Proteins



CMPDA

Cat. No.: HY-12508 CAS No.: 380607-77-2 Molecular Formula: $C_{16}H_{28}N_2O_4S_2$

Molecular Weight: 376.53 iGluR Target:

Pathway: Membrane Transporter/Ion Channel; Neuronal Signaling

-20°C Storage: Powder 3 years

In solvent

4°C 2 years -80°C 2 years

-20°C 1 year

Product Data Sheet

SOLVENT & SOLUBILITY

In Vitro

DMSO: ≥ 100 mg/mL (265.58 mM)

* "≥" means soluble, but saturation unknown.

Preparing Stock Solutions	Solvent Mass Concentration	1 mg	5 mg	10 mg
	1 mM	2.6558 mL	13.2792 mL	26.5583 mL
	5 mM	0.5312 mL	2.6558 mL	5.3117 mL
	10 mM	0.2656 mL	1.3279 mL	2.6558 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 (6.64 mM); Clear solution
- 2. Add each solvent one by one: 10% DMSO >> 90% (20% SBE-β-CD in saline) Solubility: ≥ 2.5 mg/mL (6.64 mM); Clear solution
- 3. Add each solvent one by one: 10% DMSO >> 90% corn oil Solubility: ≥ 2.5 mg/mL (6.64 mM); Clear solution

BIOLOGICAL ACTIVITY

Description

CMPDA is a positive allosteric modulator of AMPA receptors with EC50s of 45.4 ± 4.2 nM/63.4 ± 5.6 nM for GluA2i/GluA2o receptor.IC50 value: 45.4 ± 4.2 nM/63.4 ± 5.6 nM(GluA2i/GluA2o) [1]Target: AMPAR modulatorCMPDA was nearly equipotent at modulating the two isoforms of GluA2 receptors, whereas CMPDB displayed a modest preference for the flip splice variant. Similar to CX614, CMPDA slowed the rate of deactivation of GluA20 receptors approximately 2-fold but had no effect on GluA2i receptor deactivation [1].

REFERENCES 1]. Timm DE, et al. Structural and functional analysis of two new positive allosteric modulators of GluA2 desensitization and deactivation. Mol Pharmacol. 2011 Aug;80(2):267-80.				
	Caution: Product has not been fully validated for medical applications. For research use only.			
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