

**No. 46A-8 (CH<sub>3</sub>)<sub>2</sub>NH<sub>2</sub>Al(SO<sub>4</sub>)<sub>2</sub> · 6H<sub>2</sub>O, Dimethylammonium aluminum sulfate hexahydrate***(M* = 373.29; [*D*: 393.42])

1a	Ferroelectricity in (CH <sub>3</sub> ) <sub>2</sub> NH <sub>2</sub> Al(SO <sub>4</sub> ) <sub>2</sub> · 6H <sub>2</sub> O was first reported by Kirpichnikova et al. in 1988.		88Kir
b	phase	II	I 88Kir
	state	F	P
	crystal system	monoclinic	monoclinic <sup>a)</sup> 93Pie1
	space group	Pn–C <sub>s</sub> <sup>2a)</sup>	P2 <sub>1</sub> /n–C <sub>2h</sub> <sup>5b)</sup> <sup>b)</sup> 93Pie2
	Θ [K]	152	88Kir
	<i>P<sub>s</sub></i> makes an angle 43° from the <i>a</i> axis in the (101) plane. *)		88Kir
	<i>ρ<sub>X</sub></i> = 1.646 · 10 <sup>3</sup> kg m <sup>–3</sup> at 295 K.		93Pie2
	*) Here the crystal axes are chosen as adopted in [93Pie2]. Then the <i>a</i> axis corresponds to the <i>c</i> axis in [88Kir]; see Fig. 46A-8-004.		
2a	Crystal growth: evaporation of aqueous solution.		88Kir
3a	Unit cell parameters:		
	Phase I:		
	<i>a</i> = 6.403(1) Å, <i>b</i> = 10.747(2) Å, <i>c</i> = 11.282(2) Å, β = 100.47(3)°. <i>T</i> = 295 K.		93Pie2
	Kiosse et al. used different crystallographic systems.		94Kio
	The unit cell parameters given are:		
	<i>a</i> ' = 11.849(4) Å, <i>b</i> ' = 6.415(2) Å, <i>c</i> ' = 10.768(4) Å, γ' = 68.13(9)°. <i>T</i> = 295 K.		94Kio
	Relation between two axial systems: <i>a</i> ' = <i>a</i> + <i>c</i> , <i>b</i> ' = <i>c</i> , <i>c</i> ' = <i>b</i> .		
	Phase II:		
	<i>a</i> = 6.445(1) Å, <i>b</i> = 10.688(2) Å, <i>c</i> = 11.109(2) Å, β = 99.71(3)°. <i>T</i> = 95 K.		93Pie1
b	<i>Z</i> = 2 in phases I <sup>a)</sup> and II <sup>b)</sup> .		<sup>a)</sup> 93Pie1 <sup>b)</sup> 93Pie2
	Crystal structure:		
	Structure of phase I: Table 46A-8-001, Table 46A-8-002, Table 46A-8-003;		
	Fig. 46A-8-001, Fig. 46A-8-002.		
	Structure of phase II: Table 46A-8-004, Table 46A-8-005, Table 46A-8-006;		
	Fig. 46A-8-003.		
5a	Dielectric constant: Fig. 46A-8-004.		
	Dielectric dispersion: Fig. 46A-8-005, Fig. 46A-8-006, Fig. 46A-8-007, Fig. 46A-8-008.		
	Curie Weiss law: κ <sub>0</sub> = κ <sub>∞</sub> + <i>C</i> /( <i>T</i> – Θ <sub>p</sub> ), <i>T</i> > Θ <sub>I–I</sub> , where <i>C</i> = 2.81 · 10 <sup>3</sup> K, Θ <sub>p</sub> = 150.9 K, κ <sub>∞</sub> = 5. κ <sub>0</sub> : static dielectric constant.		91Sob1
	Phase diagram with regard to <i>p</i> : Fig. 46A-8-009.		
b	Nonlinear dielectric properties: Fig. 46A-8-010.		
	<i>E</i> = (1/χ <sub>p</sub> ) <i>P</i> + ξ <i>P</i> <sup>3</sup> + ζ <i>P</i> <sup>5</sup> where ξ = 4.0(5) · 10 <sup>12</sup> V m <sup>5</sup> C <sup>–3</sup> , ζ = 1.8 · 10 <sup>17</sup> V m <sup>9</sup> C <sup>–5</sup> .		89Cac
c	Spontaneous polarization and coercive field: Fig. 46A-8-011;		
	see also		95Pyk
9a	Refractive indices: <i>n</i> <sub>X'</sub> = 1.4472, <i>n</i> <sub>Y'</sub> = 1.4683, <i>n</i> <sub>Z'</sub> = 1.4706.		89Kir
	For relation between the optical axes <i>X'</i> , <i>Y'</i> , <i>Z'</i> and the crystallographic axes <i>a</i> , <i>b</i> , <i>c</i> , see Fig. 46A-8-004.		
	Birefringence: see		89Kir

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10a	Raman scattering: see	90Tor
16	Rectangular mechanical hysteresis curve (stress-strain) is observed at RT: spontaneous strain: $\approx 10^{-1}$ , coercive stress $\approx 5$ MPa. Plasticity and mechanical twin formation: see	89Kir 89Kir, 90Kir