

ERRATUM

Erratum: X-ray powder diffraction data for alaptide, 8(S)-methyl-6,9-diazaspiro/4,5/decane-7,10-dione or (cyclo(L-Alanyl-1-ami-no-1-cyclopentan carbonyl), cyclo(L-Ala-Acp) [Powder Diffr. 24, 32–34 (2009)]

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Below is the corrected table originally published on page 33 in the March 2009 issue,

TABLE I. Indexed X-ray powder diffraction data for alaptide. Only the peaks with I_{rel} of 1 or greater are given. Lattice parameters: $a=21.136(4)$ Å, $b=7.212(4)$ Å, $c=6.126(3)$ Å, $V=933.8(8)$ Å³, and $Z=4$. All lines were indexed and are consistent with the $P2_12_12_1$ space group.

$2\theta_{\text{obs}}$ (°)	d_{obs} (Å)	I_{rel}	h	k	l	$2\theta_{\text{cal}}$ (°)	d_{cal} (Å)	$\Delta 2\theta$ (°)
8.414	10.501	100	2	0	0	8.356	10.568	0.058
14.909	5.937	81	2	1	0	14.859	5.957	0.050
16.735	5.293	1	2	0	1	16.714	5.300	0.021
17.634	5.026	1	3	1	0	17.584	5.040	0.050
20.875	4.252	2	4	1	0	20.823	4.2625	0.052
22.222	3.9972	1	4	0	1	22.200	4.0012	0.022
22.841	3.8902	2	3	1	1	22.831	3.8919	0.010
24.699	3.6016	24	0	2	0	24.668	3.6061	0.031
25.609	3.4757	4	5	0	1	25.582	3.4793	0.027
26.118	3.4091	24	2	2	0	26.089	3.4129	0.021
27.803	3.2062	2	3	2	0	27.769	3.2101	0.034
28.227	3.1589	1	6	1	0	28.170	3.1653	0.057
30.008	2.9754	2	4	2	0	29.976	2.9786	0.032
31.441	2.8430	1	3	2	1	31.438	2.8433	0.003
32.655	2.7401	4	5	2	0	32.613	2.7435	0.042
33.897	2.6424	1	8	0	0	33.902	2.6420	-0.005
35.646	2.5167	2	6	2	0	35.600	2.5200	0.046
36.162	2.4820	2	8	1	0	36.179	2.4808	-0.017
36.241	2.4767	1	5	0	2	36.187	2.4803	0.054
37.059	2.4239	1	8	0	1	37.026	2.4260	0.033
38.909	2.3128	1	6	0	2	38.934	2.3114	-0.025

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