

Fig. 32A-3-001. $\alpha\text{-KIO}_3 \cdot \text{HIO}_3$. Θ_f , Θ_f^2 vs. p [87Bai].

$\alpha\text{-KIO}_3 \cdot \text{HIO}_3$

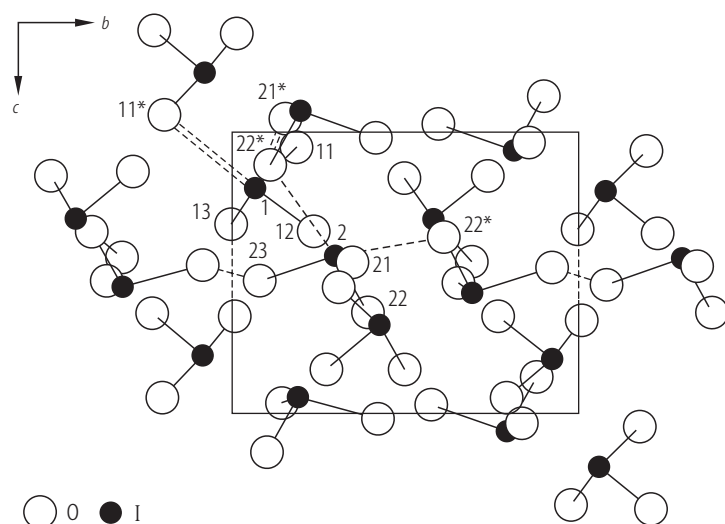


Fig. 32A-3-002. $\alpha\text{-KIO}_3 \cdot \text{HIO}_3$. Crystal structure of phase I at RT [78Tre2]. Projection along the a axis. Only the iodate groups are shown. Used crystallographic axes are (a', b', c') , the second setting of cell defined in 3b.

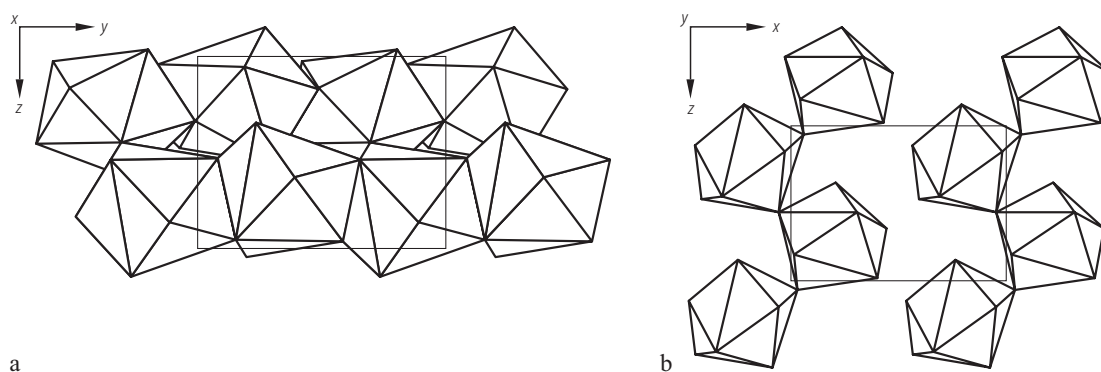
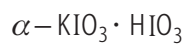


Fig. 32A-3-003. $\alpha\text{-KIO}_3 \cdot \text{HIO}_3$. Crystal structure of phase I at RT [78Tre2]. KO₉ polyhedra. (a) Projection along the *a* axis, (b) projection of polyhedra at a level *y* \approx 0.25 along the *b* axis. Used crystallographic axes are (*a'*, *b'*, *c'*), the second setting of cell defined in 3b.

