

Product Information

Amine/NHS PEG reagent, *t*-Boc-*N*-amido-PEG2-NHS ester, Purity 98%

Cat. No.: X24-09-YYX183

Size: 100 mg; 500 mg; 1 g

CAS Number: 2183440-73-3

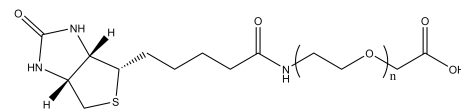
PubChem CID: 23591747

Synonym: 2183440-73-3; *t*-Boc-*N*-amido-PEG2-NHS ester; Boc-NH-PEG2-C₂-NHS

ester; 2,5-dioxopyrrolidin-1-yl

2,2-dimethyl-4-oxo-3,8,11-trioxa-5-azatetradecan-14-oate;

BocNH-PEG2-CH₂CH₂COONHS



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

Product Information

Description	<i>t</i> -Boc- <i>N</i> -amido-PEG2-NHS ester functions as another PEG linker featuring both an NHS ester and a <i>t</i> -Boc protected amino group. Its hydrophilic PEG spacer increases solubility when dissolved in aqueous media, and like before, the <i>t</i> -Boc group can be removed under mild acidic conditions to generate free amines suitable for various applications involving primary amines <i>t</i> -from proteins or modified oligonucleotides.
Molecular Weight	374.4
Molecular Formula	C ₁₆ H ₂₆ N ₂ O ₈
Functional Group 1	Boc
Functional Group 2	Amine
Functional Group 3	NHS
Reactive Group 1	Acid
Reactive Group 2	Amine
IUPAC Name	(2,5-Dioxopyrrolidin-1-yl) 3-[2-[2-[(2-methylpropan-2-yl)oxycarbonylamino]ethoxy]ethoxy]propanoate
InChI	InChI=1S/C16H26N2O8/c1-16(2,3)25-15(22)17-7-9-24-11-10-23-8-6-14(21)26-18-12(19)4-5-13(18)20/h4-11H2,1-3H3,(H,17,22)
InChI Key	FNZXIPYAXFROBS-UHFFFAOYSA-N
Canonical SMILES	CC(C)(C)OC(=O)NCCOCCOCCC(=O)ON1C(=O)CCC1=O
Form	Solid or viscous liquid
Purity	98%

Solubility	DCM
Identity	Confirmed by NMR.
Applications	This compound serves similar functions in peptide synthesis and bioconjugation as the PEG3 variant but with an optimized spacer length that may influence the flexibility and solubility of the final conjugate, enhancing its pharmacokinetic properties.
Storage	Store at -20°C.