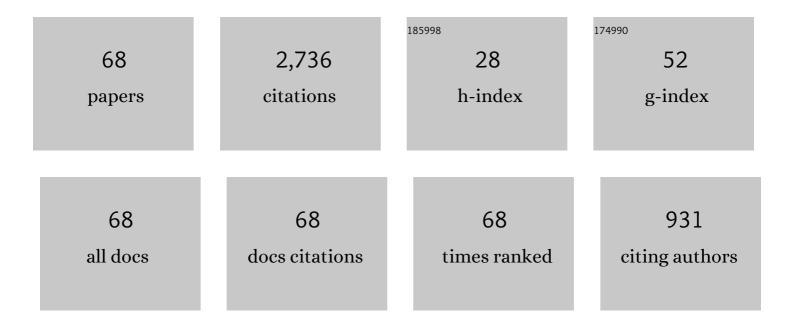
List of Publications by Year in descending order

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FREDIADO

#	Article	IF	CITATIONS
1	From square-well to Janus: Improved algorithm for integral equation theory and comparison with thermodynamic perturbation theory within the Kern-Frenkel model. Journal of Chemical Physics, 2014, 140, 094104.	1.2	19
2	Effects of patch size and number within a simple model of patchy colloids. Journal of Chemical Physics, 2010, 132, 174110.	1.2	107
3	Liquid–vapor coexistence in square-well fluids: an RHNC study. Molecular Physics, 2009, 107, 555-562.	0.8	17
4	Phase diagram and structural properties of a simple model for one-patch particles. Journal of Chemical Physics, 2009, 131, 174114.	1.2	42
5	An efficient procedure for the study of inhomogeneous liquids. Molecular Physics, 2009, 107, 301-308.	0.8	5
6	Integral equation and simulation studies of a planar nematogenic liquid in crossed external fields. Journal of Physics Condensed Matter, 2005, 17, 2801-2824.	0.7	2
7	Percolation of clusters with a residence time in the bond definition: Integral equation theory. Physical Review E, 2005, 71, 031202.	0.8	7
8	Distribution functions for a two-dimensional non-interacting quantum electron gas in an external magnetic field. Molecular Physics, 2003, 101, 1635-1639.	0.8	0
9	Structure and thermodynamics of a ferrofluid monolayer. Physical Review E, 2000, 61, 3838-3849.	0.8	47
10	Thermodynamic perturbation theory for polydisperse colloidal suspensions using orthogonal polynomial expansions. Physical Review E, 1999, 59, 6937-6945.	0.8	21
11	An integral equation study of a simple point charge model of water. Journal of Chemical Physics, 1999, 110, 1148-1153.	1.2	64
12	Static structure of polydisperse colloidal monolayers. Journal of Chemical Physics, 1998, 108, 6441-6446.	1.2	16
13	Heisenberg Spin Fluid in an External Magnetic Field. Physical Review Letters, 1998, 80, 3535-3538.	2.9	28
14	An accurate theoretical description of fluids composed of fully anisotropic molecules: Application to C2v symmetry. Journal of Chemical Physics, 1997, 107, 4642-4647.	1.2	22
15	Molecular theory of a charged particle in a polarizable nonpolar liquid. Journal of Chemical Physics, 1997, 106, 4707-4713.	1.2	15
16	Monte Carlo simulation and reference hypernetted chain equation results for structural, thermodynamic, and dielectric properties of polar heteronuclear diatomic fluids. Journal of Chemical Physics, 1996, 104, 6710-6718.	1.2	10
17	Integral equation theory of polydisperse colloidal suspensions using orthogonal polynomial expansions. Physical Review E, 1996, 54, 4411-4419.	0.8	34
18	Integral equation algorithm for fluids of fully anisotropic molecules. Journal of Chemical Physics, 1995, 103, 481-484.	1.2	37