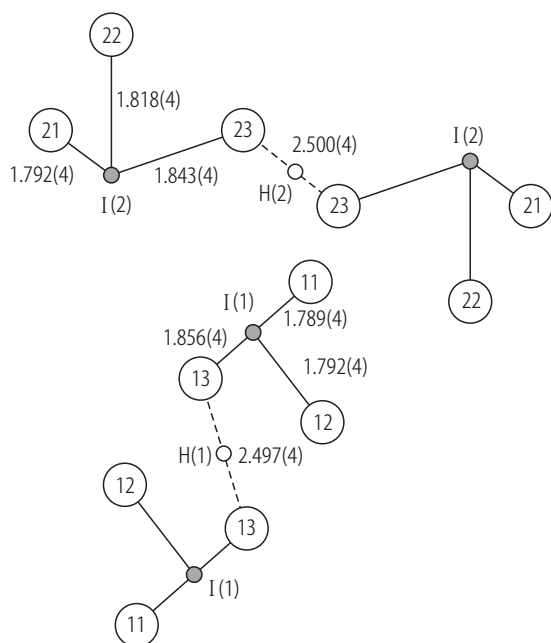


Fig. 32A-3-004. $\alpha\text{-KIO}_3 \cdot \text{HIO}_3$. Crystal structure of phase I at RT [79Tre]. Structure of non-equivalent dimers of iodate. Distances in [Å].

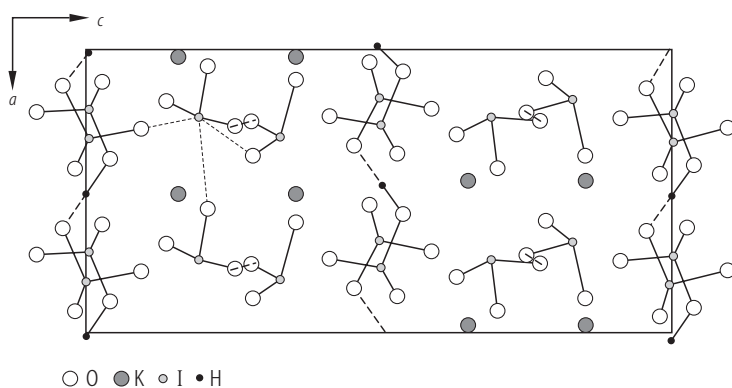
$\beta - \text{KIO}_3 \cdot \text{HIO}_3$ 

Fig. 32A-3-005. $\beta\text{-KIO}_3 \cdot \text{HIO}_3$. Crystal structure at RT [85Sor]. Projection along the b axis. Dashed lines: I–O bonds from second coordination sphere; dotted lines: hydrogen bonds.

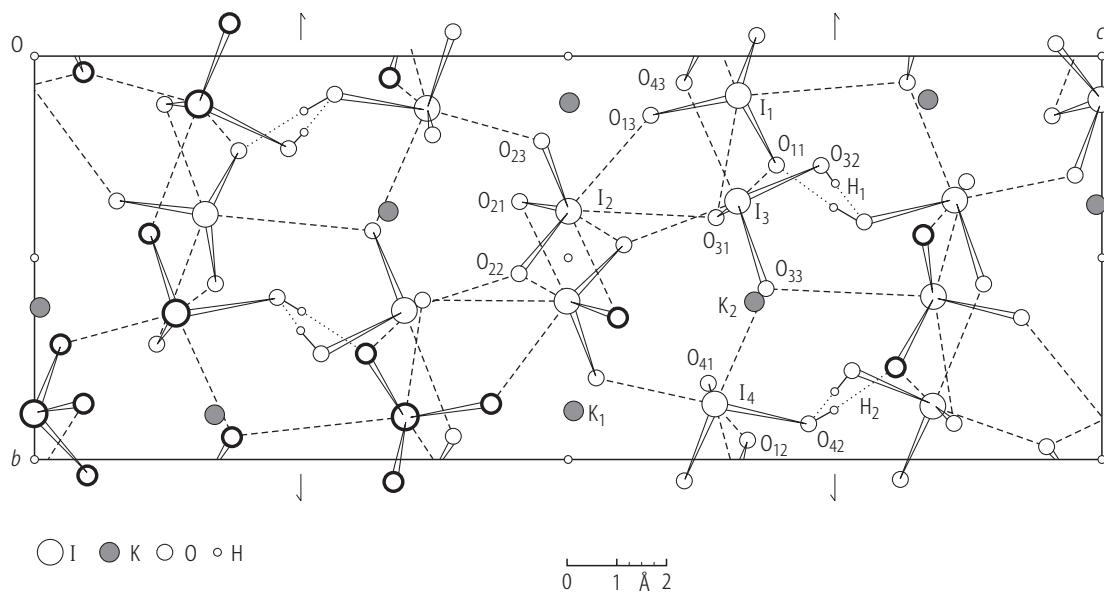
 $\gamma\text{-KIO}_3 \cdot \text{HIO}_3$ 

