

Hwankyu Lee

List of Publications by Year in descending order

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76
papers

3,386
citations

218662

26
h-index

144002

57
g-index

77
all docs

77
docs citations

77
times ranked

4682
citing authors

#	ARTICLE	IF	CITATIONS
1	The centrosomal protein nephrocystin-6 is mutated in Joubert syndrome and activates transcription factor ATF4. <i>Nature Genetics</i> , 2006, 38, 674-681.	21.4	535
2	Molecular Dynamics Studies of Polyethylene Oxide and Polyethylene Glycol: Hydrodynamic Radius and Shape Anisotropy. <i>Biophysical Journal</i> , 2008, 95, 1590-1599.	0.5	415
3	A Coarse-Grained Model for Polyethylene Oxide and Polyethylene Glycol: Conformation and Hydrodynamics. <i>Journal of Physical Chemistry B</i> , 2009, 113, 13186-13194.	2.6	338
4	Molecular Dynamics Simulations of PAMAM Dendrimer-Induced Pore Formation in DPPC Bilayers with a Coarse-Grained Model. <i>Journal of Physical Chemistry B</i> , 2006, 110, 18204-18211.	2.6	196
5	Coarse-Grained Molecular Dynamics Studies of the Concentration and Size Dependence of Fifth- and Seventh-Generation PAMAM Dendrimers on Pore Formation in DMPC Bilayer. <i>Journal of Physical Chemistry B</i> , 2008, 112, 7778-7784.	2.6	155
6	Lipid Bilayer Curvature and Pore Formation Induced by Charged Linear Polymers and Dendrimers: The Effect of Molecular Shape. <i>Journal of Physical Chemistry B</i> , 2008, 112, 12279-12285.	2.6	106
7	Coarse-Grained Model for PEGylated Lipids: Effect of PEGylation on the Size and Shape of Self-Assembled Structures. <i>Journal of Physical Chemistry B</i> , 2011, 115, 7830-7837.	2.6	104
8	Molecular Dynamics Studies of the Size, Shape, and Internal Structure of 0% and 90% Acetylated Fifth-Generation Polyamidoamine Dendrimers in Water and Methanol. <i>Journal of Physical Chemistry B</i> , 2006, 110, 4014-4019.	2.6	85
9	Multiscale Modeling of Dendrimers and Their Interactions with Bilayers and Polyelectrolytes. <i>Molecules</i> , 2009, 14, 423-438.	3.8	82
10	Effects of PEGylation on the Size and Internal Structure of Dendrimers: Self-Penetration of Long PEG Chains into the Dendrimer Core. <i>Macromolecules</i> , 2011, 44, 2291-2298.	4.8	82
11	In vitro blood cell viability profiling of polymers used in molecular assembly. <i>Scientific Reports</i> , 2017, 7, 9481.	3.3	76
12	Adsorption of Plasma Proteins onto PEGylated Lipid Bilayers: The Effect of PEG Size and Grafting Density. <i>Biomacromolecules</i> , 2016, 17, 1757-1765.	5.4	75
13	Molecular Dynamics Study of the Structure and Interparticle Interactions of Polyethylene Glycol-Conjugated PAMAM Dendrimers. <i>Journal of Physical Chemistry B</i> , 2009, 113, 13202-13207.	2.6	60
14	Cytoprotective Self-assembled RGD Peptide Nanofilms for Surface Modification of Viable Mesenchymal Stem Cells. <i>Chemistry of Materials</i> , 2017, 29, 2055-2065.	6.7	51
15	Structure and Dynamics of Helix-0 of the N-BAR Domain in Lipid Micelles and Bilayers. <i>Biophysical Journal</i> , 2008, 95, 4315-4323.	0.5	47
16	Self-Assembly of Lipids and Single-Walled Carbon Nanotubes: Effects of Lipid Structure and PEGylation. <i>Journal of Physical Chemistry C</i> , 2012, 116, 9327-9333.	3.1	41
17	Beneficial roles of H-donors as diluent and H-shuttle for asphaltenes in catalytic upgrading of vacuum residue. <i>Chemical Engineering Journal</i> , 2017, 314, 1-10.	12.7	41
18	Membrane Pore Formation Induced by Acetylated and Polyethylene Glycol-Conjugated Polyamidoamine Dendrimers. <i>Journal of Physical Chemistry C</i> , 2011, 115, 5316-5322.	3.1	39

