List of Publications by Year in descending order

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| # | Article | IF | CITATIONS |
|----|--|-----|-----------|
| 1 | Complexity of HCl and H2 molecules under q-deformed Morse potential. Indian Journal of Physics, 2022, 96, 1-7. | 0.9 | 1 |
| 2 | Multipole polarizabilities and dipole oscillator strengths of H-atom in nonideal classical plasmas. European Physical Journal Plus, 2022, 137, 1. | 1.2 | 16 |
| 3 | He atom in a quantum dot: Structural, entanglement, and information-theoretical measures. Physical Review A, 2022, 105, . | 1.0 | 4 |
| 4 | Information theory and Wigner crystallization: A model perspective. International Journal of Quantum Chemistry, 2021, 121, e26549. | 1.0 | 2 |
| 5 | Ionization of many-electron atoms by the action of two plasma models. Physical Review E, 2021, 103, 043202. | 0.8 | 10 |
| 6 | Structural modifications of two-electron systems under isotropic harmonic confinement. European Physical Journal D, 2021, 75, 1. | 0.6 | 2 |
| 7 | Excited states of the Gaussian two-electron quantum dot. European Physical Journal D, 2021, 75, 1. | 0.6 | 1 |
| 8 | Critical screening in the one- and two-electron Yukawa atoms. Physical Review A, 2018, 97, . | 1.0 | 37 |
| 9 | Exact normalized eigenfunctions for general deformed Hulthén potentials. Journal of Mathematical Physics, 2018, 59, 122103. | 0.5 | 7 |
| 10 | A new comparison theorem on Hellmann potential. Journal of Molecular Modeling, 2018, 24, 299. | 0.8 | 1 |
| 11 | Hund's rule in the (1 <i>s</i> 2 <i>s</i>)1,3 <i>S</i> states of the two-electron Debye atom. Physics of Plasmas, 2018, 25, . | 0.7 | 11 |
| 12 | Two particle system in spherically confined plasma environment. European Physical Journal D, 2017, 71, 1. | 0.6 | 8 |
| 13 | Quantum information entropies for the \$\$ell \$\$ ℓ -state Pöschl–Teller-type potential. Journal of Mathematical Chemistry, 2016, 54, 1810-1821. | 0.7 | 16 |
| 14 | Position and momentum informationâ€theoretic measures of the pseudoharmonic potential. International Journal of Quantum Chemistry, 2015, 115, 1543-1552. | 1.0 | 51 |
| 15 | Soft and hard confinement of a two-electron quantum system. European Physical Journal Plus, 2014, 129, 1. | 1.2 | 5 |
| 16 | Aneesur Rahman. Resonance, 2014, 19, 671-683. | 0.2 | 0 |
| 17 | Bound state solutions of the Deng–Fan molecular potential with the Pekeris-type approximation using the Nikiforov–Uvarov (N–U) method. Journal of Mathematical Chemistry, 2013, 51, 976-991. | 0.7 | 52 |
| 18 | Exact solutions of the SchrĶdinger equation for the pseudoharmonic potential: an application to some diatomic molecules. Journal of Mathematical Chemistry, 2012, 50, 1039-1059. | 0.7 | 69 |

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| 19 | Scaling properties of net information measures for bound states of spherical model potentials confined with finite barrier #. Journal of Chemical Sciences, 2012, 124, 241-245. | 0.7 | 7 |
| 20 | Spectral characteristics for a spherically confined â^'‹i›a‹/i›/‹i›r‹/i›+‹i›br‹/i›‹sup›2‹/sup›potential. Journal of Physics A: Mathematical and Theoretical, 2011, 44, 185307. | 0.7 | 15 |
| 21 | A generalized relative complexity measure. Journal of Statistical Mechanics: Theory and Experiment, 2011, 2011, P09016. | 0.9 | 22 |
| 22 | Static and dynamic dipole polarizabilities and electron density at origin: Ground and excited states of hydrogen atom confined in multiwalled fullerenes. International Journal of Quantum Chemistry, 2011, 111, 4425-4432. | 1.0 | 31 |
| 23 | Discrete spectra for confined and unconfined â^' <i>a</i> / <i>r</i> + <i>br</i> 2 potentials in <i>d</i> -dimensions. Journal of Mathematical Physics, 2011, 52, . | 0.5 | 11 |
