

Victor Pardo

List of Publications by Year in descending order

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86
papers

2,603
citations

249298

26
h-index

223390

49
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88
all docs

88
docs citations

88
times ranked

3939
citing authors

#	ARTICLE	IF	CITATIONS
1	Low Valence Nickelates: Launching the Nickel Age of Superconductivity. <i>Frontiers in Physics</i> , 2022, 9, .	1.0	13
2	Increasing the number of topological nodal lines in semimetals via uniaxial pressure. <i>Scientific Reports</i> , 2021, 11, 10574.	1.6	0
3	Synthesis, engineering, and theory of 2D van der Waals magnets. <i>Applied Physics Reviews</i> , 2021, 8, .	5.5	41
4	Tuning Coherent-Phonon Heat Transport in $\text{LaCoO}_3/\text{SrTiO}_3$ Superlattices. <i>Journal of Physical Chemistry Letters</i> , 2021, 12, 11878-11885.	2.1	5
5	Controlled Two-Dimensional Ferromagnetism in $1\text{T}\ddot{\text{a}}\text{CrTe}_2$: The Role of Charge Density Wave and Strain. <i>Journal of Physical Chemistry C</i> , 2020, 124, 21047-21053.	1.5	29
6	Electronic structure and magnetic exchange interactions of Cr-based van der Waals ferromagnets. A comparative study between CrBr_3 and $\text{Cr}_2\text{Ge}_2\text{Te}_6$. <i>Journal of Materials Chemistry C</i> , 2020, 8, 13582-13589.	2.7	13
7	Effects of Sr doping on the electronic and spin-state properties of infinite-layer nickelates: Nature of holes. <i>Physical Review B</i> , 2020, 102, .	1.1	26
8	Understanding the lattice thermal conductivity of SrTiO_3 from an <i>ab initio</i> perspective. <i>Physical Review Materials</i> , 2020, 4, .	1.1	11
9	Absence of Ferromagnetism in VSe_2 Caused by Its Charge Density Wave Phase. <i>Journal of Physical Chemistry C</i> , 2019, 123, 27802-27810.	1.5	88
10	Apparent auxetic to non-auxetic crossover driven by Co^{2+} redistribution in CoFe_2O_4 thin films. <i>APL Materials</i> , 2019, 7, .	2.2	11
11	Addendum: Fermiology and electron dynamics of trilayer nickelate $\text{La}_4\text{Ni}_3\text{O}_{10}$. <i>Nature Communications</i> , 2018, 9, 1952.	5.8	0
12	Electron-Hole Bilayers in CrN/MgO Heterostructures. <i>Physical Review Letters</i> , 2018, 120, 177201.	1.5	14
13	Large orbital polarization in a metallic square-planar nickelate. <i>Nature Physics</i> , 2017, 13, 864-869.	6.5	135
14	<i>Ab initio</i> study of the strain dependence of thermopower in electron-doped SrTiO_3 . <i>Journal of Physics Condensed Matter</i> , 2017, 29, 065501.	0.7	0
15	Fermiology and electron dynamics of trilayer nickelate $\text{La}_4\text{Ni}_3\text{O}_{10}$. <i>Nature Communications</i> , 2017, 8, 704.	5.8	26
16	Electron doped layered nickelates: Spanning the phase diagram of the cuprates. <i>Physical Review Materials</i> , 2017, 1, .	0.9	37
17	Ferromagnetic and insulating behavior of LaCoO_3 films grown on a (001) SrTiO_3 substrate: A simple ionic picture explained <i>ab initio</i> . <i>Physical Review Materials</i> , 2017, 1, .	0.9	3
18	Quantum spin Hall effect in rutile-based oxide multilayers. <i>Physical Review B</i> , 2016, 94, .	1.1	11

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19	Analysis of the temperature dependence of the thermal conductivity of insulating single crystal oxides. APL Materials, 2016, 4, 104815.	2.2	51
20	Efficient thermoelectric materials using nonmagnetic double perovskites with d^0 filling. Physical Review B, 2016, 94, .	1.1	2
21	Charge ordering in $NiLa_4$ Physical Review B, 2016, 94, .	1.1	40
22	Nitride multilayers as a platform for parallel two-dimensional electron-hole gases: MgO/ScN(111). Physical Review B, 2016, 93, .	1.1	6
23	Dirac topological insulator in the d_{z^2} orbital of a honeycomb oxide. Physical Review B, 2016, 94, .		
24	<i>Ab initio</i> study of nontrivial topological phases in corundum-structured d^0		