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19	A fast method of solving the hypernetted-chain equation for molecular Lennard-Jones fluids. Molecular Physics, 1995, 84, 743-755.	0.8	9
20	Numerical solution of structure integral equation theories for two-dimensional fluid mixtures. Molecular Physics, 1994, 83, 351-359.	0.8	6
21	Integral equation and simulation studies of a realistic model for liquid hydrogen chloride. Journal of Chemical Physics, 1994, 100, 1599-1605.	1.2	21
22	Trapping constant, thermal conductivity, and the microstructure of suspensions of oriented spheroids. Journal of Chemical Physics, 1991, 94, 4453-4462.	1.2	39
23	Integrals over the triplet distribution function without the triplet distribution function. Molecular Physics, 1991, 72, 1387-1395.	0.8	6
24	Integral-equation approach to the calculation of the potential distribution in a fluid. Physical Review A, 1990, 42, 7281-7288.	1.0	4
25	Twoâ€point probability function for distributions of oriented hard ellipsoids. Journal of Chemical Physics, 1990, 93, 5912-5917.	1.2	29
26	Percus-Yevick solutions for the planar dumbell fluid. Molecular Physics, 1988, 64, 1185-1193.	0.8	11
27	Referenceâ€hypernetted chain equation with anisotropic bridge function for fluids of diatomic molecules. Journal of Chemical Physics, 1988, 88, 1950-1952.	1.2	9
28	Structure, thermodynamics, and orientational correlations of the nematogenic hard ellipse fluid from the Percus-Yevick equation. Molecular Physics, 1988, 63, 623-638.	0.8	33
29	Bulk properties of composite media. I. Simplification of bounds on the shear modulus of suspensions of impenetrable spheres. Journal of Applied Physics, 1987, 62, 3503-3513.	1.1	8
30	Extension of the optimized RHNC equation to multicomponent liquids. Journal of Chemical Physics, 1987, 87, 2249-2256.	1.2	35
31	Charge distribution in plasmas with field constraint. Physical Review A, 1987, 36, 2333-2337.	1.0	15
32	Bulk properties of composite media. II. Evaluation of bounds on the shear modulus of suspensions of impenetrable spheres. Journal of Applied Physics, 1987, 62, 4135-4141.	1.1	16
33	Generalized bridge functions for the reference hypernetted-chain equation: Calculation of the electric microfield distribution in a plasma. Physical Review A, 1987, 36, 313-317.	1.0	10
34	Bulk properties of twoâ€phase disordered media. IV. Mechanical properties of suspensions of penetrable spheres at nondilute concentrations. Journal of Chemical Physics, 1987, 86, 6388-6392.	1.2	12
35	Dumbell—A program to calculate the structure and thermodynamics of a classical fluid of hard, homonuclear diatomic molecules. Computer Physics Communications, 1986, 39, 133-140.	3.0	9
36	Exact solution of the mean spherical model for the electric microfield distribution in a plasma. Physical Review A, 1986, 34, 4131-4135.	1.0	15

