# Matteo Calandra

List of Publications by Year
in descending order

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Version: 2024-02-01

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|---|---|---|---|
| 139 papers | 31,419 citations | 28274 55 h-index | 10734 138 g-index |
| 141 all docs | 141 docs citations | 141 times ranked | 27028 citing authors |

| # | Article | IF | Citations |
|---|---|---|---|
| 1 | QUANTUM ESPRESSO: a modular and open-source software project for quantum simulations of materials. Journal of Physics Condensed Matter, 2009, 21, 395502. | 1.8 | 18,183 |
| 2 | Advanced capabilities for materials modelling with Quantum ESPRESSO. Journal of Physics Condensed Matter, 2017, 29, 465901. | 1.8 | 4,303 |
| 3 | Phonon-mediated superconductivity in graphene by lithium deposition. Nature Physics, 2012, 8, 131-134. | 16.7 | 431 |
| 4 | High-Pressure Hydrogen Sulfide from First Principles: A Strongly Anharmonic Phonon-Mediated Superconductor. Physical Review Letters, 2015, 114, 157004. | 7.8 | 377 |
| 5 | Colloquium: Saturation of electrical resistivity. Reviews of Modern Physics, 2003, 75, 1085-1099. | 45.6 | 369 |
| 6 | Exponential Localization of Wannier Functions in Insulators. Physical Review Letters, 2007, 98, 046402. | 7.8 | 297 |
| 7 | Anharmonic free energies and phonon dispersions from the stochastic self-consistent harmonic approximation: Application to platinum and palladium hydrides. Physical Review B, 2014, 89, . | 3.2 | 264 |
| 8 | Chemically exfoliated single-layer<mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML"><mml:mrow><mml:msub><mml:mtext>MoS</mml:mtext><mml:mn>2</mml:mn> Stability, lattice dynamics, and catalytic adsorption from first principles. Physical Review B, 2013, 88, . | 3.2 | 240 |
| 9 | Theoretical Explanation of Superconductivity inC6Ca. Physical Review Letters, 2005, 95, 237002. | 7.8 | 223 |
| 10 | Quantum hydrogen-bond symmetrization in the superconducting hydrogen sulfide system. Nature, 2016, 532, 81-84. | 27.8 | 222 |
| 11 | Density functional perturbation theory for gated two-dimensional heterostructures: Theoretical developments and application to flexural phonons in graphene. Physical Review B, 2017, 96, . | 3.2 | 198 |
| 12 | Anomalous Highâ€Temperature Superconductivity in YH<sub>6</sub>. Advanced Materials, 2021, 33, e2006832. | 21.0 | 196 |
| 13 | Quantum crystal structure in the 250-kelvin superconducting lanthanum hydride. Nature, 2020, 578, 66-69. | 27.8 | 193 |
| 14 | Electron-phonon coupling and electron self-energy in electron-doped graphene: Calculation of angular-resolved photoemission spectra. Physical Review B, 2007, 76, . | 3.2 | 188 |
| 15 | Effect of dimensionality on the charge-density wave in few-layer<mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:mrow><mml:mn>2</mml:mn><mml:msub><mml:mrow><mml:mi>H</mml:mi><mml:mtext>-NbSe</mml:mtext> Physical Review B, 2009, 80, . | 3.2 | 184 |
| 16 | First-Principles Theory of Anharmonicity and the Inverse Isotope Effect in Superconducting Palladium-Hydride Compounds. Physical Review Letters, 2013, 111, 177002. | 7.8 | 173 |
| 17 | Charge-Density Wave and Superconducting Dome in<mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:msub><mml:mi>TiSe</mml:mi><mml:mn>2</mml:mn></mml:msub></mml:math>from Electron-Phonon Interaction. Physical Review Letters, 2011, 106, 196406. | 7.8 | 159 |
| 18 | Two-dimensional FrÃ¶hlich interaction in transition-metal dichalcogenide monolayers: Theoretical modeling and first-principles calculations. Physical Review B, 2016, 94, . | 3.2 | 155 |

| # | Article | IF | Citations |
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| 19 | Numerical study of the two-dimensional Heisenberg model using a Green function Monte Carlo technique with a fixed number of walkers. Physical Review B, 1998, 57, 11446-11456. | 3.2 | 154 |
