

Alessandro Sergi

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/9074104/publications.pdf>

Version: 2024-02-01

64
papers

1,040
citations

394421

19
h-index

454955

30
g-index

66
all docs

66
docs citations

66
times ranked

528
citing authors

#	ARTICLE	IF	CITATIONS
1	Symmetry-Induced Emergence of a Pseudo-Qutrit in the Dipolar Coupling of Two Qubits. Entropy, 2022, 24, 223.	2.2	2
2	DECIDE: A Deterministic Mixed Quantum-Classical Dynamics Approach. Applied Sciences (Switzerland), 2022, 12, 7022.	2.5	1
3	Evolution of a Non-Hermitian Quantum Single-Molecule Junction at Constant Temperature. Entropy, 2021, 23, 147.	2.2	8
4	Two-Qubit Entanglement Generation through Non-Hermitian Hamiltonians Induced by Repeated Measurements on an Ancilla. Entropy, 2020, 22, 1184.	2.2	19
5	Simulating Quantum Dynamics in Classical Nanoscale Environments. , 2020, , 515-544.		0
6	A variational mean-field study of clusterization in a zero-temperature system of soft-core bosons. EPJ Web of Conferences, 2020, 230, 00008.	0.3	3
7	Proposal of a Computational Approach for Simulating Thermal Bosonic Fields in Phase Space. Physics, 2019, 1, 402-411.	1.4	4
8	Clusterization of weakly-interacting bosons in one dimension: an analytic study at zero temperature. Journal of Physics A: Mathematical and Theoretical, 2019, 52, 015002.	2.1	9
9	Quasi-Lie Brackets and the Breaking of Time-Translation Symmetry for Quantum Systems Embedded in Classical Baths. Symmetry, 2018, 10, 518.	2.2	15
10	Freezing of soft-core bosons at zero temperature: A variational theory. Physical Review B, 2018, 98, .	3.2	18
11	Linear Quantum Entropy and Non-Hermitian Hamiltonians. Entropy, 2016, 18, 451.	2.2	20
12	On the configurational temperature NosÃˆ" Hoover thermostat. Physica A: Statistical Mechanics and Its Applications, 2016, 461, 19-35.	2.6	12
13	Quantum entropy of systems described by non-Hermitian Hamiltonians. Journal of Statistical Mechanics: Theory and Experiment, 2016, 2016, 033102.	2.3	30
14	Quantum dynamics of a plasmonic metamolecule with a time-dependent driving. Theoretical Chemistry Accounts, 2015, 134, 1.	1.4	6
15	Time correlation functions for non-Hermitian quantum systems. Physical Review A, 2015, 91, .	2.5	48
16	Embedding quantum systems with a non-conserved probability in classical environments. Theoretical Chemistry Accounts, 2015, 134, 1.	1.4	11
17	Non-Hamiltonian Modeling of Squeezing and Thermal Disorder in Driven Oscillators. Journal of Statistical Physics, 2015, 159, 255-273.	1.2	2
18	Comparison and unification of non-Hermitian and Lindblad approaches with applications to open quantum optical systems. Journal of Modern Optics, 2014, 61, 1298-1308.	1.3	54

#	ARTICLE	IF	CITATIONS
19	Computer simulation of quantum dynamics in a classical spin environment. <i>Theoretical Chemistry Accounts</i> , 2014, 133, 1.	1.4	6
20	Quantum dynamics in classical thermal baths. <i>Computer Physics Communications</i> , 2013, 184, 2474-2477.	7.5	3
21	Minimum-density anomaly and spatial ordering of softly repulsive particles in a narrow channel. <i>Soft Matter</i> , 2013, 9, 9876.	2.7	6
22	Quantum dynamics of a nano-rod under compression. <i>Physics Letters, Section A: General, Atomic and Solid State Physics</i> , 2013, 377, 1047-1051.	2.1	4
23	Transient behavior of a model fluid under applied shear. <i>Journal of Chemical Physics</i> , 2013, 138, 184501.	3.0	0
24	Quantum dynamics in the partial Wigner picture. <i>Journal of Physics A: Mathematical and Theoretical</i> , 2013, 46, 395305.	2.1	3
25	Filtering schemes in the quantum-classical Liouville approach to nonadiabatic dynamics. <i>Physical Review E</i> , 2013, 88, 033301.	2.1	16
26	Communication: Quantum dynamics in classical spin baths. <i>Journal of Chemical Physics</i> , 2013, 139, 031101.	3.0	6
27	NON-HERMITIAN QUANTUM DYNAMICS OF A TWO-LEVEL SYSTEM AND MODELS OF DISSIPATIVE ENVIRONMENTS. <i>International Journal of Modern Physics B</i> , 2013, 27, 1350163.	2.0	88
28	Stochastic simulation of long-time nonadiabatic dynamics. <i>Physica Scripta</i> , 2011, T143, 014024.	2.5	10
29	Matrix Algebras in Non-Hermitian Quantum Mechanics. <i>Communications in Theoretical Physics</i> , 2011, 56, 96-98.	2.5	7
30	MOMENTUM SHIFT IN NONADIABATIC DYNAMICS. <i>Modern Physics Letters B</i> , 2011, 25, 1271-1280.	1.9	1
31	Bulgac-Kusnezov-Nosé-Hoover thermostats. <i>Physical Review E</i> , 2010, 81, 036705.	2.1	8
32	Sampling of quantum dynamics at long time. <i>Physical Review E</i> , 2010, 81, 032101.	2.1	15
33	Numerical and analytical approach to the quantum dynamics of two coupled spins in bosonic baths. <i>Physical Review A</i> , 2009, 80, .	2.5	24
34	Nosé-Hoover dynamics in quantum phase space. <i>Journal of Physics A: Mathematical and Theoretical</i> , 2008, 41, 355304.	2.1	5
35	Quantum-classical dynamics of wave fields. <i>Journal of Chemical Physics</i> , 2007, 126, 074109.	3.0	6
36	Deterministic constant-temperature dynamics for dissipative quantum systems. <i>Journal of Physics A: Mathematical and Theoretical</i> , 2007, 40, F347-F354.	2.1	23

