

# Ard a louis

## List of Publications by Year in descending order

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137  
papers

9,250  
citations

41344

49  
h-index

45317

90  
g-index

142  
all docs

142  
docs citations

142  
times ranked

5075  
citing authors

#	ARTICLE	IF	CITATIONS
1	On the biophysics and kinetics of toehold-mediated DNA strand displacement. <i>Nucleic Acids Research</i> , 2013, 41, 10641-10658.	14.5	423
2	Structural, mechanical, and thermodynamic properties of a coarse-grained DNA model. <i>Journal of Chemical Physics</i> , 2011, 134, 085101.	3.0	379
3	Can Polymer Coils Be Modeled as "Soft Colloids"? <i>Physical Review Letters</i> , 2000, 85, 2522-2525.	7.8	374
4	Hydrodynamic interactions and Brownian forces in colloidal suspensions: Coarse-graining over time and length scales. <i>Physical Review E</i> , 2006, 74, 031402.	2.1	353
5	Accurate effective pair potentials for polymer solutions. <i>Journal of Chemical Physics</i> , 2001, 114, 4296-4311.	3.0	308
6	Introducing improved structural properties and salt dependence into a coarse-grained model of DNA. <i>Journal of Chemical Physics</i> , 2015, 142, 234901.	3.0	267
7	Sequence-dependent thermodynamics of a coarse-grained DNA model. <i>Journal of Chemical Physics</i> , 2012, 137, 135101.	3.0	265
8	Beware of density dependent pair potentials. <i>Journal of Physics Condensed Matter</i> , 2002, 14, 9187-9206.	1.8	259
9	Mean-field fluid behavior of the Gaussian core model. <i>Physical Review E</i> , 2000, 62, 7961-7972.	2.1	241
10	DNA hybridization kinetics: zippering, internal displacement and sequence dependence. <i>Nucleic Acids Research</i> , 2013, 41, 8886-8895.	14.5	203
11	Representability problems for coarse-grained water potentials. <i>Journal of Chemical Physics</i> , 2007, 126, 144509.	3.0	198
12	Influence of Polymer-Excluded Volume on the Phase-Behavior of Colloid-Polymer Mixtures. <i>Physical Review Letters</i> , 2002, 89, 128302.	7.8	182
13	Controlling crystallization and its absence: proteins, colloids and patchy models. <i>Physical Chemistry Chemical Physics</i> , 2007, 9, 2197.	2.8	179
14	Reversible self-assembly of patchy particles into monodisperse icosahedral clusters. <i>Journal of Chemical Physics</i> , 2007, 127, 085106.	3.0	176
15	Coarse-graining DNA for simulations of DNA nanotechnology. <i>Physical Chemistry Chemical Physics</i> , 2013, 15, 20395.	2.8	173
16	DNA Nanotweezers Studied with a Coarse-Grained Model of DNA. <i>Physical Review Letters</i> , 2010, 104, 178101.	7.8	162
17	Hydrodynamic and Brownian Fluctuations in Sedimenting Suspensions. <i>Physical Review Letters</i> , 2004, 93, 220601.	7.8	142
18	Polymer induced depletion potentials in polymer-colloid mixtures. <i>Journal of Chemical Physics</i> , 2002, 117, 1893-1907.	3.0	132

