

Fig. 27A-1-001. $H_{1-x}D_xCl$. Θ_f vs. x [83Cro].

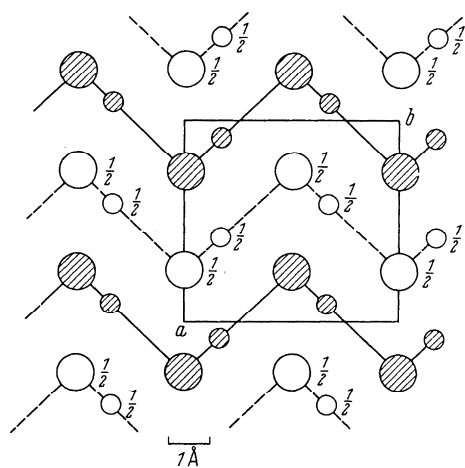


Fig. 27A-1-002. DCl. Crystal structure of phase II projected on (001) [67San1]. $T = 77.4$ K.

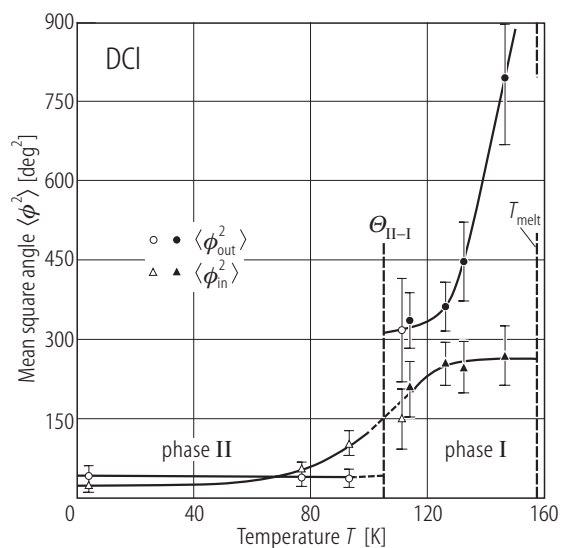


Fig. 27A-1-003. DCl. $\langle \phi_{\text{in}}^2 \rangle$, $\langle \phi_{\text{out}}^2 \rangle$ vs. T [73Nii]. $\langle \phi_{\text{in}}^2 \rangle$, $\langle \phi_{\text{out}}^2 \rangle$: mean square angular displacements within and out of the molecule plane, respectively. Open circles and triangles represent data given in [69San].

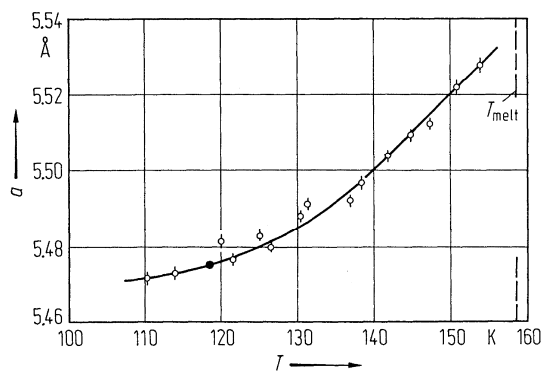


Fig. 27A-1-004. DCl. a vs. T [73Nii]. Data are normalized at 118.5 K as to coincide with the value given in [69San] shown by the solid circle.

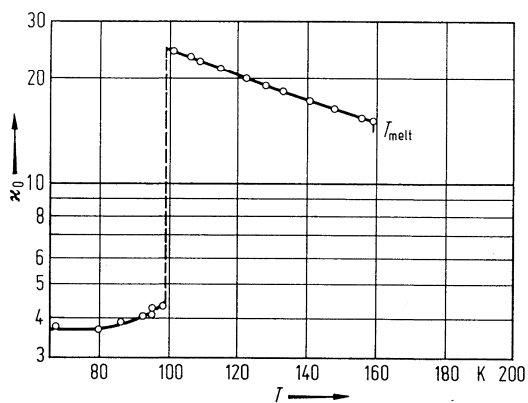


Fig. 27A-1-005. HCl (polycrystal). κ_0 vs. T [54Swe]. κ_0 : static dielectric constant.

