

## ERRATUM

**Preparation and structure of bis-tetraethylammonium(1,2-dicyanoethylene-1,2-dithiolato)(1,1-dicyanoethylene-1,1-diselenato)nickelate(II),  $[(\text{CN})_2\text{C}_2\text{Se}_2\text{NiS}_2\text{C}_2(\text{CN})_2]^{2-} [\text{N}(\text{C}_2\text{H}_5)_4]_2^+$ <sup>1</sup>**

and

**Dichloro-bis(triphenylphosphine)nickel,  $\text{Cl}_2\text{Ni}(\text{P}(\text{C}_6\text{H}_5)_3)_2$ , a redetermination<sup>2</sup>**

The cell constants are:

- for (1), page 617 (NISOMS): space group  $P\bar{1}$  ( $Z = 1$ ):  
 $a = 7.515(2)$ ,  $b = 8.717(2)$ ,  $c = 12.870(2)$  Å,  
 $\alpha = 86.84(2)$ ,  $\beta = 76.40(2)$ ,  $\gamma = 76.27(2)$ °.
- for (2), page 623 (SPAN2): space group  $P2/c$  ( $Z = 2$ ):  
 $a = 11.712(2)$ ,  $b = 8.258(1)$ ,  $c = 17.370(2)$  Å,  
 $\beta = 106.55(2)$ °.

The original manuscripts were accepted for publication as a “NOTE,” and the abstracts, containing the cell dimensions, were left out.

Fortunately, the cell dimensions are given in the same issue of *J. Crystallogr. Spectrosc. Res.*, page 601, Table 1, under the headings NISOMS and SPAN2.

<sup>1</sup>Bruins Slot, H. J., Noordik, J. H., Beurskens, P. T., Keijzers, C. P., Dietzsch, W., and Kirmse, R. (1984) *J. Crystallogr. Spectrosc. Res.* **14**, 617.

<sup>2</sup>Bruins Slot, H. J., Van Havere, W. K. L., Noordik, J. H., Beurskens, P. T., and Royo, P. (1984) *J. Crystallogr. Spectrosc. Res.* **14**, 623.