Sundiusite, a new lead sulfate oxychloride from Långban, Sweden

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Abstract

Sundiusite, $Pb_{10}(SO_4)Cl_2O_8$, is a new mineral from Långban, Sweden. It is monoclinic, C2, Cm, or C2/m, with a = 24.67(1), b = 3.781(1), c = 11.881(5)Å, $\beta = 100.07(4)^\circ$, and Z = 2. The strongest lines in the X-ray powder pattern are (Å, *I*, *hkl*) 2.981 10 510; 2.737 8 113; 3.101 6 602, $\overline{6}03$; 3.044 6 800, 403; 6.10 3 400; 3.744 3 110. Sundiusite occurs as plumose aggregates of white to colorless crystals with an adamantine luster. The Mohs hardness is about 3, and there is a perfect {100} cleavage. Optically, it appears to be biaxial (+) with all indices greater than 2.10; lath-shaped fragments are length-slow. The observed and calculated densities are 7.0 and 7.20 g/cm³, respectively. The mineral does not fluoresce in ultraviolet radiation.

The composition, as determined by electron microprobe, is PbO 93.1, FeO 0.5, SO₃ 3.5, Cl 3.0, less $O \equiv Cl 0.7$, total 99.4 weight percent, which yields the ideal formula $Pb_{10}(SO_4)Cl_2O_8$. The composition and cell geometry suggest a structural relationship to the nadorite group. Sundiusite is known only from Långban and is identical with Flink unknown #284. The name is for the late Nils Sundius.

Introduction

This new mineral species was found several years ago on a specimen in the collections of the Smithsonian Institution. The new species was later identified as one of those minerals of unknown identity noted from Långban by the mineralogist and collector, Gustav Flink (1849–1931). Flink published many lists of Långban minerals which he considered worthy of further investigation. The references to these lists are given in a report on the status of the "Flink unknowns" by Moore *et al.* (1971).

We take pleasure in naming this new mineral sundiusite after the late Swedish mineralogist Nils Sundius (1886–1976) in recognition of his numerous contributions to the study of Långban mineralogy including his classic study of the manganiferous pyroxenoids (Sundius, 1931). The pronounciation of the name is $s\bar{u}\dot{n}$ -dē- $\bar{u}s$ -ite. The name sundiusite has been previously used to denote the hypothetical sodic-calcic amphibole, NaCaNaMg₃Al₂Si₆O₂₂(OH)₂ (Phillips and Layton, 1964). However, this name was not recognized by the Subcommittee on Amphiboles, IMA, in its recent systemization of amphibole nomenclature (Leake, 1968) and indeed "sundiusite" was not even included in its list of amphibole names which should be abandoned. Since this compound is unknown as a natural mineral and since the name has not gained popular acceptance, we feel justified in using sundiusite for the new species from Långban. The mineral and name have been approved by the IMA Commission on New Minerals and Mineral Names. The type specimen is preserved in the Smithsonian Institution collection under catalog #NMNH 134984.

Physical and optical properties

Sundiusite occurs exclusively as plumose aggregates of white to colorless crystals up to 8 mm in length. The aggregates are composed of overlapping plates, which are slightly offset from one another in a diverging manner. The streak is white and the luster adamantine. There is one perfect cleavage parallel to

{100}. The mineral is quite brittle, breaking easily into lath-shaped fragments, whose longest dimension is parallel to the b crystallographic axis. A possible second cleavage may exist parallel to {001} but is very poorly developed. The Mohs hardness is approximately 3. The density, determined using a Berman balance and temperature correction, is 7.0 ± 0.2 g/cm^3 , in good agreement with the calculated value of 7.20 g/cm³. Optically, sundiusite appears to be biaxial (+) with refractive indices all >2.1. The maximum birefringence, determined with a tilting compensator, is 0.070. Cleavage laths are length-slow. There is no response to ultraviolet radiation. Sundiusite, along with blixite, occurs on fracture surfaces of an ore consisting of braunite, hausmannite, calcite, and manganoan biotite. It is the last mineral of this assemblage to form.

