

# Yoshitaka Tanimura

## List of Publications by Year in descending order

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147  
papers

7,796  
citations

53939

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58552

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152  
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152  
docs citations

152  
times ranked

2508  
citing authors

#	ARTICLE	IF	CITATIONS
1	Imaginary-time hierarchical equations of motion for thermodynamic variables. Journal of Chemical Physics, 2022, 156, 174112.	1.2	2
2	The laws of thermodynamics for quantum dissipative systems: A quasi-equilibrium Helmholtz energy approach. Journal of Chemical Physics, 2022, 157, .	1.2	4
3	Exciton transfer in organic photovoltaic cells: A role of local and nonlocal electron-phonon interactions in a donor domain. Journal of Chemical Physics, 2021, 154, 034107.	1.2	14
4	Full molecular dynamics simulations of molecular liquids for single-beam spectrally controlled two-dimensional Raman spectroscopy. Journal of Chemical Physics, 2021, 154, 124115.	1.2	0
5	Open Quantum Dynamics Theory for Non-Equilibrium Work: Hierarchical Equations of Motion Approach. Journal of the Physical Society of Japan, 2021, 90, 033001.	0.7	7
6	Probing photoinduced proton coupled electron transfer process by means of two-dimensional resonant electronic-vibrational spectroscopy. Journal of Chemical Physics, 2021, 154, 144104.	1.2	6
7	Modeling and Simulating the Excited-State Dynamics of a System with Condensed Phases: A Machine Learning Approach. Journal of Chemical Theory and Computation, 2021, 17, 3618-3628.	2.3	8
8	Open quantum dynamics theory on the basis of periodical system-bath model for discrete Wigner function. Journal of Computational Electronics, 2021, 20, 2091-2103.	1.3	3
9	Optical response of laser-driven charge-transfer complex described by Holstein-Hubbard model coupled to heat baths: Hierarchical equations of motion approach. Journal of Chemical Physics, 2021, 155, 064106.	1.2	7
10	Autobiography of Yoshitaka Tanimura. Journal of Physical Chemistry B, 2021, 125, 11787-11792.	1.2	0
11	Open quantum dynamics theory for a complex subenvironment system with a quantum thermostat: Application to a spin heat bath. Journal of Chemical Physics, 2021, 155, 244109.	1.2	2
12	Open Quantum Dynamics Theory of Spin Relaxation: Application to $^1\text{H}$ SR and Low-Field NMR Spectroscopies. Journal of the Physical Society of Japan, 2020, 89, 064710.	0.7	12
13	Numerically "exact" approach to open quantum dynamics: The hierarchical equations of motion (HEOM). Journal of Chemical Physics, 2020, 153, 020901.	1.2	219
14	Numerically "exact" simulations of entropy production in the fully quantum regime: Boltzmann entropy vs von Neumann entropy. Journal of Chemical Physics, 2020, 153, 234107.	1.2	10
15	Proton tunneling in a two-dimensional potential energy surface with a non-linear system-bath interaction: Thermal suppression of reaction rate. Journal of Chemical Physics, 2020, 152, 214114.	1.2	19
16	Modeling Intermolecular and Intramolecular Modes of Liquid Water Using Multiple Heat Baths: Machine Learning Approach. Journal of Chemical Theory and Computation, 2020, 16, 2099-2108.	2.3	13
17	Open quantum dynamics of a three-dimensional rotor calculated using a rotationally invariant system-bath Hamiltonian: Linear and two-dimensional rotational spectra. Journal of Chemical Physics, 2019, 151, 044105.	1.2	9
18	Low-Temperature Quantum Fokker-Planck and Smoluchowski Equations and Their Extension to Multistate Systems. Journal of Chemical Theory and Computation, 2019, 15, 2517-2534.	2.3	21

