

Gerardo Odriozola

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/2438956/publications.pdf>

Version: 2024-02-01

80
papers

1,308
citations

394286

19
h-index

454834

30
g-index

82
all docs

82
docs citations

82
times ranked

1084
citing authors

#	ARTICLE	IF	CITATIONS
1	Extended law of corresponding states: square-well oblates. <i>Journal of Physics Condensed Matter</i> , 2022, 34, 104002.	0.7	2
2	The role of the second virial coefficient in the vapor-liquid phase coexistence of anisotropic square-well particles. <i>Journal of Molecular Liquids</i> , 2022, 360, 119528.	2.3	1
3	Enhanced two-dimensional nematic order in slit-like pores. <i>New Journal of Physics</i> , 2021, 23, 063053.	1.2	7
4	Anisotropy-independent packing of confined hard ellipses. <i>Journal of Molecular Liquids</i> , 2021, 333, 115896.	2.3	8
5	Scaling Laws in the Diffusive Release of Neutral Cargo from Hollow Hydrogel Nanoparticles: Paclitaxel-Loaded Poly(4-vinylpyridine). <i>ACS Nano</i> , 2020, 14, 15227-15240.	7.3	15
6	A coil-to-globule transition capable coarse-grained model for poly(<i>N</i> -isopropylacrylamide). <i>Physical Chemistry Chemical Physics</i> , 2020, 22, 17913-17921.	1.3	7
7	Three-step melting of hard superdisks in two dimensions. <i>Physical Review E</i> , 2020, 102, 062603.	0.8	9
8	Ordering, clustering, and wetting of hard rods in extreme confinement. <i>Physical Review Research</i> , 2020, 2, .	1.3	13
9	Effect of Temperature on the Cononsolvency of Poly(<i>N</i> -isopropylacrylamide) (PNIPAM) in Aqueous 1-Propanol. <i>ACS Applied Polymer Materials</i> , 2019, 1, 2961-2972.	2.0	19
10	P-NIPAM in water- <i>acetone</i> mixtures: experiments and simulations. <i>Physical Chemistry Chemical Physics</i> , 2019, 21, 5106-5116.	1.3	17
11	Molecular dynamics simulations of brine-surfactant lamellas surrounded by nitrogen at different reservoir conditions. <i>Journal of Molecular Liquids</i> , 2018, 256, 480-488.	2.3	2
12	Ethane clathrates using different water- <i>ethane</i> models: Molecular dynamics. <i>Physica A: Statistical Mechanics and Its Applications</i> , 2018, 491, 89-100.	1.2	1
13	Massive replica exchange Monte Carlo algorithm: a tool to access high pressure thermodynamics of hard systems. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 27490-27500.	1.3	6
14	Phase diagram of hard squares in slit confinement. <i>Scientific Reports</i> , 2018, 8, 8886.	1.6	13
15	Competition between excluded-volume and electrostatic interactions for nanogel swelling: effects of the counterion valence and nanogel charge. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 6838-6848.	1.3	31
16	Colloid-polymer mixtures under slit confinement. <i>Journal of Chemical Physics</i> , 2017, 146, 104903.	1.2	6
17	Equivalence between particles and fields: A general statistical mechanics theory for short and long range many-body forces. <i>Fortschritte Der Physik</i> , 2017, 65, 1600072.	1.5	5
18	Conformation change of an isotactic poly (<i>N</i> -isopropylacrylamide) membrane: Molecular dynamics. <i>Journal of Chemical Physics</i> , 2017, 146, 194905.	1.2	22