Sundiusite is similar in appearance to several other minerals, the most notable of which are hydrocerussite and mendipite. The resemblance to mendipite, Pb₃O₂Cl₂, from the Mendip Hills, England is particularly striking. Although neither mineral fluoresces under ultraviolet radiation, the calcite associated with each has a similar fluorescence, which accentuates the resemblance between mendipite and sundiusite. This has led to some confusion, a matter which will be discussed in a subsequent section.

Chemistry

Sundiusite was chemically analyzed with an ARL-SEMQ electron microprobe, at an operating voltage of 15 kV and a beam current of 0.15 µA, determined using a beam current monitor. The standards used were PbO for lead, celestine for sulfur, hornblende for iron, and halite for chlorine. A wavelength-dispersive scan indicated the absence of any other elements with atomic number greater than nine. The data were corrected using a modified version of the MAGIC-4 program. The resulting analysis is presented in Table 1. Calculation of the unit-cell contents using the observed density yields Pb_{19,3}Fe_{0,3}S_{2,0}Cl_{3,9}O_{23,6} in good agreement with the theoretical composition Pb_{10} $(SO_4)Cl_2O_8$ with two formula weights per cell. Sundiusite is chemically homogeneous and is very slowly soluble in 1:1 cold HCl or HNO₃.

Crystallography

Sundiusite was studied by Weissenberg, precession, and oscillating crystal methods. The mineral is monoclinic, C2, Cm, or C2/m, with a = 24.67 esd $0.01, b = 3.781 \text{ esd } 0.001, \text{ and } c = 11.881 \text{ esd } 0.005\text{\AA},$ $\beta = 100.07 \text{ esd } 0.04^{\circ}$, and $V = 1091.1 \text{ esd } 0.8\text{\AA}^3$.

SUNDIUSITE	Theoretical Pb ₁₀ (SO ₄)Cl ₂
the second se	10 4 2

Table 1. Electron microprobe analysis of sundiusite

	3000103112	Pb10(S04)C1208
РЬО	93.1	94.30
Fe0	0.5	
50 ₃	3.5	3.38
C1	3.0	3.00
Total	100., 1	100.68
Less O = Cl	0.7	0.68
Total	99.4	100.00

Accuracy of data: ±3% of the amount present. Quantities given are in weight percent.

There is also a prominent subcell with A = 24.67, B = 3.781, C = c/3 = 3.960Å, and β = 100.07°. Cell parameters were refined by least-squares from Debye-Scherrer data utilizing NBS silicon (a =5.4309Å) as an internal standard. The X-ray powder diffraction data are given in Table 2.

Table 2. X-ray powder diffraction data for sundiusite

^I obs	d _{obs}	d _{cal}	hk1	Iobs	d _{obs}	d _{cal}	hk1
<1 1 3 <<]	12.0 9.29 7.79 6.10 5.78	12.1 9.27 7.77 6.07 (5.82	200 201 201 400 401 202	2b 2 2	2.033 1.980 1.890 1.826	2.042 2.024 2.022 1.980 1.891 1.825	913 12.0.0 713 206 020,606 11.1.3
<1 2 3	4.03 3.914 3.744	4.05 {3.917 {3.899 3.736	600 203 003 110	2 <1	1.807	{1.807 {1.805 {1.752 {1.751 {1.748}}	913 420 515 316 T16
2b 2 <1	3.553 3.428 3.349	3.578 3.537 3.532 3.426 3.356	403 203 111 310 311	Ţ	1.704	{1.710 {1.701 1.699 (1.680	116 023 14.0.3 12.0.3
6	3.101	{3.091 3.086	603 602	2	1,676	1.675 1.672 1.671	13·1·0 423 007
6 10	3.044 2.981	3.048 3.036 2.984	800 510	1	1.657 1.639	1.657 (1.643 (1.635	$\overline{7}_{16}$ 316 13.1.3
<<] 8 <]	2.892 2.737 2.628	2.911 2.735 2.629	802 113 803	2	1.608	{1.607 1.605	423 820
<1	2.559	{2.567 {2.556	711 710	<<1 1	1.551 1.490	1.546 1.492 1.491	12.0.6 10.2.0 14.0.3
<<1	2.491	{2.503 (2.484 (2.433	513 313 711	1	1.368	(1.488	15.1.0
2 1	2.428 2.274	2.429	10·0·0 713 114	<<] <<] <]	1.353 1.267 1.244		
<<1	2.186	2.195	604	<1 1	1.220		

