

Waldeci Paraguassu

List of Publications by Year in descending order

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papers

1,874
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236925

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docs citations

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times ranked

1891
citing authors

| # | ARTICLE | IF | CITATIONS |
|----|---|------|-----------|
| 1 | Phase Transitions and Coexistence of Magnetic and Electric Orders in the Methylhydrazinium Metal Formate Frameworks. <i>Chemistry of Materials</i> , 2017, 29, 2264-2275. | 6.7 | 136 |
| 2 | Pressure-induced structural phase transitions and amorphization in selected molybdates and tungstates. <i>Progress in Materials Science</i> , 2012, 57, 1335-1381. | 32.8 | 106 |
| 3 | Phonons in ferroelectric Bi ₂ WO ₆ : Raman and infrared spectra and lattice dynamics. <i>Applied Physics Letters</i> , 2008, 92, . | 3.3 | 73 |
| 4 | Temperature- and Pressure-Induced Phase Transitions in the Metal Formate Framework of [ND ₄][Zn(DCOO) ₃] and [NH ₄][Zn(HCOO) ₃]. <i>Inorganic Chemistry</i> , 2014, 53, 9615-9624. | 4.0 | 72 |
| 5 | Impedance spectroscopy analysis of BaFe ₁₂ O ₁₉ M-type hexaferrite obtained by ceramic method. <i>Ceramics International</i> , 2009, 35, 2443-2447. | 4.8 | 69 |
| 6 | Phonon calculation on olivine-like LiMPO ₄ (M = Ni, Co, Fe) and Raman scattering of the iron-containing compound. <i>Journal of Raman Spectroscopy</i> , 2005, 36, 213-220. Phonon instability driven phase transitions in ferroelectric BiW_2O_6 | 2.5 | 66 |
| 7 | Phonon instability driven phase transitions in ferroelectric BiW_2O_6 | 3.2 | 62 |
| 8 | Effect of solvent, temperature and pressure on the stability of chiral and perovskite metal formate frameworks of [NH ₂ NH ₃][M(HCOO) ₃] (M = Mn, Fe, Zn). <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 31653-31663. | 2.8 | 54 |
| 9 | Pressure-induced structural transformations in the molybdate Sc ₂ (MoO ₄) ₃ . <i>Physical Review B</i> , 2004, 69, . | 3.2 | 52 |
| 10 | Temperature-dependent Raman scattering studies of Na ₂ MoO ₄ . <i>Journal of Raman Spectroscopy</i> , 2008, 39, 937-941. | 2.5 | 52 |
| 11 | A comparative study of negative thermal expansion materials Sc ₂ (MoO ₄) ₃ and Al ₂ (WO ₄) ₃ crystals. <i>Vibrational Spectroscopy</i> , 2007, 44, 69-77. | 2.2 | 51 |
| 12 | High-pressure Raman study of Al ₂ (WO ₄) ₃ . <i>Journal of Solid State Chemistry</i> , 2004, 177, 2002-2006. | 2.9 | 50 |
| 13 | Lattice dynamics and low-temperature Raman spectroscopy studies of PMN-PT relaxors. <i>Journal of Raman Spectroscopy</i> , 2009, 40, 1144-1149. | 2.5 | 48 |
| 14 | Temperature-induced phase transformations in Na ₂ WO ₄ and Na ₂ MoO ₄ crystals. <i>Journal of Raman Spectroscopy</i> , 2011, 42, 799-802. | 2.5 | 44 |
| 15 | Structural, thermal, dielectric and phonon properties of perovskite-like imidazolium magnesium formate. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 13993-14000. | 2.8 | 43 |
| 16 | Temperature-dependent Raman scattering studies on Na ₂ Mo ₂ O ₇ disodium dimolybdate. <i>Journal of Raman Spectroscopy</i> , 2011, 42, 1114-1119. | 2.5 | 42 |
| 17 | Raman and IR Studies of Pressure- and Temperature-Induced Phase Transitions in [(CH ₂) ₃ NH ₂] ₃ [Zn(HCOO) ₃]. <i>Inorganic Chemistry</i> , 2014, 53, 12650-12657. | 4.0 | 42 |
| 18 | Raman scattering studies of pressure-induced phase transitions in perovskite formates [(CH ₃) ₂ NH ₂][Mg(HCOO) ₃] and [(CH ₃) ₂ NH ₂][Cd(HCOO) ₃]. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2016, 156, 112-117. | 3.9 | 36 |

