

## Product Information

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

**Cat. No.:** X26-02-ZQ4864

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

**Compound CID:** 52929979

**Synonym:**

PAT16(22:1(2E)(2Me,4Me[S],6Me[S])/24:0(2Me[R],3OH[R],4Me[S],6Me[S])/24:0(2Me[R],3OH[R],4Me[S],6Me[S])/22:1(2E)(2Me,4Me[S],6Me[S]))

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

#### Product Information

<b>Description</b>	Featuring hydroxylated branches at both the 6 and 2' positions, this PAT allows for the study of regiospecific hydrogen bonding on the trehalose core. It provides vital information on the molecular recognition of specialized glycolipids.
<b>Molecular Weight</b>	2123.4
<b>Molecular Formula</b>	C <sub>132</sub> H <sub>248</sub> O <sub>18</sub>
<b>IUPAC Name</b>	[(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-trimethyldocos-2-enoyl]oxyoxan-2-yl]oxy-4-[(E,4S,6S)-2,4,6-trimethyldocos-2-enoyl]oxyoxan-2-yl]methyl (2S,3S,4S,6S)-3-hydroxy-2,4,6-trimethyltetracosanoate
<b>InChI</b>	InChI=1S/C132H248O18/c1-18-23-28-33-38-43-48-53-57-59-64-68-73-78-83-88-93-106(8)98-110(12)119(135)114(16)129(141)143-103-117-121(137)124(148-128(140)113(15)101-109(11)97-105(7)92-87-82-77-72-67-63-56-51-46-41-36-31-26-21-4)126(146-118(134)95-90-85-80-75-70-61-52-47-42-37-32-27-22-5)132(145-117)150-131-125(149-130(142)115(17)120(136)111(13)99-107(9)94-89-84-79-74-69-65-60-58-54-49-44-39-34-29-24-19-2)122(138)123(116(102-133)144-131)147-127(139)112(14)100-108(10)96-104(6)91-86-81-76-71-66-62-55-50-45-40-35-30-25-20-3/h100-101,104-111,114-117,119-126,131-133,135-138H,18-99,102-103H2,1-17H3/b112-100+,113-101+/t104-,105-,106-,107-,108-,109-,110-,111-,114-,115-,116?,117?,119-,120-,121+,122?,123+,124?,125-,126+,131+,132+/m0/s1
<b>InChI Key</b>	DFHKCSYKEMQNIJ-GRIGHBBUSA-N
<b>Canonical SMILES</b>	CCCCCCCCCCCCCCCC[C@H](C)C[C@H](C)[C@@H]([C@H](C)C(=O)OCC1[C@H](C([C@H]