Fig. 32A-3-006. $\gamma\text{-KIO}_3 \cdot \text{HIO}_3$. Crystal structure at RT [72Kem]. Projection along the a axis. Dotted lines indicate hydrogen bonds; dashed lines weak I...O interactions.

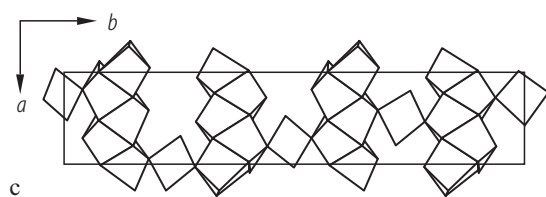
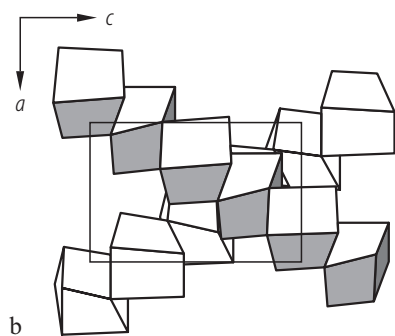
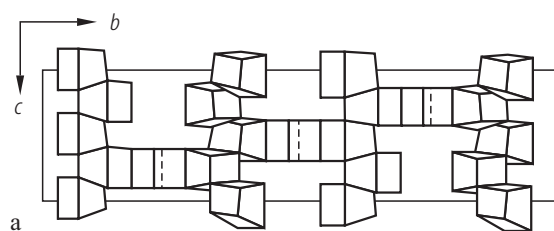
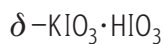


Fig. 32A-3-007. $\delta\text{-KIO}_3 \cdot \text{HIO}_3$. Crystal structure at RT [78Ily]. Fragments of three-dimensional framework of KO_8 polyhedra. (a) Chain of K(1) polyhedra and K(2) link; (b) cross section of framework parallel to (010); (c) projection along the c axis.

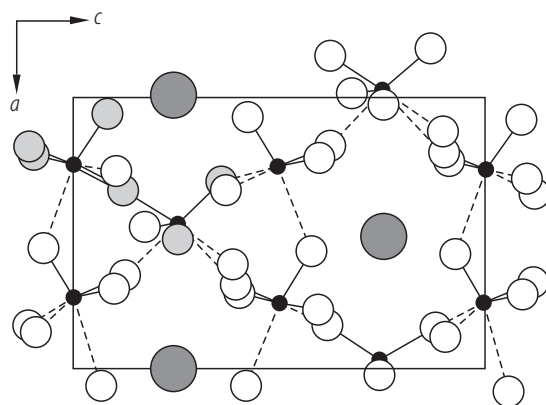
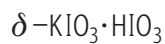


Fig. 32A-3-008. $\delta\text{-KIO}_3 \cdot \text{HIO}_3$. Crystal structure at RT [78Ily]. Projection along the b axis. A layer of structure $z \approx 0 \dots 0.125$ is shown. The large circles are K atoms. The dashed lines are “external” contacts. The dimer atoms are shaded, the contact I(1)–O(32)–I(3) is denoted by the double line.