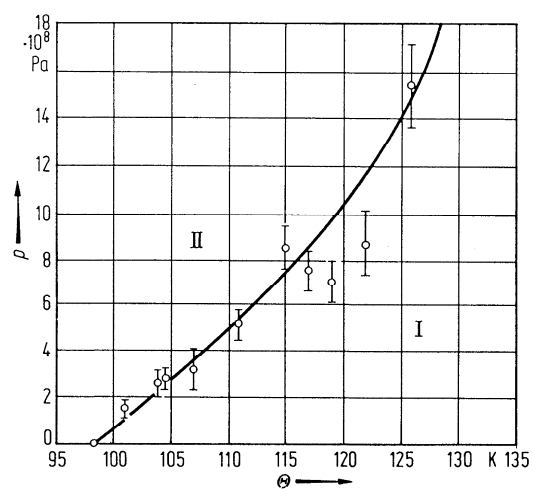


Fig. 27A-1-006. HCl. Θ vs. p [62Ste].

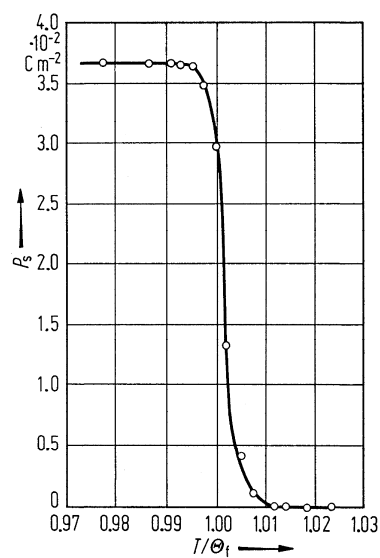


Fig. 27A-1-007. HCl. P_s vs. T/Θ_f [69Kob]. Obtained from pyroelectric measurements.

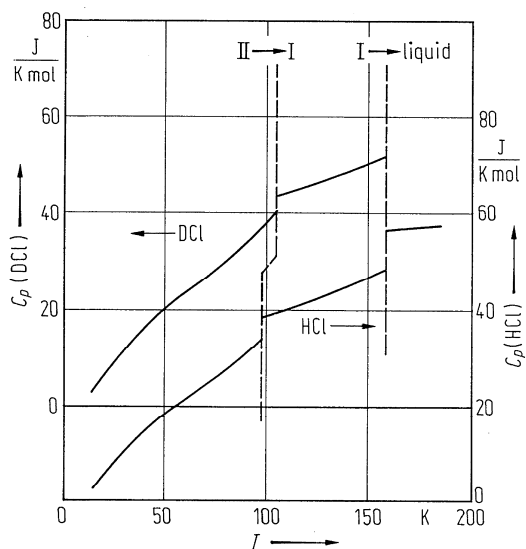


Fig. 27A-1-008. HCl, DCl. C_p vs. T [47Clu].

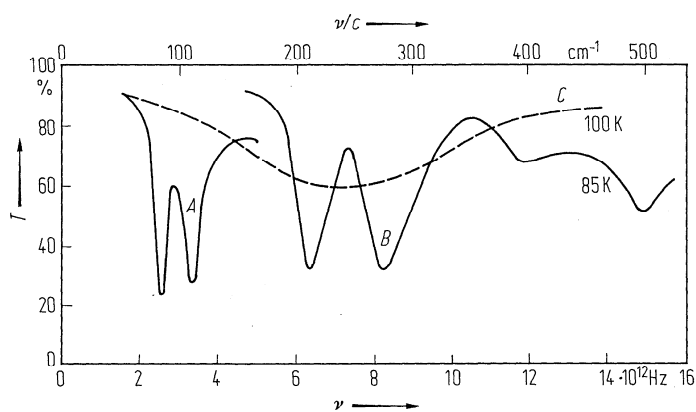


Fig. 27A-1-009. HCl. T vs. ν [67Arn]. T : transmittance, ν : frequency of incident infrared radiation. Temperature: 85 K for A, B ; 100 K for C . Specimen thickness $\approx 20 \mu\text{m}$ for A, C ; $\approx 5 \mu\text{m}$ for B .

