

# Jeremy M Hutson

## List of Publications by Citations

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231  
papers

10,353  
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55  
h-index

89  
g-index

241  
ext. papers

11,245  
ext. citations

4.4  
avg, IF

6.43  
L-index

#	Paper	IF	Citations
231	Theoretical studies of van der Waals molecules and intermolecular forces. <i>Chemical Reviews</i> , <b>1988</b> , 88, 963-988	68.1	511
230	Ultracold dense samples of dipolar RbCs molecules in the rovibrational and hyperfine ground state. <i>Physical Review Letters</i> , <b>2014</b> , 113, 205301	7.4	331
229	Creation of ultracold $^{87}\text{Rb}^{133}\text{Cs}$ molecules in the rovibrational ground state. <i>Physical Review Letters</i> , <b>2014</b> , 113, 255301	7.4	283
228	An ultracold high-density sample of rovibronic ground-state molecules in an optical lattice. <i>Nature Physics</i> , <b>2010</b> , 6, 265-270	16.2	271
227	Vibrational dependence of the anisotropic intermolecular potential of $\text{ArHF}$ . <i>Journal of Chemical Physics</i> , <b>1992</b> , 96, 6752-6767	3.9	265
226	The intermolecular potential of $\text{ArHCl}$ : Determination from high-resolution spectroscopy. <i>Journal of Chemical Physics</i> , <b>1988</b> , 89, 4550-4557	3.9	215
225	Improved potential energy surfaces for the interaction of $\text{H}_2$ with Ar, Kr, and Xe. <i>Journal of Chemical Physics</i> , <b>1987</b> , 86, 837-853	3.9	195
224	Reactions of ultracold alkali-metal dimers. <i>Physical Review A</i> , <b>2010</b> , 81,	2.6	172
223	Anisotropic intermolecular forces. <i>Molecular Physics</i> , <b>1982</b> , 45, 769-790	1.7	170
222	Vibrational dependence of the anisotropic intermolecular potential of argon-hydrogen chloride. <i>The Journal of Physical Chemistry</i> , <b>1992</b> , 96, 4237-4247		163
221	Universality of the three-body parameter for Efimov states in ultracold cesium. <i>Physical Review Letters</i> , <b>2011</b> , 107, 120401	7.4	153
220	The dynamics of open-shell Van der Waals complexes. <i>Journal of Chemical Physics</i> , <b>1991</b> , 94, 7602-7618	3.9	150
219	Precise characterization of $6\text{Li}$ Feshbach resonances using trap-sideband-resolved RF spectroscopy of weakly bound molecules. <i>Physical Review Letters</i> , <b>2013</b> , 110, 135301	7.4	149
218	Centrifugal distortion constants for diatomic molecules: an improved computational method. <i>Journal of Physics B: Atomic and Molecular Physics</i> , <b>1981</b> , 14, 851-857		129
217	Coupled channel methods for solving the bound-state Schrödinger equation. <i>Computer Physics Communications</i> , <b>1994</b> , 84, 1-18	4.2	123
216	Quantum dynamics of ultracold $\text{Na} + \text{Na}_2$ collisions. <i>Physical Review Letters</i> , <b>2002</b> , 89, 153201	7.4	120
215	Morphing ab initio potentials: A systematic study of $\text{NeHF}$ . <i>Journal of Chemical Physics</i> , <b>1999</b> , 110, 8338-8347	3.9	120

