

Mauro Rovere

List of Publications by Year in descending order

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107
papers

3,541
citations

117625

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144013

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108
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108
docs citations

108
times ranked

2054
citing authors

#	ARTICLE	IF	CITATIONS
1	Spontaneous NaCl-doped ices I _h , I _c , III, V and VI. Understanding the mechanism of ion inclusion and its dependence on the crystalline structure of ice. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 22897-22911.	2.8	7
2	Supercooled Liquids: Glass Transition and Mode Coupling Theory. <i>Soft and Biological Matter</i> , 2021, , 265-300.	0.3	0
3	Supercooled Water. <i>Soft and Biological Matter</i> , 2021, , 301-321.	0.3	0
4	Methods of Computer Simulation. <i>Soft and Biological Matter</i> , 2021, , 131-193.	0.3	0
5	Structure and dynamics of nanoconfined water and aqueous solutions. <i>European Physical Journal E</i> , 2021, 44, 136.	1.6	38
6	Advances in the study of supercooled water. <i>European Physical Journal E</i> , 2021, 44, 143.	1.6	40
7	Slow dynamics of supercooled hydration water in contact with lysozyme: examining the cage effect at different length scales. <i>Philosophical Magazine</i> , 2020, 100, 2582-2595.	1.6	6
8	Characterization of hydration water in supercooled water-trehalose solutions: The role of the hydrogen bonds network. <i>Journal of Chemical Physics</i> , 2019, 151, 044507.	3.0	8
9	Molecular dynamics simulations of freezing-point depression of TIP4P/2005 water in solution with NaCl. <i>Journal of Molecular Liquids</i> , 2018, 261, 513-519.	4.9	43
10	High density liquid structure enhancement in glass forming aqueous solution of LiCl. <i>Journal of Chemical Physics</i> , 2018, 148, 222829.	3.0	5
11	Structural Properties of Ionic Aqueous Solutions. , 2018, , 153-162.		0
12	Fragile to strong crossover and Widom line in supercooled water: A comparative study. <i>Frontiers of Physics</i> , 2018, 13, 1.	5.0	18
13	Structural properties and fragile to strong transition in confined water. <i>Journal of Chemical Physics</i> , 2017, 146, 084505.	3.0	24
14	Spontaneous NaCl-doped ice at seawater conditions: focus on the mechanisms of ion inclusion. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 9566-9574.	2.8	53
15	Microscopic origin of the fragile to strong crossover in supercooled water: The role of activated processes. <i>Journal of Chemical Physics</i> , 2017, 146, 084502.	3.0	38
16	Freezing Temperatures, Ice Nanotubes Structures, and Proton Ordering of TIP4P/ICE Water inside Single Wall Carbon Nanotubes. <i>Journal of Physical Chemistry B</i> , 2017, 121, 10371-10381.	2.6	28
17	High precision determination of the melting points of water TIP4P/2005 and water TIP4P/Ice models by the direct coexistence technique. <i>Journal of Chemical Physics</i> , 2017, 147, 244506.	3.0	60
18	Slow Dynamics and Structure of Supercooled Water in Confinement. <i>Entropy</i> , 2017, 19, 185.	2.2	5