| 20 | First-principles calculations of x-ray absorption in a scheme based on ultrasoft pseudopotentials: From<mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:mi>Î±</mml:mi></mml:math>-quartz to high-<mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:mrow><mml:msub><mml:mi>T</mml:mi><mml:mi>c</mml:mi></mml:msub></mml:mrow></mml:math>compou Physical Review B, 2009, 80, . | 3.2 | 150 |
| 21 | Electronâ€“Phonon Interactions and the Intrinsic Electrical Resistivity of Graphene. Nano Letters, 2014, 14, 1113-1119. | 9.1 | 149 |
| 22 | Phonon Dispersion and Lifetimes inMgB2. Physical Review Letters, 2003, 90, 095506. | 7.8 | 139 |
| 23 | First-principles theory of field-effect doping in transition-metal dichalcogenides: Structural properties, electronic structure, Hall coefficient, and electrical conductivity. Physical Review B, 2015, 91, . | 3.2 | 127 |
| 24 | Breakdown of Optical Phononsâ€™ Splitting in Two-Dimensional Materials. Nano Letters, 2017, 17, 3758-3763. | 9.1 | 127 |
| 25 | Wannier interpolation of the electron-phonon matrix elements in polar semiconductors: Polar-optical coupling in GaAs. Physical Review B, 2015, 92, . | 3.2 | 126 |
| 26 | Dissociation products and structures of solid<mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML"><mml:msub><mml:mi mathvariant="normal">H</mml:mi><mml:mn>2</mml:mn></mml:msub><mml:mi mathvariant="normal">S</mml:mi></mml:math>at strong compression. Physical Review B, 2016, 93, . | 3.2 | 119 |
| 27 | Effects of magnetism and doping on the electron-phonon coupling in<mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:mrow><mml:msub><mml:mrow><mml:mtext>BaFe</mml:mtext></mml:mrow><mml:mn>2</mml:mn></mml:msub></mml:mrow></mml:math> Physical Review B, 2010, 82, . | 3.2 | 112 |
| 28 | Adiabatic and nonadiabatic phonon dispersion in a Wannier function approach. Physical Review B, 2010, 82, . | 3.2 | 108 |
| 29 | Phonon-limited resistivity of graphene by first-principles calculations: Electron-phonon interactions, strain-induced gauge field, and Boltzmann equation. Physical Review B, 2014, 90, . | 3.2 | 105 |
| 30 | Second-order structural phase transitions, free energy curvature, and temperature-dependent anharmonic phonons in the self-consistent harmonic approximation: Theory and stochastic implementation. Physical Review B, 2017, 96, . | 3.2 | 100 |
| 31 | Strong anharmonicity induces quantum melting of charge density wave in<mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML"><mml:mrow><mml:mn>2</mml:mn><mml:mi>H</mml:mi><mml:mo>-</mml:mo></mml:math>TaSe2 at pressure. Physical Review B, 2015, 92, . | 3.2 | 93 |
| 32 | Evidence for Flat Bands near the Fermi Level in Epitaxial Rhombohedral Multilayer Graphene. ACS Nano, 2015, 9, 5432-5439. | 14.6 | 92 |
| 33 | Phonon Collapse and Second-Order Phase Transition in Thermoelectric SnSe. Physical Review Letters, 2019, 122, 075901. | 7.8 | 92 |
| 34 | Angular dependence of core hole screening in<mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:mrow><mml:msub><mml:mrow><mml:mtext>LiCoO</mml:mtext></mml:mrow><mml:mn>2</mml:mn></mml:r A<mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:mrow><mml:mtext>DFT</mml:mtext><mml:mo>+</mml:mo><mml:mi>U</mml:mi></mml:mrow></mml:math>c of the oxygen and cobalt<mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="i. | 3.2 | 90 |
| 35 | Anharmonic phonons in few-layer MoS<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>: Raman spectroscopy of ultralow energy compression and shear modes. Physical Review B, 2013, 87, . | 3.2 | 90 |
| 36 | Anharmonic phonon frequency shift inMgB2. Physical Review B, 2003, 68, . | 3.2 | 88 |

