

apezium faces. Each yhedra in the strucn shares edges wth one other. The corners of an hehedron shares edges. There are two Na(2) with centers $z=0.75$ and $x=0$ are vacant and the final difference er than $1.60 \text{ e}.\text{\AA}^{-3}$. sites does not give squares refinement. This polymorph is m w 460°C .

ces

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or even that there is not crystallographic order in direction in the usual sense, there is little doubt that c is nearly a multiple of 13.8 \AA . Further, there is no entering into a discussion of the elements of crystallography regarding the long pseudo-orthorhombic b axis (*ibid.*, 1967). The crystal diffraction data, some of which is given in Geller (1966), and indeed the results shown in Table 1, should suffice.

Tuinstra (1966) says that 'only in the direction of the fiber axis (c) is an ordinary indexing possible', a conclusion which is negated by the results shown in Table 1. His approach is an arbitrary one; certainly with respect to the directions perpendicular to the helix axes, he has decided arbitrarily on the disorder. Tuinstra (1966) states that the periods along the fiber axis are not indicative of c along this direction, that, for example, the ratio of the heights of the layers '3' and '1' is 2.85. The evidence is not convincing: First, note the good agreement between Q 's with the Q_0 's. Second, measurements made parallel to the rotation axis of rotation photographs cannot be considered to give very reliable spacings. Third, and most important, measurements on our photograph from layer to layer line, and the identity period calculated from them are:

Layer number	Distance (mm)	Identity period (Å)
1	3.25	13.69
2	6.58	13.78
3	10.20	13.79
4	14.47	13.67
5	not observed	
6	25.75	13.84

The average value is 13.75 \AA , but it is not better than 13.8 \AA .

Table 1. Indexing of rotation photograph of ψ -sulphur

Tuinstra Q_0	Present Work Q_0	Present Work Q_c	h k l	Tuinstra Q_0	Present Work Q_0	Present Work Q_c	h k l	Tuinstra Q_0	Present Work Q_0	Present Work Q_c	h k l	
473	478	467	002	*	2625	2634	2,12,3	3792	3774	3759	4,14,3	
613	613	610	080	*	3253	3284	245	4093	3826	3800	425	
1842	1847	1839	0,12,2	*	3713	3742	285	4093	4110	4105	465	
2050	2046	2022	044	*	4092	4137	2,20,1	4751	4719	4715	4,10,5	
2431	2422	2439	0,16,0	*	4451	4501	2,12,5	*	5639	5629	4,14,5	
2470	2475	2480	084	*	5582	5570	2,16,5	*	6472	6503	4,22,3	
2868	2894	2906	0,16,2	*	5839	5814	2,24,1	*	6957	6910	467	
3244	3242	3242	0,12,4	*	6099	6089	247					
4265	4271	4208	006	*	6610	6547	287	2250	2230	2233	660	
		4278	0,20,2	*	946	950	939	370	2707	2706	2701	
		4309	0,16,4	*		950	930	312	2800	2823	2843	6,10,0
4316	4345	4360	046	*	1032	1026	332	3284	3300	3311	6,10,2	
4815	4813	4817	086	1165	1160	1178	352	3797	3807	3798	624	
5449	5515	5587	0,24,0	1222	1224	1244	390	4172	4089	4103	664	
5618	5619	5680	0,20,4	1141	1141	1407	372	*	4762	4712	6,10,4	
*	6631	6646	0,16,6	2104	2107	2093	3,11,2	*	4923	4977	6,18,0	
					2377	2352	314	*	5484	5444	6,18,2	
962	967	941	191		2431	2428	334	*	5640	5628	6,14,4	
1146	1133	1113	113	2500	2532	2550	3,11,2	*	6171	6136	626	
1300	1318	1322	1,11,1		2578	2581	354	*	6521	6501	6,22,0	
1361	1380	1342	153		2616	3,15,0	*	6952	6968	6,22,2		
1933	1934	1876	193	3110	3099	3083	3,15,2	*	7070	7050	6,10,6	
2250	2251	2257	1,11,3		3114	394						
2978	2982	2922	1,17,1	*	3171	3226	3,17,0	2732	**	2699	711	
3113	3086	3060	135	3517	3489	3495	3,11,4	2808	**	2776	731	
				*	5547	5512	3,23,0	2950	**	2928	751	
3255	3233	3212	155	*	6295	6290	3,13,6	3470	**	3461	791	
3470	3454	3441	175	*	7916	7885	3,27,2	3712	**	3710	733	
3560	3518	3608	1,19,1	*	7962	7962	318	3867	**	3843	7,11,1	
3790	3750	3746	195	*	8263	8296	3,25,4	3863	753			
3860	3875	3857	1,17,3	*	8856	8881	3,21,6	4136	**	4092	773	
5130	5135	5118	1,15,5	*	9270	9237	3,27,4	4337	**	4397	793	
5900	5858	5865	137	1318	1316	1300	461	5414	**	5406	10,2,1	
					1942	1936	1910	4,10,1	5550	**	5520	10,4,1
6200	6096	6144	1,23,3		2250	2253	2235	423	6017	**	5978	10,6,1
	6321	6246	177		2866	2856	2824	4,14,1				
*	7005	6932	1,11,7				4,10,3					
		7058	1,25,3									

* Not reported by Tuinstra.

** Not measured in present work.

We emphasize, nevertheless, that we accept the possibility of either a very long axis or lack of order in the fiber axis direction. The nature of the reflections themselves indicates this; some appear sharper than others, and we are not sure that those that are supposed to be in the same layer are all precisely aligned. (However, the crystals are not like those with which most crystallographers usually deal.)

It is difficult to see how Tuinstra did 'index' (his quotes) his data. On page 344 of his paper (1966), he indicates a rectangular prismatic cell, then discusses a β angle of 170° , then that β is undetermined, then speaks of taking as origin for the h index in each reciprocal lattice layer, the 'point nearest to the origin in reciprocal space'. When we look at his Table 2, we find positive and negative h indices; when his $h=3$ for example, he does seem to take a β angle of 170° between his a and c axes of 8.11 and 13.8 \AA length, respectively. This means that the third layer belongs to a cell with $a=8.11$, $b=9.20$, $c=13.8 \text{ \AA}$, $\beta=170^\circ$. Other layers are indexed differently; thus, we must wonder how the intensities were calculated.

References

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