No. 1C-a20 $NaTaO_3$–$KTaO_3$

1b	Phase diagram: Fig. 1C-a20-001.	
3a	Lattice parameters: Fig. 1C-a20-002.	
5a	Dielectric constant: Fig. 1C-a20-003, Fig. 1C-a20-004. Dielectric relaxation: Fig. 1C-a20-005. Pressure effect on transition temperature: Fig. 1C-a20-006. Infrared absorption: see	69Per
c	Remanent polarization: Fig. 1C-a20-007.	
10a	Raman scattering: Fig. 1C-a20-008.	
13a	NMR: Fig. 1C-a20-009.	

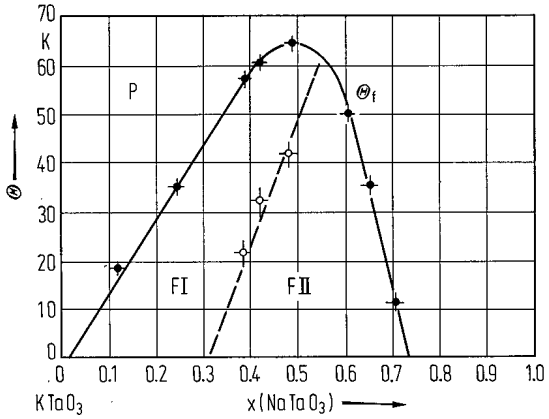

Fig. 1C-a20-001. $(K_{1-x}Na_x)TaO_3$. Θ vs. x [70Dav]. P: paraelectric cubic, F I: ferroelectric tetragonal, F II: ferroelectric orthorhombic (?).

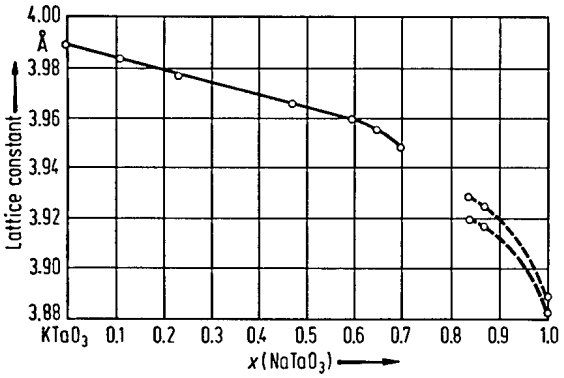

Fig. 1C-a20-002. $(K_{1-x}Na_x)TaO_3$. Lattice parameter vs. x [70Dav]. Structure for x = 0.84 appears to be tetragonal.

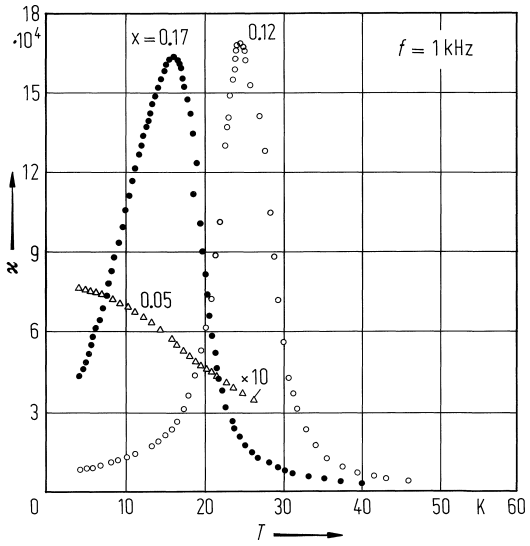

Fig. 1C-a20-003. $(K_{1-x}Na_x)TaO_3$. κ vs. T [83van]. Parameter: x. f = 1 kHz.

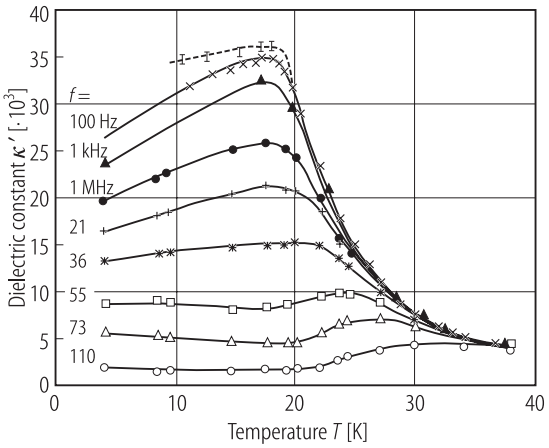

Fig. 1C-a20-004. $(K_{0.8}Na_{0.2})TaO_3$. κ' vs. T [86Mag]. Parameter: f. Broken line: static dielectric constant from Kramers-Kronig analysis.