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| 37 | Density functional theory description of hole-trapping inSiO2: A self-interaction-corrected approach. Physical Review B, 2005, 71, . | 3.2 | 86 |
| 38 | Superconductivity from doping boron icosahedra. Physical Review B, 2004, 69, . | 3.2 | 84 |
| 39 | High-<mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:msub><mml:mi>T</mml:mi><mml:mi>c</mml:mi></mml:msub></mml:math>Superconductivity in Superhard Diamondlike<mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:msub><mml:mi>BC</mml:mi><mml:mn>5</mml:mn></mml:msub></mml:math>. Physical Review Letters, 2008, 101, 016401. | 7.8 | 81 |
| 40 | Intercalant and Intermolecular Phonon Assisted Superconductivity in K-Doped Picene. Physical Review Letters, 2011, 107, 137006. | 7.8 | 79 |
| 41 | Giant Nonadiabatic Effects in Layer Metals: Raman Spectra of Intercalated Graphite Explained. Physical Review Letters, 2008, 100, 226401. | 7.8 | 75 |
| 42 | Anharmonic effects in atomic hydrogen: Superconductivity and lattice dynamical stability. Physical Review B, 2016, 93, . | 3.2 | 75 |
| 43 | Intrinsic charge transfer gap in NiO from<mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:mtext>Ni</mml:mtext><mml:mtext>â€‰</mml:mtext><mml:mi>K</mml:mi></mml:math>-edge x-ray absorption spectroscopy. Physical Review B, 2009, 79, . | 3.2 | 72 |
| 44 | The stochastic self-consistent harmonic approximation: calculating vibrational properties of materials with full quantum and anharmonic effects. Journal of Physics Condensed Matter, 2021, 33, 363001. | 1.8 | 70 |
| 45 | Anharmonic suppression of charge density waves in 2<mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:mrow><mml:mi>H</mml:mi></mml:mrow></mml:math>-NbS<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>. Physical Review B, 2012, 86, . | 3.2 | 66 |
| 46 | First-principles calculations of phonon frequencies, lifetimes, and spectral functions from weak to strong anharmonicity: The example of palladium hydrides. Physical Review B, 2015, 91, . | 3.2 | 66 |
| 47 | Electronic band structure of Two-Dimensional <mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML"><mml:msub><mml:mi mathvariant="normal">WS</mml:mi><mml:mn>2</mml:mn></mml:msub></mml:math> /Graphene van der Waals Heterostructures. Physical Review B, 2018, 97, . | 3.2 | 63 |
| 48 | Strong anharmonicity in the phonon spectra of PbTe and SnTe from first principles. Physical Review B, 2018, 97, . | 3.2 | 63 |
| 49 | Quantum Enhancement of Charge Density Wave in NbS<sub>2</sub> in the Two-Dimensional Limit. Nano Letters, 2019, 19, 3098-3103. | 9.1 | 62 |
| 50 | Projector augmented wave calculation of x-ray absorption spectra at the<mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:msub><mml:mi>L</mml:mi><mml:mrow><mml:mn>2</mml:mn><mml:mo>,</mml:mo><mml:mn>3</mml:mn></mml:mrow></mml:msub></mml:math> edges. Physical Review B, 2013, 87, . | 3.2 | 59 |
| 51 | Charge Fluctuations Close to Phase Separation in the Two-Dimensionaltâ^'JModel. Physical Review Letters, 1998, 81, 5185-5188. | 7.8 | 58 |
| 52 | Possibility of superconductivity in graphite intercalated with alkaline earths investigated with density functional theory. Physical Review B, 2006, 74, . | 3.2 | 57 |
| 53 | Electronic structure of heavily doped graphene: The role of foreign atom states. Physical Review B, 2007, 76, . | 3.2 | 57 |
| 54 | Electronic and vibrational properties of<mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML"><mml:msub><mml:mi>TiSe</mml:mi><mml:mn>2</mml:mn></mml:msub></mml:math> the charge-density-wave phase from first principles. Physical Review B, 2015, 92, . | 3.2 | 56 |

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| 55 | Pressure and stress tensor of complex anharmonic crystals within the stochastic self-consistent harmonic approximation. Physical Review B, 2018, 98, . | 3.2 | 56 |
