

Sr[S₂O₆][H₂O]₄

*hP*63

(180) *P*6₂22 – k⁵a

SrS₂O₆·4H₂O [1]

Structural features: Layers containing O₃S-SO₃ units (perpendicular to [001], two-fold disorder about [110]) alternate with puckered layers containing Sr atoms and H₂O molecules.

De Matos Gomes E. (1991) [1]

H₈O₁₀S₂Sr

a = 0.63529, *c* = 1.9218 nm, *c/a* = 3.025, *V* = 0.6717 nm³, *Z* = 3

site	Wyck.	sym.	<i>x</i>	<i>y</i>	<i>z</i>	occ.	atomic environment
O1	12 <i>k</i>	1	0.035	0.267	0.22747	0.5	
O2	12 <i>k</i>	1	0.041	0.273	0.10207	0.5	
S3	12 <i>k</i>	1	0.0654	0.3749	0.16907	0.5	
O4	12 <i>k</i>	1	0.328	0.524	0.17277	0.5	single atom S
(OH ₂)5	12 <i>k</i>	1	0.4499	0.2559	0.04247		single atom Sr
Sr6	3 <i>a</i>	222	0	0	0		

Transformation from published data (*P*6₄22): new axes -a,-b,-c

Experimental: single crystal, diffractometer, X-rays, wR = 0.040

Remarks: Short interatomic distances for partly occupied site(s). Hydrogen atoms are not taken into consideration for Pearson symbol, Wyckoff sequence and atomic environments. An ordered structure in space group (171) *P*6₂ is reported by the same author in [2].

References: [1] De Matos Gomes E. (1991), Acta Crystallogr. B 47, 12-17. [2] De Matos Gomes E. (1991), Acta Crystallogr. B 47, 7-11.