Siwaqaite, Ca₆Al₂(CrO₄)₃(OH)₁₂·26H₂O, a new mineral of the ettringite group from the pyrometamorphic Daba-Siwaqa complex, Jordan

RAFAŁ JUROSZEK^{1,*}, BILJANA KRÜGER², IRINA GALUSKINA¹, HANNES KRÜGER², YEVGENY VAPNIK³, AND EVGENY GALUSKIN¹

¹Institute of Earth Sciences, Faculty of Natural Sciences, University of Silesia, Będzińska 60, 41-205 Sosnowiec, Poland ²Institute of Mineralogy and Petrography, University of Innsbruck, Innrain 52, 6020 Innsbruck, Austria ³Department of Geological and Environmental Sciences, Ben-Gurion University of the Negev, POB 653, Beer-Sheva 84105, Israel

ABSTRACT

A new mineral, siwagaite, ideally $Ca_{6}Al_{2}(CrO_{4})_{3}(OH)_{12} \cdot 26H_{2}O[P31c, Z=2, a=11.3640(2) \text{ Å}, c=12.3640(2) \text{ Å}, c=12.3640(2)$ 21.4485(2) Å, V = 2398.78(9) Å³], a member of the ettringite group, was discovered in thin veins and small cavities within the spurrite marble at the North Siwaga complex, Lisdan-Siwaga Fault, Hashem region, Jordan. This complex belongs to the widespread pyrometamorphic rock of the Hatrurim Complex. The spurite marble is mainly composed of calcite, fluorapatite, and brownmillerite. Siwaqaite occurs with calcite and minerals of the baryte-hashemite series. It forms hexagonal prismatic crystals up to 250 µm in size, but most common are grain aggregates. Siwaqaite exhibits a canary yellow color and a yellowish-gray streak. The mineral is transparent and has a vitreous luster. It shows perfect cleavage on (10 $\overline{10}$). Parting or twinning is not observed. The calculated density of siwagaite is 1.819 g/cm³. Siwaqaite is optically uniaxial (-) with $\omega = 1.512(2)$, $\varepsilon = 1.502(2)$ (589 nm), and non-pleochroic. The empirical formula of the holotype siwagaite calculated on the basis of 8 framework cations and 26 water molecules is $Ca_{6,01}(Al_{1,87}Si_{0,12})_{\Sigma_{1,99}}[(CrO_4)_{1,71}(SO_4)_{1,13}(SeO_4)_{0,40}]_{\Sigma_{3,24}}(OH)_{11,63} \cdot 26H_2O$. X-ray diffraction (XRD), Raman, and infrared spectroscopy confirm the presence of OH⁻ groups and H₂O molecules and absence of $(CO_3)^{2-}$ groups. The crystal structure of this Cr^{6+} -analog of ettringite was solved by direct methods using single-crystal synchrotron XRD data. The structure was refined to an agreement index $R_1 = 4.54\%$. The crystal structure of siwagaite consists of $\{Ca_6[Al(OH)_6], 24H_2O\}^{6+}$ columns with the inter-column space (channels) occupied by $(CrO_4)^{2-}$, $(SO_4)^{2-}$, $(SO_4)^{2-}$, and $(SO_3)^{2-}$ groups and H₂O molecules. The tetrahedrally coordinated site occupied by different anion groups is subjected to disordering and rotation of these tetrahedra within the structure. The temperature of siwagaite formation is not higher than \sim 70–80 °C, as is evident from the mineral association and as inferred from the formation conditions of the natural and synthetic members of the ettringite group minerals, which are stable at conditions of T < 120 °C and pH = 9.5–13. The name siwagaite is derived from the name of the holotype locality-Siwaga area, where the mineral was found.

Keywords: Siwaqaite, new mineral, ettringite group, crystal structure, Raman, FTIR, Daba-Siwaqa, Jordan