

# 5«alpha»-Androstane-3«alpha»,17«beta»-diol, bis(trimethylsilyl) ether

<b>Other names:</b>	Silane, [[(3«alpha»,5«alpha»,17«beta»)-androstane-3,17-diol]bis(oxy)]bis[trimethyl-5«alpha»-Androstane-3«alpha»,17«beta»-diol, di-TMS 3,17-Bis[(trimethylsilyl)oxy]androstane, (3«alpha»,5«alpha»,17«beta»)-5«alpha»-Androstanediol-3«alpha»,17«beta», bis-TMS Androstane, 3,17-bis[(trimethylsilyl)oxy]-, (3«alpha»,5«alpha»,17«beta»)-5-«beta»-Androstan-3-«alpha»,17-«alpha»-diol, TMS 5«beta»-Androstanediol-3«alpha»,17«alpha», bis-TMS 3A,17B-Dihydroxy-5A-androstane, bis-TMS 5B-Androstane-3A,17A-diol, TMS 3A,17A-Dihydroxy-5B-androstane, TMS 5A-Androstan-3A,17B-diol, bis-TMS Androstane-3,17-diol, (3«alpha»,5«alpha»,17«beta»)-, 2tms derivative
<b>Inchi:</b>	InChI=1S/C25H48O2Si2/c1-24-15-13-19(26-28(3,4)5)17-18(24)9-10-20-21-11-12-23(27-
<b>InchiKey:</b>	KBSHKNYEUGMMDQ-LVVCFMQNSA-N
<b>Formula:</b>	C25H48O2Si2
<b>SMILES:</b>	CC12CCC(O[Si](C)(C)C)CC1CCC1C2CCC2(C)C(O[Si](C)(C)C)CCC12
<b>Mol. weight [g/mol]:</b>	436.82
<b>CAS:</b>	10426-36-5

## Physical Properties

Property code	Value	Unit	Source
log10ws	-2.92		Crippen Method
logp	7.469		Crippen Method
rinpol	2555.00		NIST Webbook
rinpol	2576.00		NIST Webbook
rinpol	2546.00		NIST Webbook
rinpol	2576.00		NIST Webbook
rinpol	2555.00		NIST Webbook
rinpol	2570.00		NIST Webbook
rinpol	2580.00		NIST Webbook

## Sources

Crippen Method:

<http://pubs.acs.org/doi/abs/10.1021/ci9903071>

**Crippen Method:**

[https://www.chemeo.com/doc/models/crippen\\_log10ws](https://www.chemeo.com/doc/models/crippen_log10ws)

**NIST Webbook:**

<http://webbook.nist.gov/cgi/cbook.cgi?ID=C10426365&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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