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19	Molecular Simulations of PEGylated Biomolecules, Liposomes, and Nanoparticles for Drug Delivery Applications. <i>Pharmaceutics</i> , 2020, 12, 533.	4.5	38
20	Quantitative Interpretation of Hydration Dynamics Enabled the Fabrication of a Zwitterionic Antifouling Surface. <i>ACS Applied Materials & Interfaces</i> , 2020, 12, 7951-7965.	8.0	38
21	Effects of Nanoparticle Electrostatics and Protein-Protein Interactions on Corona Formation: Conformation and Hydrodynamics. <i>Small</i> , 2020, 16, e1906598.	10.0	37
22	Molecular Modeling of PEGylated Peptides, Dendrimers, and Single-Walled Carbon Nanotubes for Biomedical Applications. <i>Polymers</i> , 2014, 6, 776-798.	4.5	34
23	Molecular Dynamics Studies of PEGylated Single-Walled Carbon Nanotubes: The Effect of PEG Size and Grafting Density. <i>Journal of Physical Chemistry C</i> , 2013, 117, 26334-26341.	3.1	32
24	Effects of PEGylation on the Binding Interaction of Magainin 2 and Tachyplesin I with Lipid Bilayer Surface. <i>Langmuir</i> , 2013, 29, 14214-14221.	3.5	31
25	Molecular Dynamics Studies of the Size and Internal Structure of the PAMAM Dendrimer Grafted with Arginine and Histidine. <i>Macromolecules</i> , 2011, 44, 8681-8686.	4.8	30
26	Reverse Actuation of Polyelectrolyte Effect for <i>In Vivo</i> Antifouling. <i>ACS Nano</i> , 2021, 15, 6811-6828.	14.6	30
27	Synergistic effects of magainin 2 and PGLa on their heterodimer formation, aggregation, and insertion into the bilayer. <i>RSC Advances</i> , 2015, 5, 2047-2055.	3.6	27
28	Interparticle Dispersion, Membrane Curvature, and Penetration Induced by Single-Walled Carbon Nanotubes Wrapped with Lipids and PEGylated Lipids. <i>Journal of Physical Chemistry B</i> , 2013, 117, 1337-1344.	2.6	26
29	Investigation of Ion Channel Activities of Gramicidin A in the Presence of Ionic Liquids Using Model Cell Membranes. <i>Scientific Reports</i> , 2015, 5, 11935.	3.3	24
30	Molecular Dynamics Simulations of the Anchoring and Tilting of the Lung-Surfactant Peptide SP-B1-25 in Palmitic Acid Monolayers. <i>Biophysical Journal</i> , 2005, 89, 3807-3821.	0.5	21
31	Molecular Dynamics Studies of PEGylated α -Helical Coiled Coils and Their Self-Assembled Micelles. <i>Langmuir</i> , 2014, 30, 8848-8855.	3.5	20
32	Structures, dynamics, and hydrogen-bond interactions of antifreeze proteins in TIP4P/Ice water and their dependence on force fields. <i>PLoS ONE</i> , 2018, 13, e0198887.	2.5	20
33	Formulation Optimization and In Vivo Proof-of-Concept Study of Thermosensitive Liposomes Balanced by Phospholipid, Elastin-Like Polypeptide, and Cholesterol. <i>PLoS ONE</i> , 2014, 9, e103116.	2.5	20
34	Isolating a Trimer Intermediate in the Self-Assembly of E2 Protein Cage. <i>Biomacromolecules</i> , 2012, 13, 699-705.	5.4	19
35	Effect of Arginine-Rich Peptide Length on the Structure and Binding Strength of siRNA-Peptide Complexes. <i>Journal of Physical Chemistry B</i> , 2013, 117, 6917-6926.	2.6	18
36	The binding and insertion of imidazolium-based ionic surfactants into lipid bilayers: the effects of the surfactant size and salt concentration. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 5725-5733.	2.8	18