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|----|---|-----|-----------|
| 19 | Temperature- and pressure-induced phase transitions in the niccolite-type formate framework of $[\text{H}_{3}\text{N}(\text{CH}_{2})_{4}\text{NH}_{3}][\text{Mn}_{2}(\text{HCOO})_{6}]$. Journal of Materials Chemistry C, 2016, 4, 3185-3194. | 5.5 | 36 |
| 20 | Temperature- and pressure-dependent studies of a highly flexible and compressible perovskite-like cadmium dicyanamide framework templated with protonated tetrapropylamine. Journal of Materials Chemistry C, 2019, 7, 2408-2420. | 5.5 | 32 |
| 21 | Temperature-dependent Raman spectra of $\text{Ba}_{2}\text{BiSbO}_{6}$ ceramics. Journal of Raman Spectroscopy, 2009, 40, 1205-1210. | 2.5 | 31 |
| 22 | Raman scattering study of $\text{NaAl}(\text{MoO}_{4})_{2}$ crystal under high pressures. Journal of Physics Condensed Matter, 2004, 16, 5151-5161. | 1.8 | 27 |
| 23 | Structural and optical properties of rare earth-doped $(\text{Ba}_{0.77}\text{Ca}_{0.23})_{1-x}(\text{Sm}, \text{Nd}, \text{Pr}, \text{Yb})_{x}\text{TiO}_{3}$. Journal of Applied Physics, 2011, 109, . | 2.5 | 26 |
| 24 | Temperature-dependent Raman spectra of $\text{K}_{0.2}\text{Na}_{0.8}\text{NbO}_{3}$ ceramics. Journal of Raman Spectroscopy, 2005, 36, 28-32. | 2.5 | 25 |
| 25 | High pressure effects on the structural and vibrational properties of antiferromagnetic $\text{KFe}(\text{MoO}_{4})_{2}$. Journal of Physics Condensed Matter, 2005, 17, 6285-6300. | 1.8 | 25 |
| 26 | Pressure-induced phase transitions in ferroelectric $\text{Bi}_{2}\text{MoO}_{6}$ a Raman scattering study. Journal of Physics Condensed Matter, 2010, 22, 015901. | 1.8 | 24 |
| 27 | Raman and single-crystal X-ray diffraction evidence of pressure-induced phase transitions in a perovskite-like framework of $[(\text{C}_{3}\text{H}_{7})_{4}\text{N}][\text{Mn}(\text{N}(\text{CN})_{2})_{3}]$. Dalton Transactions, 2019, 48, 9072-9078. | 3.3 | 24 |
| 28 | Raman spectroscopy study of $\text{Na}_{2}\text{MoO}_{4} \cdot 2\text{H}_{2}\text{O}$ and $\text{Na}_{2}\text{MoO}_{4}$ under hydrostatic pressure. Journal of Raman Spectroscopy, 2010, 41, 576-581. | 2.5 | 23 |
| 29 | Lattice dynamics and pressure-induced phase transitions in $\text{Bi}_{2}\text{W}_{2}\text{O}_{9}$: High-pressure Raman study. Physical Review B, 2010, 81, . | 3.2 | 23 |
| 30 | Disorder-induced symmetry lowering in the CsInMgF_{6} pyrochlore crystal. Physical Review B, 2002, 66, . | 3.2 | 22 |
| 31 | High-pressure Raman scattering of MgMoO_{4} . Vibrational Spectroscopy, 2013, 68, 34-39. | 2.2 | 22 |
| 32 | Pressure-induced phase transitions in multiferroic $\text{RbFe}(\text{MoO}_{4})_{2}$ a Raman scattering study. Journal of Solid State Chemistry, 2011, 184, 2812-2817. | 2.9 | 21 |
| 33 | Novel bimetallic MOF phosphors with an imidazolium cation: structure, phonons, high- pressure phase transitions and optical response. Dalton Transactions, 2019, 48, 242-252. | 3.3 | 21 |
| 34 | Pressure-induced crystal-amorphous transformation in $\text{Y}_{2}\text{Mo}_{3}\text{O}_{12}$. Vibrational Spectroscopy, 2013, 68, 251-256. | 2.2 | 20 |
| 35 | Temperature- and pressure-dependent studies of niccolite-type formate frameworks of $[\text{NH}_{3}(\text{CH}_{2})_{4}\text{NH}_{3}][\text{M}_{2}(\text{HCOO})_{6}]$ ($\text{M} = \text{Zn}, \text{Co}, \text{Fe}$). Physical Chemistry Chemical Physics, 2016, 18, 27613-27622. | 2.8 | 19 |
| 36 | Heterometallic perovskite-type metal-organic framework with an ammonium cation: structure, phonons, and optical response of $[\text{NH}_{4}]\text{Na}_{0.5}\text{Cr}_{x}\text{Al}_{0.5-x}(\text{HCOO})_{3}$ ($x = 0, 0.025$ and 0.5). Physical Chemistry Chemical Physics, 2018, 20, 22284-22295. | 2.8 | 19 |

