3«alpha»,16«alpha»,17«beta»-Trihydroxy-5«alpha»-androstane-tris-TMS

Other names: 5A-Androstan-3A,16A,17B-triol, tris-TMS
5«alpha»-Androstane-3«alpha», 16«alpha»,17«beta»-triol, MO TMS

Inchi: InChI=1S/C28H56O3Si3/c1-27-16-14-21(29-32(3,4)5)18-20(27)12-13-22-23(27)15-17-28

InchiKey: RQCSPBXXGJRUKU-GOWWAKEWSA-N

Formula: C28H56O3Si3

SMILES: CC12CCC(O[Si](C)(C)C)CC1CCC1C2CCC2(C)C1CC(O[Si](C)(C)C)C2O[Si](C)(C)C

Mol. weight [g/mol]: 525.00

Physical Properties

Property code	Value	Unit	Source
log10ws	-1.42		Crippen Method
logp	8.299		Crippen Method
rinpol	2774.00		NIST Webbook
rinpol	2778.00		NIST Webbook
rinpol	2785.00		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=R16616&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.