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37	Linear bands, zero-momentum Weyl semimetal, and topological transition in skutterudite-structure pnictides. <i>Physical Review B</i> , 2012, 85, .	1.1	23
38	Enhanced thermoelectric response of hole-doped LaNiO_2 . <i>Physical Review B</i> , 2012, 86, .	1.1	13
39	Ab initio calculations. Physical Review B, 2012, 86, .	0.7	0
40	Study of the influence of nanoscale doping inhomogeneities in the phase separated state of $\text{La}_{1-x}\text{Ca}_x\text{MnO}_{3-x}$. <i>Journal of Physics Condensed Matter</i> , 2012, 24, 275503.	1.1	15
41	Heteroepitaxial growth of $\text{MgO}(111)$ thin films on AlO_3 . <i>Physical Review B</i> , 2012, 86, .	1.1	26
42	Effects of applied pressure in ZnV_2O_4 and evidences for a dimerized structure. <i>Journal of Applied Physics</i> , 2011, 109, 07E158.	1.1	0
43	Electronic structure analysis of the quasi-one-dimensional oxide $\text{Sr}_6\text{Co}_5\text{O}_{15}$ within the LDA+U method. <i>Journal of Applied Physics</i> , 2011, 109, 07E114.	1.1	3
44	Dielectric properties and magnetostriction of the collinear multiferroic spinel CdVO_2 . <i>Physical Review B</i> , 2011, 84, .	1.1	73
45	Dirac Point Degenerate with Massive Bands at a Topological Quantum Critical Point. <i>Physical Review Letters</i> , 2011, 106, 056401.	2.9	44
46	Electronic structure of VO_2 . <i>Physical Review B</i> , 2011, 84, .	1.1	17
47	Charge ordering, metal-insulator transition, and magnetism. <i>Physical Review B</i> , 2011, 84, .	1.1	49
48	Evaluation of compensated magnetism in LaVO_2 . <i>Physical Review B</i> , 2011, 84, .	1.1	9
49	Electron confinement, orbital ordering, and orbital moments in VCuO . <i>Physical Review B</i> , 2010, 81, .	1.1	25
50	High-pressure magnetic and structural properties of TiOX ($X=\text{Cl}, \text{Br}$). <i>Journal of Magnetism and Magnetic Materials</i> , 2010, 322, 1069-1071.	1.0	2
51	Study of the pressure effects in TiOCl by ab initio calculations. <i>Journal of Magnetism and Magnetic Materials</i> , 2010, 322, 1072-1075.	1.0	0
52	Influence of the Ca^{2+} inhomogeneity distribution in the physical properties of $\text{La}_{0.625}\text{Ca}_{0.375}\text{MnO}_3$. <i>Physica Status Solidi C: Current Topics in Solid State Physics</i> , 2010, 7, 2620-2623.	0.8	1
53	Questionable collapse of the bulk modulus in CrN . <i>Nature Materials</i> , 2010, 9, 284-284.	13.3	2
54	Quantum Confinement Induced Molecular Correlated Insulating State in LaNi_4O_8 . <i>Physical Review Letters</i> , 2010, 105, 266402.	1.1	7