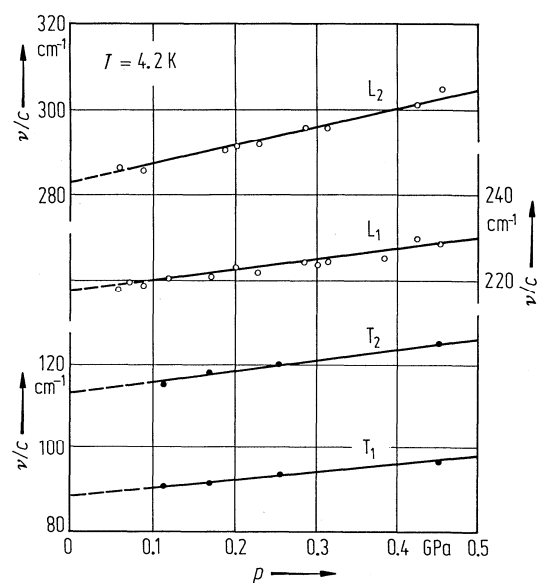


Fig. 27A-1-010. HCl. ν/c vs. p at 4.2 K [83Obr]. ν : frequency of far infrared absorption band. L and T denote librational and translational modes, respectively.

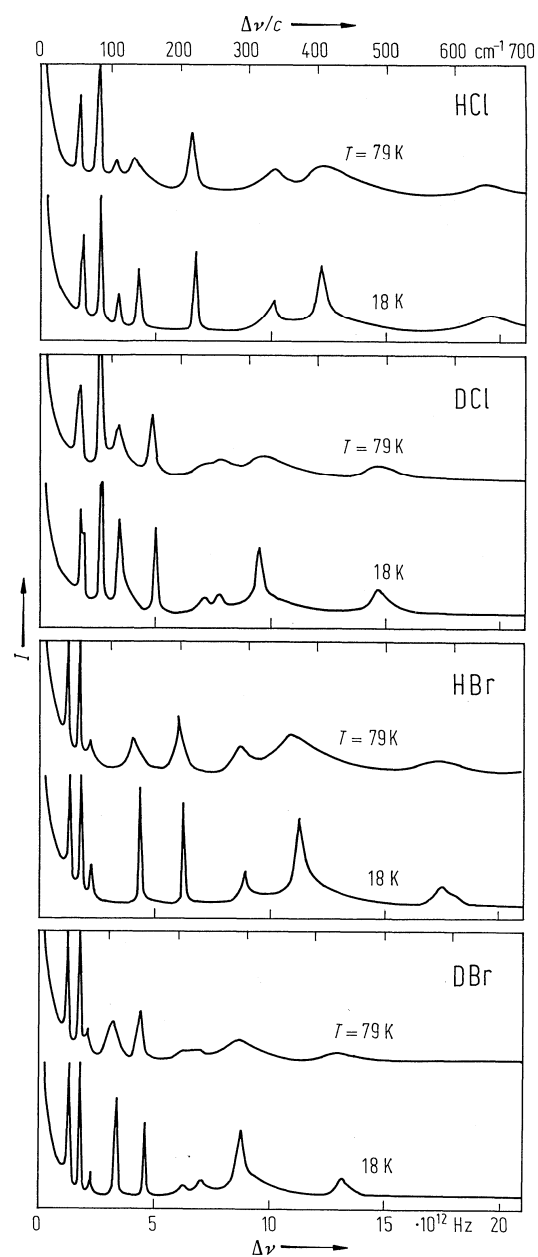


Fig. 27A-1-011. HCl, DCl, HBr, DBr. I vs. $\Delta\nu$ at 79 K and 18 K [81And]. I , $\Delta\nu$: intensity and frequency shift of Raman scattering.

