

LETTER

Crystal chemistry of a high-pressure *C2/c* clinopyroxene with six-coordinated silicon

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ABSTRACT

A $(\text{Ca}_{0.36}\text{Na}_{0.56}\text{Mg}_{0.08})(\text{Mg}_{0.73}\text{Si}_{0.27})\text{Si}_2\text{O}_6$ clinopyroxene containing both four- and six-coordinated silicon was synthesized at 15 GPa and 1600 °C and its structure determined with single-crystal X-ray diffraction. Unlike the $\text{Na}(\text{Mg}_{0.5}\text{Si}_{0.5})\text{Si}_2\text{O}_6$ clinopyroxene that exhibits an ordered *P2/n* structure with octahedrally coordinated Mg^{2+} and Si^{4+} occupying two crystallographically distinct M1 sites (Angel et al. 1988), our sample possesses a *C2/c* symmetry and shows no detectable ordering between Mg^{2+} and Si^{4+} in the M1 site. The measured unit-cell parameters are $a = 9.5792(13)$, $b = 8.7588(12)$, $c = 5.2610(6)$ Å, $\beta = 107.199(3)^\circ$, and $V = 421.7(2)$ Å³. The crystal structure of the clinopyroxene reported in this study is comparable to that for omphacites formed under high temperatures and pressures.