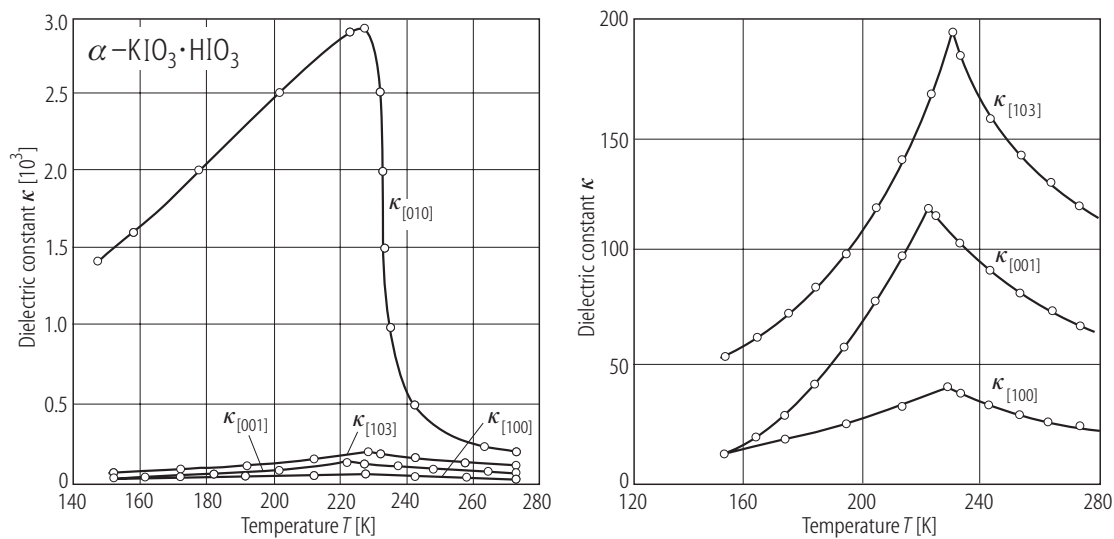


Fig. 32A-3-009. $\alpha\text{-KIO}_3 \cdot \text{HIO}_3$. κ vs. T along [010], [100], [001], [103] [78Pet1]. $f = 1$ kHz. Right panel shows details of $\kappa_{[103]}$, $\kappa_{[001]}$, $\kappa_{[100]}$ near Θ_f .

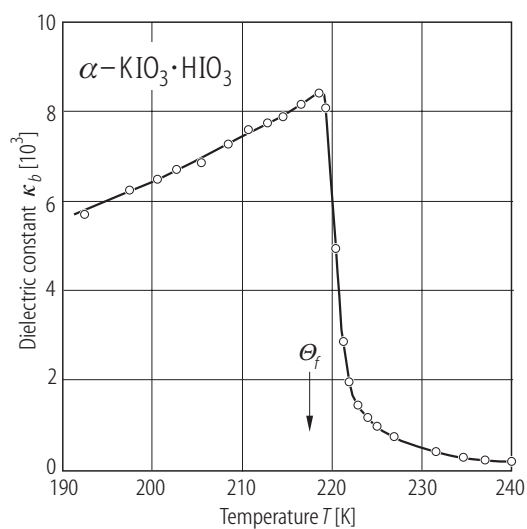


Fig. 32A-3-010. $\alpha\text{-KIO}_3 \cdot \text{HIO}_3$. κ_b vs. T [87Bai]. $f = 230$ kHz.

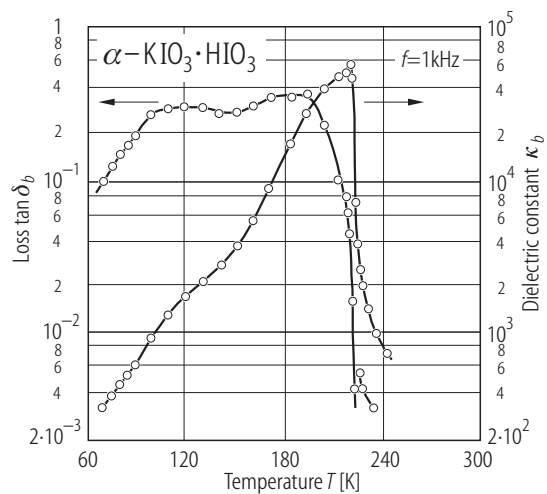


Fig. 32A-3-011. $\alpha\text{-KIO}_3 \cdot \text{HIO}_3$. κ_b , $\tan \delta_b$ vs. T [87Bai]. $f = 1$ kHz.

