Additions and Corrections

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D. J. McCabe, E. N. Duesler, and R. T. Paine*: Monodentate Coordination by a Tripodal Ligand System: Synthesis and Crystal and Molecular Structure of Bis[diisopropyl [1,2-bis(diethylcarbamoyl)-ethyl]phosphonate]erbium(III) Nitrate Monohydrate.

Page 4627. In Table II, the last 12 atom coordinates were omitted. Those atom positions are listed here.

atom	*x*	*y*	*z*
O(11)	0.1806 (5)	0.0188 (3)	0.7866 (4)
O(12)	0.1829 (5)	0.0277 (2)	0.8949 (3)
O(13)	0.1337 (6)	−0.0554 (3)	0.8402 (4)
N(6)	0.2446 (7)	0.1373 (4)	0.9874 (5)
O(14)	0.1619 (5)	0.1496 (3)	0.9409 (3)
O(15)	0.3186 (4)	0.1204 (3)	0.9636 (3)
O(16)	0.2527 (5)	0.1399 (4)	1.0502 (4)
N(7)	0.2535 (6)	0.2491 (3)	0.8383 (4)
O(17)	0.1653 (5)	0.2263 (3)	0.8226 (4)
O(18)	0.3281 (4)	0.2107 (2)	0.8522 (3)
O(19)	0.2712 (5)	0.3028 (2)	0.8393 (4)
O(20)	0.2062 (3)	0.1353 (2)	0.7201 (2)

—R. T. Paine