

# HÃœseyÄ°n YasÄ°n Uzunok

## List of Publications by Year in descending order

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29  
papers

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citations

1039406

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docs citations

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times ranked

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citing authors

#	ARTICLE	IF	CITATIONS
1	Ab initio investigation of BCS-type superconductivity in LuNi <sub>2</sub> B <sub>2</sub> . Effects of spin-orbit coupling on the electron-phonon superconductivity in the cubic Laves-phase compounds. <i>Physical Review B</i> , 2017, 96, .	1.1	33
2	Calcium and calcium compounds. <i>Physical Review B</i> , 2017, 96, .	1.1	19
3	Investigating the normal state and superconducting state properties of orthorhombic and hexagonal ZrRuP: A first-principles study. <i>Physical Review B</i> , 2019, 100, .	1.1	14
4	The effect of spin orbit interaction for superconductivity in the noncentrosymmetric superconductor CaR <sub>3</sub> Si <sub>3</sub> . <i>Journal of Alloys and Compounds</i> , 2016, 681, 205-211.	2.8	13
5	Theoretical investigation of electron-phonon interaction in the orthorhombic phase of Mo <sub>2</sub> C. <i>Journal of Alloys and Compounds</i> , 2019, 788, 842-851.	2.8	13
6	Role of spin-orbit coupling in the physical properties of LaX <sub>2</sub> Tj <sub>2</sub> ETQq000rgBT/Overlock 10 Tf 50 537 Td. <i>Physical Review B</i> , 2016, 93, .	1.1	10
7	Theoretical investigation of superconductivity in SrPd <sub>2</sub> and SrPd <sub>2</sub> . <i>Physical Review B</i> , 2016, 93, .	1.1	10
8	The influence of spin orbit interaction on phonons and superconductivity in the noncentrosymmetric superconductors LaPt <sub>3</sub> Si and LaPt <sub>3</sub> . <i>Intermetallics</i> , 2017, 86, 1-10.	1.8	10
9	Physical properties and superconductivity of Heusler compound LiGa <sub>2</sub> Rh: A first-principles calculation. <i>Solid State Communications</i> , 2020, 311, 113859.	0.9	10
10	Ab initio investigation of spin orbit coupling effect on the physical properties of IrGe superconductor. <i>Intermetallics</i> , 2019, 106, 107-114.	1.8	9
11	The effect of spin orbit interaction on the physical properties of LaTSi <sub>3</sub> (T = Ir, Pd, and Rh): First-principles calculations. <i>Journal of Applied Physics</i> , 2017, 121, 193904.	1.1	7
12	Theoretical investigation of antisymmetric spin-orbit coupling effect on the physical properties of noncentrosymmetric BaPtSb superconductor. <i>Intermetallics</i> , 2019, 108, 109-116.	1.8	7
13	Probing physical properties and superconductivity of noncentrosymmetric superconductors LaPtGe and LaPtGe <sub>3</sub> : A first-principles study. <i>Computational Materials Science</i> , 2020, 185, 109949.	1.4	7
14	The effect of martensitic phase transition from cubic to tetragonal on the physical properties of V <sub>3</sub> Si superconductor. <i>Intermetallics</i> , 2018, 96, 25-32.	1.8	5
15	Physical properties of hexagonal BaPtAs with noncentrosymmetric SrPtSb-type and centrosymmetric YPtAs-type crystal structures: Effects of spin-orbit coupling. <i>Physical Review B</i> , 2019, 100, .	1.1	5
16	Theoretical investigation of superconductivity in the non-centrosymmetric SrPtGe <sub>3</sub> and CaPtSi <sub>3</sub> compounds. <i>Philosophical Magazine</i> , 2019, 99, 198-223.	0.7	5
17	Identification of specific phonon contributions in BCS-type superconductivity of boride-carbide crystals with a layer-like structure. <i>Solid State Communications</i> , 2015, 206, 1-5.	0.9	3
18	Ab initio investigation of electron-phonon interaction in LaSn <sub>3</sub> and CaSn <sub>3</sub> . <i>Philosophical Magazine Letters</i> , 2018, 98, 375-391.	0.5	3

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19	Theoretical investigation of superconductivity mechanism in the filled skutterudites YRu <sub>4</sub> P <sub>12</sub> , YOs <sub>4</sub> P <sub>12</sub> , LaOs <sub>1</sub> P <sub>12</sub> and LaOs <sub>4</sub> As <sub>12</sub> . Journal of Physics and Chemistry of Solids, 2019, 130, 197-209.	1.9	3
20	First-principles calculations of physical properties and superconductivity of orthorhombic Mo <sub>2</sub> BC and Nb <sub>2</sub> BN. Journal of Applied Physics, 2021, 130, 153902.	1.1	3
21	Probing the physical and superconducting properties of hexagonal ZrRuAs: A first-principles calculation. Physica C: Superconductivity and Its Applications, 2020, 577, 1353715.	0.6	2
22	Probing the electron-phonon interaction in superconductivity for KSn <sub>2</sub> using the Migdal-Eliashberg theory and linear-response theory. Philosophical Magazine Letters, 2020, 100, 33-54.	0.5	2
23	Ab initio investigation of physical properties of LaT <sub>2</sub> B <sub>2</sub> C (T=Ir, Rh) compounds: A density functional theory approach. Physica C: Superconductivity and Its Applications, 2020, 568, 1353585.	0.6	1
24	Physical properties and superconductivity in the cubic A15-type Ta $\text{d}_{1e553}$ compound: A first principles study. Solid State Communications, 2022, 353, 114838.	0.9	1
25	First-principles calculations of physical properties and superconductivity of orthorhombic ScRuSi and ZrRhSi. Physical Review B, 2020, 102, .	1.1	0
26	A first-principles investigation of physical properties and superconductivity of NbPS. Solid State Sciences, 2020, 103, 106183.	1.5	0
27	The effect of spin-orbit interaction on superconductivity in the filled skutterudites MPt <sub>4</sub> Ge <sub>12</sub> (M=Ba, Sr and Th). Philosophical Magazine, 2020, 100, 2735-2758.	0.7	0
28	Impact of spin-orbit coupling on the physical properties and superconductivity of Ir-rich superconductor Mg <sub>2</sub> Ir <sub>3</sub> Si: A first-principles investigation. Journal of Physics and Chemistry of Solids, 2021, 153, 110030.	1.9	0
29	Elucidating the underlying mechanism of relatively high T <sub>c</sub> value of the orthorhombic MoRuP: a first-principles study. Philosophical Magazine, 2021, 101, 2054-2076.	0.7	0