| 24 | Phenomenological Description of a Three-Center Insertion Reaction: An Information-Theoretic Study. Journal of Physical Chemistry A, 2010, 114, 1906-1916. | 1.1 | 17 |
| 25 | Soft-core Coulomb potentials and Heun's differential equation. Journal of Mathematical Physics, 2010, 51, 022107. | 0.5 | 22 |
| 26 | Electron density and its derivatives at the nucleus for spherically confined hydrogen atom. International Journal of Quantum Chemistry, 2009, 109, 688-692. | 1.0 | 8 |
| 27 | Energies and wave functions for a soft-core Coulomb potential. Physical Review A, 2009, 80, . | 1.0 | 27 |
| 28 | Electrostatic interpretation of the force â^'â^,Vxc/â^,r connected with the exchange-correlation potential: direct relation to single-particle kinetic energy density in Be-atom. World Scientific Series in 20th Century Physics, 2009, , 689-692. | 0.0 | 0 |
| 29 | Dynamic dipole polarizabilities of the ground and excited states of confined hydrogen atom computed by means of a mapped Fourier grid method. International Journal of Quantum Chemistry, 2008, 108, 351-361. | 1.0 | 20 |
| 30 | Simultaneous degeneracy of the confined 2D hydrogen atom energy levels. Journal of Physics B: Atomic, Molecular and Optical Physics, 2008, 41, 205002. | 0.6 | 8 |
| 31 | Eigenspectrum properties of the confined 3D harmonic oscillator. Journal of Physics B: Atomic, Molecular and Optical Physics, 2008, 41, 225002. | 0.6 | 11 |
| 32 | A study of the confined 2D isotropic harmonic oscillator in terms of the annihilation and creation operators and the infinitesimal operators of the SU(2) group. Journal of Physics A: Mathematical and Theoretical, 2008, 41, 265203. | 0.7 | 10 |
| 33 | Confined hydrogen atom by the Lagrange-mesh method: Energies, mean radii, and dynamic polarizabilities. Physical Review E, 2008, 78, 026701. | 0.8 | 46 |
| 34 | Characteristic features of net information measures for constrained Coulomb potentials. Journal of Physics B: Atomic, Molecular and Optical Physics, 2007, 40, 2147-2162. | 0.6 | 37 |
| 35 | Fisher-Shannon analysis of ionization processes and isoelectronic series. Physical Review A, 2007, 76, . | 1.0 | 77 |
| 36 | Degeneracy of confinedD-dimensional harmonic oscillator. International Journal of Quantum Chemistry, 2007, 107, 798-806. | 1.0 | 42 |

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| 37 | Net information measures for modified Yukawa and Hulthén potentials. International Journal of Quantum Chemistry, 2007, 107, 1864-1874. | 1.0 | 46 |
| 38 | Information entropies for eigendensities of homogeneous potentials. Journal of Chemical Physics, 2006, 125, 074117. | 1.2 | 53 |
| 39 | DFT reactivity indices in confined many-electron atoms. Journal of Chemical Sciences, 2005, 117, 379-386. | 0.7 | 43 |
| 40 | First-order correlation-kinetic contribution to Kohn-Sham exchange charge density function in atoms, using quantal density functional theory approach. International Journal of Quantum Chemistry, 2005, 101, 231-238. | 1.0 | 4 |
| 41 | Force â^'â^,Vxc/â^,rassociated with the exchange-correlation potentialVxc(r) in the neutral Ne atom. Journal of Physics B: Atomic, Molecular and Optical Physics, 2005, 38, 1705-1717. | 0.6 | 3 |
| 42 | Characteristic features of Shannon information entropy of confined atoms. Journal of Chemical Physics, 2005, 123, 074110. | 1.2 | 153 |
| 43 | Shell-confined hydrogen atom. Journal of Chemical Physics, 2005, 122, 194324. | 1.2 | 44 |
| 44 | Alkali atoms confined to a sphere and to a fullerene cage. Canadian Journal of Physics, 2005, 83, 919-928. | 0.4 | 8 |
