plane of the central indole ring [dihedral angle 15·1 (2)°], with an average separation of 3·528 Å.

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SHORT COMMUNICATION

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Acta Cryst. (1981). B**37**, 2120

Room- and high-pressure neutron structure determination of tetrathiafulvalene–7,7,8,8-tetracyano-*p*-quinodimethane (TTF–TCNQ). Thermal expansion and isothermal compressibility. Errata.

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Abstract

Three errors in the paper by Filhol, Bravic, Gaultier, Chasseau & Vettier [*Acta Cryst.* (1981), B**37**, 1225–1235] are corrected. On page 1230, right-hand column, line 7 should read: 'In fact the C(2)···N(2) distance given in Table 6 of Kistenmacher *et al.* (1974) and wrongly used by Chasseau *et al.* (1978), is clearly not in the above directions. It is thus of importance for the discussion below to note the following point: the C(2)–H···N(2) distance referred to by these latter authors is in error (3·36 instead of 3·53 Å) and this hydrogen bond does not exist at room pressure and temperature'. In Table 6, line N(2)···D(2): the values 1·00 and 0·94 (2) should be replaced by 2·62 and 2·59 (2) respectively. In the caption of Fig. 7 the expression for α_i should read $\alpha_i = (1/l_i)(dl_i/dT)$.

All information is given in the *Abstract*.

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