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19	Modeling and analyzing a photo-driven molecular motor system: Ratchet dynamics and non-linear optical spectra. <i>Journal of Chemical Physics</i> , 2019, 150, 114103.	1.2	15
20	Hierarchical Equations of Motion Approach to Quantum Thermodynamics. <i>Fundamental Theories of Physics</i> , 2018, , 579-595.	0.1	5
21	Linear absorption spectrum of a quantum two-dimensional rotator calculated using a rotationally invariant system-bath Hamiltonian. <i>Journal of Chemical Physics</i> , 2018, 149, 084110.	1.2	6
22	Phase-space wavepacket dynamics of internal conversion via conical intersection: Multi-state quantum Fokker-Planck equation approach. <i>Chemical Physics</i> , 2018, 515, 203-213.	0.9	20
23	Hierarchical Schrödinger equations of motion for open quantum dynamics. <i>Physical Review A</i> , 2018, 98, .	1.0	32
24	Exploring complete positivity in hierarchy equations of motion. <i>New Journal of Physics</i> , 2017, 19, 013007.	1.2	8
25	Exciton-Coupled Electron Transfer Process Controlled by Non-Markovian Environments. <i>Journal of Physical Chemistry Letters</i> , 2017, 8, 5390-5394.	2.1	12
26	Probing photoisomerization processes by means of multi-dimensional electronic spectroscopy: The multi-state quantum hierarchical Fokker-Planck equation approach. <i>Journal of Chemical Physics</i> , 2017, 147, 014102.	1.2	25
27	Chapter 13 Simulating the Nonlinear Optical Response of Multichromophore Complexes. , 2017, , 467-488.		0
28	Simulating two-dimensional infrared-Raman and Raman spectroscopies for intermolecular and intramolecular modes of liquid water. <i>Journal of Chemical Physics</i> , 2016, 144, 074201.	1.2	38
29	Quantum heat current under non-perturbative and non-Markovian conditions: Applications to heat machines. <i>Journal of Chemical Physics</i> , 2016, 145, 224105.	1.2	84
30	Effects of Intermolecular Charge Transfer in Liquid Water on Raman Spectra. <i>Journal of Physical Chemistry Letters</i> , 2016, 7, 4147-4151.	2.1	22
31	Full molecular dynamics simulations of liquid water and carbon tetrachloride for two-dimensional Raman spectroscopy in the frequency domain. <i>Chemical Physics</i> , 2016, 481, 245-249.	0.9	10
32	Electron Pumping under Non-Markovian Dissipation: The Role of the Self-Consistent Field. <i>Journal of the Physical Society of Japan</i> , 2016, 85, 034803.	0.7	2
33	Fast, Accurate Simulation of Polaron Dynamics and Multidimensional Spectroscopy by Multiple Davydov Trial States. <i>Journal of Physical Chemistry A</i> , 2016, 120, 1562-1576.	1.1	60
34	Notes on simulating two-dimensional Raman and terahertz-Raman signals with a full molecular dynamics simulation approach. <i>Structural Dynamics</i> , 2015, 2, 054102.	0.9	23
35	Analysis of 2D THz-Raman spectroscopy using a non-Markovian Brownian oscillator model with nonlinear system-bath interactions. <i>Journal of Chemical Physics</i> , 2015, 142, 212421.	1.2	41
36	Real-time and imaginary-time quantum hierarchical Fokker-Planck equations. <i>Journal of Chemical Physics</i> , 2015, 142, 144110.	1.2	104

