

# Shinta Watanabe

## List of Publications by Citations

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38

papers

279

citations

10

h-index

14

g-index

42

ext. papers

325

ext. citations

2.8

avg, IF

3.03

L-index

#	Paper	IF	Citations
38	Calculations of Complete 4f and 4f-5d Energy Level Schemes of Free Trivalent Rare-Earth Ions. <i>Japanese Journal of Applied Physics</i> , <b>2004</b> , 43, L611-L613	1.4	26
37	Optical Transitions near the Fundamental Absorption Edge and Electronic Structures of YAl <sub>3</sub> (BO <sub>3</sub> ) <sub>4</sub> :Gd <sup>3+</sup> . <i>Japanese Journal of Applied Physics</i> , <b>2006</b> , 45, 146-151	1.4	22
36	First-principles relativistic calculation for 4f-5d transition energy of Ce <sup>3+</sup> in various fluoride hosts. <i>Journal of Solid State Chemistry</i> , <b>2006</b> , 179, 2438-2442	3.3	20
35	First-principles and experimental analysis of 4f-5d absorption spectra and multiplet energy levels of Pr <sup>3+</sup> , Nd <sup>3+</sup> , and U <sup>3+</sup> in LiYF <sub>4</sub> . <i>Physical Review B</i> , <b>2010</b> , 81,	3.3	16
34	[H(x)TeV <sub>9</sub> O <sub>28</sub> ](5-x)· (x=1 and 2): vanadotellurates with decavanadate structure. <i>Inorganic Chemistry</i> , <b>2011</b> , 50, 6183-8	5.1	16
33	First-principles calculation of ground and excited-state absorption spectra of ruby and alexandrite considering lattice relaxation. <i>Physical Review B</i> , <b>2009</b> , 79,	3.3	16
32	Relativistic Calculations of Complete 4f Energy Level Schemes of Free Trivalent Rare-Earth Ions. <i>Japanese Journal of Applied Physics</i> , <b>2005</b> , 44, 7488-7490	1.4	16
31	Chapter 231 First-principles calculations of transition spectra. <i>Fundamental Theories of Physics</i> , <b>2007</b> , 1-59	0.8	12
30	Relativistic many-electron calculations of Cr <sup>3+</sup> L <sub>2,3</sub> -edge x ray absorption near-edge structures for Cr <sup>3+</sup> :Al <sub>2</sub> O <sub>3</sub> and Cr <sup>3+</sup> :Cr <sub>2</sub> O <sub>3</sub> and magnetic circular dichroism of Cr <sup>3+</sup> L <sub>2,3</sub> -edge x ray absorption near-edge structures for Cr <sup>3+</sup> :Al <sub>2</sub> O <sub>3</sub> . <i>Journal of Applied Physics</i> , <b>2011</b> , 110, 123524	2.5	11
29	Experimental and Theoretical Investigations for Excitation Properties of Ba <sub>1-x</sub> Eu <sub>x</sub> MgAl <sub>10</sub> O <sub>17</sub> . <i>Journal of the Electrochemical Society</i> , <b>2007</b> , 154, J196	3.9	11
28	Experimental and First-Principles Analysis of 4f-5d Absorption Spectrum for Ce <sup>3+</sup> in LiYF <sub>4</sub> Considering Lattice Relaxation. <i>Journal of the Physical Society of Japan</i> , <b>2008</b> , 77, 084702	1.5	10
27	Control of electric, optical, thermal properties of C <sub>60</sub> films by electron-beam irradiation. <i>Carbon</i> , <b>2019</b> , 152, 882-887	10.4	9
26	Intra- and inter-atomic optical transitions of Fe, Co, and Ni ferrocyanides studied using first-principles many-electron calculations. <i>Journal of Applied Physics</i> , <b>2016</b> , 119, 235102	2.5	9
25	Shallow electron traps formed by Gd <sup>2+</sup> ions adjacent to oxygen vacancies in cerium-doped Gd <sub>3</sub> Al <sub>2</sub> Ga <sub>3</sub> O <sub>12</sub> crystals. <i>Applied Physics Letters</i> , <b>2018</b> , 113, 041906	3.4	8
24	Spectroscopic and first-principles calculation studies of the chemical forms of palladium ion in nitric acid solution for development of disposal of high-level radioactive nuclear wastes. <i>AIP Advances</i> , <b>2018</b> , 8, 045221	1.5	8
23	Energy location of Ce <sup>3+</sup> 4f level and majority carrier type in Gd <sub>3</sub> Al <sub>2</sub> Ga <sub>3</sub> O <sub>12</sub> :Ce crystals studied by surface photovoltage spectroscopy. <i>Applied Physics Letters</i> , <b>2017</b> , 110, 251101	3.4	7
22	Two-Dimensional Organometallic Kondo Lattice with Long-Range Antiferromagnetic Order. <i>Journal of Physical Chemistry C</i> , <b>2018</b> , 122, 20046-20054	3.8	7

