

# Preparation and Characterization of 2-Phenoxyethyl(thiophen-2-yl)tellane ( $\text{C}_4\text{H}_3\text{S})\text{TeCH}_2\text{CH}_2\text{OC}_6\text{H}_5$

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The synthesis and structure of  $(\text{C}_4\text{H}_3\text{S})\text{TeCH}_2\text{CH}_2\text{OC}_6\text{H}_5$  (**1**) ( $\text{C}_4\text{H}_3\text{S}$  = thiophen-2-yl) are reported and compared to those of the analogous selenium compound  $(\text{C}_4\text{H}_3\text{S})\text{SeCH}_2\text{CH}_2\text{OC}_6\text{H}_5$  previously synthesized by our group. The compound was characterized by  $^1\text{H}$ ,  $^{13}\text{C}\{^1\text{H}\}$ -, and  $^{125}\text{Te}$ -NMR spectroscopy as well as by X-ray single crystal crystallography. **1** crystallizes in the monoclinic crystal system, space group  $P2_1$ , with  $Z = 2$ , and unit cell dimensions  $a = 10.618(2)$  Å,  $b = 5.357(1)$  Å,  $c = 10.684(2)$  Å,  $\beta = 96.57(3)^\circ$ . The lattice is composed of discrete molecules that are joined together by weak hydrogen bonds into a three-dimensional network. The thiophen-2-yl ring is disordered and shows two alternative orientations with the site occupation factors of 0.70(1) and 0.30(1). All bond parameters are quite normal. The comparison of the lattices in **1** and in its selenium analogue shows that while the closest intermolecular contacts are similar, the packing of the molecules is different.

*Key words:* Telluroether, 2-Phenoxyethyl(thiophen-2-yl)tellane, X-Ray Crystallography, NMR Spectroscopy