



Product Information

Mono-sulfo-galactosylβCer (d18:1/12:0)

Cat. No.: X23-11-ZQ604

Size: 10 mg; 100 mg; 500 mg; 1 g; 10 g

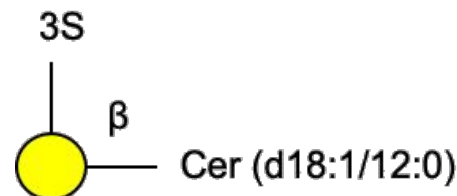
CAS Number: 852043-39-1

PubChem CID: 145710284

Synonym: 852043-39-1;

Dodecanamide-*N*-[(1*S*,2*R*,3*E*)-2-hydroxy-1-[(3-*O*-sulfo-beta-D-galactopyranosyl)oxy]methyl]-3-heptadecen-1-yl]-ammonium salt

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



Product Information

Molecular Weight	727.01
Molecular Formula	C ₃₅ H ₇₀ N ₂ O ₁₁ S
IUPAC Name	Azanium;[(2 <i>R</i> ,3 <i>R</i> ,4 <i>S</i> ,5 <i>S</i> ,6 <i>R</i>)-2-[(<i>E</i> ,2 <i>S</i> ,3 <i>R</i>)-2-(dodecanoylamino)-3-hydroxyoctadec-4-enoxy]-3,5-dihydroxy-6-(hydroxymethyl)oxan-4-yl] sulfate
InChI	InChI=1S/C36H69NO11S.H3N/c1-3-5-7-9-11-13-14-15-16-18-19-21-23-25-30(39)29(37-32(40)26-24-22-20-17-12-10-8-6-4-2)28-46-36-34(42)35(48-49(43,44)45)33(41)31(27-38)47-36;/h23,25,29-31,33-36,38-39,41-42H,3-22,24,26-28H2,1-2H3,(H,37,40)(H,43,44,45);1H3/b25-23+;/t29-,30+,31+,33-,34+,35-,36+;/m0./s1
InChI Key	ZRHCDEVQWSMWDC-CIEWKKQXSA-N
Canonical SMILES	CCCCCCCCCCCCC=CC(C(COC1C(C(C(C(O1)CO)O)OS(=O)(=O)[O-])O)NC(=O)CCCCCCCCCCC)O.[NH4+]
Isomeric SMILES	CCCCCCCCCCCCC/C=C/[C@H]([C@H](CO[C@H]1[C@@H]([C@H]([C@H]([C@H](O1)CO)O)OS(=O)(=O)[O-])O)NC(=O)CCCCCCCCCCC)O.[NH4+]
Form	Powder
Applications	Mono-sulfo-galactosylβCer (d18:1/12:0) is a non-physical sulfatidein, which can be used as a valid standard in the mass spectrometric analysis of sulfatides.
Storage	Store at -20°C.