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|----|--|-----|-----------|
| 37 | Temperature-dependent Raman scattering study of K ₂ MoO ₄ . <i>Vibrational Spectroscopy</i> , 2012, 58, 87-94. | 2.2 | 18 |
| 38 | Temperature studies of KH ₂ PO ₄ :Mn crystals using x-ray diffraction and polarized Raman scattering. <i>Physical Review B</i> , 2005, 72, . | 3.2 | 16 |
| 39 | Structural and vibrational properties of K ₃ Fe(MoO ₄) ₂ (Mo ₂ O ₇) a novel layered molybdate. <i>Journal of Physics Condensed Matter</i> , 2009, 21, 095402. | 1.8 | 16 |
| 40 | Lattice dynamics and pressure-induced phase transitions in $\hat{\pm}$ -BaTeMo ₂ O ₉ . <i>Journal of Physics Condensed Matter</i> , 2013, 25, 125404. | 1.8 | 15 |
| 41 | Temperature induced phase transformations on the Li ₂ MoO ₄ system studied by Raman spectroscopy. <i>Journal of Molecular Structure</i> , 2017, 1139, 119-124. | 3.6 | 15 |
| 42 | Pressure-enhanced ferroelectric polarisation in a polar perovskite-like [C ₂ H ₅ NH ₃] _{0.5} Cr _{0.5} (HCOO) ₃ metal-organic framework. <i>Journal of Materials Chemistry C</i> , 2019, 7, 8660-8668. | 1.8 | 14 |
| 43 | Pressure-induced phase transitions in antiferroelectric CsBi(MoO ₄) ₂ . <i>Journal of Raman Spectroscopy</i> , 2005, 36, 56-62. | 2.5 | 14 |
| 44 | Phonon properties, polymorphism, and amorphization of Dy ₂ Mo ₄ O ₁₅ under high hydrostatic pressure. <i>Physical Review B</i> , 2010, 82, . | 3.2 | 14 |
| 45 | Effect of Ni(II) doping on the structure of L-histidine hydrochloride monohydrate crystals. <i>Journal of Physics Condensed Matter</i> , 2008, 20, 275209. | 1.8 | 12 |
| 46 | A Raman scattering study of pressure-induced phase transitions in nanocrystalline Bi ₂ Mo ₆ O ₁₅ . <i>Journal of Physics Condensed Matter</i> , 2011, 23, 045401. | 1.8 | 11 |
| 47 | Room-temperature vibrational properties of the BiMn ₂ O ₅ mullite. <i>Vibrational Spectroscopy</i> , 2013, 66, 43-49. | 2.2 | 11 |
| 48 | Lattice dynamics and high-pressure Raman scattering studies of ferroelectric K ₂ MgWO ₂ (PO ₄) ₂ . <i>Physical Review B</i> , 2008, 78, . | 3.2 | 10 |
| 49 | Raman spectroscopy of d-methionine under high pressure. <i>Vibrational Spectroscopy</i> , 2014, 72, 57-61. | 2.2 | 10 |
| 50 | Near-zero thermal expansion and phase transition in In _{0.5} (ZrMg) _{0.75} Mo ₃ O ₁₂ . <i>Journal of Materials Research</i> , 2016, 31, 3240-3248. | 2.6 | 10 |
| 51 | Raman scattering studies of pressure-induced phase transitions in perovskite-like acetamidinium manganese formate. <i>Journal of Raman Spectroscopy</i> , 2018, 49, 312-316. | 2.5 | 10 |
| 52 | Stability and flexibility of heterometallic formate perovskites with the dimethylammonium cation: pressure-induced phase transitions. <i>Physical Chemistry Chemical Physics</i> , 2019, 21, 4200-4208. | 2.8 | 10 |
| 53 | Vibrational properties of Cs ₄ W ₁₁ O ₃₅ and Rb ₄ W ₁₁ O ₃₅ systems: high pressure and polarized Raman spectra. <i>Journal of Raman Spectroscopy</i> , 2011, 42, 474-481. | 2.5 | 9 |
| 54 | Pressure-induced phase transition on K ₂ MoO ₄ : A Raman scattering study and ab initio calculations. <i>Journal of Solid State Chemistry</i> , 2012, 196, 197-202. | 2.9 | 9 |

