2-O-GLYCEROL-ALPHA-D-GALACTOPYRANOSID
HEXATMS

Other names: 2-O-Glycerol-«beta»-D-galactopyranoside, MO-TMS

Inchi: InChI=1S/C27H66O8Si6/c1-36(2,3)28-19-22(20-29-37(4,5)6)31-27-26(35-41(16,17)18)25

InchiKey: ADPMNKPURASUCV-SHZNSVIDSA-N

Formula: C27H66O8Si6

SMILES: C[Si](C)(C)OCC(CO[Si](C)(C)C)O[C@@H]1O[C@@H](CO[Si](C)(C)C)[C@@H](O[Si](C)

Mol. weight [g/mol]: 687.33

Physical Properties

Property code	Value	Unit	Source
logp	7.311		Crippen Method

Sources

Cheméo Relay (1.0):

Cheméo Relay (1.0, beta):

NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?Str2File=U98090

Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l

Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws

Legend

logp: Octanol/Water partition coefficient