

3-«beta»,7-«alpha»,12-«alpha»-Trihydroxy-5-chole acid, methyl ester, TMS

InChI: InChI=1S/C34H64O5Si3/c1-23(14-17-31(35)36-4)26-15-16-27-32-28(22-30(34(26,27)3)33)/q1-3H-1/t1-3H/m1-3/s1-3
InChIKey: VIATWLYTLF7FJF-IIMSTRFDSA-N
Formula: C34H64O5Si3
SMILES: COC(=O)CCC(C)C1CCC2C3C(O[Si](C)(C)C)C=C4CC(O[Si](C)(C)C)CCC4(C)C3CC(O[Si](C)(C)C)C1
Mol. weight [g/mol]: 637.13

Physical Properties

Property code	Value	Unit	Source
log10ws	-2.41		Crippen Method
logp	9.035		Crippen Method
rinpol	3208.00		NIST Webbook
rinpol	3208.00		NIST Webbook

Sources

Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci9903071>
Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=R390062&Units=SI>

Legend

log10ws: Log10 of Water solubility in mol/l
logp: Octanol/Water partition coefficient
rinpol: Non-polar retention indices

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