

## Product Information

### MGMG(16:2(7Z,10Z)/0:0)

**Cat. No.:** X26-02-ZQ254

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

**Compound CID:** 171120561

**Synonym:** 1-(7Z,10Z-hexadecadienoyl)-3-O-β-D-galactosyl-sn-glycerol

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

#### Product Information

<b>Description</b>	A monogalactosyl monoacylglycerol (C16:2); the polyunsaturated fatty acid at sn-1 contains two cis-double bonds, providing increased chain fluidity to the glycolipid.
<b>Molecular Weight</b>	488.6
<b>Molecular Formula</b>	C <sub>25</sub> H <sub>44</sub> O <sub>9</sub>
<b>IUPAC Name</b>	[(2S)-2-hydroxy-3-[(2R,3R,4S,5R,6R)-3,4,5-trihydroxy-6-(hydroxymethyl)oxan-2-yl]oxypropyl] (7Z,10Z)-hexadeca-7,10-dienoate
<b>InChI</b>	InChI=1S/C25H44O9/c1-2-3-4-5-6-7-8-9-10-11-12-13-14-15-21(28)32-17-19(27)18-33-25-24(31)23(30)22(29)20(16-26)34-25/h6-7,9-10,19-20,22-27,29-31H,2-5,8,11-18H2,1H3/b7-6-,10-9-/t19-,20-,22+,23+,24-,25-/m1/s1
<b>InChI Key</b>	JKTGDOGAVPBQDI-NQNUALNFSA-N
<b>Canonical SMILES</b>	CCCCC/C=C\C/C=C\CCCCC(=O)OC[C@H](CO[C@H]1[C@@H]([C@H]([C@H]([C@H](O1)CO)O)O)O)O
<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>	MGMG(16:2(7Z,10Z)/0:0) can be used to study the biophysical characteristics of polyunsaturated monogalactosylmonoacylglycerols in plant lipid research.
<b>Storage</b>	Store at -20°C.