## Absolute Configuration of 2-C-Methylerythritol from Convolvulus glomeratus

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In a recent communication <sup>1</sup> we have reported the isolation of a new branched alditol (I) from Convolvulus glomeratus Choisy. The constitution and relative stereochemistry was deduced by spectroscopic methods and finally by synthesis of racemic I from eitraconic acid.

In order to elucidate the absolute stereochemistry of I, a CD study of I and the two isobutyltetrols 2 and 3 of known configurations  $^2$  has been performed. As expected, no extrema were reached in the CD measurements on aqueous solutions of I-3, and only low molecular ellipticities were obtained between 190 and 200 nm. Much more informative spectra were obtained, however, when the tetrols were investigated as their molybdate(VI) complexes (Fig. 1). From these spectra and the known erythro configuration of  $I^1$  it is obvious that I has the 2S, 3R configuration. Furthermore, since the CD band in the 330 nm region is positive for the erythritols and negative for the threitol this result suggests a method to distinguish between  $2 \cdot C$ -alkylerythritols and  $2 \cdot C$ -alkylthreitols.

Experimental. CD spectra were recorded on a Cary 60 instrument using a 1 cm cell for the 5.0 mM solutions of 1-3, and a 0.05 cm cell for the solutions of the molybdate complexes. Following the conditions used by Voelter et al., the latter solutions were 5.0 mM with respect to tetrol and 10.0 mM with respect to sodium molybdate. The pH's were 5.0-5.1, adjusted with hydrochloric acid and measured with a

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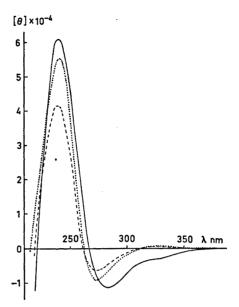


Fig. 1. CD spectra of molybdate(VI) complexes of  $1, \dots; 2, \dots;$  and  $3, \dots$ 

Beckman pH meter (accuracy 0.1 unit). The molecular ellipticities are given in degree mol<sup>-1</sup> cm<sup>2</sup>. Obtained for 1-3:  $[\theta]_{192.5} = -0.18 \times 10^3$ ,  $-0.24 \times 10^3$ ,  $+0.10 \times 10^3$ , respectively. Obtained for 1-3 as molybdate complexes: see Fig. 1.

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- Anthonsen, T., Hagen, S., Kazi, M. A., Shah, S. W. and Tagar, S. Acta Chem. Scand. B 30 (1976) 91.
- B 30 (1976) 91.
  Behr, D., Dahmén, J. and Leander, K. Acta Chem. Scand. B 30 (1976) 309.
- Voelter, W., Bayer, E., Records, R., Bunnenberg, E. and Djerassi, C. Chem. Ber. 102 (1969) 1005.

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