Product Data Sheet

Osthenol

Cat. No.: HY-N2554 CAS No.: 484-14-0 Molecular Formula: $C_{14}H_{14}O_{3}$ Molecular Weight: 230.26

Monoamine Oxidase Target: Pathway: **Neuronal Signaling**

Storage: Powder -20°C

> 4°C 2 years

3 years

In solvent -80°C 6 months

> -20°C 1 month

SOLVENT & SOLUBILITY

In Vitro

DMSO: 100 mg/mL (434.29 mM; Need ultrasonic)

Preparing Stock Solutions	Solvent Mass Concentration	1 mg	5 mg	10 mg
	1 mM	4.3429 mL	21.7146 mL	43.4292 mL
	5 mM	0.8686 mL	4.3429 mL	8.6858 mL
	10 mM	0.4343 mL	2.1715 mL	4.3429 mL

Please refer to the solubility information to select the appropriate solvent.

In Vivo

- 1. Add each solvent one by one: 10% DMSO >> 40% PEG300 >> 5% Tween-80 >> 45% saline Solubility: 2.5 mg/mL (10.86 mM); Suspended solution; Need ultrasonic
- 2. Add each solvent one by one: 10% DMSO >> 90% (20% SBE- β -CD in saline) Solubility: ≥ 2.5 mg/mL (10.86 mM); Clear solution
- 3. Add each solvent one by one: 10% DMSO >> 90% corn oil Solubility: ≥ 2.5 mg/mL (10.86 mM); Clear solution

BIOLOGICAL ACTIVITY

Description	competitive human monoami	ted coumarin isolated from the dried roots of Angelica pubescens, is selective, reversible, and ine oxidase-A (hMAO-A) inhibitor (K_i =0.26 μ M). Osthenol potently inhibits recombinant hMAO-A ows a high selectivity index for hMAO-A versus hMAO-B ^[1] .
IC ₅₀ & Target	hMAO-A 0.26 μM (Ki)	hMAO-A 0.74 μM (IC ₅₀)

l. Baek SC, et al. Osthenol, a p	
	prenylated coumarin, as a monoamine oxidase A inhibitor with high selectivity. Bioorg Med Chem Lett. 2019;29(6):839-843.
	Caution: Product has not been fully validated for medical applications. For research use only.
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REFERENCES

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