Pablo Villarreal

List of Publications by Year
in descending order

233	3,748	134610	263392
papers	citations	34 h-index	45 g-index

235	235	235	1243
all docs	docs citations	times ranked	citing authors

#	Article	IF	Citations
1	Helium structures around SF₅⁺ and SF₆⁺: novel intermolecular potential and mass spectrometry experiments. Physical Chemistry Chemical Physics, 2022, 24, 2004-2014.	1.3	5
2	A Path Integral Molecular Dynamics Simulation of a Harpoon-Type Redox Reaction in a Helium Nanodroplet. Molecules, 2021, 26, 5783.	1.7	8
3	Modelling interactions of alkali-cation dimers in He clusters. Journal of Physics: Conference Series, 2020, 1412, 202033.	0.3	0
4	Rotational-state-changing collisions between <mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML"><mml:msubsup><mml:mi mathvariant="normal">N</mml:mi><mml:mn>2</mml:mn><mml:mo>+</mml:mo></mml:msubsup></mml:math> and Rb at low energies. Physical Review A, 2020, 101, .	1.0	7
5	Theoretical Study of Cationic Alkali Dimers Interacting with He: Li₂⁺â€“He and Na₂⁺â€“He van der Waals Complexes. Journal of Physical Chemistry A, 2019, 123, 7814-7821.	1.1	9
6	A combined experimental and theoretical investigation of Cs+ ions solvated in HeN clusters. Journal of Chemical Physics, 2019, 150, 154304.	1.2	17
7	N2+(2Î£g) and Rb(2S) in a hybrid trap: modeling ion losses from radiative association paths. Physical Chemistry Chemical Physics, 2019, 21, 8342-8351.	1.3	3
8	Limitations of a Theoretical Method to Calculate the Rovibrational Spectrum of Trimers: H $$^+_3$$ 3 +. Few-Body Systems, 2018, 59, 1.	0.7	3
9	Theoretical methods for the rotationâ€“vibration spectra of triatomic molecules: distributed Gaussian functions compared with hyperspherical coordinates. International Reviews in Physical Chemistry, 2018, 37, 329-361.	0.9	4
10	Symmetry analysis of trimers rovibrational spectra: the case of Ne3. European Physical Journal D, 2018, 72, 1.	0.6	3
11	Formation of rubidium dimers on the surface of helium clusters: a first step through quantum molecular dynamics simulations. European Physical Journal D, 2018, 72, 1.	0.6	22
12	Atomic cluster collisions: ISACC-2015 (7th International Symposium)*. European Physical Journal D, 2017, 71, 1.	0.6	3
13	Comparative investigation of pure and mixed rare gas atoms on coronene molecules. Journal of Chemical Physics, 2017, 146, 034302.	1.2	16
14	Adsorption of molecular hydrogen on coronene with a new potential energy surface. Physical Chemistry Chemical Physics, 2017, 19, 26358-26368.	1.3	22
15	Rb $$_2$$ ($$^3Sigma ^+_u$$) Dimer Interacting with He Atoms: Quantum Structures of Small Clusters and Reactive Scattering Calculations Rb + RbHe $$ightarrow $$ Rb $$_2$$ ($$^3Sigma) Tj ETQq1 1 0.784314 rgBT /Overlo 2017, , 97-118.	0.2	0
16	Examination of the Feynmanâ€“Hibbs Approach in the Study of Ne_N-Coronene Clusters at Low Temperatures. Journal of Physical Chemistry A, 2016, 120, 5370-5379.	1.1	16
17	Path integral Monte Carlo investigations on doped helium clusters. International Reviews in Physical Chemistry, 2016, 35, 37-68.	0.9	22
18	Coronene molecules in helium clusters: Quantum and classical studies of energies and configurations. Journal of Chemical Physics, 2015, 143, 224306.	1.2	28

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19	A full-configuration-interaction nuclear orbital approach and application for small doped He clusters. , 2015, , .		0
20	Structural changes in water and Ar-water clusters under high pressure. Journal of Physics: Conference Series, 2015, 635, 032008.	0.3	0
21	Computational investigations of the thermodynamic properties of size-selected water and Arâ€“water clusters: high-pressure transitions. Physical Chemistry Chemical Physics, 2015, 17, 8792-8801.	1.3	22
