Inhibitors

Screening Libraries

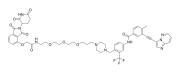
Proteins

SB1-G-187

Cat. No.: HY-137342 CAS No.: 2769753-18-4 Molecular Formula: $C_{52}H_{54}F_{3}N_{9}O_{10}$ Molecular Weight: 1022.03 **PROTACs** Target: Pathway: **PROTAC**

Storage: -20°C, stored under nitrogen

* In solvent: -80°C, 6 months; -20°C, 1 month (stored under nitrogen)



Product Data Sheet

SOLVENT & SOLUBILITY

In Vitro

DMSO: 100 mg/mL (97.84 mM; Need ultrasonic)

Preparing Stock Solutions	Solvent Mass Concentration	1 mg	5 mg	10 mg
	1 mM	0.9784 mL	4.8922 mL	9.7844 mL
	5 mM	0.1957 mL	0.9784 mL	1.9569 mL
	10 mM	0.0978 mL	0.4892 mL	0.9784 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: 5 mg/mL (4.89 mM); Suspended solution; Need ultrasonic
- 2. Add each solvent one by one: 10% DMSO >> 90% corn oil Solubility: ≥ 5 mg/mL (4.89 mM); Clear solution

BIOLOGICAL ACTIVITY

Description

SB1-G-187, a PROTAC, is a multi-kinase degrader^[1]. SB1-G-187 is a click chemistry reagent, it contains an Alkyne group and can undergo copper-catalyzed azide-alkyne cycloaddition (CuAAc) with molecules containing Azide groups.

REFERENCES

[1]. Katherine A Donovan, et al. Mapping the Degradable Kinome Provides a Resource for Expedited Degrader Development. Cell. 2020 Dec 10;183(6):1714-1731.e10.

 $\label{lem:caution:Product} \textbf{Caution: Product has not been fully validated for medical applications. For research use only.}$

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