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37	Spins Dynamics in a Dissipative Environment: Hierarchical Equations of Motion Approach Using a Graphics Processing Unit (GPU). <i>Journal of Chemical Theory and Computation</i> , 2015, 11, 3859-3865.	2.3	37
38	Dynamics of a One-Dimensional Holstein Polaron with the Hierarchical Equations of Motion Approach. <i>Journal of Physical Chemistry Letters</i> , 2015, 6, 3110-3115.	2.1	66
39	Linear and third- and fifth-order nonlinear spectroscopies of a charge transfer system coupled to an underdamped vibration. <i>Journal of Chemical Physics</i> , 2015, 142, 212423.	1.2	23
40	Quantum heat transport of a two-qubit system: Interplay between system-bath coherence and qubit-qubit coherence. <i>Journal of Chemical Physics</i> , 2015, 143, 064107.	1.2	51
41	Reduced hierarchical equations of motion in real and imaginary time: Correlated initial states and thermodynamic quantities. <i>Journal of Chemical Physics</i> , 2014, 141, 044114.	1.2	144
42	Calculating two-dimensional THz-Raman-THz and Raman-THz-THz signals for various molecular liquids: The samplers. <i>Journal of Chemical Physics</i> , 2014, 141, 124503.	1.2	26
43	Self-excited current oscillations in a resonant tunneling diode described by a model based on the Caldeira-Leggett Hamiltonian. <i>New Journal of Physics</i> , 2014, 16, 015002.	1.2	22
44	Molecular dynamics simulation for infrared spectroscopy with intramolecular forces from electronic properties of on-the-fly quantum chemical calculations. <i>International Journal of Quantum Chemistry</i> , 2013, 113, 330-335.	1.0	5
45	Quantum Suppression of Ratchet Rectification in a Brownian System Driven by a Biharmonic Force. <i>Journal of Physical Chemistry B</i> , 2013, 117, 13132-13144.	1.2	31
46	Simulation of femtosecond double-slit experiments for a chromophore in a dissipative environment. <i>Journal of Chemical Physics</i> , 2013, 139, 214302.	1.2	19
47	An Approach to Quantum Transport Based on Reduced Hierarchy Equations of Motion: Application to a Resonant Tunneling Diode. <i>Journal of the Physical Society of Japan</i> , 2013, 82, 033707.	0.7	19
48	The role of the environment time scale in light-harvesting efficiency and coherent oscillations. <i>New Journal of Physics</i> , 2012, 14, 073027.	1.2	31
49	Reduced hierarchy equations of motion approach with Drude plus Brownian spectral distribution: Probing electron transfer processes by means of two-dimensional correlation spectroscopy. <i>Journal of Chemical Physics</i> , 2012, 137, 22A550.	1.2	107
50	Dephasing by a continuous-time random walk process. <i>Physical Review E</i> , 2012, 86, 011130.	0.8	2
51	Note: Inverted time-ordering in two-dimensional-Raman-terahertz spectroscopy of water. <i>Journal of Chemical Physics</i> , 2012, 136, 236101.	1.2	24
52	System Bath Correlations and the Nonlinear Response of Qubits. <i>Journal of the Physical Society of Japan</i> , 2012, 81, 063301.	0.7	15
53	Non-Markovianity: initial correlations and nonlinear optical measurements. <i>Philosophical Transactions Series A, Mathematical, Physical, and Engineering Sciences</i> , 2012, 370, 3658-3671.	1.6	17
54	Infrared Spectral Signatures of Multilayered Surface-Fluorinated Graphene: A Molecular Dynamics Study. <i>Journal of Physical Chemistry C</i> , 2012, 116, 8343-8347.	1.5	7