22	A configurational study of helium clusters doped with Heâˆ—âˆ’ and He2âˆ—âˆ’. Journal of Chemical Physics, 2015, 142, 104303.	1.2	8
23	Quantum Features of Anionic Species He^{*â€“} and He₂^{*â€“} in Small He_N Clusters. Journal of Physical Chemistry A, 2015, 119, 11574-11582.	1.1	5
24	Variational and Path Integral Monte Carlo calculations on Helium Clusters Doped with Metastable Anions He*- and He2*-. Journal of Physics: Conference Series, 2015, 635, 072009.	0.3	0
25	Potential energy surface and bound states of the (X⁴Î£)KRbâ€K complex. International Journal of Quantum Chemistry, 2015, 115, 19-27.	1.0	1
26	Reactive scattering calculations for 87Rb+87RbHeâ†'Rb2(3Î£u+,v)+He from ultralow to intermediate energies. Journal of Chemical Physics, 2015, 142, 164304.	1.2	8
27	Path integral Monte Carlo calculations of calciumâ€doped ⁴He clusters. International Journal of Quantum Chemistry, 2014, 114, 1318-1326.	1.0	4
28	Quantum Molecular Processes Induced by Electrons, Positrons, Atoms, and Ions: From eV to Nano-eV. Journal of Physical Chemistry A, 2014, 118, 6299-6300.	1.1	0
29	Theoretical simulations of the vibrational predissociation spectra of $${m{H}} ^{+}_{5}$$ and $${m{D}}^{+}_{5}$$ clusters. Highlights in Theoretical Chemistry, 2014, , 125-131.	0.0	0
30	Quantum rotation of Rb2 (3 Î£ u +) attached to HeN droplets: a path-integral Monte Carlo study. European Physical Journal D, 2013, 67, 1.	0.6	10
31	Unravelling Coriolis temperature-dependent effects on doped helium clusters: Vib-rotational Raman spectra of (3,4He)4â€“Cl2(X). Chemical Physics Letters, 2013, 555, 12-18.	1.2	3
32	Solvent states and spectroscopy of doped helium clusters as a quantum-chemistry-like problem. Physical Chemistry Chemical Physics, 2013, 15, 10126.	1.3	11
33	Including nuclear quantum effects into highly correlated electronic structure calculations of weakly bound systems. Journal of Chemical Physics, 2013, 138, 184113.	1.2	14
34	Weakly bound finite systems: (4He)Nâ€“Rb2(3Î£u), clustering structures from a quantum Monte Carlo approach. Journal of Physics Condensed Matter, 2012, 24, 104014.	0.7	4
35	Binding energies and structures of Ca-He<mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:msub><mml:mrow /><mml:mn>2</mml:mn></mml:msub></mml:math>weakly bound triatomic complexes. Physical Review A, 2012, 86, .	1.0	9
36	Comment on â€œWeakly bound states of the He-He-Ca triatomic systemâ€• Physical Review A, 2012, 86, .	1.0	2

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37	Simulation of the infrared predissociation spectra of H5+. Physical Review A, 2012, 85, .	1.0	16
38	Direct Path-Integral Monte-Carlo simulations of H+5/D+5clusters: thermal equilibrium state properties. Journal of Physics: Conference Series, 2012, 388, 122001.	0.3	0
39	A quantum chemistry approach to energies, structures, and spectroscopy of doped helium clusters. , 2012, , .		0
40	Theoretical Investigation of the He₄Br₂ Conformers. Journal of Physical Chemistry A, 2012, 116, 7169-7176.	1.1	4
41	Spin-Polarized Rb₂ Interacting with Bosonic He Atoms: Potential Energy Surface and Quantum Structures of Small Clusters. Journal of Physical Chemistry A, 2012, 116, 2394-2404.	1.1	19
42	Helium aggregates doped with alkali dimer impurities: A finite temperature study of complexes. Computational and Theoretical Chemistry, 2012, 990, 106-111.	1.1	10
43	Binding weakly interacting partners: a study of Caâ€“He2 and its isotopomers. European Physical Journal D, 2012, 66, 1.	0.6	8
44	HeI₂ interaction potential based on an interpolation scheme. International Journal of Quantum Chemistry, 2012, 112, 2971-2975.	1.0	13
45	Theoretical simulations of the vibrational predissociation spectra of H 5 + and D 5 + clusters. Theoretical Chemistry Accounts, 2012, 131, 1.	0.5	10
