

Mitsuaki Kawamura

List of Publications by Year in descending order

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Version: 2024-02-01

27
papers

5,504
citations

687363

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552781

26
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docs citations

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times ranked

8031
citing authors

#	ARTICLE	IF	CITATIONS
1	Open-source library for the shifted Krylov subspace method of the form $\frac{1}{\alpha} \sum_{k=0}^{n-1} \beta_k z^k$ Computer Physics Communications, 2021, 258, 107536.	7.5	2
2	Diverse densest binary sphere packings and phase diagram. <i>Physical Review E</i> , 2021, 103, 023307.	2.1	12
3	The Simplest Model for Doped Poly(3,4-ethylenedioxythiophene) (PEDOT): Single-crystalline EDOT Dimer Radical Cation Salts. <i>Chemistry - A European Journal</i> , 2021, 27, 6696-6700.	3.3	6
4	The Simplest Model for Doped Poly(3,4-ethylenedioxythiophene) (PEDOT): Single-crystalline EDOT Dimer Radical Cation Salts. <i>Chemistry - A European Journal</i> , 2021, 27, 6597-6597.	3.3	0
5	RESPACK: An ab initio tool for derivation of effective low-energy model of material. <i>Computer Physics Communications</i> , 2021, 261, 107781.	7.5	40
6	DCore: Integrated DMFT software for correlated electrons. <i>SciPost Physics</i> , 2021, 10, .	4.9	7
7	Methods for constructing parameter-dependent flat-band lattices. <i>Physical Review B</i> , 2021, 103, .	3.2	5
8	GdV ₆ Sn ₆ : A Multi-carrier Metal with Non-magnetic 3d-electron Kagome Bands and 4f-electron Magnetism. <i>Journal of the Physical Society of Japan</i> , 2021, 90, .	1.6	15
9	Scaling law for Rashba-type spin splitting in quantum-well films. <i>Physical Review B</i> , 2021, 104, .	3.2	1
10	Lattice dynamics in FeSe via inelastic x-ray scattering and first-principles calculations. <i>Physical Review B</i> , 2020, 101, .	3.2	7
11	Anisotropic Triangular Lattice Realized in Rhenium Oxychlorides A ₃ ReO ₅ Cl ₂ (A = Ba, Sr). <i>Inorganic Chemistry</i> , 2020, 59, 10025-10033.	4.0	8
12	Possible Kitaev Quantum Spin Liquid State in 2D Materials with $S_{\alpha} = \frac{1}{2} \sum_{\langle ij \rangle} \tau_{ij}^{\alpha} S_i \cdot S_j$ Physical Review Letters, 2020, 124, 087205.	7.8	88
13	Microscopic characterization of the superconducting gap function in $S_{\alpha} = \frac{1}{2} \sum_{\langle ij \rangle} \tau_{ij}^{\alpha} S_i \cdot S_j$ Physical Review B, 2020, 101, .	3.2	1
14	Benchmark of density functional theory for superconductors in elemental materials. <i>Physical Review B</i> , 2020, 101, .	3.2	25
15	Vapochromism induced by intermolecular electron transfer coupled with hydrogen-bond formation in zinc dithiolene complex. <i>Journal of Materials Chemistry C</i> , 2020, 8, 14939-14947.	5.5	11
16	Effect of spin fluctuations on superconductivity in V and Nb: A first-principles study. <i>Physical Review B</i> , 2020, 102, .	3.2	6
17	Molecular beam epitaxy of superconducting $S_{\alpha} = \frac{1}{2} \sum_{\langle ij \rangle} \tau_{ij}^{\alpha} S_i \cdot S_j$ thin films. <i>Physical Review Materials</i> , 2020, 4, .	3.2	1
18	One-dimensionalization by Geometrical Frustration in the Anisotropic Triangular Lattice of the 5d Quantum Antiferromagnet Ca ₃ ReO ₅ Cl ₂ . <i>Journal of the Physical Society of Japan</i> , 2019, 88, 044708.	1.6	14

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19	FermiSurfer: Fermi-surface viewer providing multiple representation schemes. Computer Physics Communications, 2019, 239, 197-203.	7.5	181
20	mVMCâ€™ Open-source software for many-variable variational Monte Carlo method. Computer Physics Communications, 2019, 235, 447-462.	7.5	62
21	Dirac Fermions in Borophene. Physical Review Letters, 2017, 118, 096401.	7.8	353
22	Quantum lattice model solver $\langle \text{mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" id="mml23" display="inline" overflow="scroll" altimg="si23.gif"} \rangle \langle \text{mml:mi mathvariant="script"} \rangle H \langle \text{mml:mi mathvariant="normal"} \rangle \uparrow \langle \text{mml:mi} \rangle \langle \text{mml:math} \rangle$. Computer Physics Communications, 2017, 217, 180-192.	7.5	67
23	Anisotropic superconducting gaps in $\langle \text{mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML"} \rangle \langle \text{mml:mrow} \rangle \langle \text{mml:msub} \rangle \langle \text{mml:mi} \rangle \text{YNi} \langle \text{mml:mi} \rangle \langle \text{mml:mn} \rangle 2 \langle \text{mml:mn} \rangle \langle \text{mml:math} \rangle \langle \text{mml:mi} \rangle \langle \text{mml:mn} \rangle 2 \langle \text{mml:mn} \rangle \langle \text{mml:msub} \rangle \langle \text{mml:mi} \rangle \langle \text{mml:math} \rangle$: A first-principles investigation. Physical Review B, 2017, 95, .	3.2	23
24	Advanced capabilities for materials modelling with Quantum ESPRESSO. Journal of Physics Condensed Matter, 2017, 29, 465901.	1.8	4,303
25	â€™Visibleâ€™ 5d Orbital States in a Pleochroic Oxychloride. Journal of the American Chemical Society, 2017, 139, 10784-10789.	13.7	21
26	First-principles study of the pressure and crystal-structure dependences of the superconducting transition temperature in compressed sulfur hydrides. Physical Review B, 2015, 91, .	3.2	141
27	Improved tetrahedron method for the Brillouin-zone integration applicable to response functions. Physical Review B, 2014, 89, .	3.2	92