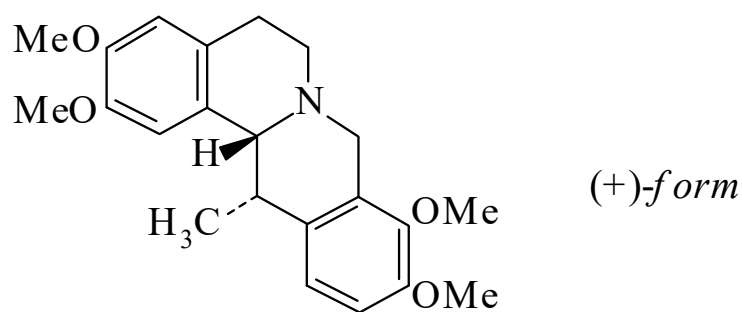




>>Entry Name: Corydaline

Synonym(s): 5,8,13,13a-Tetrahydro-2,3,9,10-tetramethoxy-13-methyl-6H-dibenzo[a,g]quinolizine, 9Cl. 2,3,9,10-Tetramethoxy-13-methylberbine, 8Cl. Corydalis A



**Molecular Formula:** C<sub>22</sub>H<sub>27</sub>NO<sub>4</sub>

**Molecular Weight:** 369.46

**Accurate Mass:** 369.194009

**Percentage Composition:** C 71.52%; H 7.37%; N 3.79%; O 17.32%

>>Variant: (+)-form

**CAS Registry Number:** 518-69-4

**Molecular Formula:** C<sub>22</sub>H<sub>27</sub>NO<sub>4</sub>

**Molecular Weight:** 369.46

**Accurate Mass:** 369.194009

**Percentage Composition:** C 71.52%; H 7.37%; N 3.79%; O 17.32%

**Biological Source:** Alkaloid from *Corydalis tuberosa* and many other *Corydalis* spp. (Fumariaceae)

**Use/Importance:** Shows analgesic and antirheumatic props.; has been evaluated clinically but with disappointing results

**Physical Description:** Prisms (EtOH)

**Melting Point:** Mp 135°

**Optical Rotation:** [ $\alpha$ ]<sub>D</sub><sup>25</sup> +295 (EtOH)

**Hazard & Toxicity:** LD<sub>50</sub> (mus, ivn) 136 mg/kg

**Rare Chemicals Library:** S44912-1

>>Variant: (±)-form

**CAS Registry Number:** 6018-35-5

**Molecular Formula:** C<sub>22</sub>H<sub>27</sub>NO<sub>4</sub>

**Molecular Weight:** 369.46

**Accurate Mass:** 369.194009

**Percentage Composition:** C 71.52%; H 7.37%; N 3.79%; O 17.32%

**Source/Synthesis:** Synthetic

**Melting Point:** Mp 135°

**References:**

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Hughes, D.W. *et al.*, *Can. J. Chem.*, 1976, **54**, 2252, (*cmr*)

Kametani, T. *et al.*, *J.C.S. Perkin I*, 1977, 1151, (*synth, pmr*)

Iwasa, K. *et al.*, *J.O.C.*, 1981, **46**, 4744, (*synth, abs config*)

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Hanaoka, M. *et al.*, *Chem. Pharm. Bull.*, 2000, **48**, 399-404, (*synth*)