214	Towards the production of ultracold ground-state RbCs molecules: Feshbach resonances, weakly bound states, and the coupled-channel model. <i>Physical Review A</i> , <b>2012</b> , 85,	2.6	115
213	Rotational predissociation of the Ar <sup>+</sup> HCl van der Waals complex: Close-coupled scattering calculations. <i>Journal of Chemical Physics</i> , <b>1983</b> , 78, 4025-4039	3.9	115
212	The intermolecular potential energy surface of Ar-HCl. <i>Molecular Physics</i> , <b>1981</b> , 43, 493-516	1.7	108
211	Molecule formation in ultracold atomic gases. <i>International Reviews in Physical Chemistry</i> , <b>2006</b> , 25, 497-526		107
210	Anisotropic intermolecular forces. <i>Molecular Physics</i> , <b>1982</b> , 45, 791-805	1.7	104
209	Atom-symmetric top van der Waals complexes: Angular momentum coupling in Ar-H <sub>2</sub> O. <i>Journal of Chemical Physics</i> , <b>1990</b> , 92, 157-168	3.9	102
208	Observation of the second triatomic resonance in Efimov's scenario. <i>Physical Review Letters</i> , <b>2014</b> , 112, 190401	7.4	100
207	Ultracold RbSr molecules can be formed by magnetoassociation. <i>Physical Review Letters</i> , <b>2010</b> , 105, 153201	7.4	91
206	Ultracold Rb-OH collisions and prospects for sympathetic cooling. <i>Physical Review Letters</i> , <b>2006</b> , 97, 183201	7.4	90
205	Ultracold Li + Li <sub>2</sub> collisions: bosonic and fermionic cases. <i>Physical Review Letters</i> , <b>2005</b> , 94, 033201	7.4	90
204	Nonadditive intermolecular forces from the spectroscopy of van der Waals trimers: Calculations on Ar <sub>2</sub> HCl. <i>Journal of Chemical Physics</i> , <b>1993</b> , 98, 5337-5351	3.9	90
203	Hyperfine energy levels of alkali-metal dimers: Ground-state polar molecules in electric and magnetic fields. <i>Physical Review A</i> , <b>2008</b> , 78,	2.6	85
202	Signatures of large amplitude motion in a weakly bound complex: High-resolution IR spectroscopy and quantum calculations for HeCO <sub>2</sub> . <i>Journal of Chemical Physics</i> , <b>1994</b> , 101, 8351-8363	3.9	85
201	Atom-molecule van der Waals complexes containing open-shell atoms. I. General theory and bending levels. <i>Journal of Chemical Physics</i> , <b>1994</b> , 101, 1939-1958	3.9	85
200	Feshbach resonances in ultracold atomic and molecular collisions: threshold behaviour and suppression of poles in scattering lengths. <i>New Journal of Physics</i> , <b>2007</b> , 9, 152-152	2.9	77
199	Ultracold atom-molecule collisions and bound states in magnetic fields: Tuning zero-energy Feshbach resonances in He-H (B). <i>Physical Review A</i> , <b>2007</b> , 75,	2.6	74
198	The intermolecular potential energy surface for CO <sub>2</sub> -Ar: Fitting to high-resolution spectroscopy of Van der Waals complexes and second virial coefficients. <i>Journal of Chemical Physics</i> , <b>1996</b> , 105, 9130-9140	3.9	72
197	Feshbach resonances, weakly bound molecular states, and coupled-channel potentials for cesium at high magnetic fields. <i>Physical Review A</i> , <b>2013</b> , 87,	2.6	69