46	New trends in atomic and molecular clusters (in honour of Gerardo Delgado-Barrio). Chemical Physics, 2012, 399, 1-3.	0.9	0
47	Path integral Monte Carlo studies of the H₅⁺/D₅⁺clusters usingab initiopotential surfaces. Physica Scripta, 2011, 84, 028109.	1.2	16
48	Quantum Features of a Barely Bound Molecular Dopant: Cs₂(³Î£_u) in Bosonic Helium Droplets of Variable Size. Journal of Physical Chemistry A, 2011, 115, 6892-6902.	1.1	19
49	Internal Proton Transfer and H₂ Rotations in the H₅⁺ Cluster: A Marked Influence on Its Thermal Equilibrium State. Journal of Physical Chemistry A, 2011, 115, 2483-2488.	1.1	30
50	Ab initio characterization of the Neâ€“I2 van der Waals complex: Intermolecular potentials and vibrational bound states. Journal of Chemical Physics, 2011, 134, 214304.	1.2	24
51	Molecular dynamics simulations of rigid and flexible water models: Temperature dependence of viscosity. Chemical Physics, 2011, 388, 9-18.	0.9	54
52	Microscopic description of small doped ³He clusters through the fullâ€configurationâ€interaction nuclear orbital approach: The (³He)_Nâ€Br₂(X) case revisited. International Journal of Quantum Chemistry, 2011, 111, 406-415.	1.0	13
53	Global potential energy surface for the ground electronic state of H: A DFT approach. International Journal of Quantum Chemistry, 2011, 111, 368-374.	1.0	12
54	Filtered stress autocorrelation functions of liquid water models. International Journal of Quantum Chemistry, 2011, 111, 375-386.	1.0	4

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55	Temperature dependence of the energetics and structure for the Ar dimer and trimer. International Journal of Quantum Chemistry, 2011, 111, 472-479.	1.0	1
56	Diffusion Monte Carlo description of Cs₂(³Î£_u)â€“(⁴He)_Nclusters: an example of weak dopantâ€“helium interaction. Physica Scripta, 2011, 84, 028107.	1.2	6
57	Theoretical characterization of intermolecular vibrational states through the multi-configuration time dependent Hartree approach: The He2.3ICl clusters. Journal of Chemical Physics, 2011, 135, 244309.	1.2	9
58	Infrared spectrum of H<mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:mrow><mml:msub><mml:mrow /><mml:mn>5</mml:mn></mml:msub><mml:msup><mml:mrow /><mml:mo>+</mml:mo></mml:msup></mml:mrow></mml:math>and D<mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:mrow><mml:msub><mml:mrow /><mml:mn>5</mml:mn></mml:msub><mml:msup	1.0	19
59	Frequency domain description of Kohlrausch response through a pair of Havriliak-Negami-type functions: An analysis of functional proximity. Physical Review E, 2011, 84, 066703.	0.8	13
60	Full-dimensional multi configuration time dependent Hartree calculations of the ground and vibrationally excited states of He2,3Br2 clusters. Journal of Chemical Physics, 2011, 135, 054303.	1.2	9
61	Quantum solvent states and rovibrational spectra of small doped H3e clusters through the full-configuration-interaction nuclear orbital approach: The (H3e)Nâ€“Cl2(X) case (Nâ‰¤4). Journal of Chemical Physics, 2010, 132, 194313.	1.2	14
62	A path-integral Monte Carlo study of a small cluster: The Ar trimer. Journal of Chemical Physics, 2010, 132, 244303.	1.2	15
63	A new accurate and full dimensional potential energy surface of H5+ based on a triatomics-in-molecules analytic functional form. Journal of Chemical Physics, 2010, 133, 024306.	1.2	53
64	Toward a realistic density functional theory potential energy surface for the H5+ cluster. Journal of Chemical Physics, 2010, 133, 054303.	1.2	24
65	Binding He atoms to hydrogen moieties: quantum features from ultraweak interactions. Molecular Physics, 2010, 108, 57-72.	0.8	2
