3«alpha»,6«beta»,7«alpha»,12«alpha»-Tetrahydro
acid, MeTMS

Other names: 3«alpha»,6«beta»,7«alpha»,12«alpha»-Tetrahydroxy-5«beta»-cholanoic acid, methyl ester-trimethylsilyl ether
3«alpha»,6«beta»,7«alpha»,12«alpha»-Tetrahydroxy-5«beta»-cholanoic acid, trimethylsilyl ether-methyl ester

Inchi: InChI=1S/C37H74O6Si4/c1-25(17-20-32(38)39-4)27-18-19-28-33-29(24-31(37(27,28)3)4

InchiKey: CAHBNWKLPVHQBG-CLTWYAMUSA-N

Formula: C37H74O6Si4

SMILES: COC(=O)CCC(C)C1CCC2C3C(O[Si](C)(C)C)C(O[Si](C)(C)C)C4CC(O[Si](C)(C)C)CCC4(

Mol. weight [g/mol]: 727.32

Physical Properties

Property code	Value	Unit	Source
log10ws	-0.82		Crippen Method
logp	9.945		Crippen Method
rinpol	3254.00		NIST Webbook
rinpol	3282.00		NIST Webbook
rinpol	3254.00		NIST Webbook

Sources

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

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

NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=R182346&Units=SI

Legend

log10ws: Log10 of Water solubility in mol/l

logp: Octanol/Water partition coefficient

rinpol: Non-polar retention indices

Cheméo (https://www.chemeo.com) is the biggest free database of chemical and physical data for the process industry.