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55	Infrared Spectral Signatures of Surface-Fluorinated Graphene: A Molecular Dynamics Study. <i>Journal of Physical Chemistry Letters</i> , 2012, 3, 246-250.	2.1	13
56	A Polarizable Water Model for Intramolecular and Intermolecular Vibrational Spectroscopies. <i>Journal of Physical Chemistry B</i> , 2011, 115, 5545-5553.	1.2	101
57	Does $\hat{\rho}$ Play a Role in Multidimensional Spectroscopy? Reduced Hierarchy Equations of Motion Approach to Molecular Vibrations. <i>Journal of Physical Chemistry A</i> , 2011, 115, 4009-4022.	1.1	62
58	Discussions on Session 6A: Applications of quantum coherence. <i>Procedia Chemistry</i> , 2011, 3, 347-351.	0.7	0
59	Non-Gaussian stochastic dynamics of spins and oscillators: A continuous-time random walk approach. <i>Physical Review E</i> , 2011, 84, 061111.	0.8	6
60	Multistate electron transfer dynamics in the condensed phase: Exact calculations from the reduced hierarchy equations of motion approach. <i>Journal of Chemical Physics</i> , 2010, 132, 214502.	1.2	85
61	Correlated fluctuations in the exciton dynamics and spectroscopy of DNA. <i>New Journal of Physics</i> , 2010, 12, 055005.	1.2	29
62	Distinct Infrared Spectral Signatures of the 1,2- and 1,4-Fluorinated Single-Walled Carbon Nanotubes: A Molecular Dynamics Study. <i>Journal of Physical Chemistry Letters</i> , 2010, 1, 1307-1311.	2.1	12
63	Non-Markovian Entanglement Dynamics in the Presence of System-Bath Coherence. <i>Physical Review Letters</i> , 2010, 104, 250401.	2.9	170
64	Coherent Multidimensional Optical Spectroscopy. <i>Accounts of Chemical Research</i> , 2009, 42, 1207-1209.	7.6	81
65	Modeling, Calculating, and Analyzing Multidimensional Vibrational Spectroscopies. <i>Accounts of Chemical Research</i> , 2009, 42, 1270-1279.	7.6	82
66	Quantum Dissipative Dynamics of Electron Transfer Reaction System: Nonperturbative Hierarchy Equations Approach. <i>Journal of the Physical Society of Japan</i> , 2009, 78, 073802.	0.7	84
67	Detecting the Dzyaloshinskii-Moriya interaction by means of pulsed EPR spectroscopy. <i>Chemical Physics Letters</i> , 2008, 457, 237-240.	1.2	4
68	Nonperturbative non-Markovian quantum master equation: Validity and limitation to calculate nonlinear response functions. <i>Chemical Physics</i> , 2008, 347, 185-193.	0.9	105
69	Ultrafast exciton transfers in DNA and its nonlinear optical spectroscopy. <i>Journal of Chemical Physics</i> , 2008, 128, 135102.	1.2	16
70	Two-dimensional fifth-order Raman spectroscopy of liquid formamide: Experiment and Theory. <i>Journal of Chemical Physics</i> , 2008, 128, 234507.	1.2	48
71	Exploring a free energy landscape by means of multidimensional infrared and terahertz spectroscopies. <i>Journal of Chemical Physics</i> , 2008, 128, 164501.	1.2	4
72	Nonequilibrium molecular dynamics simulations with a backward-forward trajectories sampling for multidimensional infrared spectroscopy of molecular vibrational modes. <i>Journal of Chemical Physics</i> , 2008, 128, 064511.	1.2	46

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73	Two-dimensional infrared surface spectroscopy for CO on Cu(100): Detection of intermolecular coupling of adsorbates. <i>Journal of Chemical Physics</i> , 2007, 126, 204703.	1.2	20
74	Ultrafast exciton-exciton coherent transfer in molecular aggregates and its application to light-harvesting systems. <i>Journal of Chemical Physics</i> , 2007, 127, 075101.	1.2	24
75	Free energy landscapes of electron transfer system in dipolar environment below and above the rotational freezing temperature. <i>Journal of Chemical Physics</i> , 2007, 126, 054504.	1.2	3
76	Dynamics of a Multimode System Coupled to Multiple Heat Baths Probed by Two-Dimensional Infrared Spectroscopy. <i>Journal of Physical Chemistry A</i> , 2007, 111, 9269-9276.	1.1	80
77	Two-dimensional Raman spectra of atomic solids and liquids. <i>Journal of Chemical Physics</i> , 2006, 124, 024508.	1.2	27
78	Calculating fifth-order Raman signals for various molecular liquids by equilibrium and nonequilibrium hybrid molecular dynamics simulation algorithms. <i>Journal of Chemical Physics</i> , 2006, 125, 074512.	1.2	57
79	Modeling vibrational dephasing and energy relaxation of intramolecular anharmonic modes for multidimensional infrared spectroscopies. <i>Journal of Chemical Physics</i> , 2006, 125, 084501.	1.2	104
80	Analyzing atomic liquids and solids by means of two-dimensional Raman spectra in frequency domain. <i>Journal of Chemical Physics</i> , 2006, 124, 194504.	1.2	19
81	Free energy landscape analysis of two-dimensional dipolar solvent model at temperatures below and above the rotational freezing point. <i>Journal of Chemical Physics</i> , 2006, 124, 124508.	1.2	5
82	Stochastic Liouville, Langevin, Fokker-Planck, and Master Equation Approaches to Quantum Dissipative Systems. <i>Journal of the Physical Society of Japan</i> , 2006, 75, 082001.	0.7	720
83	Multidimensional vibrational spectroscopy for tunneling processes in a dissipative environment. <i>Journal of Chemical Physics</i> , 2005, 123, 014503.	1.2	35
84	Multidimensional infrared spectroscopy for molecular vibrational modes with dipolar interactions, anharmonicity, and nonlinearity of dipole moments and polarizability. <i>Journal of Chemical Physics</i> , 2005, 123, 224310.	1.2	11
85	Quantum Dynamics of System Strongly Coupled to Low-Temperature Colored Noise Bath: Reduced Hierarchy Equations Approach. <i>Journal of the Physical Society of Japan</i> , 2005, 74, 3131-3134.	0.7	403
86	Two-dimensional Raman and infrared vibrational spectroscopy for a harmonic oscillator system nonlinearly coupled with a colored noise bath. <i>Journal of Chemical Physics</i> , 2004, 120, 260-271.	1.2	66
87	Energy-Level Diagrams and Their Contribution to Fifth-Order Raman and Second-Order Infrared Responses: A Distinction between Relaxation Models by Two-Dimensional Spectroscopy. <i>Journal of Physical Chemistry A</i> , 2003, 107, 8092-8105.	1.1	21
88	Two-dimensional spectroscopy for a two-dimensional rotator coupled to a Gaussian Markovian noise bath. <i>Journal of Chemical Physics</i> , 2003, 119, 1650-1660.	1.2	13
89	Two-dimensional vibrational spectroscopy of a double minimum system in a dissipative environment. <i>Journal of Chemical Physics</i> , 2003, 119, 2155-2164.	1.2	24
90	The energy landscape for solvent dynamics in electron transfer reactions: A minimalist model. <i>Journal of Chemical Physics</i> , 2002, 117, 2172-2179.	1.2	13

