

5«alpha»-Androstane-2«alpha»,3«alpha»,17«beta»

TMS

InChI=1S/C28H56O3Si3/c1-27-17-16-23-21(22(27)14-15-26(27)31-34(9,10)11)13-12-20

InchiKey:

AHPKNGPAYNSDOO-MCPUQREKSA-N

Formula:

C28H56O3Si3

SMILES:

CC12CC(O[Si](C)(C)C)C(O[Si](C)(C)C)CC1CCC1C2CCC2(C)C(O[Si](C)(C)C)CCC12

Mol. weight [g/mol]:

525.00

Physical Properties

Property code	Value	Unit	Source
log10ws	-1.42		Crippen Method
logp	8.299		Crippen Method
rinpol	2755.00		NIST Webbook

Sources

Crippen Method:

<http://pubs.acs.org/doi/abs/10.1021/ci990307l>

Crippen Method:

https://www.cheméo.com/doc/models/crippen_log10ws

NIST Webbook:

<http://webbook.nist.gov/cgi/cbook.cgi?ID=R307030&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.cheméo.com/cid/54-928-2/5-alpha-Androstane-2-alpha-3-alpha-17-beta-triol-TMS.pdf>

Generated by Cheméo on 2024-04-25 08:48:59.753817743 +0000 UTC m=+16324188.674395054.

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