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19	Mode coupling theory and fragile to strong transition in supercooled TIP4P/2005 water. <i>Journal of Chemical Physics</i> , 2016, 144, 074503.	3.0	63
20	The Widom line and dynamical crossover in supercritical water: Popular water models versus experiments. <i>Journal of Chemical Physics</i> , 2015, 143, 114502.	3.0	35
21	Relation between the two-body entropy and the relaxation time in supercooled water. <i>Physical Review E</i> , 2015, 91, 012107.	2.1	18
22	Widom line and dynamical crossovers as routes to understand supercritical water. <i>Nature Communications</i> , 2014, 5, 5806.	12.8	116
23	Do ions affect the structure of water? The case of potassium halides. <i>Journal of Molecular Liquids</i> , 2014, 189, 52-56.	4.9	31
24	Computer Simulation Study of the Structure of LiCl Aqueous Solutions: Test of Non-Standard Mixing Rules in the Ion Interaction. <i>Journal of Physical Chemistry B</i> , 2014, 118, 7680-7691.	2.6	36
25	Fragile to strong crossover at the Widom line in supercooled aqueous solutions of NaCl. <i>Journal of Chemical Physics</i> , 2013, 139, 204503.	3.0	30
26	Mode coupling and fragile to strong transition in supercooled TIP4P water. <i>Journal of Chemical Physics</i> , 2012, 137, 164503.	3.0	58
27	Water confined in MCM-41: a mode coupling theory analysis. <i>Journal of Physics Condensed Matter</i> , 2012, 24, 064109.	1.8	43
28	Ion hydration and structural properties of water in aqueous solutions at normal and supercooled conditions: a test of the structure making and breaking concept. <i>Physical Chemistry Chemical Physics</i> , 2011, 13, 19814.	2.8	75
29	Structural Properties of High and Low Density Water in a Supercooled Aqueous Solution of Salt. <i>Journal of Physical Chemistry B</i> , 2011, 115, 1461-1468.	2.6	43
30	Excess entropy of water in a supercooled solution of salt. <i>Molecular Physics</i> , 2011, 109, 2969-2979.	1.7	8
31	Structure and thermodynamics of supercooled aqueous solutions: Ionic solutes compared with water in a hydrophobic environment. <i>Journal of Molecular Liquids</i> , 2011, 159, 18-23.	4.9	4
32	Lennard-Jones binary mixture in disordered matrices: exploring the mode coupling scenario at increasing confinement. <i>Journal of Physics Condensed Matter</i> , 2011, 23, 234118.	1.8	2
33	Water at interfaces. <i>Journal of Physics Condensed Matter</i> , 2010, 22, 280301.	1.8	2
34	Anomalous dynamics of water confined in MCM-41 at different hydrations. <i>Journal of Physics Condensed Matter</i> , 2010, 22, 284102.	1.8	49
35	Dynamic Crossover in Supercooled Confined Water: Understanding Bulk Properties through Confinement. <i>Journal of Physical Chemistry Letters</i> , 2010, 1, 729-733.	4.6	148
36	A route to explain water anomalies from results on an aqueous solution of salt. <i>Journal of Chemical Physics</i> , 2010, 132, 134508.	3.0	99

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37	Molecular dynamics studies on the thermodynamics of supercooled sodium chloride aqueous solution at different concentrations. <i>Journal of Physics Condensed Matter</i> , 2010, 22, 284104.	1.8	10
38	Mode-coupling behavior of a Lennard-Jones binary mixture upon increasing confinement. <i>Physical Review E</i> , 2009, 80, 061502.	2.1	16
39	Thermodynamics of supercooled water in solutions. <i>Journal of Physics: Conference Series</i> , 2009, 177, 012003.	0.4	5
40	Effect of concentration on the thermodynamics of sodium chloride aqueous solutions in the supercooled regime. <i>Journal of Chemical Physics</i> , 2009, 130, 154511.	3.0	15
41	Thermodynamic behavior and structural properties of an aqueous sodium chloride solution upon supercooling. <i>Journal of Chemical Physics</i> , 2008, 128, 244508.	3.0	28
42	Structural properties and liquid spinodal of water confined in a hydrophobic environment. <i>Physical Review E</i> , 2007, 76, 061202.	2.1	40
43	Spinodal of supercooled polarizable water. <i>Physical Review E</i> , 2007, 75, 011201.	2.1	12
44	The phase diagram of confined fluids. <i>Journal of Molecular Liquids</i> , 2007, 134, 90-93.	4.9	8
45	Supercooled water: A molecular dynamics simulation study with a polarizable potential. <i>Journal of Molecular Liquids</i> , 2006, 127, 28-32.	4.9	3
46	Glass Transition in Confinement. <i>AIP Conference Proceedings</i> , 2006, , .	0.4	1
47	Liquid-liquid coexistence in the phase diagram of a fluid confined in fractal porous materials. <i>Europhysics Letters</i> , 2006, 75, 901-907.	2.0	8
48	Glass transition in confinement: a Lennard-Jones binary mixture study. <i>Computer Physics Communications</i> , 2005, 169, 214-217.	7.5	3
49	Local order in aqueous solutions of rare gases and the role of the solute concentration: a computer simulation study with a polarizable potential. <i>Molecular Physics</i> , 2005, 103, 501-506.	1.7	5
50	Inherent structures and Kauzmann temperature of confined liquids. <i>Physical Review E</i> , 2005, 71, 031204.	2.1	10
51	Mode coupling behavior of a Lennard-Jones binary mixture: A comparison between bulk and confined phases. <i>Journal of Chemical Physics</i> , 2005, 123, 174510.	3.0	11
52	Slow dynamics of a confined supercooled binary mixture in comparison with the bulk phase. <i>Philosophical Magazine</i> , 2004, 84, 1397-1404.	1.6	3
53	Computer simulation of the phase diagram for a fluid confined in a fractal and disordered porous material. <i>Physical Review E</i> , 2004, 70, 061505.	2.1	18
54	Effects of confinement on static and dynamical properties of water. <i>European Physical Journal E</i> , 2003, 12, 77-81.	1.6	53