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91	Vibrational spectroscopy of a harmonic oscillator system nonlinearly coupled to a heat bath. Journal of Chemical Physics, 2002, 117, 6221-6234.	1.2	56
92	On single-mode $\hat{A}$ - and V-type micromasers: quantum interference versus photon statistics. Journal of Optics B: Quantum and Semiclassical Optics, 2002, 4, 402-410.	1.4	7
93	Two-Time Correlation Function of a Two-Dimensional Quantal Rotator in a Colored Noise. Journal of the Physical Society of Japan, 2002, 71, 2414-2426.	0.7	5
94	Application of the transcorrelated Hamiltonian to the linearized coupled cluster singles and doubles model. Chemical Physics Letters, 2002, 353, 317-323.	1.2	56
95	Probing a colored-noise induced peak of a strongly damped Brownian system by one- and two-dimensional spectroscopy. Chemical Physics Letters, 2002, 358, 51-56.	1.2	6
96	Nonequilibrium initial conditions of a Brownian oscillator system observed by two-dimensional spectroscopy. Journal of Chemical Physics, 2001, 115, 2267-2281.	1.2	8
97	Biorthogonal approach for explicitly correlated calculations using the transcorrelated Hamiltonian. Journal of Chemical Physics, 2001, 115, 7865-7871.	1.2	58
98	Quantum Theory of a Two-Dimensional Rotator in a Dissipative Environment: Application to Far-Infrared Spectroscopy. Journal of the Physical Society of Japan, 2001, 70, 1167-1170.	0.7	9
99	Two-dimensional spectroscopy and harmonically coupled anharmonic oscillators. Chemical Physics, 2001, 266, 237-250.	0.9	25
100	Multi-dimensional vibrational spectroscopy measured from different phase-matching conditions. Chemical Physics Letters, 2001, 341, 329-337.	1.2	46
101	Cage Dynamics in the Third-Order Off-Resonant Response of Liquid Molecules: A Theoretical Realization. Bulletin of the Chemical Society of Japan, 2000, 73, 873-884.	2.0	3
102	Two-Dimensional Spectroscopy for Harmonic Vibrational Modes with Nonlinear System-Bath Interactions. I. Gaussian-White Case. Journal of the Physical Society of Japan, 2000, 69, 3115-3132.	0.7	52
103	Two-Dimensional Spectroscopy for Harmonic Vibrational Modes with Nonlinear System-Bath Interactions. II. Gaussian-Markovian Case. Journal of the Physical Society of Japan, 2000, 69, 4095-4106.	0.7	66
104	Electrochemistry, 2000, 68, 125-129.	0.6	0
105	Optimized perturbation approach with a Legendre transformation to a dissipative system: Correlation functions of a Morse oscillator. Physical Review E, 1999, 59, 1475-1488.	0.8	7
106	Structures and electronic phases of the bis(ethylenedithio)tetrathiafulvalene (BEDT-TTF) clusters and $\text{I}^{\ominus}$ -(BEDT-TTF) salts: A theoretical study based on ab initio molecular orbital methods. Journal of Chemical Physics, 1999, 111, 5986-5994.	1.2	31
107	Structural information from two-dimensional fifth-order Raman spectroscopy. Journal of Chemical Physics, 1999, 111, 492-503.	1.2	73
108	Polaron density matrix and effective mass at finite temperature. Physical Review B, 1999, 60, 7245-7251.	1.1	4

