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**Mengxianminite,  $\text{Ca}_2\text{Sn}_2\text{Mg}_3\text{Al}_8[(\text{BO}_3)(\text{BeO}_4)\text{O}_6]_2$ , a new borate mineral from Xianghualing skarn, Hunan Province, China**

Mengxianminite,  $\text{Ca}_2\text{Sn}_2\text{Mg}_3\text{Al}_8[(\text{BO}_3)(\text{BeO}_4)\text{O}_6]_2$ , is a new borate mineral from Xianghualing skarn, Hunan Province, southern China. It occurs in the Hsianghualite vein from this skarn, and is associated with fluorite, phlogopite, hsianghualite, magnetite, tourmaline, magnesiotaaffeite-2N2S and calcite. Mengxianminite forms subhedral to euhedral green crystals from 20 to 200  $\mu\text{m}$  long, translucent to transparent, with a vitreous luster. The crystals show perfect cleavage on {100} and good cleavage on {010}, and do not fluoresce in long- or short-wave ultraviolet light. The estimated Mohs hardness is 8, and the tenacity is brittle with irregular fracture. The calculated density is 4.17  $\text{g}/\text{cm}^3$ . Optically, mengxianminite is biaxial (-), with  $\alpha = 1.80(2)$ ,  $\beta = 1.83(2)$ ,  $\gamma = 1.84(2)$  (589 nm). Chemical analysis by electron microprobe (average of 6) gave  $\text{Al}_2\text{O}_3$  40.00,  $\text{SnO}_2$  25.96,  $\text{MgO}$  6.57,  $\text{CaO}$  8.56,  $\text{FeO}$  4.83,  $\text{B}_2\text{O}_3$  6.52,  $\text{BeO}$  4.68,  $\text{ZnO}$  1.81,  $\text{MnO}$  1.23,  $\text{Na}_2\text{O}$  1.13,  $\text{TiO}_2$  0.10,  $\text{SiO}_2$  0.04, sum 101.42 wt%. The empirical formula, calculated on the basis of 26O, 2Be and 2B atoms per formula unit, is:  $(\text{Ca}_{1.63}, \text{Na}_{0.39})_{\Sigma 2.02} (\text{Sn}_{1.84}, \text{Zn}_{0.24})_{\Sigma 2.08} (\text{Mg}_{1.74}, \text{Fe}_{0.72}, \text{Al}_{0.38}, \text{Mn}_{0.19}, \text{Ti}_{0.01})_{\Sigma 3.04} \text{Al}_8 [(\text{BO}_3)(\text{BeO}_4)\text{O}_6]_2$ . The stronger eight lines of the powder XRD pattern [d in Å (I)(hkl)] are: 3.000(35)(16 20); 2.931(100)(17 11); 2.475(29)(022); 2.430(30)(13 31); 2.375(100)(14 02/640); 2.028(52)(21 31); 1.807(35)(913); 1.530(98)(14 60/15 33). Mengxianminite is orthorhombic, space group *Fdd2*; unit-cell parameters refined from single-crystal X-ray diffraction data are:  $a = 60.689$  (3),  $b = 9.907$  (1),  $c = 5.740$  (1) Å,  $V = 3451.0$  (3) Å<sup>3</sup>,  $Z = 8$ . The structure of mengxianminite is composed of alternating O-T1-O-T2-O'-T2 layers stacked along the a axis, equal to two alternating modules: A module (O-T1-O) consists of the spinel modular and another O layer ( $\text{AlO}_6$  octahedra layer); B modular (T2-O'-T2) shows the simplified formula  $\text{CaSnAl}(\text{BeO}_4)(\text{BO}_3)$ ,  $\text{SnO}_6$  octahedra are isolated in the T2 layers, connected via  $\text{BeO}_4$  and  $\text{CaO}_{11}$  groups;  $\text{AlO}_6$  edge-sharing octahedra in the O' layer form chains running along the b axis; these chains are connected in the c direction by the  $\text{BO}_3$  triangular groups. Mengxianminite is of hydrothermal origin, crystallized during the late stage of the xianghualing skarn.

**Biography**

Can Rao has completed his PhD and Post-doctoral studies from Nanjing University, China.. He is an Associate Professor at Zhejiang University, China. He has found 3 new minerals (strontiohurlbutite, minjiangite and mengxianminite), which have been approved by IMA, and published 16 papers in reputed journals.

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