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37	Structure and thermodynamics of the dipolar hard sphere fluid from the referenceâ€hypernetted chain equation with minimized free energy. Journal of Chemical Physics, 1986, 85, 2916-2921.	1.2	33
38	Effective properties of two-phase disordered composite Media. I. Simplification of bounds on the conductivity and bulk modulus of dispersions of impenetrable spheres. Physical Review B, 1986, 33, 3370-3378.	1.1	27
39	Effective properties of two-phase disordered composite media: II. Evaluation of bounds on the conductivity and bulk modulus of dispersions of impenetrable spheres. Physical Review B, 1986, 33, 6428-6435.	1.1	102
40	Perturbation approach to the computer simulation of dipolar fluids. Journal of Chemical Physics, 1986, 85, 2913-2915.	1.2	4
41	Test of a simple analytic model for fluids of hard linear molecules. Molecular Physics, 1985, 54, 407-413.	0.8	27
42	Thermodynamic consistency conditions for the referenceâ€hypernetted chain equation with arbitrary molecular potential. Journal of Chemical Physics, 1984, 81, 4592-4593.	1.2	7
43	Choosing the reference system for liquid state perturbation theory. Molecular Physics, 1984, 52, 871-876.	0.8	82
44	Thermodynamically consistent perturbation theory for molecular fluids. Molecular Physics, 1984, 53, 363-368.	0.8	2
45	Solutions of the reference-hypernetted-chain equation with minimized free energy. Physical Review A, 1983, 28, 2374-2379.	1.0	304
46	Integral equations for fluids of linear molecules. Molecular Physics, 1982, 47, 313-317.	0.8	7
47	Integral equations for fluids of linear molecules. Molecular Physics, 1982, 47, 283-298.	0.8	93
48	Integral equations for fluids of linear molecules. Molecular Physics, 1982, 47, 299-311.	0.8	43
49	A local thermodynamic criterion for the reference-hypernetted chain equation. Physics Letters, Section A: General, Atomic and Solid State Physics, 1982, 89, 196-198.	0.9	143
50	Some topics in the molecular dynamics ensemble. Journal of Chemical Physics, 1981, 75, 5461-5463.	1.2	21
51	Hypernetted-chain solutions for the two-dimensional classical electron gas. Physical Review B, 1978, 17, 2827-2832.	1.1	72
52	Charged hard spheres in a uniform neutralizing background using â€~mixed' integral equations. Molecular Physics, 1976, 31, 1117-1127.	0.8	62
53	Calculation of NMR line shapes in calcium fluoride from modified moment expansions. Physical Review B, 1974, 9, 22-28.	1.1	21
54	Calculation of a corrected pair distribution function. Journal of Chemical Physics, 1974, 60, 1686-1687.	1.2	13

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55	Perturbation Correction for the Free Energy and Structure of Simple Fluids. Physical Review A, 1973, 8, 2548-2552.	1.0	196
56	Calculation of NMR Line Shapes Using the Memory-Function Approach. Physical Review B, 1973, 8, 3081-3092.	1.1	40
57	Perturbation correction for the free energy and structure of simple fluid mixtures. Journal of Chemical Physics, 1973, 59, 4830-4835.	1.2	16
58	Diffusive and Collective Motion in Classical Fluids. Journal of Chemical Physics, 1972, 57, 3003-3006.	1.2	5
59	Numerical Calculation of the Density Autocorrelation Function for Liquid Argon. Physical Review A, 1972, 5, 2238-2244.	1.0	3
60	Monte Carlo calculation of normal and abnormal diffusion in Ehrenfest's wind-tree model. Journal of Computational Physics, 1971, 7, 528-546.	1.9	36
61	Numerical fourier transforms in one, two, and three dimensions for liquid state calculations. Journal of Computational Physics, 1971, 8, 417-433.	1.9	141
62	Density Autocorrelation Function in a Classical Fluid from Initial Correlations. Physical Review A, 1970, 2, 1467-1473.	1.0	18
63	Equation of State of the Hardâ€Ðisk Fluid from Approximate Integral Equations. Journal of Chemical Physics, 1968, 49, 3092-3096.	1.2	162
64	N Dependence in Monte Carlo Studies of the Squareâ€Well System. Journal of Chemical Physics, 1968, 49, 4244-4245.	1.2	18
65	Effective Potential Description of the Quantum Ideal Gases. Journal of Chemical Physics, 1967, 47, 5369-5375.	1.2	80
66	Pressureâ€Consistent Integral Equation for Classical Fluids: Hardâ€6phere Solutions. Journal of Chemical Physics, 1967, 47, 4828-4833.	1.2	80
67	Approximate Methods for Obtaining Radial Distribution Functions of Fluids. Physical Review, 1965, 137, A42-A50.	2.7	23
68	Perturbation Correction to the Radial Distribution Function. Physical Review, 1964, 135, A1013-A1017.	2.7	75