

N-Boc-C1-PEG5-C3-NH2 Biochemical

N-Boc-C1-PEG5-C3-NH2

CATALOG #
AAA228958

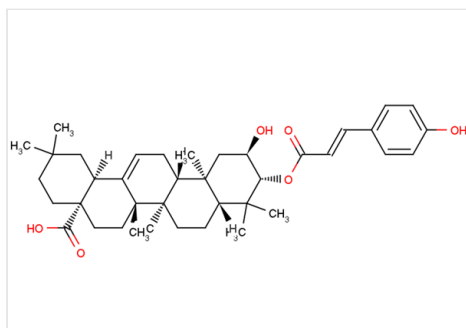
PRODUCT SYNONYMS

N-Boc-C1-PEG5-C3-NH2; N-Boc-C1-PEG5-C3-NH2 biochemical

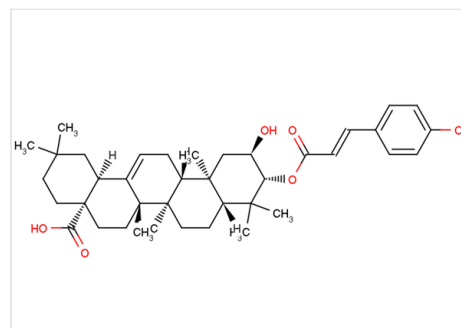
PRODUCT SPECIFICATIONS

Purity/Purification	98%
CAS	956472-77-8
Solubility	<1mg/ml refers to the product slightly soluble or insoluble
In vitro	PROTACs contain two different ligands connected by a linker; one is a ligand for an E3 ubiquitin ligase and the other is for the target protein. PROTACs exploit the intracellular ubiquitin-proteasome system to selectively degrade target proteins[1].
Preparation and Storage	Powder: -20 degree C for 3 years In solvent: -80 degree C for 2 years

IMAGES



Structure
3-O-trans-p-Coumaroylmaslinic acid, CAS 35482-91-8



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ADDITIONAL INFORMATION

Related Product Information for N-Boc-C1-PEG5-C3-NH2 biochemical

N-Boc-C1-PEG5-C3-NH2 is a PEG-based PROTAC linker that can be used in the synthesis of PROTACs.

NCBI AND UNIPROT INFORMATION

Molecular Weight	618.84
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For Research Use Only. Not for use in diagnostic or therapeutic procedures. Small volumes of product may become entrapped in the seal of the vial during shipment and storage; if necessary, briefly centrifuge the vial on a tabletop centrifuge to dislodge any liquid in the cap.