| 45 | On the importance of the "density per particle―(shape function) in the density functional theory. Journal of Chemical Physics, 2004, 120, 9969-9973. | 1.2 | 88 |
| 46 | Obtaining Kohn-Sham potential without taking the functional derivative. Bulletin of Materials Science, 2003, 26, 69-74. | 0.8 | 4 |
| 47 | Nuclear cusp of the virial exchange energy density for spherical atoms. Journal of Chemical Physics, 2002, 117, 9107-9110. | 1.2 | 1 |
| 48 | Mean excitation energy, static polarizability, and hyperpolarizability of the spherically confined hydrogen atom. Journal of Chemical Physics, 2002, 116, 4054-4057. | 1.2 | 69 |
| 49 | Improved Becke88 and PW91 exchange potentials. Journal of Physics B: Atomic, Molecular and Optical Physics, 2002, 35, 4711-4718. | 0.6 | 5 |
| 50 | A correlation of exchange energy-exchange potential at the nucleus in atoms. Journal of Chemical Physics, 2002, 116, 9570-9573. | 1.2 | 4 |
| 51 | Applications of electrostatic interpretation of components of effective Kohn–Sham potential in atoms. Journal of Chemical Physics, 2002, 117, 4684-4693. | 1.2 | 9 |
| 52 | Modified Slater potential and its application to the ground-states and excited-states of atomic systems. International Journal of Quantum Chemistry, 2002, 90, 327-333. | 1.0 | 7 |
| 53 | Static dipole and quadrupole polarizability of confined hydrogen atom withZ=N/3 (N=1-5). International Journal of Quantum Chemistry, 2002, 90, 491-496. | 1.0 | 18 |
| 54 | Ground- and excited-state cusp conditions for the electron density. Journal of Chemical Physics, 2001, 115, 6300-6308. | 1.2 | 48 |

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|----|---|-----|-----------|
| 55 | Atomic shell structure in Hartree theory. Journal of Chemical Physics, 2001, 114, 8784-8788. | 1.2 | 7 |
| 56 | Higher-order cusp of the density in certain highly excited states of atoms and molecules. Journal of Physics B: Atomic, Molecular and Optical Physics, 2000, 33, 1745-1751. | 0.6 | 41 |
| 57 | The creative Kohn. Resonance, 1999, 4, 56-63. | 0.2 | 0 |
| 58 | Characteristic features of the electrostatic potential for negative atoms within the work formalism. Journal of Chemical Physics, 1996, 104, 1025-1027. | 1.2 | 3 |
| 59 | Average local electrostatic potential and the core-valency separation in atoms. International Journal of Quantum Chemistry, 1995, 56, 399-408. | 1.0 | 16 |
| 60 | A Combination of the Work Formalism for Exchange with an Optimized Correlation Energy Functional for Atoms. Journal De Physique II, 1995, 5, 1277-1287. | 0.9 | 1 |
| 61 | Shell boundaries in rare earth atoms and tripositive ions using average local electrostatic potential. Journal of Chemical Physics, 1994, 101, 7779-7781. | 1.2 | 4 |
| 62 | Radii of monopositive atomic ions. Journal of Chemical Physics, 1993, 99, 3149-3150. | 1.2 | 8 |
| 63 | Density-functional-theory calculations of static dipole polarizability of some ions of interest in Mössbauer spectroscopy. Physical Review A, 1992, 45, 2076-2078. | 1.0 | 5 |
| 64 | Static polarizabilities for the ne-isoelectronic series using Harbola-Sahni potential. International Journal of Quantum Chemistry, 1992, 44, 1041-1044. | 1.0 | 7 |
| 65 | Upper bound to approximate ionic radii of atomic negative ions in terms of ã€^r 2〉. Journal of Chemical Physics, 1991, 95, 1421-1421. | 1.2 | 3 |
| 66 | Approximate sizes of monoatomic negative ions with fractional nuclear charge using electrostatic potentials. Theoretica Chimica Acta, 1991, 79, 373-375. | 0.9 | 0 |
| 67 | Static multipole polarizabilities ofFâ^'andClâ^'using the Harbola-Sahni potential. Physical Review A, 1991, 44, 756-757. | 1.0 | 11 |
