

# Stefano Mossa

## List of Publications by Year in descending order

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60  
papers

4,566  
citations

159585

30  
h-index

123424

61  
g-index

62  
all docs

62  
docs citations

62  
times ranked

4654  
citing authors

#	ARTICLE	IF	CITATIONS
1	The worldwide air transportation network: Anomalous centrality, community structure, and cities' global roles. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2005, 102, 7794-7799.	7.1	1,377
2	Equilibrium Cluster Phases and Low-Density Arrested Disordered States: The Role of Short-Range Attraction and Long-Range Repulsion. <i>Physical Review Letters</i> , 2004, 93, 055701.	7.8	434
3	Interplay between Time-Temperature Transformation and the Liquid-Liquid Phase Transition in Water. <i>Physical Review Letters</i> , 2002, 88, 195701.	7.8	225
4	Ground-State Clusters for Short-Range Attractive and Long-Range Repulsive Potentials. <i>Langmuir</i> , 2004, 20, 10756-10763.	3.5	187
5	Truncation of Power Law Behavior in "Scale-Free" Network Models due to Information Filtering. <i>Physical Review Letters</i> , 2002, 88, 138701.	7.8	172
6	Effective temperature of active matter. <i>Physical Review E</i> , 2008, 77, 051111.	2.1	163
7	Anomalous properties of the acoustic excitations in glasses on the mesoscopic length scale. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2009, 106, 16907-16912.	7.1	124
8	Potential energy, relaxation, vibrational dynamics and the boson peak, of hyperquenched glasses. <i>Journal of Physics Condensed Matter</i> , 2003, 15, S1051-S1068.	1.8	123
9	Disentangling density and temperature effects in the viscous slowing down of glassforming liquids. <i>Journal of Chemical Physics</i> , 2004, 120, 6135-6141.	3.0	113
10	Li <sup>+</sup> Solvation in Pure, Binary, and Ternary Mixtures of Organic Carbonate Electrolytes. <i>Journal of Physical Chemistry C</i> , 2015, 119, 4502-4515.	3.1	110
11	Measuring spatial distribution of the local elastic modulus in glasses. <i>Physical Review E</i> , 2013, 87, 042306.	2.1	104
12	Dynamics and configurational entropy in the Lewis-Wahnström model for supercooled orthoterphenyl. <i>Physical Review E</i> , 2002, 65, 041205.	2.1	98
13	Non-conservative forces and effective temperatures in active polymers. <i>Soft Matter</i> , 2011, 7, 10193.	2.7	80
14	Effective temperature of active complex matter. <i>Soft Matter</i> , 2011, 7, 3726.	2.7	79
15	Crossover (or Kovacs) Effect in an Aging Molecular Liquid. <i>Physical Review Letters</i> , 2004, 92, 045504.	7.8	78
16	Potential Energy Landscape Equation of State. <i>Physical Review Letters</i> , 2002, 88, 225701.	7.8	70
17	Acoustic excitations and elastic heterogeneities in disordered solids. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2014, 111, 11949-11954.	7.1	67
18	Molecular dynamics simulation of the fragile glass-former orthoterphenyl: A flexible molecule model. <i>Physical Review E</i> , 2000, 62, 612-630.	2.1	60

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19	Elastic heterogeneity, vibrational states, and thermal conductivity across an amorphisation transition. <i>Europhysics Letters</i> , 2013, 104, 56001.	2.0	58
20	Beating the amorphous limit in thermal conductivity by superlattices design. <i>Scientific Reports</i> , 2015, 5, 14116.	3.3	54
21	Water sub-diffusion in membranes for fuel cells. <i>Scientific Reports</i> , 2017, 7, 8326.	3.3	54
22	Routes to colloidal gel formation. <i>Computer Physics Communications</i> , 2005, 169, 166-171.	7.5	52
23	Morphology of Supported Polymer Electrolyte Ultrathin Films: A Numerical Study. <i>Journal of Physical Chemistry C</i> , 2015, 119, 1201-1216.	3.1	52
24	Inhomogeneous Transport in Model Hydrated Polymer Electrolyte Supported Ultrathin Films. <i>ACS Nano</i> , 2013, 7, 6767-6773.	14.6	50
25	Locally preferred structure in simple atomic liquids. <i>Journal of Chemical Physics</i> , 2003, 119, 8069-8074.	3.0	45
26	The Anion Effect on Li <sup>+</sup> Ion Coordination Structure in Ethylene Carbonate Solutions. <i>Journal of Physical Chemistry Letters</i> , 2016, 7, 3554-3559.	4.6	42
27	Statistical physics and liquid water at negative pressures. <i>Physica A: Statistical Mechanics and Its Applications</i> , 2002, 315, 281-289.	2.6	32
28	Aging in a Laponite colloidal suspension: A Brownian dynamics simulation study. <i>Journal of Chemical Physics</i> , 2007, 126, 014905.	3.0	32
29	Relation of vibrational excitations and thermal conductivity to elastic heterogeneities in disordered solids. <i>Physical Review B</i> , 2016, 94, .	3.2	31
30	Solvent and Salt Effect on Lithium Ion Solvation and Contact Ion Pair Formation in Organic Carbonates: A Quantum Chemical Perspective. <i>Journal of Physical Chemistry C</i> , 2018, 122, 25930-25939.	3.1	31
31	Polymer translocation through nano-pores in vibrating thin membranes. <i>Scientific Reports</i> , 2016, 6, 38558.	3.3	28
32	Aging and energy landscapes: application to liquids and glasses. <i>European Physical Journal B</i> , 2002, 30, 351-355.	1.5	25
33	Molecular dynamics simulation of the fragile glass former orthoterphenyl: A flexible molecule model. II. Collective dynamics. <i>Physical Review E</i> , 2001, 64, 021511.	2.1	23
34	Application of Statistical Physics to Understand Static and Dynamic Anomalies in Liquid Water. <i>Journal of Statistical Physics</i> , 2003, 110, 1039-1054.	1.2	23
35	Numerical evaluation of the statistical properties of a potential energy landscape. <i>Journal of Physics Condensed Matter</i> , 2003, 15, S1085-S1094.	1.8	21
36	Sub-diffusion and population dynamics of water confined in soft environments. <i>Nanoscale</i> , 2016, 8, 3314-3325.	5.6	20