66	Structuring a Quantum Solvent around a Weakly Bound Dopant: The Heâ‹¯Cs₂(³Î£_u) Complex. Journal of Physical Chemistry A, 2009, 113, 14718-14729.	1.1	23
67	An optimized full-configuration-interaction nuclear orbital approach to a â€œhard-coreâ€•interaction problem: Application to (H3e)Nâ€“Cl2(B) clusters (Nâ‰¤4). Journal of Chemical Physics, 2009, 131, 194101.	1.2	23
68	A Study of the Ar3 System at Low Temperature. Few-Body Systems, 2009, 45, 237-239.	0.7	3
69	Using a Jacobiâ€“Davidson â€œNuclear Orbitalâ€•Method for Small Doped 3 He Clusters. Few-Body Systems, 2009, 45, 233-236.	0.7	15
70	Intermolecular Ab Initio Potential and Spectroscopy of the Ground State of HeI₂ Complex Revisited. Journal of Physical Chemistry A, 2009, 113, 5754-5762.	1.1	40
71	A density-division embedding potential inversion technique. Journal of Chemical Physics, 2009, 131, 234110.	1.2	39
72	A theoretical investigation on the spectrum of the Ar trimer for high rotational excitations. Journal of Chemical Physics, 2009, 130, 154301.	1.2	18

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73	HeI[sub 2] Van der Waals Complex: Ab initio Ground and Electronic Excited Potential Surfaces for Studying Dynamics. , 2009, , .		1
74	Ab initiointeraction potentials for X and B excited states of Heâ€“I2for studying dynamics. Journal of Physics: Conference Series, 2009, 194, 122001.	0.3	0
75	Symmetry assignment in the distributed Gaussian functions method to study homonuclear rotating trimers. Chemical Physics Letters, 2008, 460, 417-422.	1.2	9
76	Viscosity of Liquid Water via Equilibrium Molecular Dynamics Simulations. Progress in Theoretical Chemistry and Physics, 2008, , 351-361.	0.2	10
77	Exact, Bornâ€“Oppenheimer, and quantum-chemistry-like calculations in helium clusters doped with light molecules: The He2N2(X) system. Journal of Chemical Physics, 2008, 128, 164313.	1.2	9
78	An inversion technique for the calculation of embedding potentials. Journal of Chemical Physics, 2008, 129, 184104.	1.2	93
79	Ab initio potential energy surface and spectrum of the B(Î3) state of the HeI2 complex. Journal of Chemical Physics, 2007, 126, 204301.	1.2	31
80	Ab initio vibrational predissociation dynamics of Heâ€“I2(B) complex. Journal of Chemical Physics, 2007, 126, 244314.	1.2	31
81	Spectral simulations of polar diatomic molecules immersed in He clusters: application to the ICl (X) molecule. Physica Scripta, 2007, 76, C96-C103.	1.2	14
82	Dynamics and Potential Energy Surfaces for small to medium size Hen-dihalogen clusters. AIP Conference Proceedings, 2007, , .	0.3	0
83	Key role of spinâ€“orbit effects in the relaxation of CO2(010) by thermal collisions with O(3Pj). Molecular Physics, 2007, 105, 1171-1181.	0.8	11
84	Vibrational and rotational bound states in floppy triatomic systems: The distributed Gaussian functions approach. Physics Reports, 2007, 452, 1-32.	10.3	22
85	Exact and quantum chemistry-like calculations in helium doped clusters: The He2Br2(X) example. International Journal of Quantum Chemistry, 2007, 107, 2756-2762.	1.0	13
86	Doped helium clusters analyzed through quantum chemistry methods. International Journal of Quantum Chemistry, 2007, 107, 2902-2921.	1.0	14
87	Additive intermolecular potentials from ab initio calculations: trends in Rg2â€“dihalogen van der Waals trimers. Theoretical Chemistry Accounts, 2007, 118, 511-517.	0.5	14
88	Potential energy surfaces and dynamics of He n Br2 van der Waals complexes. Progress in Theoretical Chemistry and Physics, 2007, , 193-202.	0.2	0
89	A theoretical study of He2ICl van der Waals cluster. Journal of Chemical Physics, 2006, 125, 014313.	1.2	34
90	Rovibrational Structures in Floppy Triatomics:Â Distributed Gaussian Functions Treatment for the Ne2H-Systemâ€. Journal of Physical Chemistry A, 2006, 110, 5487-5494.	1.1	8

