

References

- BLACKMAN, M. & LISGARTEN, N. D. (1957). *Proc. Roy. Soc.* **239**, 93.
 BLACKMAN, M. (1958). *Phil. Mag.* **3**, 831.
 BUTKOVICH, T. R. (1959). *J. Appl. Phys.* **30**, 350.
 GIBBONS, D. F. (1958). *Phys. Rev.* **112**, 136.
 JAKOB, M. & ERK, S. (1929). *Wiss. Abhandl. Phys. Tech. Reichsanstalt*, **12**, 302.
 LONSDALE, K. (1958). *Proc. Roy. Soc. A*, **247**, 424.
 MEGAW, H. D. (1934). *Nature, Lond.* **134**, 900.
 MIKSIC, M. G., SEGERMAN, E. & POST, B. (1959). *Acta Cryst.* **12**, 390.
 POWELL, R. W. (1958). *Proc. Roy. Soc. A*, **247**, 464.
 TRUBY, F. K. (1955). *Science*, **121**, 404.
 VEGARD, L. & HILLESUND, S. (1942). *Avh. Norske Vidensk. Akad.* No. 8, **1**, 24 p.

Acta Cryst. (1960). **13**, 505

X-ray data for the phosphides of aluminium, gallium, and indium. By ARRIGO ADDAMIANO, *Lamp Division, General Electric Company, Nela Park, Cleveland, Ohio, U.S.A.*

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A comparison of the intensities of the lines in powder photographs of AlP and InP obtained in this laboratory with previously reported values (Passerini, 1928; Iandelli,* 1941) has shown significant differences. We have calculated the intensities, using the relation

$$I_c = p|F|^2(1 + \cos^2 2\theta)/(\sin^2 \theta \cos \theta)$$

(p = multiplicity, θ = Bragg angle of reflection, $|F|^2$ = structure factor modulus) and atomic scattering factors for neutral atoms (*Internationale Tabellen zur Bestimmung von Kristallstrukturen*, 1935; Thomas & Umeda, 1957) and find a good agreement with the observed values.

* In private correspondence with Prof. Iandelli we learned that his calculated values for InP include an absorption correction with $\mu r = \infty$.

The relevant data and, for completeness, analogous data for GaP, are reported in Table 1.

The calculations of intensities were done on our Bendix Computer, Model G-15-D, by Mr F. W. Kuhlman, whom we wish to thank.

References

- GIESECKE, G. & PFISTER, H. (1958). *Acta Cryst.* **11**, 369.
 IANDELLI, A. (1941). *Gazz. Chim. Ital.* **71**, 58.
Internationale Tabellen zur Bestimmung von Kristallstrukturen (1935). Berlin: Borntraeger.
 PASSERINI, L. (1928). *Gazz. Chim. Ital.* **58**, 662.
 THOMAS, L. H. & UMEDA, K. (1957). *J. Chem. Phys.* **26**, 293.
 WHITE, W. E. & BUSHEY, A. H. (1944). *J. Amer. Chem. Soc.* **66**, 1666.

Table 1. X-ray data*

<i>hkl</i>	AlP		GaP		InP	
	<i>I</i> _o	<i>I</i> _c × 10 ⁻³	<i>I</i> _o	<i>I</i> _c × 10 ⁻³	<i>I</i> _o	<i>I</i> _c × 10 ⁻³
111	<i>vvs</i>	866	<i>vs</i>	2,933	<i>vs</i>	8,952
200	—	3	<i>vvw</i>	399	<i>ms to w</i>	2,236
220	<i>vs</i>	568	<i>s</i>	1,688	<i>s</i>	4,931
311	<i>s</i>	339	<i>ms</i>	1,136	<i>s</i>	4,040
222	—	53	—	138	<i>vw</i>	684
400	<i>vw</i>	88	<i>vvw</i>	281	<i>w</i>	791
331	<i>ms</i>	136	<i>w</i>	442	<i>ms</i>	1,606
420	—	<1	—	104	<i>w</i>	778
422	<i>ms</i>	250	<i>w</i>	604	<i>ms</i>	1,694
511 + 333	<i>w</i>	83 + 28	<i>vs</i>	281 + 94	<i>ms</i>	939 + 313
440	<i>vvw</i>	72	<i>vvw</i>	225	<i>w</i>	579
531	<i>ms</i>	146	<i>w</i>	523	<i>ms</i>	1,488
600 + 442	—	0 + 44	—	17 + 111	<i>w</i>	98 + 436
620	<i>ms</i>	157	<i>w</i>	520	<i>ms</i>	1,089
533	<i>vw</i>	95	<i>vw</i>	351	<i>w</i>	777
622	—	64	—	160	<i>vw</i>	471
444	<i>w</i>	143	<i>vvw</i>	425	<i>vvw</i>	459
711 + 551					<i>ms</i>	1,081 + 1,081
640					<i>vvw</i>	627
642					<i>s</i>	6,040

(*v* = very; *s* = strong; *w* = weak; *m* = moderately).

* The lattice constants are, in the order: AlP $a_0 = 5.451 \pm 0.004$ Å (present determination. Cf. White & Bushey, 1944); GaP $a_0 = 5.4505$ Å and InP $a_0 = 5.8687$ Å (Giesecke & Pfister, 1958).