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37	Effects of imidazolium-based ionic liquids on the stability and dynamics of gramicidin A and lipid bilayers at different salt concentrations. <i>Journal of Molecular Graphics and Modelling</i> , 2015, 61, 53-60.	2.4	18
38	Effect of polyelectrolyte size on multilayer conformation and dynamics at different temperatures and salt concentrations. <i>Journal of Molecular Graphics and Modelling</i> , 2016, 70, 246-252.	2.4	17
39	All-atom simulations and free-energy calculations of coiled-coil peptides with lipid bilayers: binding strength, structural transition, and effect on lipid dynamics. <i>Scientific Reports</i> , 2016, 6, 22299.	3.3	17
40	Aggregation and insertion of melittin and its analogue MelP5 into lipid bilayers at different concentrations: effects on pore size, bilayer thickness and dynamics. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 7195-7203.	2.8	17
41	Transition-Metal Dichalcogenide Artificial Antibodies with Multivalent Polymeric Recognition Phases for Rapid Detection and Inactivation of Pathogens. <i>Journal of the American Chemical Society</i> , 2021, 143, 14635-14645.	13.7	17
42	Effects of the Size, Shape, and Structural Transition of Thermosensitive Polypeptides on the Stability of Lipid Bilayers and Liposomes. <i>Macromolecules</i> , 2012, 45, 7304-7312.	4.8	16
43	Effects of the asphaltene structure and the tetralin/heptane solvent ratio on the size and shape of asphaltene aggregates. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 13931-13940.	2.8	16
44	Dynamics and stability of lipid bilayers modulated by thermosensitive polypeptides, cholesterol, and PEGylated lipids. <i>Physical Chemistry Chemical Physics</i> , 2014, 16, 3763.	2.8	15
45	Effects of temperature, salt concentration, and the protonation state on the dynamics and hydrogen-bond interactions of polyelectrolyte multilayers on lipid membranes. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 6691-6700.	2.8	15
46	Molecular Modeling of Protein Corona Formation and Its Interactions with Nanoparticles and Cell Membranes for Nanomedicine Applications. <i>Pharmaceutics</i> , 2021, 13, 637.	4.5	15
47	Effect of lipid shape on toroidal pore formation and peptide orientation in lipid bilayers. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 21340-21349.	2.8	14
48	Effect of low levels of lipid oxidation on the curvature, dynamics, and permeability of lipid bilayers and their interactions with cationic nanoparticles. <i>Journal Physics D: Applied Physics</i> , 2018, 51, 164002.	2.8	13
49	Supramolecular Functionalization for Improving Thermoelectric Properties of Single-Walled Carbon Nanotubes—Small Organic Molecule Hybrids. <i>ACS Applied Materials & Interfaces</i> , 2020, 12, 51387-51396.	8.0	13
50	Membrane penetration and curvature induced by single-walled carbon nanotubes: the effect of diameter, length, and concentration. <i>Physical Chemistry Chemical Physics</i> , 2013, 15, 16334.	2.8	11
51	Effect of peptide conformation on TiO ₂ biomineralization. <i>Dalton Transactions</i> , 2013, 42, 13817.	3.3	11
52	Structural effects of tachyplesin I and its linear derivative on their aggregation and mobility in lipid bilayers. <i>Journal of Molecular Graphics and Modelling</i> , 2015, 59, 123-128.	2.4	11
53	Adsorption of plasma proteins onto PEGylated single-walled carbon nanotubes: The effects of protein shape, PEG size and grafting density. <i>Journal of Molecular Graphics and Modelling</i> , 2017, 75, 1-8.	2.4	11
54	Alpha-Hederin Nanopore for Single Nucleotide Discrimination. <i>ACS Nano</i> , 2019, 13, 1719-1727.	14.6	11