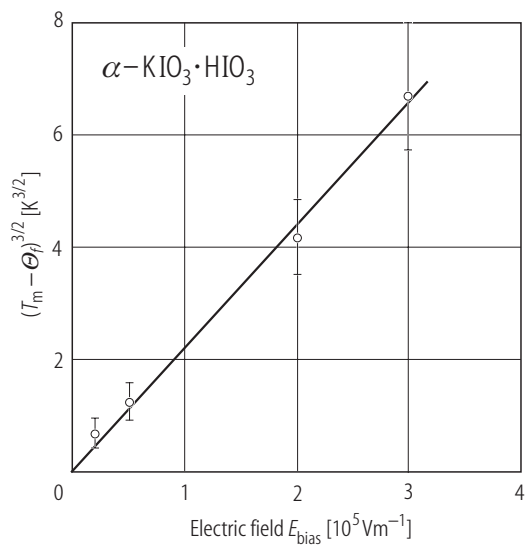


Fig. 32A-3-012. $\alpha\text{-KIO}_3 \cdot \text{HIO}_3$. $(T_m - \Theta_f)^{3/2}$ vs. E_{bias} [91Bai]. T_m : temperature of maximum κ_b .

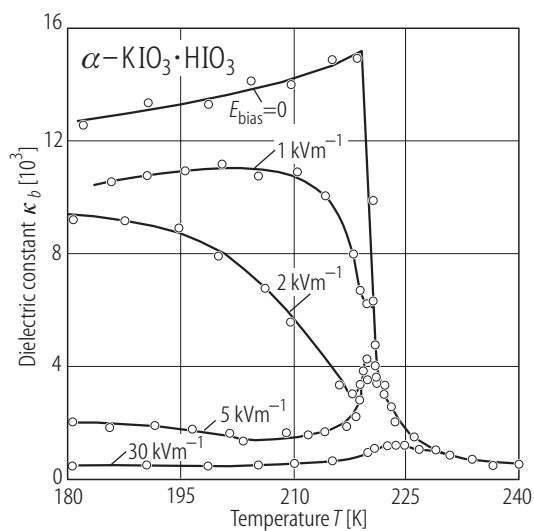


Fig. 32A-3-013. $\alpha\text{-KIO}_3 \cdot \text{HIO}_3$. κ_b vs. T [91Bai]. Parameter: $E_{\text{bias}}, f = 1$ kHz.

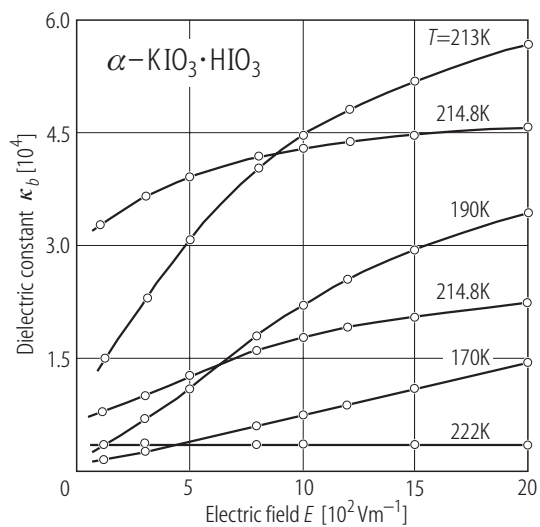


Fig. 32A-3-014. $\alpha\text{-KIO}_3 \cdot \text{HIO}_3$. κ_b vs. E [91Bai]. Parameter: $T, f = 1$ kHz. E : intensity of measuring ac electric field. The sequence of the experiment was 222 K \rightarrow 214.8 K (upper line) \rightarrow 213 K \rightarrow 190 K \rightarrow 170 K \rightarrow 214.8 K (lower line).

