

2«alpha»,3«alpha»-cyclopropane-5«alpha»-andro

bisTMS

InChI: InChI=1S/C26H48O2Si2/c1-25-12-11-21-19(20(25)9-10-24(25)28-30(6,7)8)15-23(27-29(28)24)26
InChIKey: XHXIUUDIZDQNZJE-CQRJRCTISA-N
Formula: C26H48O2Si2
SMILES: CC12CCC3C(CC(O[Si](C)(C)C)C4CC5CC5CC43C)C1CCC2O[Si](C)(C)C
Mol. weight [g/mol]: 448.83

Physical Properties

Property code	Value	Unit	Source
log10ws	-2.75		Crippen Method
logp	7.325		Crippen Method
rinpol	2654.00		NIST Webbook

Sources

Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci9903071>
Crippen Method: https://www.cheméo.com/doc/models/crippen_log10ws
NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=R385887&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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