

No. 41A-9 CsLiWO₄, Cesium lithium tungstate*(M* = 387.69)

1a	Ferroelectricity in CsLiWO ₄ was reported by Aleksandrov et al. in 1981.			81Ale
b	phase	III	II	I
	state	F	F	P
	crystal system	monoclinic	trigonal	cubic
	space group	Bm – C _s ³	R3m – C _{3v} ⁵	F $\bar{4}$ 3m – T _d ²
	Θ [K]	191	221	
	$P_s \parallel [111]$.			81Ale
	$\rho_X = 4.42 \cdot 10^3 \text{ kg m}^{-3}$.			80Oka
3a	Unit cell parameters:			
	Phase I (at 298 K): $a = 8.350(1) \text{ \AA}$.			80Oka
	Phase II: trigonal system: $a = 5.91 \text{ \AA}$, $\alpha = 59.5^\circ$;			81Ale
	hexagonal system: $a = b = 5.836(3) \text{ \AA}$, $c = 14.322(8) \text{ \AA}$.			81Ale
	Phase III: $a = 5.85 \text{ \AA}$, $b = 5.86 \text{ \AA}$, $c = 10.25 \text{ \AA}$, $\beta = 125.5^\circ$.			81Ale
b	$Z = 4$ (in phase I), $Z = 3$ (in phase II), $Z = 2$ (in phase III).			85All
	Crystal structure of phase I: Fig. 41A-9-001, Fig. 41A-9-002, Fig. 41A-9-003.			
	Fractional coordinates and isotropic temperature parameters: Table 41A-9-001, Table 41A-9-002.			
	Interatomic distances: Table 41A-9-003, Table 41A-9-004.			
	Crystal structure of phase II: Fig. 41A-9-004.			
5a	Dielectric constant at 180...250 K: Fig. 41A-9-005.			
c	Spontaneous polarization at 160...230 K: Fig. 41A-9-006.			
6a	Specific heat capacity: Fig. 41A-9-007.			
	See also			84Ale
	Transition entropy ΔS_m :			83Fle
	at II–I transition: $0.3 R$, at III–II transition: $0.334 R$.			
	Transition heat ΔQ_m :			83Fle
	at II–I transition: 600 J mol^{-1} , at III–II transition: 450 J mol^{-1} .			
9b	Electrooptic coefficient: Fig. 41A-9-008.			
e	Nonlinear optical properties: $I_{2\omega} / I_{2\omega}^{\text{SiO}_2} = 35$. $I_{2\omega}$: SHG intensity.			84Kab