

# Sho Hayakawa

## List of Publications by Year in descending order

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11  
papers

136  
citations

1307594

7  
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1281871

11  
g-index

11  
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docs citations

11  
times ranked

77  
citing authors

#	ARTICLE	IF	CITATIONS
1	Temperature-dependent mechanisms of dislocation–twin boundary interactions in Ni-based equiatomic alloys. <i>Acta Materialia</i> , 2021, 211, 116886.	7.9	28
2	Atomistic simulations of grain boundary energies in austenitic steel. <i>Journal of Materials Science</i> , 2019, 54, 5570-5583.	3.7	20
3	Screw dislocation–spherical void interactions in fcc metals and their dependence on stacking fault energy. <i>Journal of Materials Science</i> , 2019, 54, 11509-11525.	3.7	19
4	Effects of stacking fault energies on the interaction between an edge dislocation and an 8.0-nm-diameter Frank loop of self-interstitial atoms. <i>Nuclear Materials and Energy</i> , 2016, 9, 581-586.	1.3	16
5	Atomistic simulations for the effects of stacking fault energy on defect formations by displacement cascades in FCC metals under Poisson’s deformation. <i>Journal of Materials Science</i> , 2019, 54, 11096-11110.	3.7	12
6	Molecular dynamic simulations evaluating the effect of the stacking fault energy on defect formations in face-centered cubic metals subjected to high-energy particle irradiation. <i>Computational Materials Science</i> , 2021, 195, 110479.	3.0	10
7	Interactions between clusters of self-interstitial atoms via a conservative climb in BCC–Fe. <i>Philosophical Magazine</i> , 2018, 98, 2311-2325.	1.6	8
8	Atomistic modeling of meso-timescale processes with SEAKMC: A perspective and recent developments. <i>Computational Materials Science</i> , 2021, 194, 110390.	3.0	8
9	Behavior of a self-interstitial-atom type dislocation loop in the periphery of an edge dislocation in BCC-Fe. <i>Nuclear Materials and Energy</i> , 2016, 9, 592-597.	1.3	7
10	Saddle point sampling using scaled normal coordinates. <i>Computational Materials Science</i> , 2021, 200, 110785.	3.0	4
11	Interaction between a dislocation and nanotwin–hcp lamella in Ni-based concentrated alloys from atomistic simulations. <i>Scripta Materialia</i> , 2022, 218, 114810.	5.2	4