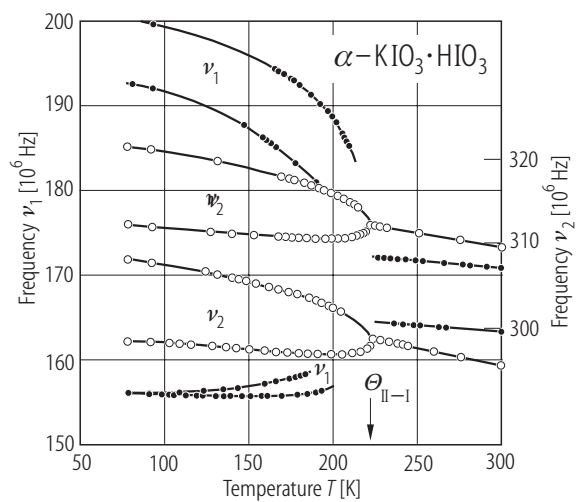


Fig. 32A-3-015. $\alpha\text{-KIO}_3 \cdot \text{HIO}_3$. ν_1 , ν_2 vs. T [78Pet1]. ν_1 , ν_2 : ^{127}I NQR frequency of $(1/2 \leftrightarrow 3/2)$ and $(3/2 \leftrightarrow 5/2)$ transition, respectively.