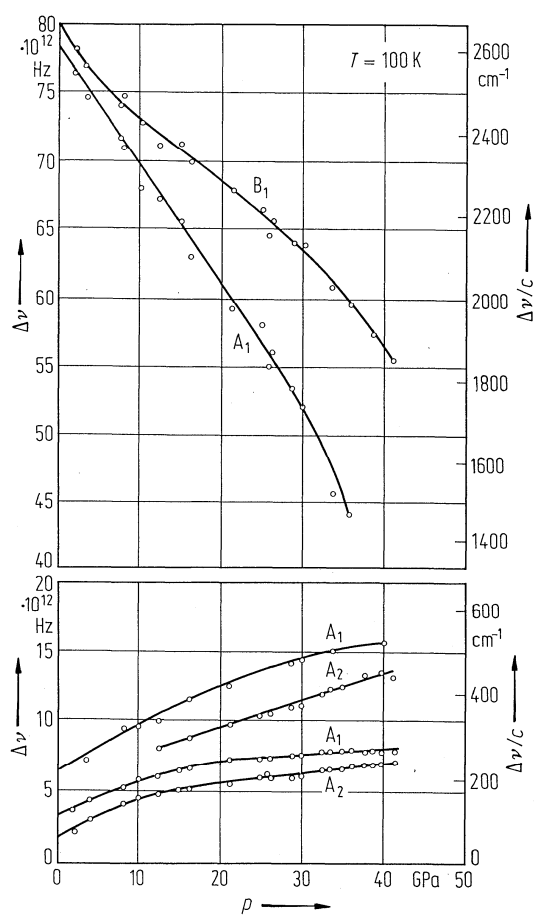


Fig. 27A-1-012. HCl. $\Delta\nu$ vs. p at 100 K [84Joh]. $\Delta\nu$: Raman frequency shift.