| 68 | Isoelectronic changes in energy of quark atoms and molecules via the Levy equation. Theoretica Chimica Acta, 1990, 77, 57-60. | 0.9 | 2 |
| 69 | Quadrupole polarizabilities ofFâ^',Clâ^', andBrâ^'using Ξatheory. Physical Review A, 1990, 42, 3115-3116. | 1.0 | 1 |
| 70 | Generalized Zâ€ŧransition state calculations of isoelectronic change in energy of atoms. Journal of Chemical Physics, 1990, 92, 2113-2114. | 1.2 | 1 |
| 71 | Zâ€ŧransition state calculations of energy changes and electrostatic potentials in isoelectronic atoms and molecules. Journal of Chemical Physics, 1989, 90, 4373-4378. | 1.2 | 12 |
| 72 | Approximate radii for singly negative ions of 3d, 4d, and 5d metal atoms. Journal of Chemical Physics, 1989, 91, 5123-5124. | 1.2 | 33 |

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|----|--|-----|-----------|
| 73 | Density functional theory calculations of one electron Rydberg states in Li atom. Theoretica Chimica Acta, 1989, 76, 373-375. | 0.9 | 12 |
| 74 | Electronegativities and isoelectronic energy and electronegativity differences for monatomic systems with nonintegral nuclear charges: Local-spin-density-functional calculations. Physical Review A, 1989, 40, 2260-2264. | 1.0 | 6 |
| 75 | Characteristic features of the electrostatic potentials of singly negative monoatomic ions. Journal of Chemical Physics, 1989, 90, 4370-4372. | 1.2 | 125 |
| 76 | Quadrupole antishielding factors forFâ^',Clâ^', andBrâ^'usingΞatheory. Physical Review A, 1988, 38, 4330-4332. | 1.0 | 2 |
| 77 | Two-electron one-photon x-ray transition energy in the Kr isoelectronic series. Physical Review A, 1987, 36, 2971-2972. | 1.0 | 2 |
| 78 | Sternheimer Antishielding Functions ß(r) and γ(r) for Rare Earth Atoms. Zeitschrift Fur Naturforschung - Section A Journal of Physical Sciences, 1986, 41, 37-46. | 0.7 | 7 |
| 79 | Approximate upper bounds to the momentum expectation value ratios ?p 2?/?p ?1? and ?p?/?p ?1? in atoms. Theoretica Chimica Acta, 1986, 70, 115-117. | 0.9 | 0 |
| 80 | A simple bound to average electronic density and energy for atoms within an isoelectronic series. International Journal of Quantum Chemistry, 1986, 29, 165-167. | 1.0 | 5 |
| 81 | Electric-field-gradient sternheimer function for closed shell ions. Hyperfine Interactions, 1986, 30, 253-264. | 0.2 | 8 |
| 82 | Empirical estimates of electrostatic potentials at the nuclei in molecules. International Journal of Quantum Chemistry, 1985, 27, 231-232. | 1.0 | 2 |
| 83 | Nuclear quadrupole Sternheimer shielding-antishielding function for N3-, Cl-and Cd2+ions in crystals. Journal of Physics C: Solid State Physics, 1985, 18, 3153-3156. | 1.5 | 2 |
| 84 | Bounds to average radial electron density in atoms using generalized Hölder inequality. Journal of Chemical Physics, 1985, 83, 3709-3709. | 1.2 | 7 |
| 85 | 3d contribution to Sternheimer antishielding in Sc-Ni. Journal of Physics F: Metal Physics, 1984, 14, 521-523. | 1.6 | 0 |
| 86 | Empirical static quadrupole polarisability for closed shell ions in solids. Journal of Physics C: Solid State Physics, 1984, 17, L227-L228. | 1.5 | 2 |
| 87 | Empirical static quadrupole polarizability for some closed shell free atoms and ions. Journal of Chemical Physics, 1984, 80, 584-585. | 1.2 | 6 |
| 88 | Approximate upper bounds to average radial electron density expectation value ã€^ Ï〉 within an isoelectronic series of atoms. Journal of Chemical Physics, 1984, 81, 5213-5214. | 1.2 | 6 |
| 89 | A simple Vne–i̇́0 relationship for atomic isoelectronic series. Journal of Chemical Physics, 1984, 81, 2861-2861. | 1.2 | 2 |
| 90 | Semiempirical estimates of static multipole polarizabilities for free atoms. International Journal of Quantum Chemistry, 1984, 26, 1051-1052. | 1.0 | 2 |