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91	An overview on potential energy surfaces of rare-gas dihalogen van der Waals clusters. Physica Scripta, 2006, 73, C57-C63.	1.2	13
92	A full-configuration interaction â€œnuclear orbitalâ€•method to study doped HeN3 clusters (Nâ©½4). Journal of Chemical Physics, 2006, 125, 221101.	1.2	26
93	Vibrational quenching of CO2(010) by collisions with O(P3) at thermal energies: A quantum-mechanical study. Journal of Chemical Physics, 2006, 124, 164302.	1.2	15
94	Polar di-halogen molecules solvated in bosonic helium clusters: The paradigm ofICl(X). Physical Review A, 2006, 74, .	1.0	33
95	STRUCTURE AND DYNAMICS OF VAN DER WAALS COMPLEXES: FROM TRIATOMIC TO MEDIUM SIZE CLUSTERS. , 2006, , .		0
96	ABINITIO VAN DERWAALS POTENTIAL ENERGYSURFACES APPLICATION TO COMPLEXES OF BROMINE MOLECULE WITH HELIUM ATOMS. , 2006, , 347-369.		0
97	Br2(X) Microsolvation in Helium Clusters: Effect of the Interaction on the Quantum Solvent Density Distribution. ChemPhysChem, 2005, 6, 1348-1356.	1.0	19
98	Ab initio calculations, potential representation and vibrational dynamics of He2Br2 van der Waals complex. Journal of Chemical Physics, 2005, 122, 044305.	1.2	36
99	Energies and density distributions of(He4)Nclusters doped withBr2(X): A Hartree-like approach. Physical Review A, 2005, 71, .	1.0	35
100	The binding of He4 and He3 to a hydrogen molecule: A computational study for pH2 and oH2. Journal of Chemical Physics, 2005, 122, 084308.	1.2	20
101	A complete configurational study for the bound states of Ne trimers. Journal of Chemical Physics, 2005, 122, 084313.	1.2	21
102	Bound-state energies in argon trimers via a variational expansion: The effects from many-body corrections. Journal of Chemical Physics, 2005, 122, 144319.	1.2	23
103	Role of Boson-Fermion Statistics on the Raman Spectra ofBr2(X)in Helium Clusters. Physical Review Letters, 2004, 93, 053401.	2.9	46
104	The open-shell interaction of He with the Bâ€Š3Îu(0+) state of Br2: Anab initiostudy and its comparison with a diatomics-in-molecule perturbation model. Journal of Chemical Physics, 2004, 120, 2182-2192.	1.2	20
105	HeBr2complex: ground-state potential and vibrational dynamics fromab initiocalculations. Molecular Physics, 2004, 102, 2277-2283.	0.8	37
106	Raman spectra of (He)N-Br2(X) clusters: The role of boson/fermion statistics in a quantum solvent. Journal of Chemical Physics, 2004, 121, 2975-2984.	1.2	25
107	Three-Dimensional ab Initio Potential and Ground State Dynamics of the HeI2Complex. Journal of Physical Chemistry A, 2004, 108, 6065-6071.	1.1	36
108	Structures and Energetics of Hn+ Clusters (n = 5â€”11). ChemInform, 2003, 34, no.	0.1	1

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109	CCSD(T) potential energy surface and bound rovibrational level calculations for the Arâ€“ICl(X) complex. Chemical Physics Letters, 2003, 375, 328-336.	1.2	31
110	Structure and bonding of ArClF: Intermolecular potential surface. Israel Journal of Chemistry, 2003, 43, 279-286.	1.0	9
111	Structures and Energetics of Clusters (n = 5â˜11). Journal of Physical Chemistry A, 2003, 107, 4768-4772.	1.1	42
112	CCSD(T) intermolecular potential between He atom and ClF molecule: Comparison with experiment. Journal of Chemical Physics, 2003, 119, 4216-4222.	1.2	36
113	The 4He Trimer: Structure and Energetics of a Very Unusual Molecule. Collection of Czechoslovak Chemical Communications, 2003, 68, 1-22.	1.0	6
114	Vibrational predissociation of NeBr2 (X,â€‰v=1) using anab initiopotential energy surface. Journal of Chemical Physics, 2002, 117, 10019-10025.	1.2	26
115	He79Br2 B,v=8â†X,vâ€³=0 excitation spectrum: Ab initio prediction and spectroscopic manifestation of a linear isomer. Journal of Chemical Physics, 2002, 117, 6117-6120.	1.2	37