| 56 | Critical Role of the Exchange Interaction for the Electronic Structure and Charge-Density-Wave Formation in <mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:mrow><mml:msub><mml:mrow><mml:mi>TiSe</mml:mi></mml:mrow><mml:mrow><mml:mn>2</mml:mn></m Physical Review Letters, 2017, 119, 176401. | 7.8 | 55 |
| 57 | Electron-phonon coupling and phonon self-energy inMgB2: Interpretation ofMgB2Raman spectra. Physical Review B, 2005, 71, . | 3.2 | 53 |
| 58 | Phonon-Assisted Magnetic Mott-Insulating State in the Charge Density Wave Phase of Single-Layer <mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:mrow><mml:mn>1</mml:mn><mml:mi>T</mml:mi><mml:mtext>â^'</mml:mtext><mml:msub><mml:mrow><m Physical Review Letters, 2018, 121, 026401. | 7.8 | 52 |
| 59 | Magnetic gap opening in rhombohedral-stacked multilayer graphene from first principles. Physical Review B, 2017, 95, . | 3.2 | 48 |
| 60 | Anharmonic and non-adiabatic effects in MgB2: Implications for the isotope effect and interpretation of Raman spectra. Physica C: Superconductivity and Its Applications, 2007, 456, 38-44. | 1.2 | 47 |
| 61 | Electrical resistivity at large temperatures: Saturation and lack thereof. Physical Review B, 2002, 66, . | 3.2 | 46 |
| 62 | Dynamical properties of a strongly correlated model for quarter-filled layered organic molecular crystals. Physical Review B, 2003, 68, . | 3.2 | 46 |
| 63 | Electrochemical doping of few-layer ZrNCl from first principles: Electronic and structural properties in field-effect configuration. Physical Review B, 2014, 89, . | 3.2 | 46 |
| 64 | Flat electronic bands in long sequences of rhombohedral-stacked graphene. Physical Review B, 2018, 97, . | 3.2 | 46 |
| 65 | Weak anharmonic effects inMgB2: A comparative inelastic x-ray scattering and Raman study. Physical Review B, 2007, 75, . | 3.2 | 41 |
| 66 | Metal-insulator transition and charge ordering in the extended Hubbard model at one-quarter filling. Physical Review B, 2002, 66, . | 3.2 | 40 |
| 67 | Black metal hydrogen above 360â€‰GPa driven by proton quantum fluctuations. Nature Physics, 2021, 17, 63-67. | 16.7 | 40 |
| 68 | High-pressure phase diagram of hydrogen and deuterium sulfides from first principles: Structural and vibrational properties including quantum and anharmonic effects. Physical Review B, 2018, 97, . | 3.2 | 38 |
| 69 | Formation of hot-electron ensembles quasiequilibrated in momentum space by ultrafast momentum scattering of highly excited hot electrons photoinjected into the Î“ valley of GaAs. Physical Review B, 2016, 93, . | 3.2 | 37 |
| 70 | Weak Dimensionality Dependence and Dominant Role of Ionic Fluctuations in the Charge-Density-Wave Transition of <mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:mrow><mml:msub><mml:mrow><mml:mi>NbSe</mml:mi></mml:mrow><mml:mrow><mml:mn>2</mml:mn></m Physical Review Letters, 2020, 125, 106101. | 7.8 | 37 |
| 71 | Electronic structure and magnetic properties of few-layer Cr <sub>2</sub> Ge <sub>2</sub> Te <sub>6</sub> : the key role of nonlocal electronâ€“electron interaction effects. 2D Materials, 2019, 6, 045042. | 4.4 | 36 |
| 72 | Search for highTcin layered structures: The case of LiB. Physical Review B, 2007, 75, . | 3.2 | 35 |

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| 73 | Two-Dimensional Analysis of the Double-Resonant 2D Raman Mode in Bilayer Graphene. Physical Review Letters, 2014, 113, 187401. | 7.8 | 35 |
| 74 | Effect of electron doping on lattice instabilities in single-layer <mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML"><mml:mrow><mml:mn>1</mml:mn><mml:mi>H</mml:mi><mml:mo>’</mml:m Physical Review B, 2017, 95, . | 3.2 | 35 |