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| 91 | Radial sternheimer shielding-antishielding for He, Ne, and Ar isoelectronic series. Hyperfine Interactions, 1983, 14, 243-247. | 0.2 | 0 |
| 92 | Lithiated ammonia, amide anions, and ammonium ions. An ab initio study of structures, bonding, and energetic relationships. Inorganic Chemistry, 1983, 22, 496-503. | 1.9 | 60 |
| 93 | Isoelectronic changes in the Hartree–Fock properties of atoms. Journal of Chemical Physics, 1983, 78, 4779-4780. | 1.2 | 0 |
| 94 | Sternheimer antishielding functions for 3d transition metal atoms. Journal of Physics F: Metal Physics, 1982, 12, L97-L99. | 1.6 | 1 |
| 95 | Radial Sternheimer shielding-antishielding for rare gas atoms. Hyperfine Interactions, 1982, 12, 113-118. | 0.2 | 3 |
| 96 | Slater transition state calculations of valence electron spin–orbit splitting in atoms. Journal of Chemical Physics, 1981, 75, 5971-5971. | 1.2 | 2 |
| 97 | Static dipole and quadrupole polarizabilities of alkaline earth atoms. International Journal of Quantum Chemistry, 1981, 19, 373-376. | 1.0 | 14 |
| 98 | Overlap sternheimer shielding-antishielding functions for halogen ions in ionic solids. Hyperfine Interactions, 1981, 9, 355-359. | 0.2 | 3 |
| 99 | Polarizabilities and Sternheimer antishielding factor for ferriclike ions. Physical Review A, 1981, 23, 1026-1029. | 1.0 | 10 |
| 100 | On the geometry of NaPO3. Journal of Chemical Physics, 1981, 75, 1043-1044. | 1.2 | 9 |
| 101 | On deviations in Ruedenberg energy formula for molecules. Journal of Chemical Physics, 1981, 74, 1500-1501. | 1.2 | 1 |
| 102 | Slater transition state calculations of electron affinity of heavy atoms. Journal of Chemical Physics, 1981, 75, 1037-1038. | 1.2 | 21 |
| 103 | The Spherical Crystal Potential and Physical Properties of Ions in Crystals. A HF alculation for the 10, 18, and 36 Electron Closed Shell Ions of Xâ€Ray Scattering Factors, Diamagnetic Susceptibilities, and Dipole Polarizabilities. Zeitschrift Fur Elektrotechnik Und Elektrochemie, 1980, 84, 1240-1251. | 0.9 | 10 |
| 104 | Isoelectronic changes in total Hartree-Fock energy of atoms. Theoretica Chimica Acta, 1980, 57, 179-180. | 0.9 | 1 |
| 105 | Approximate relationship for isoelectronic changes in atomic average electron density. International Journal of Quantum Chemistry, 1980, 17, 1241-1242. | 1.0 | 4 |
| 106 | Energy-eigenvalue sum relationship in molecules. International Journal of Quantum Chemistry, 1980, 18, 907-909. | 1.0 | 3 |
| 107 | Sternheimer quadrupole shielding-antishielding function for F-, Cl-, Br-and I-in ionic solids. Journal of Physics C: Solid State Physics, 1980, 13, L423-L426. | 1.5 | 1 |
| 108 | Dependence on distance of the Sternheimer shielding-antishielding factor for 3d elements. Journal of Physics F: Metal Physics, 1980, 10, L67-L70. | 1.6 | 5 |

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| 109 | Hartree–Fock–Slater transition state calculations of KLL Auger energies in Na and Mg. Journal of Chemical Physics, 1980, 73, 4704-4705. | 1.2 | 3 |
| 110 | Effect of self-consistency and crystalline potential in the solid state on nuclear quadrupole Sternheimer antishielding factors in closed-shell ions. Physical Review B, 1980, 22, 4167-4179. | 1.1 | 111 |
| 111 | Quadrupole polarizability and shielding-antishielding factors forCe3+in ionic solids. Physical Review B, 1980, 22, 5542-5543. | 1.1 | 0 |
| 112 | Quadrupole polarizabilities ofFe3+- andRu3+-like ions. Physical Review A, 1980, 21, 693-694. | 1.0 | 2 |
| 113 | Isoelectronic energy changes from energy–density functionals. Journal of Chemical Physics, 1980, 73, 5403-5404. | 1.2 | 2 |