116	The van der Waals potential energy surfaces and structures of Heâ€“ICl and Neâ€“ICl clusters. Journal of Chemical Physics, 2002, 117, 7017-7023.	1.2	51
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127	Effective resolvent applied to interacting resonances. European Physical Journal D, 2001, 15, 215-219.	0.6	5
128	Rotationâ€“vibration interaction in 4He trimers. Chemical Physics Letters, 2001, 335, 105-110.	1.2	10
129	The structure of a weakly bound ionic trimer: Calculations for the 4He2Hâ¯ complex. Journal of Chemical Physics, 2001, 114, 5520-5530.	1.2	18
130	Photodissociation of NeBr2(B) below and above the dissociation limit of Br2(B). Journal of Chemical Physics, 2001, 115, 2566-2575.	1.2	23
131	GonzÃ¡lez-Lezanaet al.Reply:. Physical Review Letters, 2001, 86, 4190-4190.	2.9	4
132	Application of a time-dependent Hartree approach on several surfaces to the vibrational predissociation of Ne₂I₂. Molecular Physics, 2000, 98, 1783-1791.	0.8	3
133	Blueshifts of the Bâ†X excitation spectra of He79Br2 using a DIM-based potential. Chemical Physics Letters, 2000, 318, 578-584.	1.2	14
134	Searching for Efimov states in triatomic systems: The case of LiHe 2. Europhysics Letters, 2000, 50, 567-573.	0.7	30
135	Theoretical simulations of the He79Br2â€ŠB, v=8â†X, vâ€³=0 excitation spectrum: Spectroscopic manifestation of a linear isomer?. Journal of Chemical Physics, 2000, 113, 4620-4628.	1.2	25
136	The weakly bound ground state of the LiHe2 triatomic system. Physical Chemistry Chemical Physics, 2000, 2, 4067-4073.	1.3	41
137	Diffusion Monte-Carlo Calculations of Quasi-Bound States of Rare Gas-Halogen Clusters: a Diabatic Approach. , 2000, , 93-102.		0
138	Comparative configurational study for He, Ne, and Ar trimers. Journal of Chemical Physics, 1999, 110, 9000-9010.	1.2	70
139	Efimov States forH4eTrimers?. Physical Review Letters, 1999, 82, 1648-1651.	2.9	69
140	Diabatic diffusion Monte Carlo calculations of the energy and structure of the Cl 2 He n (n =1â€“10) clusters. Computational and Theoretical Chemistry, 1999, 493, 125-132.	1.5	6
141	A combined experimental-theoretical study of the vibrational predissociation and product rotational distributions for high vibrational levels of He79Br2. Journal of Chemical Physics, 1999, 110, 256-266.	1.2	35
142	Quantum-mechanical study of the resonances of the SN2 reaction Cl-+CH3Clâ†'ClCH3+Cl-. Physical Chemistry Chemical Physics, 1999, 1, 1197-1203.	1.3	35
143	An adiabatic model for rare gasâ€“halogen van der Waals complexes: application to HeBr2(B).. Computational and Theoretical Chemistry, 1998, 433, 107-111.	1.5	4
144	Decision support systems generator for industrial companies Module IV: Forecasting support system. Computers and Industrial Engineering, 1998, 35, 315-318.	3.4	3

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145	Theoretical study of VRT energy transfer in Ne+I2(B) collisions using a spectroscopic interaction potential. Journal of the Chemical Society, Faraday Transactions, 1998, 94, 2307-2313.	1.7	6
146	Potential energy surface and spectroscopy of clusters of rare-gas atoms with cyclopropane. Journal of Chemical Physics, 1998, 109, 9288-9299.	1.2	5
147	A hybrid classical/quantum approach to cluster fragmentation dynamics: Application to the vibrational predissociation of He2Cl2. Journal of Chemical Physics, 1998, 108, 1989-1996.	1.2	17
148	Vibronic shifts for LiH in X and A states. Journal of Physics B: Atomic, Molecular and Optical Physics, 1997, 30, L427-L434.	0.6	16
149	Strongly mixed resonances in the photofragmentation of HeBr2 near Br2(B) dissociation: Stabilization and close-coupling studies. Journal of Chemical Physics, 1997, 106, 3216-3226.	1.2	21