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19	Colloid-Polymer Mixtures in the Protein Limit. <i>Physical Review Letters</i> , 2003, 90, 068304.	7.8	127
20	Effective forces in colloidal mixtures: From depletion attraction to accumulation repulsion. <i>Physical Review E</i> , 2002, 65, 061407.	2.1	125
21	A nucleotide-level coarse-grained model of RNA. <i>Journal of Chemical Physics</i> , 2014, 140, 235102.	3.0	117
22	Phase separation in binary hard-core mixtures: An exact result. <i>Physical Review Letters</i> , 1992, 68, 3363-3365.	7.8	108
23	The stability of a crystal with diamond structure for patchy particles with tetrahedral symmetry. <i>Journal of Chemical Physics</i> , 2010, 132, 234511.	3.0	93
24	How To Derive and Parameterize Effective Potentials in Colloid~Polymer Mixtures. <i>Macromolecules</i> , 2002, 35, 1860-1869.	4.8	90
25	Optimizing DNA Nanotechnology through Coarse-Grained Modeling: A Two-Footed DNA Walker. <i>ACS Nano</i> , 2013, 7, 2479-2490.	14.6	88
26	Force-Induced Rupture of a DNA Duplex: From Fundamentals to Force Sensors. <i>ACS Nano</i> , 2015, 9, 11993-12003.	14.6	86
27	Direct Simulation of the Self-Assembly of a Small DNA Origami. <i>ACS Nano</i> , 2016, 10, 1724-1737.	14.6	86
28	Effective potentials for polymers and colloids: beyond the van der Waals picture of fluids?. <i>Philosophical Transactions Series A, Mathematical, Physical, and Engineering Sciences</i> , 2001, 359, 939-960.	3.4	85
29	Plectoneme tip bubbles: Coupled denaturation and writhing in supercoiled DNA. <i>Scientific Reports</i> , 2015, 5, 7655.	3.3	84
30	Theory of asymmetric nonadditive binary hard-sphere mixtures. <i>Physical Review E</i> , 2001, 64, 051202.	2.1	81
31	Phase diagram of model anisotropic particles with octahedral symmetry. <i>Journal of Chemical Physics</i> , 2007, 127, 054501.	3.0	80
32	The Arrival of the Frequent: How Bias in Genotype-Phenotype Maps Can Steer Populations to Local Optima. <i>PLoS ONE</i> , 2014, 9, e86635.	2.5	78
33	Inhibition of protein crystallization by evolutionary negative design. <i>Physical Biology</i> , 2004, 1, P9-P13.	1.8	77
34	Coarse-grained modelling of the structural properties of DNA origami. <i>Nucleic Acids Research</i> , 2019, 47, 1585-1597.	14.5	75
35	Many-body interactions and correlations in coarse-grained descriptions of polymer solutions. <i>Physical Review E</i> , 2001, 64, 021801.	2.1	73
36	Measuring colloidal interactions with confocal microscopy. <i>Journal of Chemical Physics</i> , 2007, 127, 044507.	3.0	73

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37	Knot-Controlled Ejection of a Polymer from a Virus Capsid. <i>Physical Review Letters</i> , 2009, 102, 088101.	7.8	72
38	Modelling the self-assembly of virus capsids. <i>Journal of Physics Condensed Matter</i> , 2010, 22, 104101.	1.8	71
39	Density profiles and surface tension of polymers near colloidal surfaces. <i>Journal of Chemical Physics</i> , 2002, 116, 10547-10556.	3.0	70
40	Self-Assembly and Evolution of Homomeric Protein Complexes. <i>Physical Review Letters</i> , 2009, 102, 118106.	7.8	68
41	Coarse-grained simulations of DNA overstretching. <i>Journal of Chemical Physics</i> , 2013, 138, 085101.	3.0	66
42	Design of hidden thermodynamic driving for non-equilibrium systems via mismatch elimination during DNA strand displacement. <i>Nature Communications</i> , 2020, 11, 2562.	12.8	66
43	A tractable genotype-phenotype map modelling the self-assembly of protein quaternary structure. <i>Journal of the Royal Society Interface</i> , 2014, 11, 20140249.	3.4	62
44	Monodisperse self-assembly in a model with protein-like interactions. <i>Journal of Chemical Physics</i> , 2009, 131, 175102.	3.0	61
45	Genetic Correlations Greatly Increase Mutational Robustness and Can Both Reduce and Enhance Evolvability. <i>PLoS Computational Biology</i> , 2016, 12, e1004773.	3.2	61
46	Exotic fluids and crystals of soft polymeric colloids. <i>Journal of Physics Condensed Matter</i> , 2002, 14, 7681-7698.	1.8	60
47	The structure of the genotype-phenotype map strongly constrains the evolution of non-coding RNA. <i>Interface Focus</i> , 2015, 5, 20150053.	3.0	58
48	Interplay between hydrodynamic and Brownian fluctuations in sedimenting colloidal suspensions. <i>Physical Review E</i> , 2008, 77, 011402.	2.1	57
49	Dynamic Colloidal Stabilization by Nanoparticle Halos. <i>Physical Review Letters</i> , 2004, 93, 248303.	7.8	55
50	Force-Induced Unravelling of DNA Origami. <i>ACS Nano</i> , 2018, 12, 6734-6747.	14.6	55
51	Influence of solvent quality on effective pair potentials between polymers in solution. <i>Physical Review E</i> , 2003, 67, 041801.	2.1	54
52	Self-assembly of monodisperse clusters: Dependence on target geometry. <i>Journal of Chemical Physics</i> , 2009, 131, 175101.	3.0	54
53	DNA hairpins destabilize duplexes primarily by promoting melting rather than by inhibiting hybridization. <i>Nucleic Acids Research</i> , 2015, 43, 6181-6190.	14.5	54
54	Modelling Toehold-Mediated RNA Strand Displacement. <i>Biophysical Journal</i> , 2015, 108, 1238-1247.	0.5	54

