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ProductInformation

Aconitine

Product Number **A8001** Store at Room Temperature

Product Description

Molecular Formula: C₃₄H₄₇NO₁₁ Molecular Weight: 645.7 CAS Number: 302-27-2 Melting Point: 204 °C¹

Specific Rotation: $+17.3^{\circ}$ (chloroform)¹ Synonym: $(1\alpha,3\alpha,6\alpha,14\alpha,15\alpha,16\beta)$ -20-ethyl-1,6,16-trimethoxy-4-(methoxymethyl)aconitane-3,8,13,14,15-pentol 8-acetate 14-benzoate¹

Aconitine is an alkaloid with a diterpenoid structure that occurs naturally in the roots of *Aconitum napellus* (*Ranunculaceae*), or the monkshood plant. Being the principal alkaloid in the *A. napellus* root, aconitine has been studied for its pharmacological properties. One mechanism of action is suppression of the inactivation of voltage-dependent Na $^+$ channels via the binding to neurotoxin binding site 2 of the α -subunit of the channel protein. A review of the use of aconitine and other compounds to target voltage-gated sodium channels has been published.

The interaction of aconitine (10-100 nM) and 6-benzoylheteratisine (3 μ M) in the rat hippocampus has been studied.⁵ The effect of aconitine on the expression of MHC class II in macrophages has been investigated.⁶ Aconitine (1 μ M) has been shown to suppress both normal and epileptiform activity in rat hippocampal slices as an anticonvulsive agent.⁷

HPLC/SPE/MS and capillary LC/FAB-MS methods for the analysis of aconitine from biological samples have been published.^{8,9}

Precautions and Disclaimer

For Laboratory Use Only. Not for drug, household or other uses.

Preparation Instructions

This product is soluble in chloroform (50 mg/ml), yielding a clear, colorless to light yellow solution. Its solubility in chloroform has been reported to be 500 mg/ml. It is also soluble in absolute alcohol (35 mg/ml), and ether (20 mg/ml).

References

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