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55	Insulator transition through a semi-Dirac point in oxide nanostructures: layers confined within VO_2 TiO_2 Physical Review B, 2010, 81, .	1.1	54
56	Fermiology and magnetism in weak itinerant ferromagnet CoS_2 : an <i>ab initio</i> study. Journal of Physics Condensed Matter, 2010, 22, 505602.	0.7	13
57	Half-Metallic Semi-Dirac-Point Generated by Quantum Confinement in TiO_2 Physical Review Letters, 2009, 102, 166803.	2.9	213
58	Reduction of the bulk modulus at high pressure CrN . Nature Materials, 2009, 8, 947-951.	13.3	154
59	Electronic structure of dimerized spinel ZnV_2O_4 . Journal of Magnetism and Magnetic Materials, 2009, 321, 679-681.	1.0	3
60	Enhanced Dimerization of TiOCl under Pressure: Spin-Peierls to Peierls Transition. Physical Review Letters, 2009, 102, 056406.	2.9	23
61	Compensated magnetism by design in double perovskite oxides. Physical Review B, 2009, 80, .	1.1	82
62	Tight-Binding Modeling and Low-Energy Behavior of the Semi-Dirac Point. Physical Review Letters, 2009, 103, 016402.	2.9	165
63	Pressure-induced metal-insulator transition in. Physica B: Condensed Matter, 2008, 403, 1639-1641.	1.3	13
64	Electronic structure of the antiferromagnetic phase of. Physica B: Condensed Matter, 2008, 403, 1636-1638.	1.3	13
65	Magnetism driven by anion vacancies in superconducting FeSe Physical Review B, 2008, 78, .	1.1	95
66	Homopolar Bond Formation in ZnV_2O_4 Close to a Metal-Insulator Transition. Physical Review Letters, 2008, 101, 256403.	2.9	60
67	Magnetism from p states in alkaline earth monoxides: Trends with varying N impurity concentration. Physical Review B, 2008, 78, .	1.1	77
68	Crystallographic and magnetic structure of $\text{SrCoO}_{2.5}$ Neutron study coupled with band-structure calculations. Physical Review B, 2008, 78, .	1.1	162
69	Effect of spin fluctuations on the thermodynamic and transport properties of the itinerant ferromagnet CoS_2 Physical Review B, 2008, 78, .	1.1	24
70	Ising-type behavior in the antiferromagnetic phase of BaCo_2O_7 from first-principles BaCo_2O_7 Physical Review B, 2008, 78, .	1.1	9
71	A_2VO_4 A_2VO_4 stretchy="false" V_2O_5 stretchy="false" V_2O_5 Physical Review B, 2008, 78, .	2.9	62
72	Ferromagnetic clusters in polycrystalline BaCoO_3 . Journal of Magnetism and Magnetic Materials, 2007, 316, e670-e673.	1.0	15

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73	Ising-type anisotropy and spin state transitions in from first-principles calculations. Journal of Magnetism and Magnetic Materials, 2007, 316, e656-e658.	1.0	0
74	First principle calculations of charge ordering in manganites. Journal of Magnetism and Magnetic Materials, 2007, 316, e734-e737.	1.0	3
75	Dynamic magnetic behavior of BaCoO ₃ quasi-one-dimensional perovskite. Physical Review B, 2006, 74, .	1.1	18
76	Possible non-collinear magnetic configurations in BaCoO ₃ . Physica B: Condensed Matter, 2006, 378-380, 556-557.	1.3	2
77	Orbital and magnetic structure of quasi-1D cobaltites: Ab initio calculations and experiment. Journal of Magnetism and Magnetic Materials, 2006, 300, 48-52.	1.0	3
78	First-principles study of the spin-state transitions in GdBaCo ₂ O _{5.5} . Physical Review B, 2006, 73, .	1.1	19
79	Bandstructure study of magnetic and orbital order in $\text{CaMn}_2\text{Si}_2\text{O}_{10}$. <small>xmlns:xocs="http://www.elsevier.com/xml/xocs/dtd" xmlns:xs="http://www.w3.org/2001/XMLSchema" xmlns:xsi="http://www.w3.org/2001/XMLSchema-instance" xmlns="http://www.elsevier.com/xml/ja/dtd" xmlns:ja="http://www.elsevier.com/xml/ja/dtd" xmlns:mml="http://www.w3.org/1998/Math/MathML" xmlns:tb="http://www.elsevier.com/xml/common/table/dtd"</small>	1.0	2
80	Influence of interfacial scattering and surface roughness on giant magnetoresistance in Fe/Cr trilayers using ab initio layer potentials. Journal of Magnetism and Magnetic Materials, 2005, 290-291, 392-395.	1.0	3
81	Ab initio electronic structure of rare earth orthoferrites. Journal of Magnetism and Magnetic Materials, 2005, 290-291, 396-399.	1.0	19
82	Effects of Sr doping on the magnetic properties of Sr _x Ba _{1-x} CoO ₃ . Applied Physics Letters, 2005, 86, 202507.	1.5	7
83	Evidence for magnetic clusters in BaCoO ₃ . Physical Review B, 2004, 70, .	1.1	25
84	Magnetic structure and orbital ordering in BaCoO ₃ from first-principles calculations. Physical Review B, 2004, 70, .	1.1	33
85	Geometry optimization and electronic structure of BaCoO ₃ . Solid State Communications, 2003, 128, 101-106.	0.9	13
86	Possible spin configurations and magnetism in BaCoO ₃ perovskite. International Journal of Quantum Chemistry, 2003, 91, 252-256.	1.0	9