

## Product Information

### 2-O-hexadecanoyl-3-O-(2,4S,6S-trimethyl-2E-pentacosenoyl)-6-O-(2,4S,6S-trimethyl-2E-pentacosenoyl)-2'-O-(2R,4S,6S-trimethyl-3R-hydroxy-tetracosanoyl)-4'-O

Cat. No.: X26-02-ZQ5065

Size: 1 mg; 10 mg; 25 mg; 100 mg

Compound CID: 52930180

Synonym:

PAT16(25:1(2E)(2Me,4Me[S],6Me[S])/25:1(2E)(2Me,4Me[S],6Me[S])/24:0(2Me[R],3OH[R],4Me[S],6Me[S])/24:1(2E)(2Me,4Me[S],6Me[S]))

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

#### Product Information

Description	This variant combines a hydroxylated branch with dual C25 and a C24 pthienoic acid to examine how functionalized sites modify bilayer organization. It is utilized in structural biology to characterize the lipid antigen-binding specificity of host immune proteins.
Molecular Weight	2189.5
Molecular Formula	$C_{138}H_{258}O_{17}$
IUPAC Name	[(3R,5R,6R)-5-hexadecanoyloxy-3-hydroxy-6-[(2R,3S,5S)-4-hydroxy-6-(hydroxymethyl)-3-[(2S,3S,4S,6S)-3-hydroxy-2,4,6-trimethyltetracosanoyl]oxy-5-[(E,4S,6S)-2,4,6-trimethyltetracos-2-enoyl]oxyoxan-2-yl]oxy-4-[(E,4S,6S)-2,4,6-trimethylpentacos-2-enoyl]oxyoxan-2-yl)methyl (E,4S,6S)-2,4,6-trimethylpentacos-2-enoate
InChI	InChI=1S/C138H258O17/c1-18-23-28-33-38-43-48-53-57-61-65-69-73-77-82-87-92-97-111(6)102-115(10)106-119(14)133(144)148-110-124-127(142)130(153-135(146)121(16)108-117(12)104-113(8)99-94-89-84-79-74-70-66-62-58-54-49-44-39-34-29-24-19-2)132(151-125(140)101-96-91-86-81-76-67-52-47-42-37-32-27-22-5)138(150-124)155-137-131(154-136(147)122(17)126(141)118(13)105-114(9)100-95-90-85-80-75-71-64-60-56-51-46-41-36-31-26-21-4)128(143)129(123(109-139)149-137)152-134(145)120(15)107-116(11)103-112(7)98-93-88-83-78-72-68-63-59-55-50-45-40-35-30-25-20-3/h106-108,111-118,122-124,126-132,137-139,141-143H,18-105,109-110H2,1-17H3/b119-106+,120-107+,121-108+/t111-,112-,113-,114-,115-,116-,117-,118-,122-,123?,124?,126-,127+,128?,129+,130?,131-,132+,137+,138+/m0/s1
InChI Key	LZPXFZCXHWQDE-OPBNFDCFSA-N



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<b>Canonical SMILES</b>	<chem>CCCCCCCCCCCCCCCCCCCC[C@H](C)C[C@H](C)/C=C(C)/C(=O)OCC1[C@H](C([C@H]([C@H](O1)O[C@@H]2[C@H](C([C@@H](C(O2)CO)OC(=O)/C=C/[C@@H](C)C[C@@H](C)CCCCCCCCCCCCCCCCC)/C)O)OC(=O)[C@@H](C)[C@H]([C@@H](C)C[C@@H](C)CCCCCCCCCCCCCCCC)O)OC(=O)CCCCCCCCCCCCCCCC)OC(=O)/C=C/[C@@H](C)C[C@@H](C)CCCCCCCCCCCCCCCC)/C)O</chem>
<b>Source</b>	Chemical synthesis
<b>Form</b>	Solid or liquid
<b>Purity</b>	≥90%
<b>Identity</b>	Confirmed by NMR/HPLC/MS.
<b>Stability</b>	The product is stable for one year when stored at the recommended temperature in lyophilized powder.
<b>Quality Level</b>	Research level
<b>Applications</b>	Used to examine how positional isomerism of hydroxyl groups affects the recognition of glycolipids by the immune system.
<b>Storage</b>	Store at -20°C.

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