

# Sedigheh Bigdeli

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/6815764/publications.pdf>

Version: 2024-02-01

11  
papers

260  
citations

1163117

8  
h-index

1281871

11  
g-index

11  
all docs

11  
docs citations

11  
times ranked

179  
citing authors

#	ARTICLE	IF	CITATIONS
1	On the third-generation Calphad databases: An updated description of Mn. <i>Physica Status Solidi (B): Basic Research</i> , 2015, 252, 2199-2208.	1.5	52
2	New description of metastable hcp phase for unaries Fe and Mn: Coupling between first-principles calculations and CALPHAD modeling. <i>Physica Status Solidi (B): Basic Research</i> , 2016, 253, 1830-1836.	1.5	34
3	Thermodynamic evaluation of pure Co for the third generation of thermodynamic databases. <i>Physica Status Solidi (B): Basic Research</i> , 2017, 254, 1600231.	1.5	32
4	A New Description of Pure C in Developing the Third Generation of Calphad Databases. <i>Journal of Phase Equilibria and Diffusion</i> , 2018, 39, 832-840.	1.4	29
5	An insight into using DFT data for Calphad modeling of solid phases in the third generation of Calphad databases, a case study for Al. <i>Calphad: Computer Coupling of Phase Diagrams and Thermochemistry</i> , 2019, 65, 79-85.	1.6	29
6	A thermodynamic assessment of the binary Fe-Mn system for the third generation of Calphad databases. <i>Calphad: Computer Coupling of Phase Diagrams and Thermochemistry</i> , 2019, 64, 185-195.	1.6	26
7	High-temperature oxidation behaviour of Al <sub>x</sub> FeCrCoNi and AlTiVCr compositionally complex alloys. <i>Npj Materials Degradation</i> , 2020, 4, .	5.8	21
8	A method for handling the extrapolation of solid crystalline phases to temperatures far above their melting point. <i>Calphad: Computer Coupling of Phase Diagrams and Thermochemistry</i> , 2020, 68, 101737.	1.6	19
9	Thermodynamic assessment of binary erythritol-xylitol phase diagram for phase change materials design. <i>Calphad: Computer Coupling of Phase Diagrams and Thermochemistry</i> , 2018, 60, 29-36.	1.6	8
10	Strategies for High-Temperature Corrosion Simulations of Fe-Based Alloys Using the Calphad Approach: Part I. <i>Journal of Phase Equilibria and Diffusion</i> , 2021, 42, 403-418.	1.4	7
11	A Generalized Approach Obeying the Third Law of Thermodynamics for the Expression of Lattice Stability and Compound Energy: A Case Study of Unary Aluminum. <i>Journal of Phase Equilibria and Diffusion</i> , 2018, 39, 519-531.	1.4	3