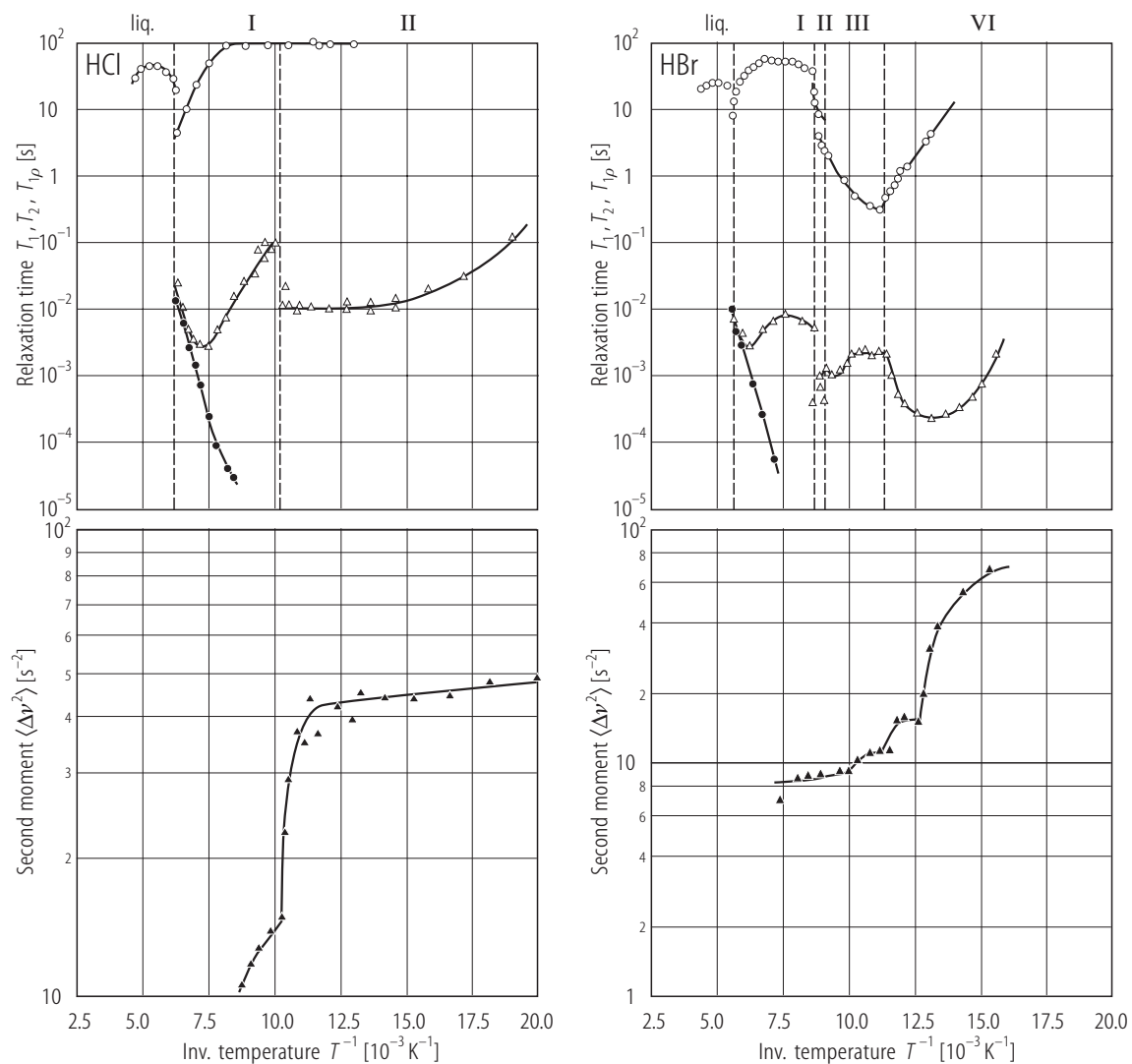


Fig. 27A-1-013. HCl, HBr. T_1 , T_2 , $T_{1\rho}$, $\langle \Delta\nu^2 \rangle$ vs. $1/T$ [68Gen]. Open circle: proton spin lattice relaxation time T_1 , full circles: spin-spin relaxation time T_2 , open triangles: rotating frame relaxation time $T_{1\rho}$, full triangles: second moment $\langle \Delta\nu^2 \rangle$. rf magnetic field strength $H_1 = 1273.2 \text{ A m}^{-1}$.

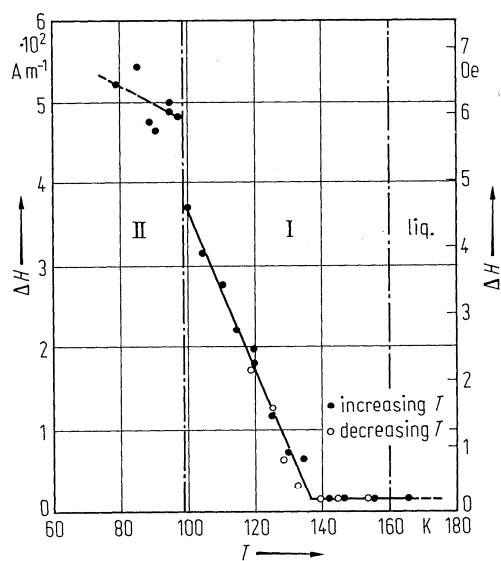


Fig. 27A-1-014. HCl. ΔH vs. T [49Alp]. ΔH : NMR line width of proton.

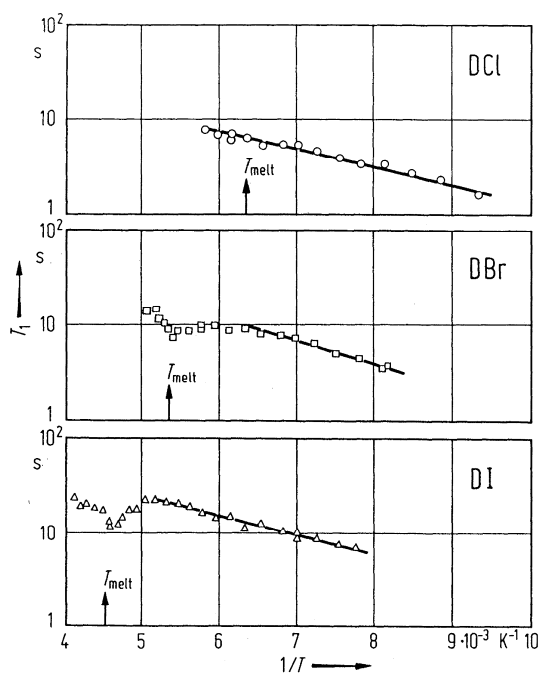


Fig. 27A-1-015. DCl, DBr, DI. T_1 vs. $1/T$ [68Gen]. T_1 : deuteron spin lattice relaxation time.

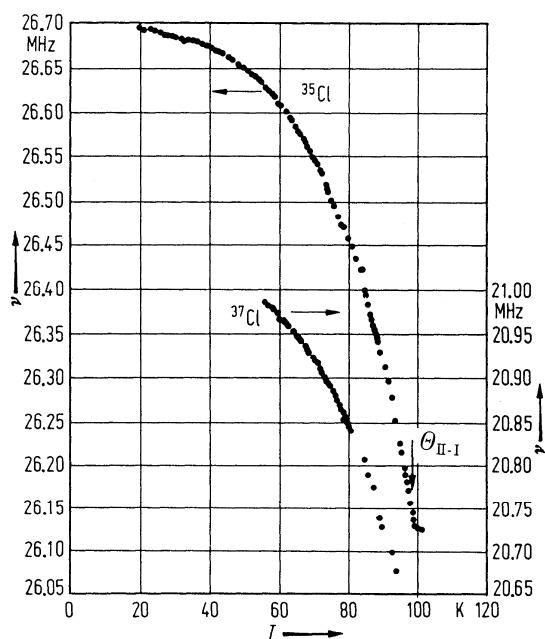


Fig. 27A-1-016. HCl. ν vs. T [68Oku]. ν : NQR frequency. See also [77Mac].