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109	Femtochemistry. Proceedings of the National Academy of Sciences of the United States of America, 1999, 96, 8823-8824.	3.3	4
110	A thermal bath induced new resonance on the linear and nonlinear spectra of a two-level system. Chemical Physics, 1999, 242, 367-385.	0.9	3
111	Two-dimensional line-shape analysis of photon-echo signal. Chemical Physics Letters, 1999, 314, 488-495.	1.2	56
112	Ab initio MO studies of DCNQI molecules. Synthetic Metals, 1999, 103, 2099-2100.	2.1	1
113	Ab Initio MO Studies on Electronic States of DCNQI Molecules. Journal of Physical Chemistry B, 1999, 103, 266-270.	1.2	3
114	Absorption line shape of impurity molecule driven by a fractal noise. Chemical Physics Letters, 1998, 289, 97-104.	1.2	10
115	Pump-probe spectra and nuclear dynamics for a dissipative molecular system in a strong laser field: predissociation dynamics. Chemical Physics Letters, 1998, 292, 28-34.	1.2	19
116	Theoretical study on electron correlation of 1-D (DCNQI) <sub>2</sub> M (M=Li, Ag) salts. Chemical Physics Letters, 1998, 298, 15-20.	1.2	8
117	Two-dimensional THz spectroscopy of liquids: non-linear vibrational response to a series of THz laser pulses. Chemical Physics Letters, 1998, 295, 298-304.	1.2	36
118	Fifth-order two-dimensional vibrational spectroscopy of a Morse potential system in condensed phases. Chemical Physics, 1998, 233, 217-229.	0.9	55
119	Coherent two-dimensional Raman scattering: Frequency-domain measurement of the intra- and intermolecular vibrational interactions. Journal of Chemical Physics, 1998, 108, 1326-1334.	1.2	71
120	Spectral random walks and line broadening of impurity molecules in an Ising spin glass environment. Journal of Chemical Physics, 1998, 108, 1851-1858.	1.2	16
121	The (2n+1)th-order off-resonant spectroscopy from the (n+1)th-order anharmonicities of molecular vibrational modes in the condensed phase. Journal of Chemical Physics, 1997, 106, 1687-1698.	1.2	100
122	Gaussian- $\epsilon$ Markovian quantum Fokker- $\epsilon$ Planck approach to nonlinear spectroscopy of a displaced Morse potentials system: Dissociation, predissociation, and optical Stark effects. Journal of Chemical Physics, 1997, 107, 1779-1793.	1.2	71
123	Two-time correlation functions of a harmonic system nonbilinearly coupled to a heat bath: Spontaneous Raman spectroscopy. Physical Review E, 1997, 56, 2747-2750.	0.8	45
124	First-, third-, and fifth-order resonant spectroscopy of an anharmonic displaced oscillators system in the condensed phase. Journal of Chemical Physics, 1997, 106, 2078-2095.	1.2	54
125	Femtosecond two-dimensional spectroscopy from anharmonic vibrational modes of molecules in the condensed phase. Journal of Chemical Physics, 1997, 107, 2267-2283.	1.2	140
126	Interplay of inhomogeneity and anharmonicity in 2D Raman spectroscopy of liquids. Chemical Physics Letters, 1997, 277, 159-166.	1.2	39



