SACADA Database Code: 258

Topology: 3^6T4

# of independent nodes (IN): 6
Transitivity: [6996]
Space Group: Pm-3m
Pearson: cP200
Coordination Number (CN): 3

Year: 2010

Data

Name	Pressure, GPa	Density, g/cm³	Gap, eV	Relative energy, eV/atom	Bulk, GPa	Shear, GPa	Vickers, GPa	Refs
3^6T4 (SACADA #258)		.960		0.888	74.0	23.1	3.5	SACADA[1]
c200		0.96	Metal		76.4			doi: 10.1088/0953-8984/22/33/334220
c200								doi: 10.3103/s1063457610020012

Elasticity tensor (kBar)[1]

1160.7072	529.3844	529.3844	-0.0000	0.0000	-0.0000
529.3844	1160.7072	529.3844	0.0000	-0.0000	0.0000
529.3844	529.3844	1160.7072	0.0000	0.0000	-0.0000
-0.0000	0.0000	0.0000	187.9857	-0.0000	0.0000
0.0000	-0.0000	-0.0000	-0.0000	187.9857	0.0000
-0.0000	0.0000	0.0000	0.0000	0.0000	187.9857

[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].