#	ARTICLE	IF	CITATIONS
37	On the geometry and entropy of non-Hamiltonian phase space. <i>Journal of Statistical Mechanics: Theory and Experiment</i> , 2007, 2007, P02013-P02013.	2.3	13
38	On Computational Strategies within Molecular Dynamics Simulation. <i>Physics Essays</i> , 2007, 20, 629-640.	0.4	4
39	Freezing a Single Distal Motion in Dihydrofolate Reductase. <i>Journal of Physical Chemistry B</i> , 2006, 110, 2435-2441.	2.6	15
40	Statistical mechanics of quantum-classical systems with holonomic constraints. <i>Journal of Chemical Physics</i> , 2006, 124, 024110.	3.0	12
41	Nonadiabatic chemical reactions. <i>Computer Physics Communications</i> , 2005, 169, 400-403.	7.5	4
42	Quantum-Classical Wigner-Liouville Equation. <i>Ukrainian Mathematical Journal</i> , 2005, 57, 891-899.	0.5	0
43	Phase space flows for non-Hamiltonian systems with constraints. <i>Physical Review E</i> , 2005, 72, 031104.	2.1	14
44	Non-Hamiltonian commutators in quantum mechanics. <i>Physical Review E</i> , 2005, 72, 066125.	2.1	38
45	Experimental and Simulative Dissociation of Dimeric Cu,Zn Superoxide Dismutase Doubly Mutated at the Intersubunit Surface. <i>Biophysical Journal</i> , 2005, 88, 2875-2882.	0.5	3
46	Non-Equilibrium Molecular Dynamics. , 2005, , 745-761.		10
47	Simulating Reactions That Occur Once in a Blue Moon. , 2005, , 1597-1611.		3
48	Non-Equilibrium Molecular Dynamics. , 2005, , 745-761.		1
49	Simulating Reactions That Occur Once in a Blue Moon. , 2005, , 1597-1611.		0
50	Generalized bracket formulation of constrained dynamics in phase space. <i>Physical Review E</i> , 2004, 69, 021109.	2.1	19
51	Quantum-classical limit of quantum correlation functions. <i>Journal of Chemical Physics</i> , 2004, 121, 7565.	3.0	26
52	Simulating quantum dynamics in classical environments. <i>Theoretical Chemistry Accounts</i> , 2003, 110, 49-58.	1.4	60
53	Nonadiabatic reaction rates for dissipative quantum-classical systems. <i>Journal of Chemical Physics</i> , 2003, 119, 12776-12783.	3.0	21
54	Non-Hamiltonian equilibrium statistical mechanics. <i>Physical Review E</i> , 2003, 67, 021101.	2.1	42

#	ARTICLE	IF	CITATIONS
55	Quantum-classical dynamics of nonadiabatic chemical reactions. <i>Journal of Chemical Physics</i> , 2003, 118, 8566-8575.	3.0	38
56	Effective binding force calculation in a dimeric protein by molecular dynamics simulation. <i>Journal of Chemical Physics</i> , 2002, 116, 6329-6338.	3.0	27
57	Dynamics-Function Correlation in Cu, Zn Superoxide Dismutase: A Spectroscopic and Molecular Dynamics Simulation Study. <i>Biophysical Journal</i> , 2001, 80, 2556-2567.	0.5	26
58	Density Functional Study of the Photoactive Yellow Protein's Chromophore. <i>Journal of Physical Chemistry B</i> , 2001, 105, 4386-4391.	2.6	59
59	Non-Hamiltonian equations of motion with a conserved energy. <i>Physical Review E</i> , 2001, 64, 056125.	2.1	53
60	Structure of phosphorus-selenium glasses: results from ab initio molecular dynamics simulations. <i>Molecular Physics</i> , 2000, 98, 701-707.	1.7	5
61	Reversible integrators for basic extended system molecular dynamics. <i>Molecular Physics</i> , 1999, 97, 825-832.	1.7	15
62	First-principles simulation of phosphorus-selenium systems. <i>Chemical Physics Letters</i> , 1996, 259, 301-306.	2.6	3
63	Molecular dynamics study of the plastic-crystalline phase transition of tetraphosphorus triselenide. <i>Molecular Physics</i> , 1995, 84, 727-742.	1.7	6
64	Simulation of superoxide-superoxide dismutase association rate for six natural variants. Comparison with the experimental catalytic rate. <i>The Journal of Physical Chemistry</i> , 1994, 98, 10554-10557.	2.9	28