



## Product Information

### E3 ligase ligand-linker conjugate inhibitor, (S,R,S)-AHPC-PEG4-NH<sub>2</sub>, Purity ≥98%

**Cat. No.:** X24-09-YM716

**Size:** 100 mg; 200 mg; 500 mg; 1 g

**CAS Number:** 2010159-57-4

**Compound CID:** 124177628

**Synonym:** 2010159-57-4; VH032-PEG4-NH<sub>2</sub>; VHL ligand-linker conjugate 4; E3

ligase ligand-linker conjugate 7 free base;

(2S,4R)-1-[(2S

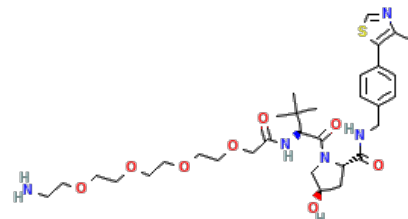
)-2-[[2-[2-[2-[2-(2-Aminoethoxy)ethoxy]ethoxy]ethoxy]acetyl]amino]-3,3-dimethylbutan

oyl]-4-hydroxy-N

-[[4-(4-methyl-1,3-thiazol-5-yl)phenyl]methyl]pyrrolidine-2-carboxamide; E3 ligase

ligand-linker conjugate inhibitor

**This product is for research use only and is not intended for diagnostic use.**



#### Product Information

|                          |                                                                                                                                                                                                                                                                                                       |
|--------------------------|-------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|
| <b>Description</b>       | (S,R,S)-AHPC-PEG4-NH <sub>2</sub> is an effective PROTAC inhibitor, that has the ability to prevent the pathway of E3 ligase ligand-linker conjugation. The molecular weight of the compound is 663.83, and its molecular formula is C <sub>32</sub> H <sub>49</sub> N <sub>5</sub> O <sub>8</sub> S. |
| <b>Molecular Weight</b>  | 663.83                                                                                                                                                                                                                                                                                                |
| <b>Molecular Formula</b> | C <sub>32</sub> H <sub>49</sub> N <sub>5</sub> O <sub>8</sub> S                                                                                                                                                                                                                                       |
| <b>IUPAC Name</b>        | (2S,4R)-1-[(2S<br><br>N-[[4-(4-methyl-1,3-thiazol-5-yl)phenyl]methyl]pyrrolidine-2-carboxamide                                                                                                                                                                                                        |
| <b>InChI</b>             | InChI=1S/C32H49N5O8S/c1-22-28(46-21-35-22)24-7-5-23(6-8-24)18-34-30(40)26-17-25(38)19-37(26)31(41)29(32(2,3)4)36-27(39)20-45-16-15-44-14-13-43-12-11-42-10-9-33/h5-8,21,25-26,29,38H,9-20,33H2,1-4H3,(H,34,40)(H,36,39)/t25-,26+,29-/m1/s1                                                            |
| <b>InChI Key</b>         | VSMWDQCRYIOFMX-UWPQUIUOSA-N                                                                                                                                                                                                                                                                           |
| <b>Canonical SMILES</b>  | CC1=C(SC=N1)C2=CC=C(C=C2)CNC(=O)C3CC(CN3C(=O)C(C(C)(C)C)NC(=O)COCCOCCOCCOCCN)O                                                                                                                                                                                                                        |
| <b>Form</b>              | Lyophilized powder                                                                                                                                                                                                                                                                                    |



|                      |                                                                                                                 |
|----------------------|-----------------------------------------------------------------------------------------------------------------|
| <b>Purity</b>        | ≥98%                                                                                                            |
| <b>Impurities</b>    | Free from inappropriate visible particulates, foreign matter, discoloration, or other defects.                  |
| <b>Identity</b>      | Confirmed by NMR/HPLC/MS.                                                                                       |
| <b>Stability</b>     | In its lyophilized form, the chemical remains stable for 36 months.                                             |
| <b>Quality Level</b> | Research grade                                                                                                  |
| <b>Applications</b>  | (S,R,S)-AHPC-PEG4-NH <sub>2</sub> plays a key role in enhancing the pharmacological properties of linked drugs. |
| <b>Storage</b>       | Store at -20°C, and keep desiccated.                                                                            |