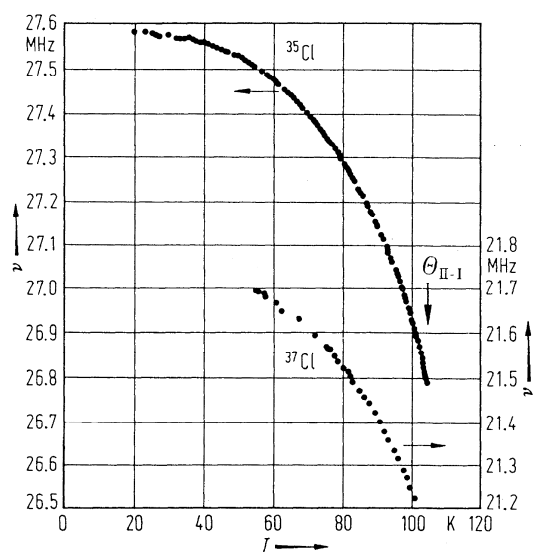


Fig. 27A-1-017. DCl. ν vs. T [68Oku]. ν : NQR frequency.

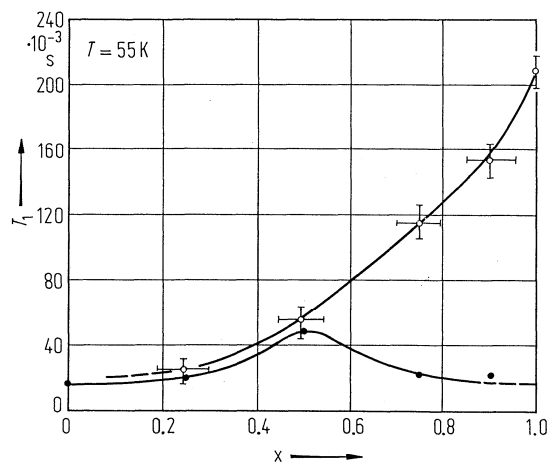


Fig. 27A-1-018. $\text{H}_{1-x}\text{D}_x\text{Cl}$. T_1 vs. x at 55 K [83Cro]. T_1 : spin-lattice relaxation time for ^{35}Cl belonging to H-Cl dipoles (full circles) and to D-Cl dipoles (open circles).

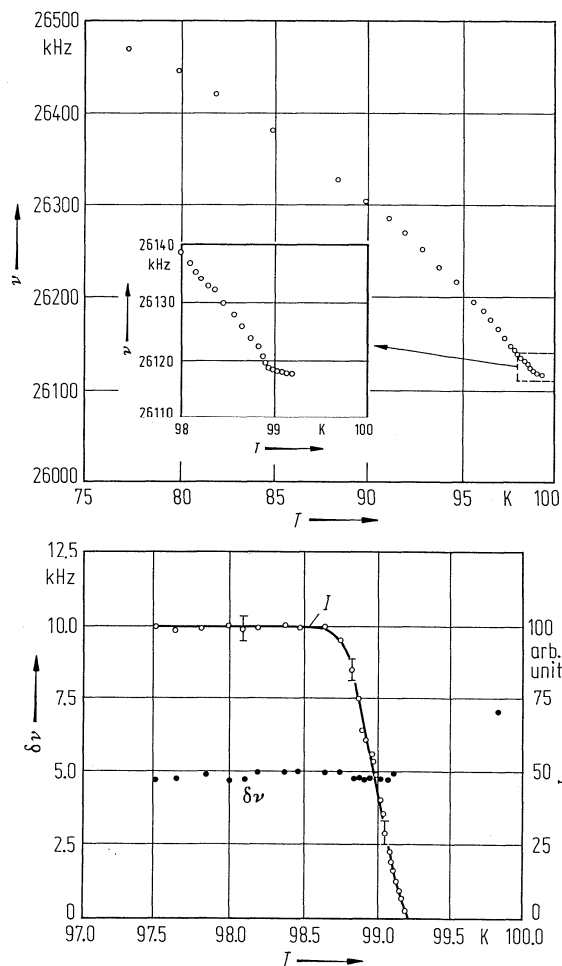


Fig. 27A-1-019. HCl. ν , $\delta\nu$, I vs. T [80Rig]. ν , $\delta\nu$, I : frequency, linewidth (half width at half maximum) and intensity of ^{35}Cl NQR line. See also Fig. 27A-1-016.