| 114 | Isoelectronic Changes in the Behaviour of Nonrelativistic Electron Density Near the Nucleus in Free Ions. Zeitschrift Fur Naturforschung - Section A Journal of Physical Sciences, 1979, 34, 525-526. | 0.7 | 0 |
| 115 | Isoelectronic Changes in the Core Radius of Fe3+ and Ru3+ Like Ions. Zeitschrift Fur Naturforschung - Section A Journal of Physical Sciences, 1979, 34, 1253-1254. | 0.7 | 0 |
| 116 | Variation of shellwise s electron density at the nucleus in ionic solids. Journal of Physics C: Solid State Physics, 1979, 12, L75-L77. | 1.5 | 1 |
| 117 | Effect of a crystalline spherical potential on the Fermi-contact term inMn2+andFe3+ions. Physical Review B, 1979, 20, 2901-2904. | 1.1 | 0 |
| 118 | An approximate density equation for isoelectronic changes in atoms. Journal of Chemical Physics, 1979, 70, 5334-5334. | 1.2 | 6 |
| 119 | Effect of ionic size on the Koopmans ionization potentials in positive ions. Journal of Chemical Physics, 1979, 71, 559-560. | 1.2 | 4 |
| 120 | Sternheimer valence shielding and antishielding factors for several2pand3patoms and ions. Physical Review A, 1979, 20, 1323-1326. | 1.0 | 4 |
| 121 | Variation of core-electron contributions to Mössbauer isomer shift with ionic size. Physical Review B, 1979, 19, 2447-2448. | 1.1 | 1 |
| 122 | Atomic quadrupole moments:np2(n+1)sexcited-state configuration of N, P, As, and Sb. Physical Review A, 1979, 20, 2276-2277. | 1.0 | 6 |
| 123 | The isoelectronic energy–density relationship in atoms:2â€36 electron series. Journal of Chemical Physics, 1979, 71, 3551-3551. | 1.2 | 4 |
| 124 | Virial scaling corrections to Koopmans' ionization potential in positive ions of He, Ne, and Ar series. Journal of Chemical Physics, 1979, 70, 2025-2026. | 1.2 | 1 |
| 125 | Sternheimer antishielding factors of F?, Cl?, Br? and I?. Theoretica Chimica Acta, 1979, 52, 181-187. | 0.9 | 9 |
| 126 | The total Hartree–Fock SCF energy in atoms and ions. Journal of Chemical Physics, 1979, 70, 586. | 1.2 | 9 |

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| 127 | Slater transitionâ€state calculations of shakeâ€up energies in Ne, Ar, and Kr. Journal of Chemical Physics, 1979, 71, 1035-1036. | 1.2 | 9 |
| 128 | Notizen: Orbital Binding Energies within the Statistical Exchange Approximation: Ne and Ar Isoelectronic Series. Zeitschrift Fur Naturforschung - Section A Journal of Physical Sciences, 1979, 34, 901-902. | 0.7 | 2 |
| 129 | Atomic quadrupole moments of Al and In. Physical Review A, 1978, 18, 1327-1328. | 1.0 | 6 |
| 130 | Sternheimer shielding factors for atomic quadrupole moments:np(n+1)sconfiguration of C, Si, Ge, Sn, and Pb. Physical Review A, 1978, 18, 26-28. | 1.0 | 4 |
| 131 | Nuclear hexadecapole antishielding factors: Tripositive lanthanide and tetrapositive actinide ions. Physical Review A, 1978, 18, 2450-2452. | 1.0 | 2 |
| 132 | Sternheimer valence shielding and antishielding factors for some ions of interest in Mössbauer spectroscopy. Physical Review B, 1977, 16, 107-114. | 1.1 | 12 |
| 133 | Quadrupole antishielding factors and polarizabilities in ionic crystals. Physical Review B, 1977, 15, 95-102. | 1.1 | 35 |
| 134 | Sternheimer antishielding factors for core electrons in metals: Comparison with free-ion results. Physical Review A, 1977, 16, 1786-1788. | 1.0 | 9 |
| 135 | Quadrupole antishielding factors for actinide ions. Physical Review A, 1976, 14, 880-881. | 1.0 | 4 |
| 136 | Quadrupole antishielding factors for some3d5,4d5, and5d5ions. Physical Review A, 1976, 14, 539-542. | 1.0 | 13 |
| 137 | Nuclear hexadecapole antishielding factors. Physical Review A, 1975, 11, 1162-1167. | 1.0 | 8 |