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Product Data Sheet

5-HT1A modulator 2 hydrochloride

Cat. No.:	HY-136621	
CAS No.:	3880-76-0	Ŷ
Molecular Formula:	C ₁₁ H ₁₆ CINO	\downarrow \land \backslash NH ₂
Molecular Weight:	213.7	
Target:	5-HT Receptor	
Pathway:	GPCR/G Protein; Neuronal Signaling	\checkmark \checkmark
Storage:	4°C, sealed storage, away from moisture * In solvent : -80°C, 6 months; -20°C, 1 month (sealed storage, away from moisture)	H-CI

SOLVENT & SOLUBILITY

In Vitro	DMSO : 100 mg/mL (467.95 mM; Need ultrasonic)					
	Preparing Stock Solutions	Solvent Mass Concentration	1 mg	5 mg	10 mg	
		1 mM	4.6795 mL	23.3973 mL	46.7946 mL	
		5 mM	0.9359 mL	4.6795 mL	9.3589 mL	
		10 mM	0.4679 mL	2.3397 mL	4.6795 mL	
	Please refer to the solubility information to select the appropriate solvent.					
In Vivo	1. Add each solvent Solubility: ≥ 2.5 m	one by one: 10% DMSO >> 40% PE g/mL (11.70 mM); Clear solution	G300 >> 5% Tween-80) >> 45% saline		
	2. Add each solvent one by one: 10% DMSO >> 90% (20% SBE-β-CD in saline) Solubility: ≥ 2.5 mg/mL (11.70 mM); Clear solution					

Diological Activity				
Description	5-HT1A modulator 2 hydrochloride, a derivative of 8-OH-DPAT (HY-112061), is a modulator of 5-HT _{1A} with a K _i of 53 nM for 5- HT _{1A} binding ^[1] .			
IC ₅₀ & Target	Ki: 53 nM (5-HT _{1A}) ^[1]			
In Vitro	5-HT1A modulator 2 hydrochloride (compound 3) binds to 5-HT _{1A} with a K _i of 53 nM ^[1] . MCE has not independently confirmed the accuracy of these methods. They are for reference only.			

REFERENCES

[1]. Naiman, N., et al. 2-(Alkylamino) tetralin derivatives: interaction with 5-HT1A serotonin binding sites. Journal of Medicinal Chemistry, 1989;32(1), 253–256.

Caution: Product has not been fully validated for medical applications. For research use only.

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