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55	Long-range correlations in the mechanics of small DNA circles under topological stress revealed by multi-scale simulation. <i>Nucleic Acids Research</i> , 2016, 44, gkw815.	14.5	54
56	TacoxDNA: A user-friendly web server for simulations of complex DNA structures, from single strands to origami. <i>Journal of Computational Chemistry</i> , 2019, 40, 2586-2595.	3.3	54
57	Formation of dodecagonal quasicrystals in two-dimensional systems of patchy particles. <i>Journal of Chemical Physics</i> , 2012, 136, 054904.	3.0	52
58	Correlation of automorphism group size and topological properties with program-size complexity evaluations of graphs and complex networks. <i>Physica A: Statistical Mechanics and Its Applications</i> , 2014, 404, 341-358.	2.6	51
59	Characterizing the Motion of Jointed DNA Nanostructures Using a Coarse-Grained Model. <i>ACS Nano</i> , 2017, 11, 12426-12435.	14.6	51
60	Symmetry and simplicity spontaneously emerge from the algorithmic nature of evolution. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2022, 119, e2113883119.	7.1	50
61	Nonmonotonic variation with salt concentration of the second virial coefficient in protein solutions. <i>Physical Review E</i> , 2003, 67, 051404.	2.1	49
62	From genotypes to organisms: State-of-the-art and perspectives of a cornerstone in evolutionary dynamics. <i>Physics of Life Reviews</i> , 2021, 38, 55-106.	2.8	49
63	Crystallization and phase separation in nonadditive binary hard-sphere mixtures. <i>Physical Review E</i> , 2000, 61, R1028-R1031.	2.1	48
64	Extracting bulk properties of self-assembling systems from small simulations. <i>Journal of Physics Condensed Matter</i> , 2010, 22, 104102.	1.8	47
65	Stick boundary conditions and rotational velocity auto-correlation functions for colloidal particles in a coarse-grained representation of the solvent. <i>Journal of Physics Condensed Matter</i> , 2005, 17, S3393-S3399.	1.8	46
66	Self-assembly, modularity, and physical complexity. <i>Physical Review E</i> , 2010, 82, 026117.	2.1	46
67	Relating monomer to centre-of-mass distribution functions in polymer solutions. <i>Europhysics Letters</i> , 2002, 58, 53-59.	2.0	45
68	Metallization of fluid hydrogen. <i>Philosophical Transactions Series A, Mathematical, Physical, and Engineering Sciences</i> , 1998, 356, 119-138.	3.4	44
69	Templated self-assembly of patchy particles. <i>Soft Matter</i> , 2011, 7, 3423.	2.7	44
70	Input-output maps are strongly biased towards simple outputs. <i>Nature Communications</i> , 2018, 9, 761.	12.8	43
71	Discrete charge patterns, Coulomb correlations and interactions in protein solutions. <i>Europhysics Letters</i> , 2002, 57, 731-737.	2.0	42
72	Evolutionary dynamics in a simple model of self-assembly. <i>Physical Review E</i> , 2011, 83, 066105.	2.1	42

