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**Ethanolamine hydrogen *d*-tartrate: optical properties and X-ray diffraction data.\*** By E. G. STEWARD, *Research Laboratories of The General Electric Company Limited, Wembley, England*

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Ethanolamine hydrogen *d*-tartrate,  $C_6H_{13}NO_7$ , may be crystallized from an aqueous solution containing the appropriate proportions of ethanolamine and *d*-tartaric acid.

The crystals, which are piezoelectric, belong to the monoclinic sphenoidal class and have the appearance shown in Fig. 1(a), the sphenoids {011} and {1 $\bar{1}$ 0} defining

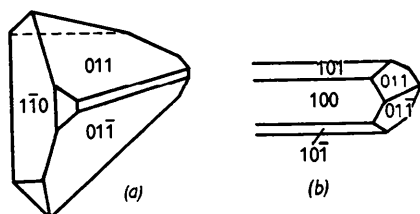


Fig. 1. Crystal habit.

the general shape of the crystals. When grown rapidly from a highly supersaturated solution, however, needles with pronounced pinacoids {100}, {101} and {10 $\bar{1}$ } are formed (Fig. 1(b)).

The principal X-ray diffraction data and optical properties of these crystals are given below and in Table 1.

$C_6H_{13}NO_7$ . Molecular weight = 211.2.  
Density (18° C.) = 5.51 g.cm.<sup>-3</sup>.

\* Communication No. 488 from the Staff of the Research Laboratories of The General Electric Company Limited, Wembley, England.

Table 1. Principal 'powder' lines.

(Intensities visually estimated; *d* not corrected for absorption).

<i>d</i> (Å)	Intensity	<i>d</i> (Å)	Intensity
7.64	<i>w</i>	3.34	<i>s</i>
5.92	<i>s</i>	3.22	<i>m-s</i>
5.75	<i>w</i>	3.12	<i>w</i>
5.59	<i>w</i>	2.96	<i>w</i>
5.34	<i>m</i>	2.86	<i>w-m</i>
4.65	<i>m</i>	2.81	<i>w-m</i>
4.40	<i>m</i>	2.67	<i>m</i>
3.90	<i>m</i>	2.63	<i>w</i>
3.81	<i>vs</i>	2.40	<i>w-m</i>
3.73	<i>m</i>	2.38	<i>m</i>
3.56	<i>w</i>	2.11	<i>w-m</i>
3.46	<i>m</i>		

(*w* = weak; *m* = medium; *s* = strong; *vs* = very strong)

Dimensions ( $\pm 0.03$  Å) of selected monoclinic structure cell:

$$a = 8.83, b = 7.51, c = 7.60 \text{ Å}; \beta = 92^\circ.$$

Molecules per unit cell = 8 (calculated density = 5.57 g.cm.<sup>-3</sup>).

Probable space group:  $P2_1-C_2^2$ ; 010 halved.

Cleavage: 001 (excellent).

Biaxial negative: optical axial angle ( $2V$ ) approximately  $19\frac{1}{2}^\circ$  at 18° C. (sodium light).

Refractive indices ( $\pm 0.001$ ); sodium light; 18° C.:

$$\gamma = 1.551 \text{ parallel to } b.$$

$$\beta = 1.549 \text{ parallel to } a \text{ (to within } \frac{1}{2}^\circ).$$

$$\alpha = 1.485.$$

Dispersion: too weak for positive description.

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**A simple mechanical structure-factor computing aid.** By V. VAND,\* *Chemistry Department, The University, Glasgow W. 2, Scotland*

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A calculating analogue machine has been constructed which has proved useful in structure-factor calculations. It calculates simultaneously eleven values

$$C_n = \cos 2\pi(nu + m), \quad n = 0, 1, 2, \dots, 10,$$

for any value of  $0 \leq u \leq 1$  and  $0 \leq m \leq 1$ .

These two variables are fed into the machine by means of two hand-wheels. In order to reach a three-digit accuracy of setting, each variable is displayed on a pair of dials geared in a 1:10 ratio, the first dial reading the first digit, the second dial the second and third digits. The machine essentially consists of a 0:1:2:...:10 gear-box, in which the rotation *u* of the first hand-wheel is multiplied by *n* by a train of gears (see Fig. 1). The

rotation *m* is then added to each product, for example by means of a differential. The sum  $nu + m$  is then displayed as the angular displacement of the pointers on the eleven

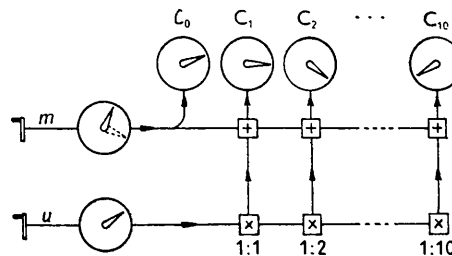


Fig. 1. Principle of the machine.

\* Imperial Chemical Industries Research Fellow.