- 196 Ultracold collisions involving heteronuclear alkali metal dimers. *Physical Review Letters*, **2005**, 94, 200407.4 69
- 195 Anisotropic intermolecular forces from Hartree-Fock plus damped dispersion (HFD) calculations. *Molecular Physics*, **1984**, 52, 763-781 1.7 68
- 194 Vibrational predissociation and infrared spectrum of the ArHCl van der Waals molecule. *Journal of Chemical Physics*, **1984**, 81, 2357-2362 3.9 68
- 193 Ultracold quantum dynamics: Spin-polarized K+K2 collisions with three identical bosons or fermions. *Physical Review A*, **2005**, 71, 2.6 67
- 192 The ArHF intermolecular potential: Overtone spectroscopy and ab initio calculations. *Journal of Chemical Physics*, **1993**, 99, 9337-9349 3.9 67
- 191 Selective adsorption resonances in the scattering of helium atoms from xenon coated graphite: Close-coupling calculations and potential dependence. *Journal of Chemical Physics*, **1983**, 79, 5179-5187 3.9 67
- 190 Production of ultracold NH molecules by sympathetic cooling with Mg. *Physical Review Letters*, **2009**, 103, 183201 7.4 64
- 189 Molecular collisions in ultracold atomic gases. *International Reviews in Physical Chemistry*, **2007**, 26, 1-28 7 64
- 188 Nonadditive intermolecular forces from the spectroscopy of van der Waals trimers: A theoretical study of Ar<sub>2</sub>-HF. *Physical Review A*, **1995**, 51, 239-250 2.6 64
- 187 Atom-Molecule van der Waals Complexes Containing Open-Shell Atoms. 2. The Bound States of Cl-HCl. *The Journal of Physical Chemistry*, **1994**, 98, 5844-5854 64
- 186 On the long-range and short-range behavior of potentials from reproducing kernel Hilbert space interpolation. *Journal of Chemical Physics*, **2000**, 112, 4415-4416 3.9 63
- 185 Microwave spectroscopy and interaction potential of the long-range He...Ar<sup>+</sup> ion. *Journal of Chemical Physics*, **1995**, 102, 2379-2403 3.9 62
- 184 Magnetically tunable Feshbach resonances in ultracold Li-Yb mixtures. *Physical Review Letters*, **2012**, 108, 043201 7.4 61
- 183 Three-body nonadditive forces between spin-polarized alkali-metal atoms. *Physical Review A*, **2003**, 67, 2.6 61
- 182 The intermolecular potential of NeHCl: Determination from high-resolution spectroscopy. *Journal of Chemical Physics*, **1989**, 91, 4448-4454 3.9 60
- 181 Spectroscopic properties and potential surfaces for atom-diatom van der Waals molecules. *Molecular Physics*, **1980**, 41, 1123-1141 1.7 59
- 180 A new approach to perturbation theory for breakdown of the Born-Oppenheimer approximation. *Molecular Physics*, **1980**, 41, 1113-1122 1.7 58
- 179 Predissociation of weak-anisotropy Van der Waals molecules. Theory, approximations and practical predictions. *Faraday Discussions of the Chemical Society*, **1982**, 73, 339 57

178	Large effects of electric fields on atom-molecule collisions at millikelvin temperatures. <i>Physical Review Letters</i> , <b>2011</b> , 106, 193201	7.4	56
177	Interactions and dynamics in Li+Li <sub>2</sub> ultracold collisions. <i>Journal of Chemical Physics</i> , <b>2007</b> , 127, 074302	3.9	56
176	Prospects of forming ultracold molecules in 2 $\Sigma$ states by magnetoassociation of alkali-metal atoms with Yb. <i>Physical Review A</i> , <b>2013</b> , 87,	2.6	55
175	I-NoLLS: A program for interactive nonlinear least-squares fitting of the parameters of physical models. <i>Computer Physics Communications</i> , <b>1997</b> , 102, 252-268	4.2	55
174	Anisotropic intermolecular potentials. III. Rare-gas-Hydrogen bromide systems. <i>Journal of Chemical Physics</i> , <b>1989</b> , 91, 4455-4461	3.9	54
173	A theoretical study of the Ar <sub>2</sub> HCl van der Waals cluster. <i>Journal of Chemical Physics</i> , <b>1989</b> , 90, 1337-1344	3.9	54
172	Sticky collisions of ultracold RbCs molecules. <i>Nature Communications</i> , <b>2019</b> , 10, 3104	17.4	53
171	Interaction of NH(X <sup>3</sup> $\Sigma^-$ ) molecules with rubidium atoms: implications for sympathetic cooling and the formation of extremely polar molecules. <i>Physical Review Letters</i> , <b>2004</b> , 92, 163202	7.4	53
170	An evaluation of existing potential energy surfaces for CO <sub>2</sub> -Ar: Pressure broadening and high-resolution spectroscopy of van der Waals complexes. <i>Journal of Chemical Physics</i> , <b>1996</b> , 104, 2156-2166	3.9	53
169	Methods for calculating the bound state energies of van der Waals trimers: Applications to Ar <sub>3</sub> . <i>Journal of Chemical Physics</i> , <b>1993</b> , 98, 2160-2169	3.9	53
168	Ultracold molecules for quantum simulation: rotational coherences in CaF and RbCs. <i>Quantum Science and Technology</i> , <b>2019</b> , 4, 014010	5.5	53
167	Low-energy collisions of NH <sub>3</sub> and ND <sub>3</sub> with ultracold Rb atoms. <i>Physical Review A</i> , <b>2009</b> , 79,	2.6	51
166	Non-additive intermolecular forces from the spectroscopy of Van der Waals trimers: A comparison of Ar <sub>2</sub> HF and Ar <sub>2</sub> HCl, including H/D isotope effects. <i>Journal of Chemical Physics</i> , <b>1997</b> , 106, 6288-6301	3.9	51
165	Spectral line shape parameters for HF in a bath of Ar are accurately predicted by a potential inferred from spectra of the van der Waals dimer. <i>Journal of Chemical Physics</i> , <b>1994</b> , 100, 891-898	3.9	51
164	Close-coupling calculations of transport and relaxation cross sections for H <sub>2</sub> in Ar. <i>Journal of Chemical Physics</i> , <b>1984</b> , 80, 1135-1149	3.9	51
163	Morphing the He-DCS intermolecular potential. <i>Journal of Chemical Physics</i> , <b>2001</b> , 115, 5059-5065	3.9	50
162	Pairwise-additive models for atom-surface interaction potentials: An ab initio study of He-LiF. <i>Physical Review B</i> , <b>1986</b> , 33, 3724-3735	3.3	49
161	New vibration-rotation code for tetraatomic molecules exhibiting wide-amplitude motion: WAVR4. <i>Computer Physics Communications</i> , <b>2004</b> , 163, 117-131	4.2	47

