Screening Libraries

Product Data Sheet



Oxatomide

Cat. No.: HY-123205 CAS No.: 60607-34-3 Molecular Formula: $C_{27}H_{30}N_4O$ Molecular Weight: 426.55

Target: Histamine Receptor; P2X Receptor; 5-HT Receptor

Pathway: GPCR/G Protein; Immunology/Inflammation; Neuronal Signaling; Membrane

Transporter/Ion Channel

Storage: Powder -20°C 3 years

> 4°C 2 years

-80°C 2 years In solvent

1 year -20°C

SOLVENT & SOLUBILITY

In Vitro

DMSO: 250 mg/mL (586.10 mM; Need ultrasonic)

	Solvent Mass Concentration	1 mg	5 mg	10 mg
Preparing Stock Solutions	1 mM	2.3444 mL	11.7220 mL	23.4439 mL
	5 mM	0.4689 mL	2.3444 mL	4.6888 mL
	10 mM	0.2344 mL	1.1722 mL	2.3444 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: ≥ 2.08 mg/mL (4.88 mM); Clear solution
- 2. Add each solvent one by one: 10% DMSO >> 90% (20% SBE-β-CD in saline) Solubility: ≥ 2.08 mg/mL (4.88 mM); Clear solution
- 3. Add each solvent one by one: 10% DMSO >> 90% corn oil Solubility: ≥ 2.08 mg/mL (4.88 mM); Clear solution

BIOLOGICAL ACTIVITY

Description

Oxatomide is a potent and orally active dual H1-histamine receptor and P2X7 receptor antagonist with antihistamine and anti-allergic activity. Oxatomide almost completely blocks the ATP-induced current in human P2X7 receptors (IC $_{50}$ of 0.95 μ M). Oxatomide inhibits ATP-induced Ca^{2+} influx with an IC_{50} value of 0.43 μ M and also inhibits serotonin^{[1][2]}.

IC₅₀ & Target P2X7 H₁ Receptor serotonin

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〈Ohmori, et al. Pharmac				
	cological studies on oxatomide	e (KW-4354). (7) Antagonistic effe	ts on chemical mediators. Nihon Yakurigaku Zasshi. 1983	May;81(5):399-4
	Caution: Product has n	ot been fully validated for m	edical applications. For research use only.	
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REFERENCES

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