150	Nonradiative lifetimes for LiH in the A state using adiabatic and diabatic schemes. Journal of Chemical Physics, 1997, 107, 10515-10522.	1.2	52
151	Theoretical spectroscopy and dynamics of fragmentation of the He79Br2 complex. , 1997, , .		1
152	Collisional heating in ionic argon clusters The Ar+Ar2+ case. Journal of the Chemical Society, Faraday Transactions, 1997, 93, 901-907.	1.7	1
153	Mode excitation dynamics in the fragmentation of Ar3+: An helicity decoupling study. Journal of Chemical Physics, 1997, 106, 1718-1728.	1.2	6
154	Title is missing!. Molecular Engineering, 1997, 7, 219-230.	0.2	5
155	A full quantum study of the vibrational predissociation mechanisms in Ar3+ cluster. Chemical Physics, 1997, 218, 71-81.	0.9	6
156	Half- and full-collision VT energy transfer in the Heî—¸Br2(B) system. Chemical Physics Letters, 1997, 269, 448-454.	1.2	8
157	Sampling the Initial Conditions for Quasiclassical Trajectory Studies of Vibrational Predissociation Dynamics. Topics in Molecular Organization and Engineering, 1997, , 219-230.	0.1	0
158	Computed energy curves for modelling the dissociation of helium trimer ions. Chemical Physics Letters, 1996, 259, 641-646.	1.2	6
159	Quasiclassical dynamics of the I2â€“Ne2vibrational predissociation: A comparison with experiment. Journal of Chemical Physics, 1996, 104, 8405-8412.	1.2	46
160	Vibrational predissociation dynamics of the He79Br2van der Waals molecule: A quantum mechanical study. Journal of Chemical Physics, 1996, 105, 7454-7463.	1.2	46
161	Fragmentation of Ar3+ clusters via vibrational predissociation. Chemical Physics Letters, 1995, 242, 336-342.	1.2	19
162	Vibrational predissociation of I2î—¸Ne. A quasiclassical dynamical study. Chemical Physics Letters, 1995, 243, 236-242.	1.2	32

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163	Time-dependent Hartree study of lifetimes for the Ne2I2 van der Waals cluster. Chemical Physics Letters, 1995, 246, 197-203.	1.2	21
164	Wave packet study of the Ne2I2 fragmentation dynamics: a four degrees of freedom model. Chemical Physics Letters, 1995, 246, 187-196.	1.2	27
165	Fragmentation dynamics of ionized argon clusters: an effective potential model. Zeitschrift FÃ¼r Physik D-Atoms Molecules and Clusters, 1995, 35, 115-124.	1.0	13
166	Vector Correlations in the Photopredissociation of van der Waals Molecules. The Journal of Physical Chemistry, 1994, 98, 3307-3316.	2.9	12
167	Energy levels and structure of tetraâ€atomic van der Waals clusters. Journal of Chemical Physics, 1994, 101, 2217-2230.	1.2	64
168	Competitive internal transfers in metastable decay of cluster ions. Journal of Chemical Physics, 1994, 100, 6472-6486.	1.2	12
169	Metastable decay of rare gas cluster ions ? the (Ar3)+ system. Zeitschrift FÃ¼r Physik D-Atoms Molecules and Clusters, 1994, 30, 77-77.	1.0	2
170	Theoretical Spectroscopy and Dynamics of Tetra-Atomic Van Der Waals Clusters. Jerusalem Symposia on Quantum Chemistry and Biochemistry, 1994, , 57-72.	0.2	1
171	Metastable decay of rare gas cluster ions ? the (Ar3)+ system. Zeitschrift FÃ¼r Physik D-Atoms Molecules and Clusters, 1993, 27, 357-364.	1.0	4
172	Selective adsorption resonances at rainbow conditions in the scattering of atoms by stepped surfaces: application to the 4He/Cu(117) system. Surface Science, 1993, 290, L693-L698.	0.8	12
173	Coherence effects between intramolecular vibrational relaxation and dissociation in triatomic van der Waals systems. Journal of Chemical Physics, 1993, 99, 1035-1049.	1.2	36
174	Study of the selective adsorption phenomenon in the 4He/Cu(11Î±) (with Î± = 0, 3, 5, 7) elastic scattering: the critical kinematic effect. Surface Science, 1992, 274, 21-34.	0.8	30
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