

5B-Cholestane-3A,7A,26-triol, TMS

Inchi: InChI=1S/C36H72O3Si3/c1-26(25-37-40(5,6)7)15-14-16-27(2)30-17-18-31-34-32(20-22-
InchiKey: NSAXRTMXFXEPGS-FIDJAZAJSA-N
Formula: C36H72O3Si3
SMILES: CC(CCCC(C)C1CCC2C3C(O[Si](C)(C)C)CC4CC(O[Si](C)(C)C)CCC4(C)C3CCC12C)CC
Mol. weight [g/mol]: 637.21

Physical Properties

Property code	Value	Unit	Source
log10ws	-3.93		Crippen Method
logp	10.989		Crippen Method
rinpol	3491.00		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=R585205&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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<https://www.chemeo.com/cid/42-478-5/5B-Cholestane-3A-7A-26-triol-TMS.pdf>

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