160	Potential energy surfaces for ArDH (X 2) obtained by fitting to high-resolution spectroscopy. <i>Journal of Chemical Physics</i> , <b>1993</b> , 99, 7477-7486	3.9	46
159	A spectroscopic puzzle in ArHF solved: The test of a new potential. <i>Journal of Chemical Physics</i> , <b>1992</b> , 97, 8009-8018	3.9	46
158	Parity doubling in open-shell van der Waals complexes. <i>Chemical Physics Letters</i> , <b>1992</b> , 193, 355-363	2.5	46
157	Calculating energy levels of isomerizing tetra-atomic molecules. II. The vibrational states of acetylene and vinylidene. <i>Journal of Chemical Physics</i> , <b>2005</b> , 122, 064309	3.9	45
156	Cold collisions between OH and Rb: The field-free case. <i>Physical Review A</i> , <b>2007</b> , 75,	2.6	44
155	On the rotational constants of floppy molecules. <i>Chemical Physics Letters</i> , <b>1994</b> , 222, 257-262	2.5	44
154	Vibrational predissociation of hydrogen, deuterium, and hydrogen deuteride-argon van der Waals molecules. <i>The Journal of Physical Chemistry</i> , <b>1983</b> , 87, 2713-2720		44
153	Vibrational predissociation of the NeD <sub>2</sub> H <sub>4</sub> and ArD <sub>2</sub> H <sub>4</sub> van der Waals complexes. <i>Journal of Chemical Physics</i> , <b>1984</b> , 81, 4474-4480	3.9	44
152	Production of optically trapped RbCs <sup>87</sup> Feshbach molecules. <i>Physical Review A</i> , <b>2014</b> , 89,	2.6	43
151	The atom-surface interaction potential for He-NaCl: A model based on pairwise additivity. <i>Surface Science</i> , <b>1986</b> , 173, 337-350	1.8	43
150	High resolution radiofrequency spectroscopy of ArHCl. <i>Journal of Chemical Physics</i> , <b>1981</b> , 74, 6520-6521	3.1	43
149	Observation of Feshbach resonances between alkali and closed-shell atoms. <i>Nature Physics</i> , <b>2018</b> , 14, 881-884	16.2	42
148	Modeling sympathetic cooling of molecules by ultracold atoms. <i>Physical Review A</i> , <b>2015</b> , 92,	2.6	41
147	Non-additive intermolecular forces from the spectroscopy of van der Waals trimers: far-infrared spectra and calculations on Ar <sub>2</sub> -DCl. <i>Molecular Physics</i> , <b>1994</b> , 81, 579-598	1.7	41
146	Predissociation of HD/Ar van der Waals molecules by internal rotation. <i>Journal of Chemical Physics</i> , <b>1983</b> , 78, 4040-4043	3.9	41
145	The potential energy surface of HeHCN determined by fitting to high-resolution spectroscopic data. <i>Journal of Chemical Physics</i> , <b>1996</b> , 105, 440-450	3.9	40
144	Ultracold polar molecules as qubits. <i>New Journal of Physics</i> , <b>2020</b> , 22, 013027	2.9	40
143	Rotationally inelastic scattering in CH <sub>4</sub> +He, Ne, and Ar: State-to-state cross sections via direct infrared laser absorption in crossed supersonic jets. <i>Journal of Chemical Physics</i> , <b>1996</b> , 105, 3497-3516	3.9	39