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55	Dispersion and bilayer interaction of single-walled carbon nanotubes modulated by covalent and noncovalent PEGylation. <i>Molecular Simulation</i> , 2015, 41, 1254-1263.	2.0	10
56	Design of a reversible inverted pH-responsive caged protein. <i>Biomaterials Science</i> , 2015, 3, 627-635.	5.4	9
57	Effects of hydrophobic and hydrogen-bond interactions on the binding affinity of antifreeze proteins to specific ice planes. <i>Journal of Molecular Graphics and Modelling</i> , 2019, 87, 48-55.	2.4	9
58	Multilayer Nanofilms via Inkjet Printing for Stabilizing Growth Factor and Designing Desired Cell Developments. <i>Advanced Healthcare Materials</i> , 2017, 6, 1700216.	7.6	8
59	Modulation of the Vault Protein-Protein Interaction for Tuning of Molecular Release. <i>Scientific Reports</i> , 2017, 7, 14816.	3.3	8
60	Effect of Protein Corona on Nanoparticle-Lipid Membrane Binding: The Binding Strength and Dynamics. <i>Langmuir</i> , 2021, 37, 3751-3760.	3.5	8
61	Multivalent Nanosheet Antibody Mimics for Selective Microbial Recognition and Inactivation. <i>Advanced Materials</i> , 2021, 33, e2101376.	21.0	8
62	Effects of imidazolium-based ionic surfactants on the size and dynamics of phosphatidylcholine bilayers with saturated and unsaturated chains. <i>Journal of Molecular Graphics and Modelling</i> , 2015, 60, 162-168.	2.4	7
63	Mechanistic Pathway of Lipid Phase-Dependent Lipid Corona Formation on Phenylalanine-Functionalized Gold Nanoparticles: A Combined Experimental and Molecular Dynamics Simulation Study. <i>Journal of Physical Chemistry B</i> , 2022, 126, 2241-2255.	2.6	7
64	Effect of the structural difference between Bax- $\hat{L}\pm 5$ and Bcl-xL- $\hat{L}\pm 5$ on their interactions with lipid bilayers. <i>Physical Chemistry Chemical Physics</i> , 2014, 16, 981-988.	2.8	6
65	Heterodimer and pore formation of magainin 2 and PGLa: The anchoring and tilting of peptides in lipid bilayers. <i>Biochimica Et Biophysica Acta - Biomembranes</i> , 2020, 1862, 183305.	2.6	6
66	Self-assembly of mixtures of a dendrimer and lipids: effects of hydrophobicity and electrostatics. <i>Molecular Simulation</i> , 2012, 38, 534-539.	2.0	5
67	Corona Formation: Effects of Nanoparticle Electrostatics and Protein-Protein Interactions on Corona Formation: Conformation and Hydrodynamics (Small 10/2020). <i>Small</i> , 2020, 16, 2070054.	10.0	5
68	Effects of salt on the size and internal structure of PAMAM dendrimers at different pH. <i>Molecular Simulation</i> , 2012, 38, 589-594.	2.0	4
69	Supramolecular Protein Assembly Retains Its Structural Integrity at Liquid-Liquid Interface. <i>Advanced Materials Interfaces</i> , 2020, 7, 1901674.	3.7	4
70	Self-assembled DNA hollow spheres from microsponges. <i>Biofabrication</i> , 2019, 11, 025016.	7.1	3
71	Topological analysis of single-stranded DNA with an alpha-hederin nanopore. <i>Biosensors and Bioelectronics</i> , 2021, 171, 112711.	10.1	3
72	All-Atom Simulations and Free Energy Calculations of Antibodies Bound to the Spike Protein of SARS-CoV-2: The Binding Strength and Multivalent Hydrogen-Bond Interactions. <i>Advanced Theory and Simulations</i> , 2021, 4, 2100012.	2.8	2

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73	Supramolecular Assemblies: Supramolecular Protein Assembly Retains Its Structural Integrity at Liquid–Liquid Interface (Adv. Mater. Interfaces 4/2020). Advanced Materials Interfaces, 2020, 7, 2070021.	3.7	1
74	A simple strategy for signal enhancement in lateral flow assays using superabsorbent polymers. Mikrochimica Acta, 2021, 188, 364.	5.0	1
75	Disassembly and trimer formation of E2 protein cage: the effects of C-terminus, salt, and protonation state. Journal Physics D: Applied Physics, 2018, 51, 365402.	2.8	0
76	Antibacterial Strategies: Multivalent Nanosheet Antibody Mimics for Selective Microbial Recognition and Inactivation (Adv. Mater. 22/2021). Advanced Materials, 2021, 33, 2170173.	21.0	0