

3«beta»,26-dihydroxy-5-cholestene, TMS

Inchi: InChI=1S/C33H62O2Si2/c1-24(23-34-36(5,6)7)12-11-13-25(2)29-16-17-30-28-15-14-26-
InchiKey: XPOWNWJGSYYWLA-LHVSONKFSA-N
Formula: C33H62O2Si2
SMILES: CC(CCCC(C)C1CCC2C3CC=C4CC(O[Si](C)(C)C)CCC4(C)C3CCC12C)CO[Si](C)(C)C
Mol. weight [g/mol]: 547.02

Physical Properties

Property code	Value	Unit	Source
log10ws	-5.52		Crippen Method
logp	10.079		Crippen Method
rinpol	3465.00		NIST Webbook

Sources

Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=R493780&Units=SI>
Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci9903071>

Legend

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

Latest version available from:

<https://www.chemeo.com/cid/64-275-6/3-beta-26-dihydroxy-5-cholestene-TMS.pdf>

Generated by Cheméo on 2024-04-24 16:42:35.673247228 +0000 UTC m=+16266204.593824544.

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