142	Atom-spherical top van der Waals complexes: A theoretical study. <i>Journal of Chemical Physics</i> , <b>1994</b> , 100, 2505-2521	3.9	39
141	Prospects for sympathetic cooling of polar molecules: NH with alkali-metal and alkaline-earth atoms--a new hope. <i>Faraday Discussions</i> , <b>2009</b> , 142, 191-201; discussion 221-55	3.6	38
140	Manipulating ultracold polar molecules with microwave radiation: The influence of hyperfine structure. <i>Physical Review A</i> , <b>2009</b> , 80,	2.6	37
139	The potential energy surface and near-dissociation states of He-H <sub>2</sub> <sup>+</sup> . <i>Journal of Chemical Physics</i> , <b>1999</b> , 110, 3418-3427	3.9	37
138	Cold and ultracold NH-NH collisions in magnetic fields. <i>Physical Review A</i> , <b>2011</b> , 83,	2.6	36
137	The prospects of sympathetic cooling of NH molecules with Li atoms. <i>European Physical Journal D</i> , <b>2011</b> , 65, 151-160	1.3	36
136	Stimulating the production of deeply bound RbCs molecules with laser pulses: the role of spin-orbit coupling in forming ultracold molecules. <i>New Journal of Physics</i> , <b>2009</b> , 11, 055011	2.9	35
135	Interaction between LiH molecule and Li atom from state-of-the-art electronic structure calculations. <i>Journal of Chemical Physics</i> , <b>2011</b> , 134, 114109	3.9	35
134	Prospects for producing ultracold NH <sub>3</sub> molecules by sympathetic cooling: A survey of interaction potentials. <i>Physical Review A</i> , <b>2008</b> , 78,	2.6	34
133	molscat: A program for non-reactive quantum scattering calculations on atomic and molecular collisions. <i>Computer Physics Communications</i> , <b>2019</b> , 241, 9-18	4.2	33
132	Cold collisions of N (4S) atoms and NH (3 $\Sigma$ ) molecules in magnetic fields. <i>Physical Chemistry Chemical Physics</i> , <b>2011</b> , 13, 3669-80	3.6	33
131	Avoided crossings between bound states of ultracold cesium dimers. <i>Physical Review A</i> , <b>2008</b> , 78,	2.6	33
130	Multichannel quantum defect theory for cold molecular collisions. <i>Physical Review A</i> , <b>2011</b> , 84,	2.6	32
129	Calculating nuclear quadrupole coupling constants for van der Waals complexes. <i>Molecular Physics</i> , <b>1995</b> , 84, 185-199	1.7	32
128	Controlling the rotational and hyperfine state of ultracold Rb <sup>87</sup> Cs <sup>133</sup> molecules. <i>Physical Review A</i> , <b>2016</b> , 94,	2.6	31
127	Observation of a microwave spectrum of the long-range He [H <sub>2</sub> <sup>+</sup> complex. <i>Chemical Physics Letters</i> , <b>1996</b> , 260, 395-405	2.5	31
126	Hyperfine energy levels of alkali-metal dimers: Ground-state homonuclear molecules in magnetic fields. <i>Physical Review A</i> , <b>2009</b> , 79,	2.6	30
125	Long range intermolecular forces in triatomic systems: connecting the atom-atom and atom-atom-atom representations. <i>Molecular Physics</i> , <b>2006</b> , 104, 23-31	1.7	30

