NDM-1 inhibitor-3

Cat. No.:	HY-150758				
Molecular Formula:	C ₁₆ H ₁₂ O ₄				
Molecular Weight:	268.26				
Target:	Bacterial; Beta-lactamase				
Pathway:	Anti-infection				
Storage:	Powder	-20°C	3 years		
		4°C	2 years		
	In solvent	-80°C	6 months		
		-20°C	1 month		

SOLVENT & SOLUBILITY

In Vitro	DMSO : 100 mg/mL (372.77 mM; Need ultrasonic)						
Prep Stoc		Solvent Mass Concentration	1 mg	5 mg	10 mg		
	Preparing Stock Solutions	1 mM	3.7277 mL	18.6386 mL	37.2773 mL		
		5 mM	0.7455 mL	3.7277 mL	7.4555 mL		
		10 mM	0.3728 mL	1.8639 mL	3.7277 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 (9.32 mM); Clear solution						
	2. Add each solvent one by one: 10% DMSO >> 90% (20% SBE-B-CD in saline) Solubility: \geq 2.5 mg/mL (9.32 mM); Clear solution						

Description	NDM-1 inhibitor-3 (Compound 89) is a New Delhi Metallo- β -lactamase-1 (NDM-1) inhibitor with a K _i of 4 μ M ^[1] .				
IC ₅₀ & Target	NDM-1 ^[1]				
In Vitro	New Delhi Metallo-β-lactamase-1 (NDM-1), a metallo-β-lactamase (MBL) first discovered at the end of 2008, is now considered as the most clinically relevant target for antibiotic resistance due to its worldwide prevalence ^[1] . MCE has not independently confirmed the accuracy of these methods. They are for reference only.				

REFERENCES

Product Data Sheet

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[1]. Caburet J, et al. A fragment-based drug discovery strategy applied to the identification of NDM-1 β-lactamase inhibitors. Eur J Med Chem. 2022 Jul 11;240:114599.

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

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