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37	Glassy Potts model: A disordered Potts model without a ferromagnetic phase. <i>Physical Review B</i> , 1999, 59, 8401-8404.	3.2	17
38	Vibrational dynamics and thermodynamics, ideal glass transitions and folding transitions, in liquids and biopolymers. <i>AIP Conference Proceedings</i> , 2004, , .	0.4	15
39	Sound damping in glasses: Interplay between anharmonicities and elastic heterogeneities. <i>Physical Review B</i> , 2020, 101, .	3.2	15
40	An operational scheme to determine the locally preferred structure of model liquids. <i>Journal of Non-Crystalline Solids</i> , 2006, 352, 4847-4850.	3.1	14
41	Cutoff nonlinearities in the low-temperature vibrations of glasses and crystals. <i>Physical Review E</i> , 2016, 93, 043314.	2.1	14
42	Anharmonic thermodynamics of vacancies using a neural network potential. <i>Physical Review Materials</i> , 2019, 3, .	2.4	14
43	Water confined in self-assembled ionic surfactant nano-structures. <i>Soft Matter</i> , 2015, 11, 2469-2478.	2.7	13
44	The effect of polymorphism on the structural, dynamic and dielectric properties of plastic crystal water: A molecular dynamics simulation perspective. <i>Journal of Chemical Physics</i> , 2019, 150, 124506.	3.0	13
45	Impact of elastic heterogeneity on the propagation of vibrations at finite temperatures in glasses. <i>Condensed Matter Physics</i> , 2019, 22, 43604.	0.7	13
46	Liquid stability in a model for ortho-terphenyl. <i>Journal of Chemical Physics</i> , 2004, 120, 6128-6134.	3.0	11
47	Molecular dynamics simulation study of the high frequency sound waves in the fragile glass former orthoterphenyl. <i>Journal of Chemical Physics</i> , 2002, 116, 1077-1084.	3.0	9
48	Response to: "Comment on "Disentangling density and temperature effects in the viscous slowing down of glassforming liquids" [J. Chem. Phys. 121, 11503 (2004)]. <i>Journal of Chemical Physics</i> , 2004, 121, 11505.	3.0	9
49	Orientational and induced contributions to the depolarized Rayleigh spectra of liquid and supercooled ortho-terphenyl. <i>Journal of Chemical Physics</i> , 2002, 117, 3289-3295.	3.0	8
50	Equilibrium and out-of-equilibrium thermodynamics in supercooled liquids and glasses. <i>Journal of Physics Condensed Matter</i> , 2003, 15, S351-S357.	1.8	8
51	Effect of Surface Hydrophilicity on the Formation of Nafion Thin Films Inside PEMFC Catalyst Layers: A Computational Study. <i>ECS Transactions</i> , 2013, 45, 101-108.	0.5	8
52	Structure and dynamics of liquid CS <sub>2</sub> : Going from ambient to elevated pressure conditions. <i>Journal of Chemical Physics</i> , 2016, 145, 154505.	3.0	8
53	Vibrational origin of the fast relaxation processes in molecular glass formers. <i>Europhysics Letters</i> , 2002, 60, 92-98.	2.0	7
54	From Ionic Surfactants to Nafion through Convolutional Neural Networks. <i>Journal of Physical Chemistry B</i> , 2020, 124, 8918-8927.	2.6	7

#	ARTICLE	IF	CITATIONS
55	Quenches and crunches: Does the system explore in ageing the same part of the configuration space explored in equilibrium?. The Philosophical Magazine: Physics of Condensed Matter B, Statistical Mechanics, Electronic, Optical and Magnetic Properties, 2002, 82, 695-705.	0.6	6
56	Quenches and crunches: does the system explore in ageing the same part of the configuration space explored in equilibrium?. The Philosophical Magazine: Physics of Condensed Matter B, Statistical Mechanics, Electronic, Optical and Magnetic Properties, 2002, 82, 695-705.	0.6	6
57	Re-entrant Phase Transitions and Dynamics of a Nanoconfined Ionic Liquid. Physical Review X, 2018, 8, .	8.9	5
58	Intermittent rearrangements accompanying thermal fluctuations distinguish glasses from crystals. Journal of Chemical Physics, 2020, 153, 154501.	3.0	5
59	Time correlation functions for quantum systems: Validating Bayesian approaches for harmonic oscillators and beyond. Journal of Chemical Physics, 2021, 155, 134108.	3.0	1
60	Water at Positive and Negative Pressures. , 2002, , 59-67.		1