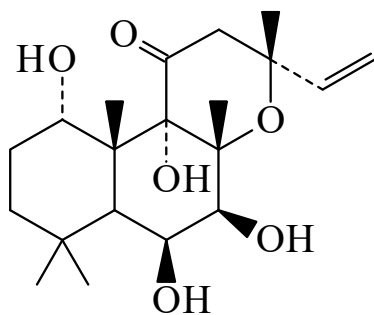




>>Entry Name: 8,13-Epoxy-1,6,7,9-tetrahydroxy-14-labden-11-one



**CAS Registry Number:** 66428-84-0

**Related CAS Registry Numbers(s):** 64657-11-0 112420-42-5 114298-68-9

**Molecular Formula:** C<sub>20</sub>H<sub>32</sub>O<sub>6</sub>

**Molecular Weight:** 368.469

**Accurate Mass:** 368.21989

**Percentage Composition:** C 65.19%; H 8.75%; O 26.05%

>>Variant: (1 $\alpha$ ,6 $\beta$ ,7 $\beta$ ,8 $\alpha$ ,9 $\alpha$ ,13*R*)-form

**CAS Registry Number:** 64657-20-1

**Molecular Formula:** C<sub>20</sub>H<sub>32</sub>O<sub>6</sub>

**Molecular Weight:** 368.469

**Accurate Mass:** 368.21989

**Percentage Composition:** C 65.19%; H 8.75%; O 26.05%

**Biological Source:** Constit. of *Coleus forskohlii*

**Physical Description:** Cryst.

**Melting Point:** Mp 177-180°

**Sigma:** D3533

>>Derivative: 6-Ac

**Synonym(s):** Coleonol B

**CAS Registry Number:** 64657-21-2

**Molecular Formula:** C<sub>22</sub>H<sub>34</sub>O<sub>7</sub>

**Molecular Weight:** 410.506

**Accurate Mass:** 410.230455

**Percentage Composition:** C 64.37%; H 8.35%; O 27.28%

**Biological Source:** Constit. of *Coleus forskohlii*

**Physical Description:** Cryst.

**Melting Point:** Mp 208-210°

**Optical Rotation:** [ $\alpha$ ]<sub>D</sub> +2.92 (c, 1 in CHCl<sub>3</sub>)

**Sigma:** A9420

>>Derivative: 7-Ac

**Synonym(s):** Coleonol. Forskolin. **Colforsin**, INN, USAN. Boforsin. HL 362

**CAS Registry Number:** 66575-29-9

**Molecular Formula:** C<sub>22</sub>H<sub>34</sub>O<sub>7</sub>

**Molecular Weight:** 410.506

**Accurate Mass:** 410.230455

**Percentage Composition:** C 64.37%; H 8.35%; O 27.28%

**Biological Source:** Constit. of *Coleus forskohlii*

**Biological Use/Importance:** Potassium channel (K<sub>v</sub>) blocker. Antihypertensive agent used to treat glaucoma

**Physical Description:** Cryst.

**Melting Point:** Mp 230-232°

**Optical Rotation:** [ $\alpha$ ]<sub>D</sub><sup>25</sup> -26.2 (c, 1.68 in CHCl<sub>3</sub>)

**Solubility:** BERDY SOL: Sol. MeOH, C<sub>6</sub>H<sub>6</sub>; poorly sol. H<sub>2</sub>O, hexane

**Calculated Log P Data:** Log P 1.08 (calc)

**UV:** [neutral]  $\lambda_{\max}$  0 (end) ( $\epsilon$ ) (MeOH) (Derep)

**Fluka:** 47735

**Sigma:** F3917