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127	Sensitivity of two-dimensional fifth-order Raman response to the mechanism of vibrational mode-mode coupling in liquid molecules. <i>Chemical Physics Letters</i> , 1997, 278, 175-183.	1.2	57
128	Ab initio molecular orbital calculations by the resonating Hartree-Fock approach: superposition of non-orthogonal Slater determinants. <i>Chemical Physics Letters</i> , 1996, 263, 687-690.	1.2	20
129	Unified time-path approach to the generating functional of the Brownian oscillator system: The bilinearly corrected Feynman rule for nonequilibrium processes. <i>Physical Review E</i> , 1996, 53, 214-227.	0.8	16
130	Unified time-path approach to the effect of anharmonicity on the molecular vibrational spectroscopy in solution. <i>Journal of Chemical Physics</i> , 1996, 105, 7294-7309.	1.2	31
131	Femtosecond pump-probe spectroscopy of intermolecular vibrations in molecular dimers. <i>Journal of Chemical Physics</i> , 1995, 103, 1981-1984.	1.2	14
132	Multistate quantum Fokker-Planck approach to nonadiabatic wave packet dynamics in pump-probe spectroscopy. <i>Journal of Chemical Physics</i> , 1994, 101, 3049-3061.	1.2	93
133	Femtosecond Two-Dimensional Raman Spectroscopy of Liquid Water. <i>The Journal of Physical Chemistry</i> , 1994, 98, 12466-12470.	2.9	64
134	Quantum Brownian Oscillator Analysis of Pump-Probe Spectroscopy in the Condensed Phase. , 1994, , 327-343.		1
135	Optical Stark Spectroscopy of a Brownian Oscillator in Intense Fields. <i>Journal of the Physical Society of Japan</i> , 1994, 63, 66-77.	0.7	85
136	Temperature dependence and non-Condon effects in pump-probe spectroscopy in the condensed phase. <i>Journal of the Optical Society of America B: Optical Physics</i> , 1993, 10, 2263.	0.9	36
137	Description of nonlinear optical response using phase space wave packets. <i>The Journal of Physical Chemistry</i> , 1993, 97, 12596-12601.	2.9	14
138	Real-time path-integral approach to quantum coherence and dephasing in nonadiabatic transitions and nonlinear optical response. <i>Physical Review E</i> , 1993, 47, 118-136.	0.8	119
139	Two-dimensional femtosecond vibrational spectroscopy of liquids. <i>Journal of Chemical Physics</i> , 1993, 99, 9496-9511.	1.2	559
140	The interplay of tunneling, resonance, and dissipation in quantum barrier crossing: A numerical study. <i>Journal of Chemical Physics</i> , 1992, 96, 8485-8496.	1.2	102
141	Quantum and classical Fokker-Planck equations for a Gaussian-Markovian noise bath. <i>Physical Review A</i> , 1991, 43, 4131-4142.	1.0	189
142	Nonperturbative expansion method for a quantum system coupled to a harmonic-oscillator bath. <i>Physical Review A</i> , 1990, 41, 6676-6687.	1.0	256
143	Time Evolution of a Quantum System in Contact with a Nearly Gaussian-Markoffian Noise Bath. <i>Journal of the Physical Society of Japan</i> , 1989, 58, 101-114.	0.7	706
144	Time-Dependent Spectrum of a Two-Level System Coupled to a Heat Bath Driven by Pulsed Laser. <i>Journal of the Physical Society of Japan</i> , 1989, 58, 3001-3012.	0.7	5

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145	Second Order Optical Process of a Three-Level System in Contact with a Nearly Gaussian-Markoffian Noise Bath. Journal of the Physical Society of Japan, 1989, 58, 1850-1859.	0.7	13
146	Two-Time Correlation Functions of a System Coupled to a Heat Bath with a Gaussian-Markoffian Interaction. Journal of the Physical Society of Japan, 1989, 58, 1199-1206.	0.7	58
147	Second Order Optical Process of A Randomly Modulated Multi-Level Atom. Journal of the Physical Society of Japan, 1986, 55, 4550-4565.	0.7	7