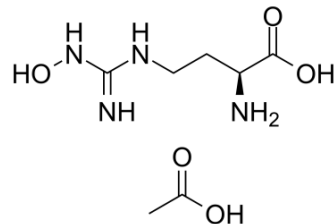


nor-NOHA monoacetate

Cat. No.:	HY-112885B		
CAS No.:	2250019-93-1		
Molecular Formula:	C ₇ H ₁₆ N ₄ O ₅		
Molecular Weight:	236.23		
Target:	Arginase		
Pathway:	Immunology/Inflammation; Metabolic Enzyme/Protease		
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 (423.32 mM; Need ultrasonic)
 H₂O : 50 mg/mL (211.66 mM; Need ultrasonic)

Preparing Stock Solutions	Solvent		Mass		
	Concentration		1 mg	5 mg	10 mg
	1 mM		4.2332 mL	21.1658 mL	42.3316 mL
	5 mM		0.8466 mL	4.2332 mL	8.4663 mL
	10 mM		0.4233 mL	2.1166 mL	4.2332 mL

Please refer to the solubility information to select the appropriate solvent.

In Vivo

- Add each solvent one by one: 10% DMSO >> 40% PEG300 >> 5% Tween-80 >> 45% saline
 Solubility: ≥ 2.5 mg/mL (10.58 mM); Clear solution
- Add each solvent one by one: 10% DMSO >> 90% (20% SBE-β-CD in saline)
 Solubility: ≥ 2.5 mg/mL (10.58 mM); Clear solution
- Add each solvent one by one: 10% DMSO >> 90% corn oil
 Solubility: ≥ 2.5 mg/mL (10.58 mM); Clear solution

BIOLOGICAL ACTIVITY

Description

nor-NOHA (Nω-Hydroxy-nor-L-arginine) monoacetate is a potent and selective arginase inhibitor. nor-NOHA monoacetate inhibits rat liver arginase with a K_i of 0.5 μM^[1].

REFERENCES

[1]. Tenu JP, et al. Effects of the new arginase inhibitor N(omega)-hydroxy-nor-L-arginine on NO synthase activity in murine macrophages. Nitric Oxide. 1999;3(6):427-438.

[2]. Custot J, et al. A new α -amino acid N w -hydroxy-nor-L-arginine: A highly-affinity inhibitor of arginase well adapted to bind to its manganese cluster. Journal of the American Chemical Society 119, 4086-4087(1997).

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

Tel: 609-228-6898

Fax: 609-228-5909

E-mail: tech@MedChemExpress.com

Address: 1 Deer Park Dr, Suite Q, Monmouth Junction, NJ 08852, USA