Secondary and Tertiary Structural Features of Aptamer–Ligand Interaction Shared by RNA Aptamers with Different Primary Sequences. Molecules, 2019, 24, 4535.	3.8	6
3	Pre-engineering Thinking and the Engineering Habits of Mind in Preschool Classroom. Early Childhood Education Journal, 2019, 47, 187-198.	2.7	25
4	Study strategies and beliefs about learning as a function of academic achievement and achievement goals. Memory, 2018, 26, 683-690.	1.7	69
5	Students' Use of Optional Online Reviews and Its Relationship to Summative Assessment Outcomes in Introductory Biology. CBE Life Sciences Education, 2017, 16, ar23.	2.3	18
6	Engineering Thinking in Prekindergarten Children: A Systematic Literature Review. Journal of Engineering Education, 2017, 106, 454-474.	3.0	25
7	Prequestions do not enhance the benefits of retrieval in a STEM classroom. Cognitive Research: Principles and Implications, 2017, 2, 42.	2.0	17
8	A Classroom Study on the Relationship Between Student Achievement and Retrieval-Enhanced Learning. Educational Psychology Review, 2016, 28, 353-375.	8.4	55
9	Measuring the Effectiveness of Team-Based Leaning Outcomes in a Human Factors Course. Proceedings of the Human Factors and Ergonomics Society, 2015, 59, 337-341.	0.3	1
10	Estimating Error in Diffusion Coefficients Derived from Molecular Dynamics Simulations. Journal of Chemical Theory and Computation, 2015, 11, 4586-4592.	5. 3	107
11	A coarse-graining approach for molecular simulation that retains the dynamics of the all-atom reference system by implementing hydrodynamic interactions. Journal of Chemical Physics, 2014, 141, 174107.	3.0	23
12	An adaptable pentaloop defines a robust neomycin-B RNA aptamer with conditional ligand-bound structures. Rna, 2014, 20, 815-824.	3. 5	12
13	Ab Initio Study of Molecular Interactions in Cellulose lÎ \pm . Journal of Physical Chemistry B, 2013, 117, 10430-10443.	2.6	22
14	Molecular dynamics simulations on the Tre1 G protein-coupled receptor: exploring the role of the arginine of the NRY motif in Tre1 structure. BMC Structural Biology, 2013, 13, 15.	2.3	6
15	Investigating the malleability of RNA aptamers. Methods, 2013, 63, 178-187.	3.8	14
16	Evaluation of coarse-grained mapping schemes for polysaccharide chains in cellulose. Journal of Chemical Physics, 2013, 138, 214108.	3.0	27
17	Multiscale Modeling for Host-Guest Chemistry of Dendrimers in Solution. Polymers, 2012, 4, 463-485.	4. 5	14
18	Dendrimer–Fullerenol Soft-Condensed Nanoassembly. Journal of Physical Chemistry C, 2012, 116, 15775-15781.	3.1	16

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19	A coarse-grained model for \hat{l}^2 -d-glucose based on force matching. Theoretical Chemistry Accounts, 2012, 131, 1.	1.4	10
20	A biophysical perspective of understanding nanoparticles at large. Physical Chemistry Chemical Physics, 2011, 13, 7273.	2.8	63
21	Reintroducing explicit solvent to a solvent-free coarse-grained model. Physical Review E, 2011, 84, 025701.	2.1	5
22	Molecular dynamics simulation of fractal aggregate diffusion. Physical Review E, 2010, 82, 051402.	2.1	14
23	Fluorescence resonance energy transfer between phenanthrene and PAMAM dendrimers. Physical Chemistry Chemical Physics, 2010, 12, 9285.	2.8	35
24	Cell Trafficking of Carbon Nanotubes Based on Fluorescence Detection. Methods in Molecular Biology, 2010, 625, 135-151.	0.9	13
25	Molecular dynamics simulation of mixed matrix nanocomposites containing polyimide and polyhedral oligomeric silsesquioxane (POSS). Polymer, 2009, 50, 1324-1332.	3.8	65
26	Computational and Experimental Analyses Converge to Reveal a Coherent Yet Malleable Aptamer Structure That Controls Chemical Reactivity. Journal of the American Chemical Society, 2009, 131, 14747-14755.	13.7	14
27	Coarse-Grained Intermolecular Potentials Derived From The Effective Fragment Potential: Application To Water, Benzene, And Carbon Tetrachloride. Challenges and Advances in Computational Chemistry and Physics, 2009, , 197-218.	0.6	3
28	Nanoparticle Ordering <i>via</i> Functionalized Block Copolymers in Solution. ACS Nano, 2008, 2, 1259-1265.	14.6	44
29	Computer Simulation of Architectural and Molecular Weight Effects on the Assembly of Amphiphilic Linearâ "Dendritic Block Copolymers in Solution. Langmuir, 2008, 24, 3030-3036.	3.5	38
30	Modeling Styreneâ^'Styrene Interactionsâ€. Journal of Physical Chemistry A, 2006, 110, 519-525.	2.5	36
31	Effect of Terminal Group Modification on the Solution Properties of Dendrimers:  A Molecular Dynamics Simulation Study. Macromolecules, 2006, 39, 4247-4255.	4.8	34
32	Biomolecule-directed assembly of nanoscale building blocks studied via lattice Monte Carlo simulation. Journal of Chemical Physics, 2004, 121, 3919-3929.	3.0	17
33	Effect of pressure on the complete phase behavior of binary mixtures. AICHE Journal, 2004, 50, 215-225.	3.6	13
34	Tethered Nano Building Blocks:Â Toward a Conceptual Framework for Nanoparticle Self-Assembly. Nano Letters, 2003, 3, 1341-1346.	9.1	311
35	Simulated Assembly of Nanostructured Organic/Inorganic Networks. Nano Letters, 2003, 3, 989-994.	9.1	69
36	Equilibria between solid, liquid, and vapor phases in binary Lennard–Jones mixtures. Fluid Phase Equilibria, 2002, 194-197, 197-206.	2.5	22

#	Article	IF	CITATION
37	Monte Carlo simulations of complete phase diagrams for binary Lennard–Jones mixtures. Fluid Phase Equilibria, 2001, 182, 37-46.	2.5	37
38	Molecular simulation of complete phase diagrams for binary mixtures. AICHE Journal, 2001, 47, 1664-1675.	3.6	37
39	Solid–liquid phase equilibrium for binary Lennard-Jones mixtures. Journal of Chemical Physics, 1999, 110, 11433-11444.	3.0	73
40	Team-Based Learning in Engineering Classrooms: Feedback Form and Content Adds Value to the Learning Experience. , 0, , .		8