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73	Effects of Interparticle Attractions on Colloidal Sedimentation. <i>Physical Review Letters</i> , 2010, 104, 068301.	7.8	40
74	Atomic Tunneling from a Scanning-Tunneling or Atomic-Force Microscope Tip: Dissipative Quantum Effects from Phonons. <i>Physical Review Letters</i> , 1995, 74, 1363-1366.	7.8	39
75	The Asakura-Oosawa model in the protein limit: the role of many-body interactions. <i>Journal of Physics Condensed Matter</i> , 2003, 15, S3429-S3442.	1.8	38
76	Density profiles and solvation forces for a Yukawa fluid in a slit pore. <i>Journal of Chemical Physics</i> , 2008, 128, 204704.	3.0	38
77	Invited article: Thermodynamic perturbation theory of the phase behaviour of colloid/interacting polymer mixtures. <i>Molecular Physics</i> , 2004, 102, 1-11.	1.7	37
78	The self-assembly of DNA Holliday junctions studied with a minimal model. <i>Journal of Chemical Physics</i> , 2009, 130, 065101.	3.0	36
79	Epistasis can lead to fragmented neutral spaces and contingency in evolution. <i>Proceedings of the Royal Society B: Biological Sciences</i> , 2012, 279, 1777-1783.	2.6	36
80	Effect of Bending Rigidity on the Knotting of a Polymer under Tension. <i>ACS Macro Letters</i> , 2012, 1, 1352-1356.	4.8	36
81	DNA Cruciform Arms Nucleate through a Correlated but Asynchronous Cooperative Mechanism. <i>Journal of Physical Chemistry B</i> , 2012, 116, 11616-11625.	2.6	36
82	The structure of colloid-polymer mixtures. <i>Europhysics Letters</i> , 1999, 46, 741-747.	2.0	35
83	Characterizing DNA Star-Tile-Based Nanostructures Using a Coarse-Grained Model. <i>ACS Nano</i> , 2016, 10, 4236-4247.	14.6	35
84	Probing ion-ion and electron-ion correlations in liquid metals within the quantum hypernetted chain approximation. <i>Physical Review B</i> , 2000, 61, 11400-11410.	3.2	34
85	Coarse-graining diblock copolymer solutions: a macromolecular version of the Widom-Rowlinson model. <i>Molecular Physics</i> , 2005, 103, 3045-3054.	1.7	34
86	Effect of topology on dynamics of knots in polymers under tension. <i>Europhysics Letters</i> , 2010, 89, 20001.	2.0	34
87	Re-entrant phase behavior for systems with competition between phase separation and self-assembly. <i>Journal of Chemical Physics</i> , 2011, 134, 104905.	3.0	34
88	Generalized depletion potentials. <i>Journal of Physics Condensed Matter</i> , 2001, 13, L777-L784.	1.8	32
89	The effect of scale-free topology on the robustness and evolvability of genetic regulatory networks. <i>Journal of Theoretical Biology</i> , 2010, 267, 48-61.	1.7	32
90	Phenotype Bias Determines How Natural RNA Structures Occupy the Morphospace of All Possible Shapes. <i>Molecular Biology and Evolution</i> , 2022, 39, .	8.9	32

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91	Coarse-graining polymers as soft colloids. <i>Physica A: Statistical Mechanics and Its Applications</i> , 2002, 306, 251-261.	2.6	31
92	Simulating a burnt-bridges DNA motor with a coarse-grained DNA model. <i>Natural Computing</i> , 2014, 13, 535-547.	3.0	30
93	The crossover from single file to Fickian diffusion. <i>Faraday Discussions</i> , 2010, 144, 285-299.	3.2	28
94	The effect of topology on the structure and free energy landscape of DNA kissing complexes. <i>Journal of Chemical Physics</i> , 2012, 136, 215102.	3.0	28
95	The Role of Loop Stacking in the Dynamics of DNA Hairpin Formation. <i>Journal of Physical Chemistry B</i> , 2014, 118, 14326-14335.	2.6	27
96	Contingency, convergence and hyper-astronomical numbers in biological evolution. <i>Studies in History and Philosophy of Science Part C: Studies in History and Philosophy of Biological and Biomedical Sciences</i> , 2016, 58, 107-116.	1.3	25
97	Extending Linear Response: Inferences from Electron-Ion Structure Factors. <i>Physical Review Letters</i> , 1998, 81, 4456-4459.	7.8	24
98	Viscous fingering at ultralow interfacial tension. <i>Soft Matter</i> , 2013, 9, 10599.	2.7	24
99	Zigzag transitions and nonequilibrium pattern formation in colloidal chains. <i>Journal of Chemical Physics</i> , 2013, 139, 134908.	3.0	22
100	Dynamics of solutes with hydrodynamic interactions: Comparison between Brownian dynamics and stochastic rotation dynamics simulations. <i>Physical Review E</i> , 2013, 88, 043304.	2.1	22
101	Identifying Physical Causes of Apparent Enhanced Cyclization of Short DNA Molecules with a Coarse-Grained Model. <i>Journal of Chemical Theory and Computation</i> , 2019, 15, 4660-4672.	5.3	22
102	Phase Separation of Penetrable Core Mixtures. <i>Journal of Statistical Physics</i> , 2003, 110, 1015-1037.	1.2	21
103	Pattern formation in colloidal explosions. <i>Europhysics Letters</i> , 2011, 94, 48008.	2.0	21
104	Confinement of knotted polymers in a slit. <i>Molecular Physics</i> , 2011, 109, 1289-1295.	1.7	20
105	Phase separation of a multiple occupancy lattice gas. <i>Journal of Physics A</i> , 2004, 37, 577-590.	1.6	18
106	Multi-scale coarse-graining for the study of assembly pathways in DNA-brick self-assembly. <i>Journal of Chemical Physics</i> , 2018, 148, 134910.	3.0	18
107	Effect of polymer-polymer interactions on the surface tension of colloid-polymer mixtures. <i>Journal of Chemical Physics</i> , 2003, 119, 12667-12672.	3.0	17
108	Influence of solvent quality on polymer solutions: A Monte Carlo study of bulk and interfacial properties. <i>Journal of Chemical Physics</i> , 2004, 121, 612.	3.0	17