| 75 | Strong anharmonicity and high thermoelectric efficiency in high-temperature SnS from first principles. Physical Review B, 2019, 100, . | 3.2 | 35 |
| 76 | Polarized Resonant Inelastic X-Ray Scattering as an Ultrafine Probe of Excited States ofLa2CuO4. Physical Review Letters, 2006, 96, 077006. | 7.8 | 34 |
| 77 | From antiferromagnetism tod-wave superconductivity in the two-dimensionalt-Jmodel. Physical Review B, 2000, 61, R11894-R11897. | 3.2 | 30 |
| 78 | K-edge x-ray absorption spectra in transition-metal oxides beyond the single-particle approximation: Shake-up many-body effects. Physical Review B, 2012, 86, . | 3.2 | 30 |
| 79 | Thermodynamic stabilities of ternary metal borides: Anab initioguide for synthesizing layered superconductors. Physical Review B, 2008, 78, . | 3.2 | 29 |
| 80 | Possible phase separation and weak localization in the absence of a charge-density wave in single-phase<mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML"><mml:mrow><mml:mn>1</mml:mn><mml:mi>T</mml:mi></mml:mrow></mml:n xmlns:mml="http://www.w3.org/1998/Math/MathML"><mml:msub><mml:mi>VS</mml:mi><mml:mn>2</mml:mn></mml:msub></mml: Physical Review B, 2014, 89, . | 3.2 | 28 |
| 81 | van der Waals driven anharmonic melting of the 3D charge density wave in VSe2. Nature Communications, 2021, 12, 598. | 12.8 | 28 |
| 82 | Charge density wave and spin 1/2 insulating state in single layer 1T-NbS <sub>2</sub>. 2D Materials, 2019, 6, 035041. | 4.4 | 27 |
| 83 | Anharmonic enhancement of superconductivity in metallic molecularCmcaâ€‰â€‰â^'â€‰â€‰4 hydrogen at high pressure: a first-principles study. Journal of Physics Condensed Matter, 2016, 28, 494001. | 1.8 | 26 |
| 84 | Violation of Ioffe-Regel condition but saturation of resistivity of the high- T c cuprates. Europhysics Letters, 2003, 61, 88-94. | 2.0 | 25 |
| 85 | Local and nonlocal electron-phonon couplings in K<mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:msub><mml:mrow /><mml:mn>3</mml:mn></mml:msub></mml:math>picene and the effect of metallic screening. Physical Review B, 2012, 86, . | 3.2 | 25 |
| 86 | Anharmonicity and Doping Melt the Charge Density Wave in Single-Layer TiSe<sub>2</sub>. Nano Letters, 2020, 20, 4809-4815. | 9.1 | 24 |
| 87 | Understanding the Photomagnetic Behavior in Copper Octacyanomolybdates. Journal of Physical Chemistry A, 2012, 116, 8678-8683. | 2.5 | 23 |
| 88 | Density-functional calculation of static screening in two-dimensional materials: The long-wavelength dielectric function of graphene. Physical Review B, 2015, 91, . | 3.2 | 21 |
| 89 | Saturation of Electrical Resistivity in Metals at Large Temperatures. Physical Review Letters, 2001, 87, 266601. | 7.8 | 20 |
| 90 | Phonon softening inNaxCoO2â^™yH2O: Implications for the Fermi surface topology and the superconducting state. Physical Review B, 2006, 74, . | 3.2 | 19 |

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| 91 | First-principles determination of the Raman fingerprint of rhombohedral graphite. Physical Review Materials, 2017, 1, . | 2.4 | 19 |
| 92 | X-ray magnetic and natural circular dichroism from first principles: Calculation of <mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML"><mml:mi>K</mml:mi></mml:math> - and <mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML"><mml:msub><mml:mi>L</mml:mi><mml:mn>1</mml:mn></mml:msub></mml:m -edge spectra. Physical Review B, 2017, 96, . | 3.2 | 18 |
| 93 | Misfit Layer Compounds: A Platform for Heavily Doped 2D Transition Metal Dichalcogenides. Advanced Functional Materials, 2021, 31, 2007706. | 14.9 | 17 |
| 94 | Theory of the thickness dependence of the charge density wave transition in 1 T-TiTe<sub>2</sub>. 2D Materials, 2020, 7, 045032. | 4.4 | 17 |