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|----|---|-----|-----------|
| 55 | Structural and vibrational properties of carbonophosphates: Na ₃ MCO ₃ PO ₄ (M = Mn, Fe, Co and Ni). Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy, 2019, 223, 117298. | 3.9 | 9 |
| 56 | Mechanism of Pressure-Induced Phase Transitions and Structure-Property Relations in Methylhydrazinium Manganese Hypophosphite Perovskites. Journal of Physical Chemistry C, 2021, 125, 10121-10129. | 3.1 | 9 |
| 57 | Temperature-dependent Raman scattering of KDP:Mn (0.9% weight of Mn) crystal. Journal of Raman Spectroscopy, 2010, 41, 1318-1322. | 2.5 | 7 |
| 58 | Temperature-dependent Raman scattering study on Cs ₄ W ₁₁ O ₃₅ and Rb ₄ W ₁₁ O ₃₅ systems. Journal of Solid State Chemistry, 2013, 199, 7-14. | 2.9 | 7 |
| 59 | Vibrational spectroscopy study and ab initio calculation on ZnMoO ₄ system. Journal of Molecular Structure, 2020, 1206, 127776. | 3.6 | 7 |
| 60 | Lattice dynamics calculations and high-pressure Raman spectra of the ZnMoO ₄ . Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy, 2020, 239, 118501. | 3.9 | 7 |
| 61 | Vibrational studies of hexagonal bronze systems: phonon calculation and high pressure induced phase transformation. Journal of Raman Spectroscopy, 2009, 40, 1150-1157. | 2.5 | 6 |
| 62 | High pressure Raman scattering study on Sm ₂ Mo ₄ O ₁₅ system. Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy, 2017, 174, 80-85. | 3.9 | 6 |
| 63 | Effect of Alkali and Trivalent Metal Ions on the High-Pressure Phase Transition of [C ₂ H ₅ NH ₃] ^I _{0.5} M ^{III} _{0.5} (HCOO) ₆ (M ^I = Na, K and M ^{III} = Cr, Al) Heterometallic Perovskites. Journal of Physical Chemistry C, 2020, 124, 6337-6348. | 3.1 | 6 |
| 64 | Investigation of phase transitions in LiK _{1-x} (NH ₄) _x SO ₄ mixed crystal. Solid State Communications, 1999, 109, 507-512. | 1.9 | 5 |
| 65 | Evaluating Al _{2-x} Ga _x W ₃ O ₁₂ system for thermal shock resistance. Journal of Solid State Chemistry, 2019, 277, 149-158. | 2.9 | 5 |
| 66 | A temperature-dependent Raman scattering and X-ray diffraction study of K ₂ Mo ₂ O ₇ ·H ₂ O and ab initio calculations of K ₂ Mo ₂ O ₇ . Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy, 2021, 263, 120184. | 3.9 | 5 |
| 67 | Vibrational properties of RbNd(WO ₄) ₂ : high pressure Raman study, structural and phonon calculations. Journal of Physics Condensed Matter, 2011, 23, 405901. | 1.8 | 4 |
| 68 | Lattice dynamics and high-pressure Raman scattering studies of CoTeMoO ₆ crystal. Vibrational Spectroscopy, 2016, 84, 153-158. | 2.2 | 4 |
| 69 | Local impurity-phase generation in laser irradiated Li _x Co _{0.9} Ga _{0.1} O ₂ . Chemical Physics Letters, 2004, 397, 520-526. | 2.6 | 3 |
| 70 | Ionic properties of an organic-inorganic sol-gel hybrid based on polydimethylsiloxane and tetraethoxysilane doped with sodium dodecyl sulfate. Journal of Applied Polymer Science, 2010, 115, 851-854. | 2.6 | 3 |
| 71 | Pressure-induced structural transformations in In _{2-x} Y _x (MoO ₄) ₂ . Journal of Applied Physics, 2003, 94, 014301. | 2.5 | 3 |
| 72 | Raman characterization of single-crystalline Ga _{0.96} Mn _{0.04} As:Zn nanowires realized by ion-implantation. Nanotechnology, 2019, 30, 335202. | 2.6 | 3 |

