

Songming Wan

List of Publications by Year in descending order

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papers

303
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933447

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16
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docs citations

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times ranked

227
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#	ARTICLE	IF	CITATIONS
1	Room temperature Raman spectroscopy and ^{29}Si MAS NMR combined with high temperature Raman spectroscopy and DFT calculation of $x\text{MgO}-(1-x)\text{CaO}$ - SiO_2 glasses and melts. <i>Ceramics International</i> , 2022, 48, 4911-4920.	4.8	5
2	The structural origin of the $15\frac{1}{4}\mu\text{m}$ residual absorption in the BaGa_4Se_7 crystal. <i>Journal of Materials Chemistry C</i> , 2022, 10, 649-654.	5.5	3
3	Effect of MgO on the structure of SiO_2 -poor/rich MgO - CaO - SiO_2 melts by in situ high temperature time-gated Raman spectroscopy and theoretical calculation. <i>Journal of Raman Spectroscopy</i> , 2022, 53, 1635-1646.	2.5	2
4	Fine structures and their impacts on the characteristic Raman spectra of molten binary alkali tungstates. <i>Journal of Raman Spectroscopy</i> , 2021, 52, 1452-1461.	2.5	3
5	Quantitative analysis on the microstructure of molten binary KF - AlF_3 system by in situ Raman spectroscopy assisted with first principles method. <i>Journal of Raman Spectroscopy</i> , 2020, 51, 187-192.	2.5	8
6	Structural investigations on two typical lithium germanate melts by in situ Raman spectroscopy and density functional theory calculations. <i>CrystEngComm</i> , 2020, 22, 701-707.	2.6	9
7	Refractive properties of the $\pm\text{-BaGeO}_3$ crystal and their origins: a density functional theory study. <i>CrystEngComm</i> , 2020, 22, 6620-6625.	2.6	2
8	Quantitative studies on the structure of $x\text{CaO}$ -($1-x$) SiO_2 glasses and melts by in-situ Raman spectroscopy, ^{29}Si MAS NMR and quantum chemistry ab initio calculation. <i>Journal of Non-Crystalline Solids</i> , 2020, 546, 120252.	3.1	10
9	BaGeO_3 : A Mid-IR Transparent Crystal with Superstrong Raman Response. <i>Inorganic Chemistry</i> , 2020, 59, 3542-3545.	4.0	3
10	In Situ Raman Spectroscopy and DFT Studies of the Li_2GeO_3 Melt Structure. <i>Inorganic Chemistry</i> , 2019, 58, 5025-5030.	4.0	19
11	Temperature Dependent Micro-Structure of KAlF_4 from Solid to Molten States. <i>Materials</i> , 2018, 11, 1846.	2.9	4
12	Synthesis of CuInTe_2 nanowires: A polycrystalline-to-single-crystalline transformation process. <i>Journal of the American Ceramic Society</i> , 2018, 101, 5358-5362.	3.8	0
13	Quantitative Studies on the Structure of Molten Binary Potassium Molybdates by in Situ Raman Spectroscopy and Quantum Chemistry ab Initio Calculations. <i>Analytical Chemistry</i> , 2018, 90, 9085-9092.	6.5	11
14	In situ high-temperature Raman spectroscopic studies of the vibrational characteristics and microstructure evolution of sodium tungstate dihydrate crystal during heating and melting. <i>Journal of Raman Spectroscopy</i> , 2018, 49, 1693-1705.	2.5	15
15	Micro-structure studies of the molten binary K_3AlF_6 - Al_2O_3 system by in situ high temperature Raman spectroscopy and theoretical simulation. <i>Inorganic Chemistry Frontiers</i> , 2018, 5, 1861-1868.	6.0	37
16	Investigation on the Structure of a LiB_3O_5 - $\text{Li}_2\text{Mo}_3\text{O}_{10}$ High-Temperature Solution for Understanding the $\text{Li}_2\text{Mo}_3\text{O}_{10}$ Flux Behavior. <i>Inorganic Chemistry</i> , 2017, 56, 3623-3630.	4.0	9
17	In-situ studies on the micro-structure evolution of $\text{A}_2\text{W}_2\text{O}_7$ ($\text{A} = \text{Li, Na, K}$) during melting by high temperature Raman spectroscopy and density functional theory. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2017, 185, 188-196.	3.9	18
18	Structural studies of a Li_2O - $4\text{B}_2\text{O}_3$ melt by high-temperature Raman spectroscopy and density functional theory. <i>CrystEngComm</i> , 2017, 19, 5721-5726.	2.6	4

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19	Raman and Density Functional Theory Studies of $\text{Li}_2\text{Mo}_4\text{O}_{13}$ Structures in Crystalline and Molten States. <i>Inorganic Chemistry</i> , 2017, 56, 14129-14134.	4.0	11
20	Raman Spectral and Density Functional Theory Analyses of the CsB_3O_5 Melt Structure. <i>Inorganic Chemistry</i> , 2016, 55, 7098-7102.	4.0	10
21	Structural analyses of a K_2O -rich KNbO_3 melt and the mechanism of KNbO_3 crystal growth. <i>CrystEngComm</i> , 2015, 17, 2636-2641.	2.6	15
22	In situ Raman investigation of a LiB_3O_5 melt toward understanding the structural memory phenomena. <i>CrystEngComm</i> , 2015, 17, 9357-9362.	2.6	6
23	Raman spectroscopy and density functional theory analyses of the melt structure in a $\text{Li}_2\text{B}_4\text{O}_7$ crystal growth system. <i>CrystEngComm</i> , 2014, 16, 3086-3090.	2.6	18
24	New insights into the BiB_3O_6 melt structure. <i>CrystEngComm</i> , 2013, 15, 995-1000.	2.6	14
25	Structural investigation of Li_2O - B_2O_3 - MoO_3 glasses and high-temperature solutions: toward understanding the mechanism of flux-induced growth of lithium triborate crystal. <i>CrystEngComm</i> , 2013, 15, 356-364.	2.6	11
26	In situ investigation of the microstructure of $\text{KGd}(\text{WO}_4)_2$ crystal growth boundary layer by confocal laser Raman microscopy. <i>CrystEngComm</i> , 2012, 14, 8722.	2.6	2
27	High temperature Raman spectroscopy study on the microstructure of the boundary layer around a growing LiB_3O_5 crystal. <i>CrystEngComm</i> , 2011, 13, 5239.	2.6	18
28	Segregation during crystal growth from melt and absorption cross section determination by optical absorption method. <i>Science in China Series G: Physics, Mechanics and Astronomy</i> , 2008, 51, 481-491.	0.2	13
29	Raman Spectroscopy Study on CsB_3O_5 Crystal Melt Boundary Layer Structure. <i>Crystal Growth and Design</i> , 2008, 8, 412-414.	3.0	23
30	A thermodynamic criterion for the choice of flux and its validity in NaBO_2 -fluxed $\text{Li}_2\text{-BaB}_2\text{O}_4$ crystal growth. <i>CrystEngComm</i> , 0, , .	2.6	0