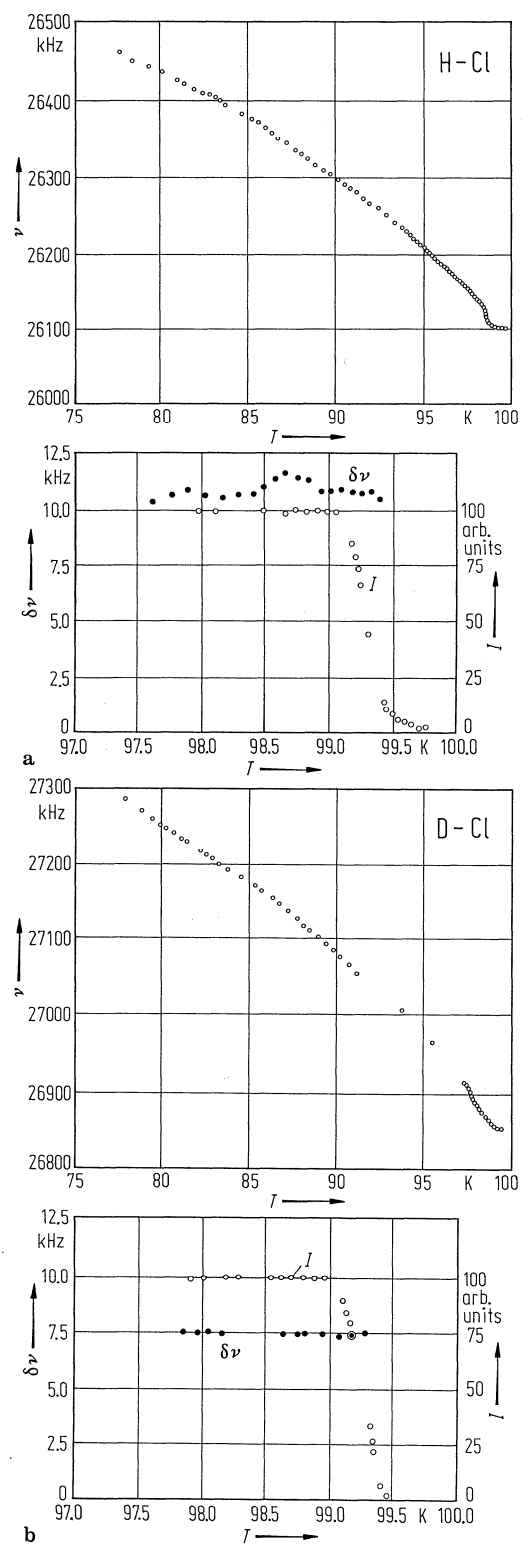


Fig. 27A-1-020. $\text{H}_{0.75}\text{D}_{0.25}\text{Cl}$. ν , $\delta\nu$, I vs. T [80Rig]. ν , $\delta\nu$, I : frequency, linewidth (half width at half maximum) and intensity of ^{35}Cl NQR line (a) for H-Cl molecules and (b) for D-Cl molecules.

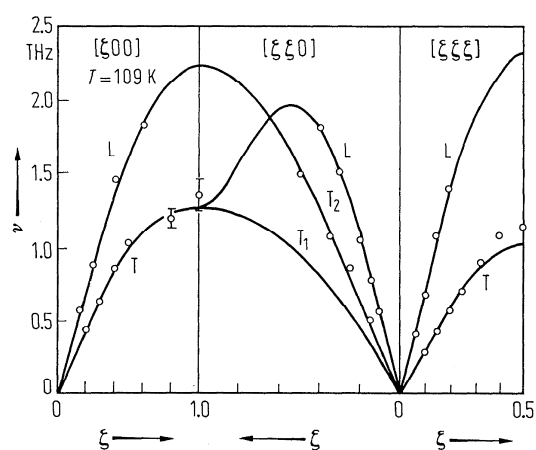


Fig. 27A-1-021. DCl. ν vs. ζ . [75Pre]. ν : phonon frequency, ζ : reduced wave vector coordinate.

