C-7280948

Cat. No.:	HY-15890		
CAS No.:	587850-67-7		
Molecular Formula:	C ₁₄ H ₁₆ N ₂ O ₂ S		
Molecular Weight:	276.35		
Target:	Histone Methyltransferase		
Pathway:	Epigenetics		
Storage:	Powder	-20°C	3 years
		4°C	2 years
	In solvent	-80°C	2 years
		-20°C	1 year

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SOLVENT & SOLUBILITY

In Vitro	DMSO : ≥ 38 mg/mL (137.51 mM) * "≥" means soluble, but saturation unknown.					
	Preparing Stock Solutions	Solvent Mass Concentration	1 mg	5 mg	10 mg	
		1 mM	3.6186 mL	18.0930 mL	36.1860 mL	
		5 mM	0.7237 mL	3.6186 mL	7.2372 mL	
		10 mM	0.3619 mL	1.8093 mL	3.6186 mL	
	Please refer to the so	lubility information to select the app	propriate solvent.			
In Vivo	 Add each solvent one by one: 10% DMSO >> 40% PEG300 >> 5% Tween-80 >> 45% saline Solubility: ≥ 2.5 mg/mL (9.05 mM); Clear solution Add each solvent one by one: 10% DMSO >> 90% corn oil Solubility: ≥ 2.5 mg/mL (9.05 mM); Clear solution 					
	50tability. ± 2.5 m	5/ m2 (3.03 mm), elear 3010101				

BIOLOGICALACTIVITY				
Description	C-7280948 is a selective and potent protein methyltransferase1 (PRMT1) inhibitor with an IC $_{50}$ value of 12.75 μ M ^[1] .			
IC ₅₀ & Target	IC50: 12.75 μM (PRMT1) ^[1]			
In Vitro	C-7280948 is compound 6a in this article, it against hPRMT1 with an IC ₅₀ value of 12.75 μM. Especially arginine methyltransferases also target non-histone protein substrates and are therefore often called protein methyltransferases (PRMTs). The subtype PRMT1 has been linked to the activation of estrogen and androgen receptors and therefore may represent a new treatment option for hormone-dependent cancer.			

Product Data Sheet

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H₂N.

MCE has not independently confirmed the accuracy of these methods. They are for reference only.

REFERENCES

[1]. Bissinger EM, et al. Acyl derivatives of p-aminosulfonamides and dapsone as new inhibitors of the arginine methyltransferase hPRMT1. Bioorg Med Chem. 2011 Jun 15;19(12):3717-3731.

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

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