

3«alpha»,16«beta»,17«alpha»-Trihydroxy-5«alpha»-androstane-tris-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: RQCSPBXXGJRUKE-KBHMRCOTSA-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	2678.00		NIST Webbook
rinpol	2678.00		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=R16634&Units=SI>

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/41-338-1/3-alpha-16-beta-17-alpha-Trihydroxy-5-alpha-androstane-tris-TMS.pdf>

Generated by Cheméo on 2025-12-05 13:40:54.71036035 +0000 UTC m=+4690252.240401006.

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