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55	Computer simulation of structural properties of dilute aqueous solutions of argon at supercritical conditions. <i>Journal of Chemical Physics</i> , 2003, 118, 3646-3650.	3.0	10
56	Slow dynamics of a confined supercooled binary mixture: Direct space analysis. <i>Physical Review E</i> , 2003, 67, 041202.	2.1	36
57	Slow dynamics of a confined supercooled binary mixture. II. Qspace analysis. <i>Physical Review E</i> , 2003, 68, 061209.	2.1	17
58	Strong layering effects and anomalous dynamical behaviour in confined water at low hydration. <i>Journal of Physics Condensed Matter</i> , 2003, 15, S145-S150.	1.8	12
59	Anomalous dynamics of confined water at low hydration. <i>Journal of Physics Condensed Matter</i> , 2003, 15, 7625-7633.	1.8	50
60	Double dynamical regime of confined water. <i>Journal of Physics Condensed Matter</i> , 2003, 15, 1521-1529.	1.8	14
61	Microscopic two-dimensional lattice model of dimer granular compaction with friction. <i>Physical Review E</i> , 2002, 66, 031301.	2.1	7
62	Stretched exponential relaxation in a diffusive lattice model. <i>Physical Review E</i> , 2002, 65, 026127.	2.1	5
63	Slow dynamics of k-mers on a square lattice. <i>The Philosophical Magazine: Physics of Condensed Matter B, Statistical Mechanics, Electronic, Optical and Magnetic Properties</i> , 2002, 82, 375-381.	0.6	2
64	Mode Coupling relaxation scenario in a confined glass former. <i>Europhysics Letters</i> , 2002, 57, 212-218.	2.0	29
65	Confined water in the low hydration regime. <i>Journal of Chemical Physics</i> , 2002, 117, 369-375.	3.0	86
66	Layer analysis of the structure of water confined in vycor glass. <i>Journal of Chemical Physics</i> , 2002, 116, 342.	3.0	143
67	Supercooled confined water and the mode coupling scenario. <i>Physica A: Statistical Mechanics and Its Applications</i> , 2002, 304, 53-58.	2.6	3
68	Single particle dynamics of a confined Lennard-Jones mixture in the supercooled regime. <i>Physica A: Statistical Mechanics and Its Applications</i> , 2002, 314, 530-538.	2.6	1
69	Random sequential adsorption and diffusion of dimers and k-mers on a square lattice. <i>Journal of Chemical Physics</i> , 2001, 114, 7563-7569.	3.0	19
70	Modifications of the hydrogen bond network of liquid water in a cylindrical SiO ₂ pore. <i>Journal of Molecular Liquids</i> , 2000, 85, 127-137.	4.9	92
71	Molecular dynamics study of the glass transition in confined water. <i>European Physical Journal Special Topics</i> , 2000, 10, Pr7-203-Pr7-206.	0.2	2
72	Non-exponential kinetic behaviour of confined water. <i>Europhysics Letters</i> , 2000, 49, 183-188.	2.0	52

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73	Supercooled Confined Water and the Mode Coupling Crossover Temperature. <i>Physical Review Letters</i> , 2000, 85, 4317-4320.	7.8	142
74	Glass transition and layering effects in confined water: A computer simulation study. <i>Journal of Chemical Physics</i> , 2000, 113, 11324-11335.	3.0	172
75	Water in confined geometries: experiments and simulations. <i>Journal of Physics Condensed Matter</i> , 2000, 12, A345-A350.	1.8	74
76	Studies of water in confinement by experiments and simulations. <i>European Physical Journal Special Topics</i> , 2000, 10, Pr7-187-Pr7-193.	0.2	13
77	Water in porous glasses. A computer simulation study. <i>Journal of Molecular Liquids</i> , 1999, 80, 165-178.	4.9	65
78	Evidence of glassy behaviour of water molecules in confined states. <i>The Philosophical Magazine: Physics of Condensed Matter B, Statistical Mechanics, Electronic, Optical and Magnetic Properties</i> , 1999, 79, 1923-1930.	0.6	23
79	Water in porous glasses. A computer simulation Study. <i>Journal of Molecular Liquids</i> , 1999, 80, 165-178.	4.9	33
80	A molecular dynamics simulation of water confined in a cylindrical SiO ₂ pore. <i>Journal of Chemical Physics</i> , 1998, 108, 9859-9867.	3.0	127
81	Theory of gas-gas phase transition in rare-gas binary mixtures. <i>Journal of Chemical Physics</i> , 1996, 105, 2020-2027.	3.0	4
82	Inelasticity effects in the neutron diffraction measurements from water steam using pulsed sources. <i>Journal of Molecular Liquids</i> , 1995, 64, 221-240.	4.9	11
83	Translational absorption band in low density mixtures of noble gases: the He-Xe case. <i>Molecular Physics</i> , 1995, 84, 1065-1075.	1.7	3
84	Fluid-fluid phase separation in binary mixtures of asymmetric non-additive hard spheres. <i>Journal of Physics Condensed Matter</i> , 1994, 6, A163-A166.	1.8	31
85	Simulation studies of gas-liquid transitions in two dimensions via a subsystem-block-density distribution analysis. <i>European Physical Journal B</i> , 1993, 90, 215-228.	1.5	58
86	A mean-field study of the temperature dependence of the layer magnetizations in a semi-infinite Ising model. <i>Nuovo Cimento Della Societa Italiana Di Fisica D - Condensed Matter, Atomic, Molecular and Chemical Physics, Biophysics</i> , 1993, 15, 541-546.	0.4	1
87	Computer simulation of critical phenomena in fluids. <i>Journal of Physics Condensed Matter</i> , 1993, 5, B193-B200.	1.8	10
88	Liquid-solid transition in the bond-particle model for elemental semiconductors. <i>The Philosophical Magazine: Physics of Condensed Matter B, Statistical Mechanics, Electronic, Optical and Magnetic Properties</i> , 1992, 65, 921-932.	0.6	2
89	Magnetic-Phase Transitions of Ising Surfaces with Modified Surface-Bulk Coupling: a Monte Carlo Study. <i>Europhysics Letters</i> , 1992, 20, 547-552.	2.0	14
90	Phase equilibrium in liquid binary mixtures of non-additive hard spheres. <i>Physica Scripta</i> , 1992, T45, 251-252.	2.5	7