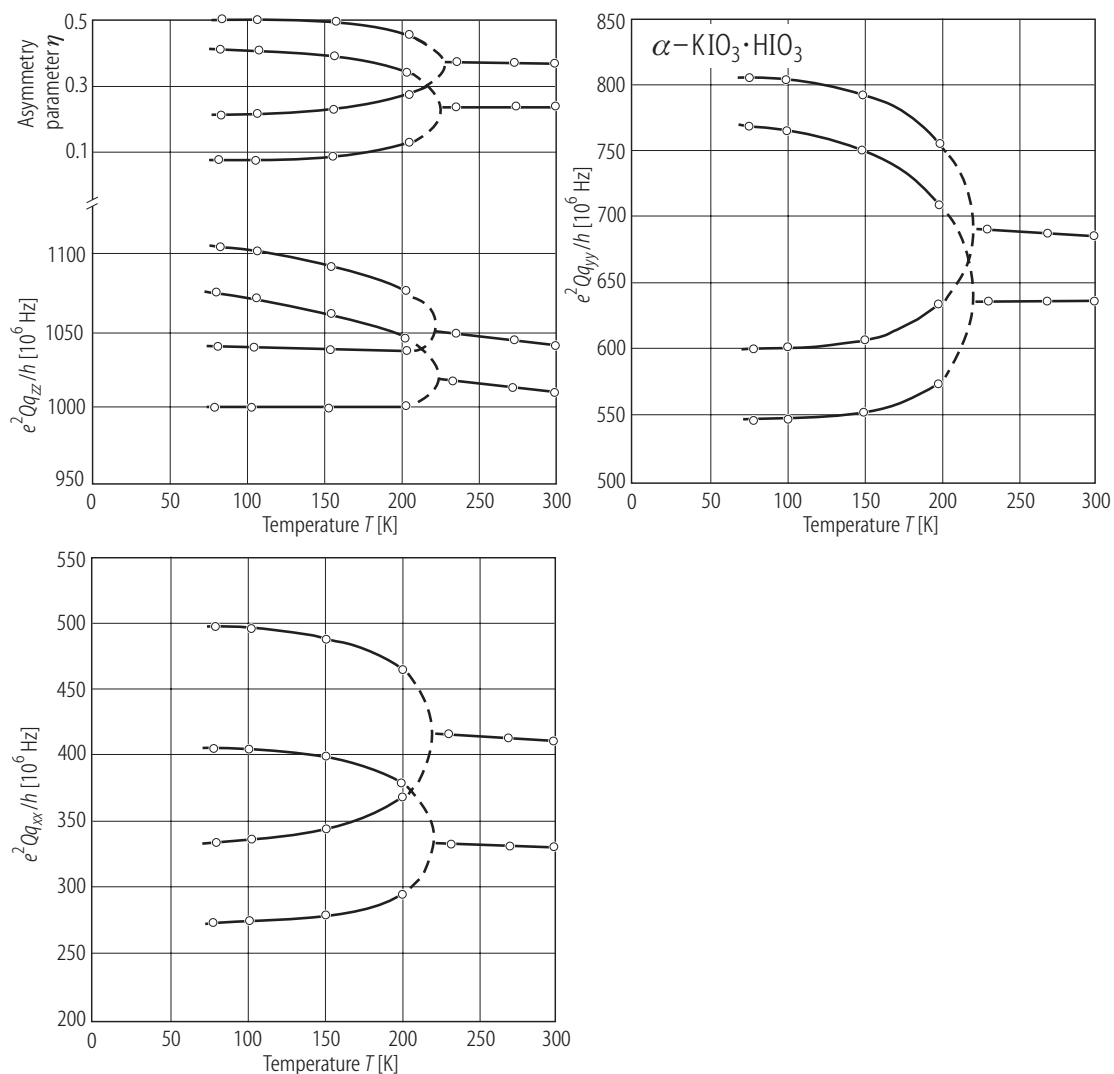


Fig. 32A-3-016. $\alpha\text{-KIO}_3 \cdot \text{HIO}_3$, e^2Qq_{zz}/h , e^2Qq_{yy}/h , e^2Qq_{xx}/h and η vs. T [83Pet]. e^2Qq_{zz}/h , e^2Qq_{yy}/h , e^2Qq_{xx}/h : components of nuclear quadrupole coupling tensor of ^{127}I , η : asymmetry parameter.

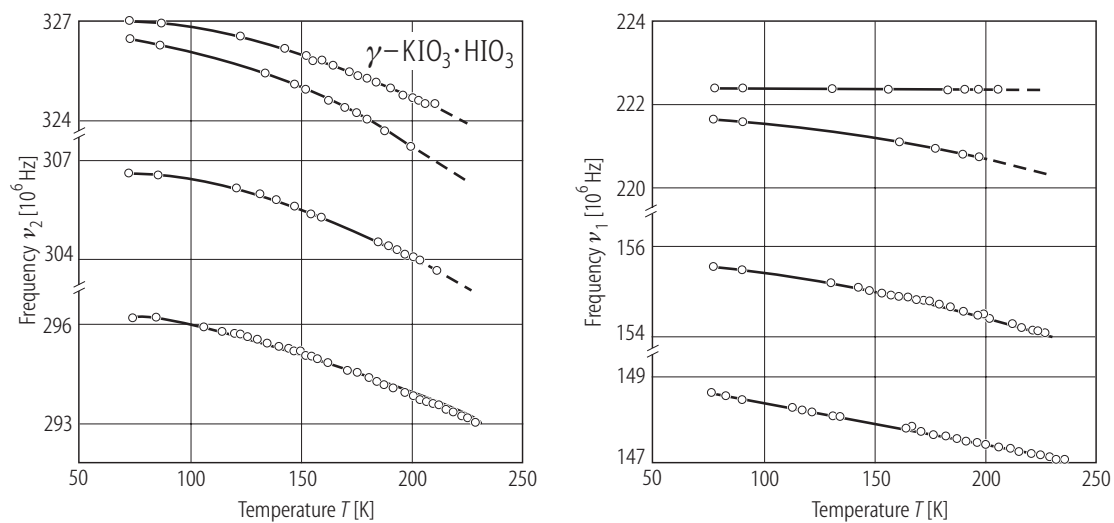


Fig. 32A-3-017. $\gamma\text{-KIO}_3 \cdot \text{HIO}_3$. ν_1 , ν_2 vs. T [78Pet2]. ν_1 , ν_2 : ¹²⁷I NQR frequency of (1/2 \leftrightarrow 3/2) and (3/2 \leftrightarrow 5/2) transition, respectively.