ERRATUM

Structure of tetra(*n*-butyl)ammonium tetraiodo- μ, μ' -diiododiplatinate(II), [(*n*-C₄H₉)₄N]₂ [Pt₂I₆]¹

A computer programming error resulted in the omission of several minus signs from the published coordinates (Table 3, page 386) and thermal parameters (Table 4, page 387). The corrected tables are listed below. We thank E. C. Lingafelter for calling our attention to this problem.

¹R. D. Rogers, H. Isci, and W. R. Mason (1984) *J. Cryst. Spectrosc. Res.* **14**, 383.

Table 3. Final fractional coordinates for
 $[(n\text{-C}_4\text{H}_9)_4\text{N}]_2[\text{Pt}_2\text{I}_6]$

Atom	x/a	y/b	z/c
Pt(1)	0.3705(2)	0.7825(2)	0.0996(1)
Pt(2)	0.1310(2)	0.8060(2)	0.1414(1)
I(1)	0.1985(4)	0.7449(4)	0.0558(2)
I(2)	0.3028(4)	0.8465(4)	0.1841(2)
I(3)	0.4266(4)	0.7141(4)	0.0110(2)
I(4)	0.5373(4)	0.8291(4)	0.1511(2)
I(5)	-0.0379(4)	0.7650(4)	0.0900(2)
I(6)	0.0790(4)	0.8669(4)	0.2325(2)
N(1)	0.206(4)	-0.067(3)	0.862(2)
N(2)	0.277(4)	0.450(4)	0.886(2)
C(1)	0.249(4)	0.015(4)	0.899(2)
C(2)	0.303(5)	-0.011(5)	0.960(3)
C(3)	0.336(5)	0.080(5)	0.989(3)
C(4)	0.386(6)	0.058(6)	1.054(3)
C(5)	0.282(5)	-0.144(5)	0.861(3)
C(6)	0.362(5)	-0.110(5)	0.830(3)
C(7)	0.424(5)	-0.196(5)	0.834(3)
C(8)	0.510(6)	-0.173(6)	0.806(3)
C(9)	0.131(5)	-0.118(5)	0.890(3)
C(10)	0.047(5)	-0.049(5)	0.896(3)
C(11)	-0.016(5)	-0.116(5)	0.927(3)
C(12)	-0.102(6)	-0.055(6)	0.933(4)
C(13)	0.165(4)	-0.019(4)	0.806(2)
C(14)	0.123(5)	-0.093(5)	0.766(3)
C(15)	0.090(5)	-0.043(5)	0.701(3)
C(16)	0.180(5)	-0.027(5)	0.679(3)
C(17)	0.176(5)	0.497(5)	0.880(3)
C(18)	0.182(5)	0.589(5)	0.914(3)
C(19)	0.080(5)	0.627(5)	0.910(3)
C(20)	0.028(5)	0.566(5)	0.948(3)
C(21)	0.338(5)	0.520(5)	0.857(3)
C(22)	0.428(7)	0.480(6)	0.853(4)
C(23)	0.489(6)	0.547(6)	0.824(3)
C(24)	0.581(7)	0.506(6)	0.819(4)
C(25)	0.320(5)	0.433(4)	0.945(3)
C(26)	0.264(5)	0.382(5)	0.980(3)
C(27)	0.318(6)	0.373(6)	1.044(4)
C(28)	0.265(6)	0.313(6)	1.077(3)
C(29)	0.257(5)	0.359(5)	0.849(3)
C(30)	0.210(5)	0.366(5)	0.788(3)
C(31)	0.198(6)	0.262(5)	0.767(3)
C(32)	0.152(8)	0.264(8)	0.698(5)

Table 4. Thermal parameters for $[(n\text{-C}_4\text{H}_9)_4\text{N}]_2[\text{Pt}_2\text{I}_6]$

Atom	U_{11}	U_{22}	U_{33}	U_{12}	U_{13}	U_{23}
Pt(1)	0.021(3)	0.041(2)	0.038(2)	0.004(2)	0.004(2)	0.001(2)
Pt(2)	0.021(2)	0.039(2)	0.033(2)	0.000(2)	0.005(2)	-0.004(2)
I(1)	0.031(4)	0.077(4)	0.038(3)	-0.004(3)	0.004(3)	-0.018(3)
I(2)	0.029(4)	0.099(5)	0.049(3)	-0.006(3)	-0.001(3)	-0.031(3)
I(3)	0.040(4)	0.078(4)	0.055(3)	0.014(3)	0.018(3)	-0.004(3)
I(4)	0.021(4)	0.058(4)	0.082(4)	-0.001(3)	0.001(3)	-0.001(3)
I(5)	0.026(4)	0.055(4)	0.068(4)	-0.008(3)	0.005(3)	-0.002(3)
I(6)	0.049(4)	0.061(4)	0.048(3)	-0.010(3)	0.018(3)	0.010(3)
N(1)	2(1)					
N(2)	3(1)					
C(1)	2(1)					
C(2)	4(2)					
C(3)	4(2)					
C(4)	6(2)					
C(5)	4(2)					
C(6)	4(2)					
C(7)	4(2)					
C(8)	6(2)					
C(9)	3(2)					
C(10)	4(2)					
C(11)	4(2)					
C(12)	7(2)					
C(13)	2(1)					
C(14)	4(2)					
C(15)	4(2)					
C(16)	4(2)					
C(17)	4(2)					
C(18)	4(2)					
C(19)	4(2)					
C(20)	5(2)					
C(21)	4(2)					
C(22)	7(2)					
C(23)	6(2)					
C(24)	7(2)					
C(25)	3(2)					
C(26)	4(2)					
C(27)	7(2)					
C(28)	6(2)					
C(29)	4(2)					
C(30)	4(2)					
C(31)	6(2)					
C(32)	10(3)					