

16-Dehydropregnenolone, MO-TMS

Inchi: InChI=1S/C25H41NO2Si/c1-17(26-27-4)21-10-11-22-20-9-8-18-16-19(28-29(5,6)7)12-14
InchiKey: CISBSUGRAJLSLW-ALNVHOQCSA-N
Formula: C25H41NO2Si
SMILES: CON=C(C)C1=CCC2C3CC=C4CC(O[Si](C)(C)C)CCC4(C)C3CCC12C
Mol. weight [g/mol]: 415.68

Physical Properties

Property code	Value	Unit	Source
log10ws	-4.86		Crippen Method
logp	6.728		Crippen Method
rinpol	2802.00		NIST Webbook

Sources

NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=R512702&Units=SI>
Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci9903071>
Crippen Method: 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

Latest version available from:

<https://www.chemeo.com/cid/63-084-9/16-Dehydropregnenolone-MO-TMS.pdf>

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