SACADA Database Code: 574

Topology: 4^3T149-CA

# of independent nodes (IN): 3
Transitivity: [39(10)5]
Space Group: I2/m
Pearson: mS24
Coordination Number (CN): 4

Year: 2021

Data

Name	Pressure, GPa	Density, g/cm³	Gap, eV	Relative energy, eV/atom	Bulk, GPa	Shear, GPa	Vickers, GPa	Refs
4^3T149-CA (SACADA #574)		3.246		0.287	390.0	379.9	67.1	SACADA[1]
4^3T149-CA								doi: 10.1038/s41524-021-00491-y

Elasticity tensor (kBar)[1]

12004.0236	523.3074	770.9832	-0.0000	0.0000	-162.2944
523.3074	10177.3311	942.9098	0.0000	0.0000	-135.5444
770.9832	942.9098	8674.0866	-0.0000	-0.0000	70.9359
-0.0000	0.0000	-0.0000	4226.1766	-67.8130	0.0000
0.0000	-0.0000	-0.0000	-67.8130	2769.2523	0.0000
-162.2944	-135.5444	70.9359	0.0000	0.0000	3029.4570

[1] We apply the density functional theory (DFT) approach by using the Vienna Ab Initio Simulation Package (VASP) to calculate the total energy and properties of carbon allotropes.

DFT calculations

We apply the density functional theory (DFT) approach by using the Vienna Ab Initio Simulation Package (VASP) package [6] to calculate the total energy of carbon allotropes. The Generalized Gradient Approximation [7] (GGA) for exchange-correlational functional is used everywhere. The energy cutoff set to 600 eV. Fully automatic Γ-centered k-points mesh with a reciprocal-space resolution of $2\pi \times 0.025$ Å^{-1} is applied. We used tetrahedron method with Blöchl corrections to perform the k-point integration. The convergence thresholds are set at 10^{-6} eV for energy and 10^{-5} eV Å^{-1} for ionic forces. Polycrystalline elastic moduli — the bulk modulus, the shear modulus, Young's modulus, and the Poisson's ratio ν — have been calculated within the Voigt–Reuss–Hill [8] approximation. The Vicker's hardness H_v has been estimated according to Oganov's model [9].