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109	Polymer solutions: from hard monomers to soft polymers. <i>Journal of Physics Condensed Matter</i> , 2005, 17, S3185-S3193.	1.8	17
110	Characterizing the bending and flexibility induced by bulges in DNA duplexes. <i>Journal of Chemical Physics</i> , 2015, 142, 165101.	3.0	16
111	Density-Functional Study of Interfacial Properties of Colloid~Polymer Mixtures. <i>Journal of Physical Chemistry B</i> , 2005, 109, 6640-6649.	2.6	15
112	Hydrodynamics of confined colloidal fluids in two dimensions. <i>Physical Review E</i> , 2009, 79, 051402.	2.1	14
113	Coarse-grained modelling of supercoiled RNA. <i>Journal of Chemical Physics</i> , 2015, 143, 243122.	3.0	13
114	Generic predictions of output probability based on complexities of inputs and outputs. <i>Scientific Reports</i> , 2020, 10, 4415.	3.3	12
115	Complex dynamics of knotted filaments in shear flow. <i>Europhysics Letters</i> , 2010, 92, 34003.	2.0	11
116	From Concentration Profiles to Polymer Osmotic Equations of State. <i>ChemPhysChem</i> , 2005, 6, 1760-1764.	2.1	10
117	Measuring Internal Forces in Single-Stranded DNA: Application to a DNA Force Clamp. <i>Journal of Chemical Theory and Computation</i> , 2020, 16, 7764-7775.	5.3	10
118	How Péclet number affects microstructure and transient cluster aggregation in sedimenting colloidal suspensions. <i>Journal of Chemical Physics</i> , 2012, 136, 064517.	3.0	9
119	Combining quantum and classical density functional theory for ion~electron mixtures. <i>Journal of Non-Crystalline Solids</i> , 2002, 312-314, 60-68.	3.1	8
120	Boolean Threshold Networks as Models of Genotype-Phenotype Maps. <i>Springer Proceedings in Complexity</i> , 2020, , 143-155.	0.3	8
121	Phase Separation of a Model Binary Polymer Solution in an External Field. <i>Journal of Physical Chemistry B</i> , 2006, 110, 3661-3665.	2.6	6
122	Taylor dispersion of colloidal particles in narrow channels. <i>Molecular Physics</i> , 2015, 113, 2538-2545.	1.7	6
123	The interplay of supercoiling and thymine dimers in DNA. <i>Nucleic Acids Research</i> , 2022, 50, 2480-2492.	14.5	6
124	Electron~ion structure factors and the general accuracy of linear response. <i>Journal of Non-Crystalline Solids</i> , 1999, 250-252, 9-14.	3.1	5
125	Extracting short-ranged interactions from structure factors. <i>Molecular Physics</i> , 2011, 109, 2945-2951.	1.7	5
126	Fluid-Solid Phase Separation in Hard-Sphere Mixtures is Unrelated to Bond Percolation. <i>Physical Review Letters</i> , 2000, 84, 1840-1840.	7.8	4



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127	Preliminary Evaluation of a Kinetic Parameter Estimator with Application to Direct Parametric Reconstruction. , 0, , .		3
128	Reply to Ocklenburg and Mundorf: The interplay of developmental bias and natural selection. Proceedings of the National Academy of Sciences of the United States of America, 2022, 119, .	7.1	3
129	Coarse-Grained Modelling of Extreme DNA Bending. Biophysical Journal, 2014, 106, 66a.	0.5	2
130	Coarse-Grained Modeling of RNA for Biology and Nanotechnology. Biophysical Journal, 2017, 112, 369a.	0.5	1
131	Reconfigurable Tâ€junction DNA Origami. Angewandte Chemie - International Edition, 2020, 59, 15942-15946.	13.8	1
132	Is Water an Amniotic Eden or a Corrosive Hell? Emerging Perspectives on the Strangest Fluid in the Universe. , 2010, , 3-9.		0
133	Precision control of DNA-based molecular reactions. , 2016, , .		0
134	A touch of awe: crafting meaning from the wonder of the cosmos. Annals of the New York Academy of Sciences, 2018, 1432, 46-62.	3.8	0
135	Reconfigurable Tâ€junction DNA Origami. Angewandte Chemie, 2020, 132, 16076-16080.	2.0	0
136	The long and winding road to understanding organismal construction. Physics of Life Reviews, 2022, 42, 19-24.	2.8	0
137	Free energy landscapes of DNA and its assemblies: perspectives from coarse-grained modelling. Frontiers of Nanoscience, 2022, , 195-210.	0.6	0