124	Vibrational relaxation of CO ( $v=1$ ) by inelastic collisions with 3He and 4He. <i>Journal of Chemical Physics</i> , <b>1995</b> , 103, 2528-2537	3.9	29
123	Prospects for sympathetic cooling of molecules in electrostatic, ac and microwave traps. <i>European Physical Journal D</i> , <b>2011</b> , 65, 141-149	1.3	28
122	Regular and irregular vibrational states: Localized anharmonic modes in Ar <sub>3</sub> . <i>Journal of Chemical Physics</i> , <b>1999</b> , 110, 902-911	3.9	28
121	The intermolecular potential of Ar-acetylene. Information from infrared and microwave spectroscopy. <i>Chemical Physics Letters</i> , <b>1992</b> , 198, 1-8	2.5	28
120	Microwave Shielding of Ultracold Polar Molecules. <i>Physical Review Letters</i> , <b>2018</b> , 121, 163401	7.4	28
119	Nonadditive intermolecular forces in Ar <sub>n</sub> HF van der Waals clusters: Effects on the HF vibrational frequency shift. <i>Journal of Chemical Physics</i> , <b>1999</b> , 111, 8378-8383	3.9	27
118	Calculations of the spectra of rare gas dimers and trimers: Implications for additive and nonadditive intermolecular forces in Ne <sub>2</sub> Ar, Ne <sub>2</sub> Kr, Ne <sub>2</sub> Xe, Ar <sub>2</sub> Ne, Ar <sub>3</sub> , Ar <sub>2</sub> Kr and Ar <sub>2</sub> Xe. <i>Journal of Chemical Physics</i> , <b>1995</b> , 103, 3386-3391	3.9	27
117	Microwave spectroscopy and interaction potential of the long-range He <sup>+</sup> Kr <sup>+</sup> ion: An example of Hund's case (e). <i>Journal of Chemical Physics</i> , <b>1996</b> , 105, 8602-8614	3.9	27
116	Dramatic reductions in inelastic cross sections for ultracold collisions near Feshbach resonances. <i>Physical Review Letters</i> , <b>2009</b> , 103, 163201	7.4	26
115	Line shape, transport and relaxation properties from intermolecular potential energy surfaces: The test case of CO <sub>2</sub> Ar. <i>Journal of Chemical Physics</i> , <b>1997</b> , 107, 1824-1834	3.9	26
114	Microwave electronic spectrum of the Ne <sup>+</sup> Ne <sup>+</sup> long-range complex: The interaction potential. <i>Journal of Chemical Physics</i> , <b>2002</b> , 116, 3662-3669	3.9	26
113	Coupled channel bound state calculations: Calculating expectation values without wavefunctions. <i>Chemical Physics Letters</i> , <b>1988</b> , 151, 565-569	2.5	26
112	Use of calculated centrifugal distortion constants ( $D_{JK}$ and $M_J$ ) in the analysis of the B $\leftarrow$ X system of I <sub>2</sub> . <i>Journal of Molecular Spectroscopy</i> , <b>1982</b> , 96, 266-278	1.3	26
111	Feshbach resonances in ultracold 85Rb. <i>Physical Review A</i> , <b>2013</b> , 87,	2.6	25
110	Chemistry. Ultracold chemistry. <i>Science</i> , <b>2010</b> , 327, 788-9	33.3	25
109	Intermolecular potential energy surfaces and bound states in FHF. <i>Journal of Chemical Physics</i> , <b>2000</b> , 112, 592-600	3.9	25
108	Bound-state wave functions from coupled channel calculations using log-derivative propagators: Application to spectroscopic intensities in ArHF. <i>Journal of Chemical Physics</i> , <b>1994</b> , 101, 5578-5584	3.9	25
107	The secular equation/perturbation theory method for calculating spectra of van der Waals complexes. <i>Journal of Chemical Physics</i> , <b>1985</b> , 83, 1197-1203	3.9	25