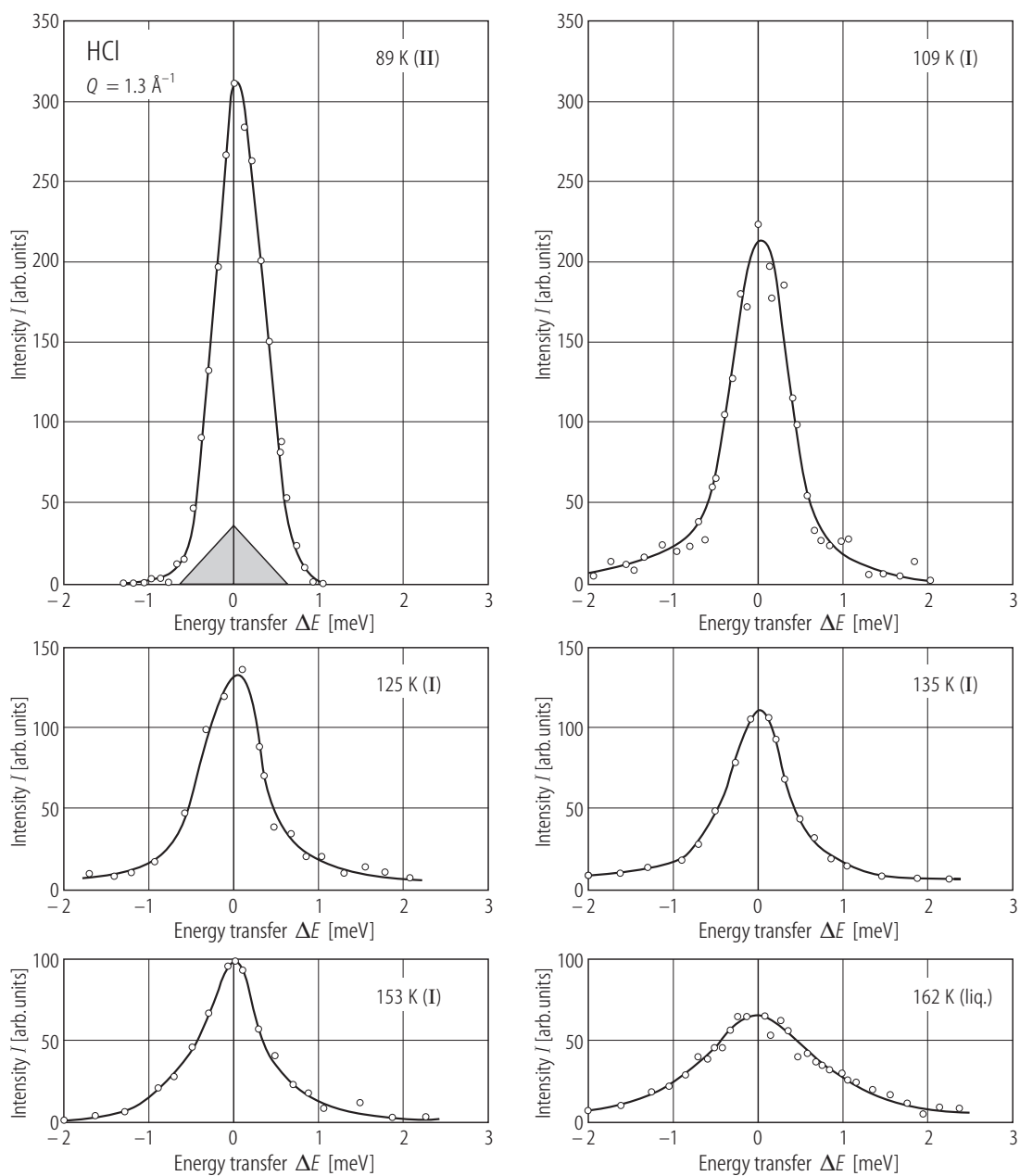


Fig. 27A-1-022. HCl. I vs. ΔE [78Fuj]. Parameter: T . I : incoherent neutron scattering intensity. ΔE : energy transfer of neutrons. Wave number coordinate $Q = 1.3 \text{ \AA}^{-1}$.