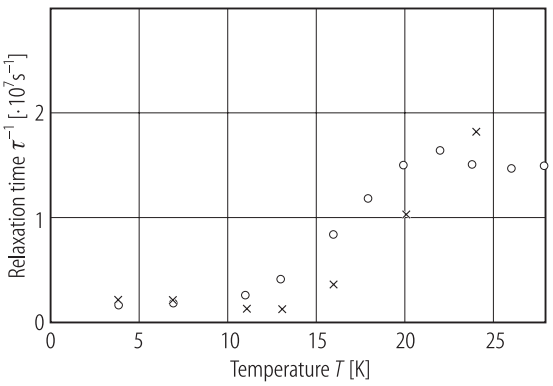

Fig. 1C-a20-005. $(K_{0.8}Na_{0.2})TaO_3$. τ^{-1} vs. T [88Mag]. τ : relaxation time. Circles: dielectric, crosses: acoustic.

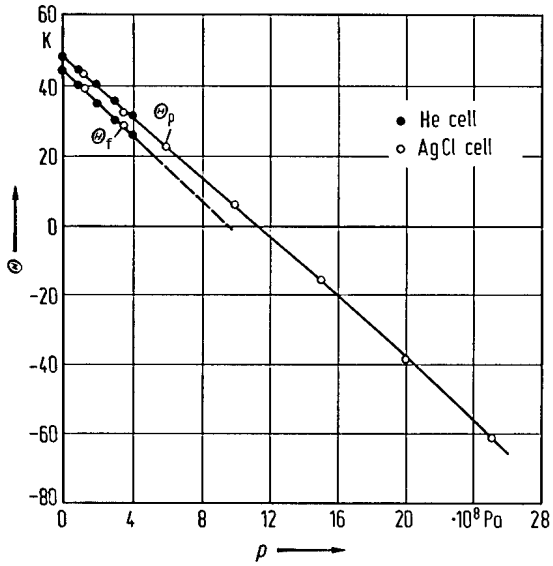

Fig. 1C-a20-006. $(K_{0.72}Na_{0.28})TaO_3$. Θ vs. p [74Sam]. Θ_p: paraelectric Curie temperature.

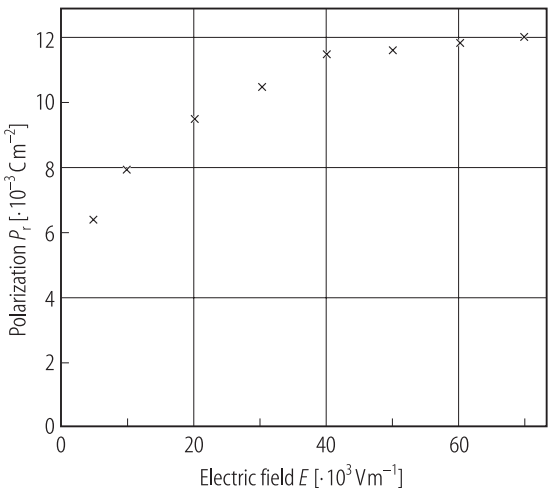

Fig. 1C-a20-007. $(K_{0.82}Na_{0.18})TaO_3$. P_r vs. E [87Lan]. E: applied electric field.

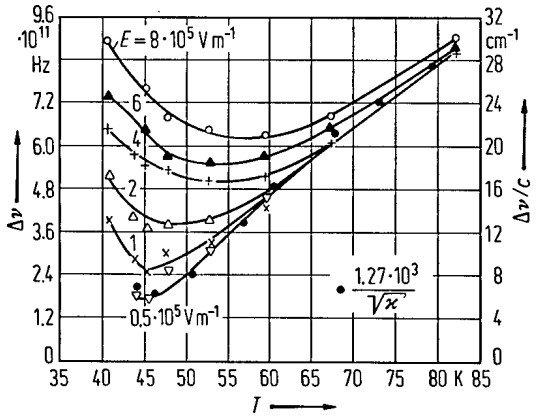

Fig. 1C-a20-008. $(K_{0.7}Na_{0.3})TaO_3$. $\Delta\nu$ vs. T [72Dav]. $\Delta\nu$: Raman frequency shift of the soft phonon mode. Parameter: electric field along [001] axis. Full circles indicate best fit of low-field data to $\kappa^{-1/2}$.

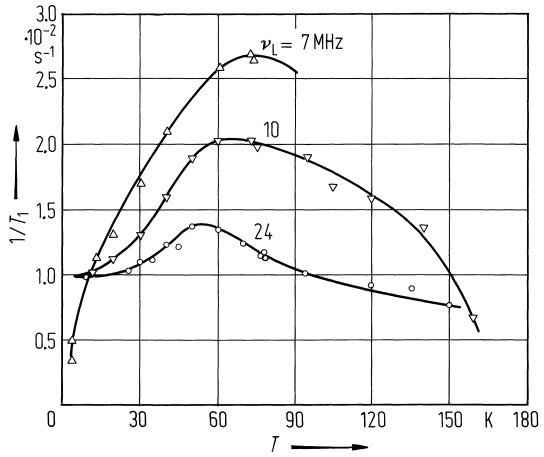

Fig. 1C-a20-009. $(K_{0.82}Na_{0.18})TaO_3$. $1/T_1$ vs. T [83Hoc]. Parameter: ν_L (Larmor frequency). T_1: spin-lattice relaxation time of ^{23}Na nuclei.

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