#	ARTICLE	IF	CITATIONS
19	Critical behavior of hard squares in strong confinement. <i>Physical Review E</i> , 2017, 95, 042610.	0.8	6
20	Phase behaviour of short range triangle well fluids: A comparison with lysozyme suspensions. <i>Journal of Molecular Liquids</i> , 2017, 225, 723-729.	2.3	4
21	Anomalous structural transition of confined hard squares. <i>Physical Review E</i> , 2016, 94, 050603.	0.8	9
22	Thermodynamic properties of triangle-well fluids in two dimensions: MC and MD simulations. <i>Journal of Chemical Physics</i> , 2016, 145, 174505.	1.2	8
23	Entropy Driven Self-Assembly in Charged Locking Key Particles. <i>Journal of Physical Chemistry B</i> , 2016, 120, 5966-5974.	1.2	8
24	Parallel Replica Exchange Monte Carlo Applied to Hard Systems. <i>Communications in Computer and Information Science</i> , 2016, , 392-418.	0.4	0
25	Effect of orientational restriction on monolayers of hard ellipsoids. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 4547-4556.	1.3	10
26	Using a Parallel Genetic Algorithm to Fit a Pulsed Townsend Discharge Simulation to Experiments. <i>Communications in Computer and Information Science</i> , 2016, , 343-355.	0.4	2
27	Corresponding states law for a generalized Lennard-Jones potential. <i>Journal of Chemical Physics</i> , 2015, 143, 024504.	1.2	17
28	A heuristic rule for classification of classical fluids: Master curves for Mie, Yukawa and square-well potentials. <i>Chemical Physics Letters</i> , 2015, 631-632, 26-29.	1.2	18
29	Wall particle interactions and depletion forces in narrow slits. <i>Current Opinion in Colloid and Interface Science</i> , 2015, 20, 24-31.	3.4	5
30	Coexistence and interfacial properties of a triangle-well mimicking the Lennard-Jones fluid and a comparison with noble gases. <i>Journal of Chemical Physics</i> , 2015, 142, 074706.	1.2	6
31	Phase diagram of two-dimensional hard ellipses. <i>Journal of Chemical Physics</i> , 2014, 140, 204502.	1.2	45
32	Empty liquid phase of colloidal ellipsoids: The role of shape and interaction anisotropy. <i>Journal of Chemical Physics</i> , 2014, 140, 134905.	1.2	10
33	Coexistence and interfacial properties of triangle-well fluids. <i>Molecular Physics</i> , 2014, 112, 2114-2121.	0.8	13
34	Simple effective rule to estimate the jamming packing fraction of polydisperse hard spheres. <i>Physical Review E</i> , 2014, 89, 040302.	0.8	34
35	Towards understanding the empty liquid of colloidal platelets: vapour liquid phase coexistence of square-well oblate ellipsoids. <i>Soft Matter</i> , 2013, 9, 5277.	1.2	18
36	Empty liquid state and self-assembly of high valence non-spherical colloidal systems. <i>Soft Matter</i> , 2013, 9, 11178.	1.2	10

#	ARTICLE	IF	CITATIONS
37	Structure and coexistence properties of shoulder-“square well fluids. <i>Journal of Molecular Liquids</i> , 2013, 185, 70-75.	2.3	2
38	Statistical Mechanics Approach to Lock-Key Supramolecular Chemistry Interactions. <i>Physical Review Letters</i> , 2013, 110, 105701.	2.9	30
39	Constant-force approach to discontinuous potentials. <i>Journal of Chemical Physics</i> , 2013, 138, 214105.	1.2	18
40	Further details on the phase diagram of hard ellipsoids of revolution. <i>Journal of Chemical Physics</i> , 2013, 138, 064501.	1.2	41
41	Expansion of Natural Na ⁺ and Ca ²⁺ Montmorillonites in the Presence of NaCl and Surfactant Solutions. <i>Energy & Fuels</i> , 2012, 26, 2578-2584.	2.5	6
42	Revisiting the phase diagram of hard ellipsoids. <i>Journal of Chemical Physics</i> , 2012, 136, 134505.	1.2	60
43	Ion-specific colloidal aggregation: Population balance equations and potential of mean force. <i>Journal of Chemical Physics</i> , 2011, 135, 134704.	1.2	4
44	Entropy effects in self-assembling mechanisms: Also a view from the information theory. <i>Journal of Molecular Liquids</i> , 2011, 164, 87-100.	2.3	11
45	Assisted crystal growing by tempering metastable vapor-“liquid fluids. <i>Chemical Physics Letters</i> , 2011, 501, 466-469.	1.2	5
46	Equilibrium equation of state of a hard sphere binary mixture at very large densities using replica exchange Monte Carlo simulations. <i>Journal of Chemical Physics</i> , 2011, 134, 054504.	1.2	30
47	Vapor-“liquid surface tension of strong short-range Yukawa fluid. <i>Journal of Chemical Physics</i> , 2011, 134, 154702.	1.2	12
48	Communication: Equation of state of hard oblate ellipsoids by replica exchange Monte Carlo. <i>Journal of Chemical Physics</i> , 2011, 134, 201103.	1.2	8
49	Hard ellipsoids: Analytically approaching the exact overlap distance. <i>Journal of Chemical Physics</i> , 2011, 135, 084508.	1.2	11
50	Phase behaviour and separation kinetics of symmetric non-additive hard discs. <i>Molecular Simulation</i> , 2010, 36, 175-185.	0.9	6
51	Ion-induced reversibility in the aggregation of hydrophobic colloids. <i>Soft Matter</i> , 2010, 6, 1114.	1.2	12
52	Stability mechanisms for plate-like nanoparticles immersed in a macroion dispersion. <i>Journal of Physics Condensed Matter</i> , 2009, 21, 424107.	0.7	10
53	Replica exchange Monte Carlo applied to hard spheres. <i>Journal of Chemical Physics</i> , 2009, 131, 144107.	1.2	24
54	Population Inversion of a NAHS Mixture Adsorbed into a Cylindrical Pore. <i>Journal of Physical Chemistry C</i> , 2008, 112, 18028-18033.	1.5	16

