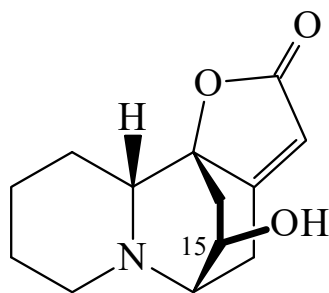




>>Entry Name: Securinol A



**CAS Registry Number:** 5008-48-0

**Molecular Formula:** C<sub>13</sub>H<sub>17</sub>NO<sub>3</sub>

**Molecular Weight:** 235.282

**Accurate Mass:** 235.120844

**Percentage Composition:** C 66.36%; H 7.28%; N 5.95%; O 20.40%

**Biological Source:** Alkaloid from the leaves of *Securinega suffruticosa* and from *Margaritaria indica* (Euphorbiaceae)

**Physical Description:** Cryst. (C<sub>6</sub>H<sub>6</sub>)

**Melting Point:** Mp 135-136°

**Optical Rotation:** [ $\alpha$ ]<sub>D</sub><sup>17</sup>+58.2 (c, 0.14 in CHCl<sub>3</sub>)

>>Derivative: Picrate

**Physical Description:** Cryst. + 1H<sub>2</sub>O (EtOH)

**Melting Point:** Mp 186-187 dec.°

>>Derivative: 15-Epimer

**Synonym(s):** Securinol B

**CAS Registry Number:** 30155-10-3

**Molecular Formula:** C<sub>13</sub>H<sub>17</sub>NO<sub>3</sub>

**Molecular Weight:** 235.282

**Accurate Mass:** 235.120844

**Percentage Composition:** C 66.36%; H 7.28%; N 5.95%; O 20.40%

**Biological Source:** Alkaloid from leaves of *Securinega suffruticosa* (Euphorbiaceae)

**Physical Description:** Needles (petrol)

**Melting Point:** Mp 159-160°

**Optical Rotation:** [ $\alpha$ ]<sub>D</sub> +120 (c, 0.15 in EtOH)

**Other Data:** Probable struct. based on revised struct. for Securinol A. Provisional numbering

>>Derivative: 15-Epimer, methanesulfonyl

**Melting Point:** Mp 176-177°

**References:**

Horii, Z. *et al.*, *Chem. Pharm. Bull.*, 1965, **13**, 1307; 1970, **18**, 2009, (*isol, uv, ord, ir, pmr, ms*)

Arbain, D. *et al.*, *J.C.S. Perkin 1*, 1991, 1863, (*isol, pmr, cmr, cryst struct, abs config*)