| 95 | Electronic structure of pristine and Ni-substituted <mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML"><mml:mrow><mml:mi>La</mml:mi><mml:mi>Fe</mml:mi><mml:msub><mml:mi mathvariant="normal">O</mml:mi><mml:mn>3</mml:mn></mml:msub></mml:mrow></mml:math> from near edge x-ray absorption fine structure experiments and first-principles simulations. Physical | 3.6 | 17 |
| 96 | Maximum<mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:mrow><mml:msub><mml:mi>T</mml:mi><mml:mi>c</mml:mi></mml:msub></mml:mrow></mml:math>at the verge of a simultaneous order-disorder and lattice-softening transition in superconducting<mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:mrow><mml:msub><mml:mrow><mml:mtext>CaC</mml:mtext></mml:mrow><mml:mn>6</mml:mn></mml:ms Physical Review B, 2008, 78, . | 3.2 | 16 |
| 97 | Charge density waves go nano. Nature Nanotechnology, 2015, 10, 737-738. | 31.5 | 16 |
| 98 | Atomic and electronic structure of trilayer graphene/SiC(0001): Evidence of Strong Dependence on Stacking Sequence and charge transfer. Scientific Reports, 2016, 6, 33487. | 3.3 | 16 |
| 99 | Energy relaxation mechanism of hot-electron ensembles in GaAs: Theoretical and experimental study of its temperature dependence. Physical Review B, 2018, 97, . | 3.2 | 16 |
| 100 | Chiral Spin Texture in the Charge-Density-Wave Phase of the Correlated Metallic <mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:mrow><mml:mi>Pb</mml:mi><mml:mrow><mml:mo>/</mml:mo><mml:mi>Si</mml:mi><mml:mo stretchy="false">(</mml:mo><mml:mn>111</mml:mn><mml:mo) Tj ETQq0 0 0 rgBT /Overlock 10 Tf 50 372 Td (stretchy="false">)</mm | 7.8 | 15 |
| 101 | Neutron scattering study of the high-energy graphitic phonons in superconducting<mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:mrow><mml:msub><mml:mrow><mml:mtext>CaC</mml:mtext></mml:mrow><mml:mn>6</mml:mn></mml:msub></mml:mrow></mml:math> Physical Review B, 2011, 84, . | 3.2 | 14 |
| 102 | Comparative study of the phonons in nonsuperconducting BaC<mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:msub><mml:mrow /><mml:mn>6</mml:mn></mml:msub></mml:math>and superconducting CaC<mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:msub><mml:mrow /><mml:mn>6</mml:mn></mml:msub></mml:math>using inelastic x-ray scattering. Physical Review B, 2011, 84, . | 3.2 | 14 |
| 103 | Universal Increase in the Superconducting Critical Temperature of Two-Dimensional Semiconductors at Low Doping by the Electron-Electron Interaction. Physical Review Letters, 2015, 114, 077001. | 7.8 | 14 |
| 104 | Quantum effects in muon spin spectroscopy within the stochastic self-consistent harmonic approximation. Physical Review Materials, 2019, 3, . | 2.4 | 13 |
| 105 | Origin of superconductivity of CaC6 and of other intercalated graphites. Physica Status Solidi (B): Basic Research, 2006, 243, 3458-3463. | 1.5 | 12 |
| 106 | Phonon dispersion and low-energy anomaly in<mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:mrow><mml:msub><mml:mrow><mml:mtext>CaC</mml:mtext></mml:mrow><mml:mn>6</mml:mn></mml:ms inelastic neutron and x-ray scattering experiments. Physical Review B, 2010, 81, . | 3.2 | 12 |
| 107 | Superconductivity in metalâ€coated graphene. Physica Status Solidi (B): Basic Research, 2012, 249, 2544-2548. | 1.5 | 12 |
| 108 | Long-Range Rhombohedral-Stacked Graphene through Shear. Nano Letters, 2020, 20, 5017-5023. | 9.1 | 12 |

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| 111 | Multi-scale theoretical approach to X-ray absorption spectra in disordered systems: an application to the study of Zn(ii) in water. Physical Chemistry Chemical Physics, 2018, 20, 24775-24782. | 2.8 | 10 |
| 112 | Field-effect-driven half-metallic multilayer graphene. Physical Review B, 2018, 98, . | 3.2 | 10 |
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