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91	Freezing of liquid alkali metals as screened ionic plasmas. <i>Journal of Physics Condensed Matter</i> , 1991, 3, 1627-1636.	1.8	5
92	Statistical mechanical models for liquid and amorphous structure in covalently bonded systems. <i>Nuovo Cimento Della Societa Italiana Di Fisica D - Condensed Matter, Atomic, Molecular and Chemical Physics, Biophysics</i> , 1990, 12, 619-632.	0.4	1
93	Hybrid molecular dynamics. <i>Computer Physics Communications</i> , 1990, 60, 311-318.	7.5	29
94	The gas-liquid transition of the two-dimensional Lennard-Jones fluid. <i>Journal of Physics Condensed Matter</i> , 1990, 2, 7009-7032.	1.8	147
95	Ordering Transitions Induced by Coulomb Interactions. <i>Physics and Chemistry of Materials With Low-dimensional Structures</i> , 1989, , 221-238.	1.0	5
96	Liquid Structure and Freezing of Metals and Salts*. <i>Zeitschrift Fur Physikalische Chemie</i> , 1988, 156, 411-424.	2.8	4
97	Block Density Distribution Function Analysis of Two-Dimensional Lennard-Jones Fluids. <i>Europhysics Letters</i> , 1988, 6, 585-590.	2.0	104
98	Thermodynamic Consistency and Entropy Change in the Density-Wave Theory of Freezing. <i>Physics and Chemistry of Liquids</i> , 1987, 16, 157-162.	1.2	1
99	Liquid Dichlorides and Dichloride-Monochloride Mixtures. <i>ECS Proceedings Volumes</i> , 1987, 1987-7, 195-199.	0.1	0
100	Structure and dynamics of molten salts. <i>Reports on Progress in Physics</i> , 1986, 49, 1001-1081.	20.1	218
101	Freezing of a modulated liquid: The superionic-to-normal transition of strontium chloride. <i>Solid State Communications</i> , 1985, 55, 1109-1111.	1.9	12
102	On the density-wave theory of classical Wigner crystallisation. <i>Journal of Physics C: Solid State Physics</i> , 1985, 18, 3445-3455.	1.5	50
103	Liquid structure and freezing of the two-dimensional classical electron fluid. <i>Journal of Physics C: Solid State Physics</i> , 1985, 18, 4011-4019.	1.5	9
104	Freezing of Ionic Melts into Normal and Superionic Phases. <i>Physics and Chemistry of Liquids</i> , 1983, 13, 113-122.	1.2	20
105	Dynamic corrections for neutron scattering of water in the free molecule approximation. <i>Journal of Chemical Physics</i> , 1982, 77, 2647-2655.	3.0	19
106	Liquid alkali metals and alloys as electron-ion plasmas. <i>Physica B: Physics of Condensed Matter & C: Atomic, Molecular and Plasma Physics, Optics</i> , 1981, 111, 11-23.	0.9	48
107	The inelasticity correction for heavy water. <i>Journal of Chemical Physics</i> , 1980, 73, 3729-3734.	3.0	17