106	Hyperfine structure of alkali-metal diatomic molecules. <i>Physical Review A</i> , <b>2017</b> , 96,	2.6	24
105	Calculating energy levels of isomerizing tetra-atomic molecules. I. The rovibrational bound states of Ar <sub>2</sub> HF. <i>Journal of Chemical Physics</i> , <b>2003</b> , 118, 4896-4904	3.9	24
104	Physical origin of oscillations in the three-dimensional collision amplitudes of heavy-light-heavy systems. Semiclassical quantization of chaotic scattering. <i>Journal of Chemical Physics</i> , <b>1993</b> , 98, 3929-3944	2.9	24
103	The augmented secular equation method for calculating spectra of van der Waals complexes. Application to the infrared spectrum of Ar <sub>2</sub> Cl. <i>Journal of the Chemical Society, Faraday Transactions 2</i> , <b>1986</b> , 82, 1163-1171		24
102	Robust entangling gate for polar molecules using magnetic and microwave fields. <i>Physical Review A</i> , <b>2020</b> , 101,	2.6	23
101	ac Stark effect in ultracold polar Rb <sup>87</sup> Cs <sup>133</sup> molecules. <i>Physical Review A</i> , <b>2017</b> , 96,	2.6	23
100	Contrasting the wide Feshbach resonances in Li <sub>6</sub> and Li <sub>7</sub> . <i>Physical Review A</i> , <b>2014</b> , 89,	2.6	23
99	On the choice of inertial axes for interpreting spectroscopic properties of van der Waals complexes. <i>Journal of Chemical Physics</i> , <b>1994</b> , 101, 5438-5440	3.9	23
98	Non-additive intermolecular forces from the spectroscopy of Van der Waals trimers: the effect of monomer vibrational excitation in Ar <sub>2</sub> HF and Ar <sub>2</sub> HCl. <i>Faraday Discussions</i> , <b>1994</b> , 97, 119-129	3.6	23
97	Effective-range approximations for resonant scattering of cold atoms. <i>Physical Review A</i> , <b>2014</b> , 89,	2.6	22
96	Large-amplitude quantum mechanics in polyatomic hydrides. II. A particle-on-a-sphere model for XH(n) (n=4,5). <i>Journal of Chemical Physics</i> , <b>2008</b> , 128, 094306	3.9	22
95	Potential energy surfaces and bound states for the open-shell van der Waals cluster BrHF. <i>Journal of Chemical Physics</i> , <b>2003</b> , 119, 8873-8881	3.9	22
94	A semiempirical model for atom-surface dispersion coefficients. <i>Surface Science</i> , <b>1986</b> , 165, 289-302	1.8	22
93	Feshbach spectroscopy of an ultracold mixture of <sup>85</sup> Rb and <sup>133</sup> Cs. <i>Physical Review A</i> , <b>2013</b> , 87,	2.6	21
92	Quadrupolar contributions to the atom-surface Van Der Waals interaction. <i>Surface Science</i> , <b>1986</b> , 175, L775-L781	1.8	21
91	Cold atomic and molecular collisions: approaching the universal loss regime. <i>New Journal of Physics</i> , <b>2015</b> , 17, 045019	2.9	20
90	Three-body parameter for Efimov states in Li <sub>6</sub> . <i>Physical Review A</i> , <b>2014</b> , 90,	2.6	20
89	Magnetically tunable Feshbach resonances in Li+Yb(3PJ). <i>Physical Review A</i> , <b>2013</b> , 88,	2.6	20

88	Effect of hyperfine interactions on ultracold molecular collisions: NH(3) with Mg(1S) in magnetic fields. <i>Physical Review A</i> , <b>2011</b> , 84,	2.6	20
87	On the coupled-channel calculation of bound states for trimeric systems using hyperspherical coordinates. <i>Journal of Chemical Physics</i> , <b>1989</b> , 91, 4197-4203	3.9	20
86	Coupled channel calculations on the vibrational predissociation of the ethylene dimer. <i>Chemical Physics Letters</i> , <b>1986</b> , 125, 477-480	2.5	20
85	Close-coupling calculations of transport and relaxation cross sections for H2 in Ar. <i>Journal of Chemical Physics</i> , <b>1987</b> , 86, 854-857	3.9	20
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