21	Chemical forms of molybdenum ion in nitric acid solution studied using liquid-phase X-ray absorption fine structure, Ultraviolet-Visible absorption spectroscopy and first-principles calculations. <i>Chemical Physics Letters</i> , <b>2019</b> , 723, 76-81	2.5	6
20	Spectroscopic and theoretical studies on the structural, electronic, and optical properties of zinc octaethylporphyrin/C co-deposited films. <i>Journal of Chemical Physics</i> , <b>2017</b> , 147, 214701	3.9	6
19	Morphological and optical properties of $\alpha$ and $\beta$ phase zinc (II) phthalocyanine thin films for application to organic photovoltaic cells. <i>Journal of Chemical Physics</i> , <b>2020</b> , 153, 144704	3.9	6
18	Current Situation and Future Development of Discrete Variational Multielectron Method. <i>Advances in Quantum Chemistry</i> , <b>2008</b> , 54, 297-314	1.4	5
17	Visualizing cation vacancies in Ce:Gd <sub>3</sub> Al <sub>2</sub> Ga <sub>3</sub> O <sub>12</sub> scintillators by gamma-ray-induced positron annihilation lifetime spectroscopy. <i>Applied Physics Express</i> , <b>2020</b> , 13, 085505	2.4	4
16	Visualizing hidden electron trap levels in Gd <sub>3</sub> Al <sub>2</sub> Ga <sub>3</sub> O <sub>12</sub> :Ce crystals using a mid-infrared free-electron laser. <i>Applied Physics Letters</i> , <b>2018</b> , 112, 031112	3.4	4
15	Molecular orbital calculations of Eu-doped SrAl <sub>2</sub> O <sub>4</sub> clusters. <i>Solid State Communications</i> , <b>2015</b> , 206, 42-45	4.6	4
14	Luminescence properties of YAl <sub>3</sub> (BO <sub>3</sub> ) <sub>4</sub> substituted with Sc <sup>3+</sup> ions. <i>Physica Status Solidi (A) Applications and Materials Science</i> , <b>2006</b> , 203, 2701-2704	1.6	3
13	Correlational analysis of Eu <sup>3+</sup> charge transfer state using La effective charge in La-based mixed-anion host compounds. <i>Japanese Journal of Applied Physics</i> , <b>2017</b> , 56, 032601	1.4	2
12	Comparative study of Auger-free luminescence of Rb <sub>2</sub> ZnCl <sub>4</sub> crystals between experiment and calculation. <i>Physica Status Solidi C: Current Topics in Solid State Physics</i> , <b>2013</b> , 10, 993-996		2
11	Sorption Properties of Aluminum Hexacyanoferrate for Platinum Group Elements. <i>Chemistry Letters</i> , <b>2020</b> , 49, 83-86	1.7	2
10	Electron-beam irradiation of photopolymerized C <sub>60</sub> film studied using in situ scanning tunneling microscope, in situ Fourier-transform infrared spectroscopy, and first-principles calculations. <i>AIP Advances</i> , <b>2020</b> , 10, 085212	1.5	2
9	Local environment of W and Mo atoms in CaW <sub>1-x</sub> Mo <sub>x</sub> O <sub>4</sub> (x = 0.12) solid solution studied by X-ray structural analyzes. <i>Japanese Journal of Applied Physics</i> , <b>2019</b> , 58, 120602	1.4	2
8	Theoretical and Experimental Consideration of Valence Band X-ray Photoelectron Spectroscopy Spectra of Cr-Deficient MgCr <sub>2-x</sub> O <sub>4</sub> . <i>Japanese Journal of Applied Physics</i> , <b>2007</b> , 46, 4175-4178	1.4	1
7	Chemical forms of rhodium ion in pure water and nitric acid solution studied using ultraviolet-visible spectroscopy and first-principles calculations. <i>IOP Conference Series: Materials Science and Engineering</i> , <b>2020</b> , 835, 012001	0.4	1
6	Structural analyses of Gd <sub>3</sub> (Al,Ga) <sub>5</sub> O <sub>12</sub> garnet solid solutions via X-ray and UV absorption spectroscopy experiments for Gd atoms. <i>Journal of Alloys and Compounds</i> , <b>2021</b> , 867, 159055	5.7	1
5	Photopolymerization effects on the external quantum efficiency of fullerene/zinc phthalocyanine heterojunction solar cells. <i>AIP Advances</i> , <b>2021</b> , 11, 075227	1.5	1
4	The uptake mechanism of palladium ions into Prussian-blue nanoparticles in a nitric acid solution toward application for the recycling of precious metals from electronic and nuclear wastes.. <i>RSC Advances</i> , <b>2021</b> , 11, 20701-20707	3.7	1

- 3 The uptake characteristics of Prussian-blue nanoparticles for rare metal ions for recycling of precious metals from nuclear and electronic wastes.. *Scientific Reports*, **2022**, 12, 5135 4.9 1
- 2 Local structure analysis of Sb, Bi, and Ag dopant atoms in Mg<sub>2</sub>Si semiconductor by x-ray absorption spectroscopy and first-principles calculation. *Journal of Applied Physics*, **2021**, 130, 245105 2.5 0
- 1 Comparative Study on Optical Properties of YPO<sub>4</sub>: Mn, Zr Phosphor by Experiment and Calculation **2015**, 217-235