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Erratum

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Erratum

Relationship Between IR C=O Frequenceis and their Solid State Structures, V. M. Kolb, T. E. Janota, C. L. Dantzman, M. L. Kozenski, D. P. Strommen, J. P. Snyder and G. E. Tipsword, *Mol. Cryst., Liq. Cryst.*, **211**, pp. 403-406 (1992).

Additional deuteration experiments and analysis of data suggest that the bands assigned as C=N for *o*-OCH_3-APSC, *m*-OCH_3-APSC, *o*-NO_2-APSC and *o*-F-APSC (p. 404) are actually C=C aromatic bands. The corresponding C=N bands are found at 1628, 1627, 1613 and 1621 cm^{-1}, respectively.