Product Data Sheet

Dacisteine

Cat. No.: HY-121765 CAS No.: 18725-37-6 Molecular Formula: $C_7H_{11}NO_4S$ Molecular Weight: 205.23

Target: **Endogenous Metabolite** Pathway: Metabolic Enzyme/Protease

Storage: Powder -20°C

> -80°C In solvent 6 months -20°C

3 years 4°C 2 years

SOLVENT & SOLUBILITY

DMSO: ≥ 250 mg/mL (1218.15 mM) In Vitro

* "≥" means soluble, but saturation unknown.

1 month

Preparing Stock Solutions	Solvent Mass Concentration	1 mg	5 mg	10 mg
	1 mM	4.8726 mL	24.3629 mL	48.7258 mL
	5 mM	0.9745 mL	4.8726 mL	9.7452 mL
	10 mM	0.4873 mL	2.4363 mL	4.8726 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 (10.13 mM); Clear solution
- 2. Add each solvent one by one: 10% DMSO >> 90% (20% SBE-β-CD in saline) Solubility: ≥ 2.08 mg/mL (10.13 mM); Clear solution
- 3. Add each solvent one by one: 10% DMSO >> 90% corn oil Solubility: ≥ 2.08 mg/mL (10.13 mM); Clear solution

BIOLOGICAL ACTIVITY

Description Dacisteine (N,S-Diacetyl-L-cysteine) is a cysteine derivative and displays a less New Delhi metallo-beta-lactamase-1 (NDM-1) inhibitor with an IC $_{50}$ value of 1000 $\mu M^{[1]}$. Dacisteine can be used for the treatment of cardiovascular and cerebrovascular diseases caused by platelet aggregation^[2].

IC₅₀ & Target IC50: 1000 μM (NDM-1)^[1]

In Vivo

Dacisteine is a preparing agent for treating or preventing cardiovascular and cerebrovascular diseases caused by platelet aggregation.

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

REFERENCES

[1]. Cui-Gai Bail, et al. Cysteine and Its Derivatives as New Delhi Metallo-beta-lactamase-1 Inhibitors. Current Enzyme Inhibition, 2015, 11, 46-57

[2]. FanKe, et al. Application of compound of thioether acid structure in preparing anti-platelet aggregation drugs. Patent WO2019007015A1.

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

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