

Abstract Submitted
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Low-energy phonon dispersion in $\text{LaFe}_4\text{Sb}_{12}$ ANDREAS LEITHE-JASPER, Max Planck Institute for Chemical Physics of Solids, Dresden, Germany, MARTIN BOEHM, HANNU MUTKA, MICHAEL M. KOZA, Institut Laue-Langevin, Grenoble, France — We studied the vibrational dynamics of a single crystal of $\text{LaFe}_4\text{Sb}_{12}$ by three-axis inelastic neutron spectroscopy. The dispersion of phonons with wave vectors q along $[xx0]$ and $[xxx]$ directions in the energy range of eigenmodes with high amplitudes of lanthanum vibrations, i.e., at $\hbar\omega < 12$ meV is identified. Symmetry-avoided anticrossing dispersion of phonons is established in both monitored directions and distinct eigenstates at high-symmetry points and at the Brillouin-zone center are discriminated. The experimentally derived phonon dispersion and intensities are compared with and backed up by *ab initio* lattice dynamics calculations. results of the computer model match well with the experimental data.

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