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|----|--|-----|-----------|
| 73 | Temperature and Pressure Dependent Phonon Dynamics Properties of Gallium Selenide Telluride. Journal of Raman Spectroscopy, 0, , . | 2.5 | 3 |
| 74 | High-pressure Raman scattering study of ferroelectric K_3 Physical Review B, 2010, 82, . | 3.2 | 2 |
| 75 | Evaluation of carrier density and mobility in Mn ion-implanted GaAs:Zn nanowires by Raman spectroscopy. Nanotechnology, 2020, 31, 205705. | 2.6 | 2 |
| 76 | Vibrational spectroscopy and lattice dynamic calculation on the MnMoO4 system. Journal of Solid State Chemistry, 2022, 311, 123105. | 2.9 | 2 |
| 77 | Structural, vibrational and magnetic properties of monoclinic La2FeMnO6 double perovskite. Vacuum, 2022, 202, 111140. | 3.5 | 2 |
| 78 | Lattice dynamics and high pressure properties of K ⁺ ionic conducting system KNbTeO ₆ . Journal of Raman Spectroscopy, 2020, 51, 2517-2524. | 2.5 | 1 |
| 79 | Two new low-temperature phase transitions in the Li(NH4)1-xNaxSO4 system. Phase Transitions, 2004, 77, 921-928. | 1.3 | 0 |
| 80 | Computer simulation of Na2ThF6 single crystals: prediction of a phase transition under hydrostatic pressures. Journal of Physics Condensed Matter, 2008, 20, 165202. | 1.8 | 0 |