

# **Product** Data Sheet

## **Rhodionin**

Cat. No.:HY-N0241CAS No.:85571-15-9Molecular Formula: $C_{21}H_{20}O_{11}$ Molecular Weight:448.38

Target: Cholinesterase (ChE)

Pathway: Neuronal Signaling

Storage: 4°C, protect from light

\* In solvent: -80°C, 6 months; -20°C, 1 month (protect from light)

### **SOLVENT & SOLUBILITY**

In Vitro

DMSO: 16.67 mg/mL (37.18 mM; Need ultrasonic)

Preparing Stock Solutions	Solvent Mass Concentration	1 mg	5 mg	10 mg
	1 mM	2.2303 mL	11.1513 mL	22.3025 mL
	5 mM	0.4461 mL	2.2303 mL	4.4605 mL
	10 mM	0.2230 mL	1.1151 mL	2.2303 mL

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

In Vivo

1. Add each solvent one by one: 10% DMSO >> 90% (20% SBE-β-CD in saline) Solubility: ≥ 1.67 mg/mL (3.72 mM); Clear solution

#### **BIOLOGICAL ACTIVITY**

**Description** Rhodionin, isolated from the root of Rhodiola crenulata, is a specific non-competitive cytochrome P450 2D6 inhibitor with

an IC  $_{50}$  of 0.761  $\mu$ M and a Ki of 0.769  $\mu$ M  $^{[1]}$ . Rhodionin exhibits potent, dose-dependent inhibitory effects on acetylcholinesterase (AChE) with IC  $_{50}$  ranged from 57.50 to 2.43  $\mu$ g/mL  $^{[2]}$ . Rhodionin exhibits potent DPPH free radical

scavenging activities, with an IC $_{50}$  of 19.49  $\mu M^{[3]}$ .

IC<sub>50</sub> & Target AChE

#### **REFERENCES**

[1]. Xu W, et al. Two potent cytochrome P450 2D6 inhibitors found in Rhodiola rosea. Pharmazie. 2013 Dec;68(12):974-6.

[2]. Li FJ, et al. Molecular interaction studies of acetylcholinesterase with potential acetylcholinesterase inhibitors from the root of Rhodiola crenulata using molecular docking and isothermal titration calorimetry methods. Int J Biol Macromol. 2017 Nov;104(Pt A):527-532.

[3]. Choe KI, et al. The antioxid 27;17(10):11484-94.	dant and anti-inflammatory effects of phenolic compounds isolated from the root of Rhodiola sachalinensis A. BOR. Molecules. 20	012 Sep
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