114.6 mm Debye-Scherrer camera, CuK α radiation, NBS silicon internal standard, b = broad line

The subcell dimensions of sundiusite suggest a relationship between the new mineral and members of the nadorite group. As shown in Table 3, the lattice parameters of orthorhombic nadorite, perite, and blixite are related to those of synthetic tetragonal nadorite (SN) such that $a \simeq c \simeq a_{\rm SN} \sqrt{2}$ and $b \simeq c_{\rm SN}$ or $2c_{\rm SN}$. In other words the basic geometrical unit common to all of them is either tetragonal or pseudotetragonal with parameters $A \simeq 4$ and $C \simeq 12$ Å. This corresponds to the sundiusite subcell with its A parameter halved. It may also be noted that nadorite and its congeners possess a single perfect cleavage perpendicular to their 12 (or 25) Å translation, reflecting the layered character of their crystal structures (Giuseppetti and Tadini, 1973). Sundiusite also possesses a single perfect cleavage nearly perpendicular to its 25 Å translation. From the foregoing similarities we infer that sundiusite is a lead oxychloride with a layered structure related to those of the nadorite group.

If sundiusite is indeed related to nadorite as proposed, the presence of the sulfate ion is somewhat incongruous, as the nadorite structure contains no tetrahedral anions. An explanation may be that the nadorite minerals are a part of a larger family of bismuth, lead, and antimony oxyhalides (*Structure Reports*, 1947–1948, p. 305–320), some of which contain complex anions, *e.g.* CO_3^{2-} and MoO_4^{2-} , which play the same structural role as the halide ions. This may also be the role of SO_4^{2-} in the sundiusite structure.

Identity of sundiusite and Flink unknown #284

The original label that accompanied the specimen gives very little information; it simply states "#284, 285, Långban." In one of his lists of new or incompletely described minerals from Långban (Flink, 1924), the number 284 is assigned to a lath-like, snow white, highly lustrous mineral from the Amerika stope. This description is in good agreement with the appearance of sundiusite, and it would seem that the

Table 3. Crystal chemical data for minerals of the nadorite group

	Formula	Space Group	Unit Cell		(A)
			а	Ь	С
Nadorite ^a	PbSb0 ₂ C1	Cmem	5.60	12.24	5.45
Perite ^b	PbBi0 ₂ C1	Cmem	5.63	12.42	5.58
Blixite ^C	Pb ₁₆ (0,0H) _{16-x} C1 ₈	Orthorhombic	5.83	25.47	5.69
Nadorite ^d (synthetic)	PbSb02C1	I4/mmm	a = 3.90, c = 12.29		

^aGiuseppetti and Tadini (1973), ^bGillberg (1961), ^CGabrielson <u>et al.</u> (1960), ^dSillén and Melander (1941).

two minerals are identical (Flink #285 is blixite). However, we note that Moore et al. (1971) attribute #284 to mendipite on the basis of density and singlecrystal X-ray photographs. Subsequent to our characterization of sundiusite, a number of specimens of Flink #284 were located in the Flink collection at Harvard University. Both sundiusite and mendipite were identified by powder X-ray photography on these specimens, but the two minerals do not occur together on any one specimen. Hence, Flink number 284 was probably assigned to two separate minerals, which would have been difficult for Flink to distinguish by sight or density measurements, the densities of sundiusite and mendipite being 7.20 and 7.22 g/ cm³, respectively. Mendipite has also been reported from Långban by Goñi and Guillemin (1953).

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