#	ARTICLE	IF	CITATIONS
55	Entropy driven key-lock assembly. <i>Journal of Chemical Physics</i> , 2008, 129, 111101.	1.2	43
56	Linking Phase Behavior and Reversible Colloidal Aggregation at Low Concentrations: Simulations and Stochastic Mean Field Theory. <i>Journal of Physical Chemistry B</i> , 2007, 111, 5564-5572.	1.2	17
57	Aggregation kinetics of latex microspheres in alcohol-water media. <i>Journal of Colloid and Interface Science</i> , 2007, 310, 471-480.	5.0	12
58	Electrolyte distribution around two like-charged rods: Their effective attractive interaction and angular dependent charge reversal. <i>Journal of Chemical Physics</i> , 2006, 124, 134902.	1.2	31
59	Two rods confined by positive plates: effective forces and charge distribution profiles. <i>Journal of Physics Condensed Matter</i> , 2006, 18, S2335-S2352.	0.7	6
60	Effect of Confinement on the Interaction between Two Like-Charged Rods. <i>Physical Review Letters</i> , 2006, 97, 018102.	2.9	20
61	Stability of Ca-montmorillonite hydrates: A computer simulation study. <i>Journal of Chemical Physics</i> , 2005, 123, 174708.	1.2	15
62	Stability of K-Montmorillonite Hydrates: Hybrid MC Simulations. <i>Journal of Chemical Theory and Computation</i> , 2005, 1, 1211-1220.	2.3	15
63	Brownian dynamics simulations of Laponite colloid suspensions. <i>Physical Review E</i> , 2004, 70, 021405.	0.8	42
64	Irreversible versus reversible aggregation: Mean field theory and experiments. <i>Journal of Chemical Physics</i> , 2004, 121, 5468-5481.	1.2	17
65	Na-montmorillonite hydrates under ethane rich reservoirs: NPzT and $\hat{1}/4$ PzT simulations. <i>Journal of Chemical Physics</i> , 2004, 121, 4266-4275.	1.2	17
66	Colloidal aggregation with sedimentation: concentration effects. <i>European Physical Journal E</i> , 2004, 13, 165-178.	0.7	19
67	Coupled aggregation and sedimentation processes: stochastic mean field theory. <i>Physica A: Statistical Mechanics and Its Applications</i> , 2004, 335, 35-46.	1.2	7
68	Na-Montmorillonite Hydrates under Basin Conditions: Hybrid Monte Carlo and Molecular Dynamics Simulations. <i>Langmuir</i> , 2004, 20, 2010-2016.	1.6	29
69	Simulated Reversible Aggregation Processes for Different Interparticle Potentials: The Cluster Aging Phenomenon. <i>Journal of Physical Chemistry B</i> , 2003, 107, 2180-2188.	1.2	17
70	Modeling the aggregation of partially covered particles: Theory and simulation. <i>Physical Review E</i> , 2003, 68, 011404.	0.8	15
71	Coupled aggregation and sedimentation processes: The sticking probability effect. <i>Physical Review E</i> , 2003, 67, 031401.	0.8	14
72	Constant bond breakup probability model for reversible aggregation processes. <i>Physical Review E</i> , 2002, 65, 031405.	0.8	40

#	ARTICLE	IF	CITATIONS
73	Coupled aggregation and sedimentation processes: Three-dimensional off-lattice simulations. European Physical Journal E, 2002, 7, 153-161.	0.7	1
74	Title is missing!. European Physical Journal E, 2002, 7, 153-161.	0.7	12
75	The DLCA-RLCA transition arising in 2D-aggregation: simulations and mean field theory. European Physical Journal E, 2001, 5, 471-480.	0.7	25
76	A Light Scattering Study of the Transition Region between Diffusion- and Reaction-Limited Cluster Aggregation. Journal of Colloid and Interface Science, 2001, 240, 90-96.	5.0	49
77	A probabilistic aggregation kernel for the computer-simulated transition from DLCA to RLCA. Europhysics Letters, 2001, 53, 797-803.	0.7	58
78	The kinetics of irreversible aggregation processes. , 2001, , 87-90.		1
79	Multiple contact kernel for diffusionlike aggregation. Physical Review E, 2000, 62, 8335-8343.	0.8	39
80	Dynamic scaling concepts applied to numerical solutions of Smoluchowski's rate equation. Journal of Chemical Physics, 1999, 111, 7657-7667.	1.2	31