

# Silane,

# [(3«alpha»,5«alpha»,17«beta»)-androstane-3,17-diyloxy]bis(1,1-dimethylethyl)dimethylsilane

Other names:

Silane,  
[(3«alpha»,5«alpha»,17«beta»)-androstane-3,17-diyloxy]bis(1,1-dimethylethyl)dimethylsilane  
Androstane; silane deriv.

3-«alpha»,17-«beta»-Bis(tert-butyldimethylsilyloxy)-5-«alpha»-androstane

5«alpha»-Androstane-3«alpha»,17«beta»-diol, bis(tert-butyldimethylsilyl) ether

InChI:

InChI=1S/C31H60O2Si2/c1-28(2,3)34(9,10)32-23-17-19-30(7)22(21-23)13-14-24-25-15-

InChIKey:

VXQXLCZNGSUBJR-QHDZDOOTSA-N

Formula:

C31H60O2Si2

SMILES:

CC1CCC(O[Si](C)(C)C(C)(C)C)CC1CCC1C2CCC2(C)C(O[Si](C)(C)C(C)(C)C)CCC12

Mol. weight [g/mol]:

520.98

CAS:

57711-48-5

## Physical Properties

Property code	Value	Unit	Source
log10ws	-5.43		Crippen Method
logP	9.810		Crippen Method
rinpol	3070.80		NIST Webbook

## Sources

NIST Webbook:

<http://webbook.nist.gov/cgi/cbook.cgi?ID=C57711485&Units=SI>

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)

## Legend

log10ws: Log10 of Water solubility in mol/l

logP: Octanol/Water partition coefficient

rinpol: Non-polar retention indices

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[https://www.chemeo.com/cid/76-538-1/Silane-3-alpha-5-alpha-17-beta-androstane-3-17-diyloxy\]bis\(1,1-dimethylethyl\)dimethylsilane](https://www.chemeo.com/cid/76-538-1/Silane-3-alpha-5-alpha-17-beta-androstane-3-17-diyloxy]bis(1,1-dimethylethyl)dimethylsilane)

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