

17-«alpha»-Hydroxy-7-«alpha»,17-«alpha»-dimethyl-17-TMS

InChI: InChI=1S/C24H42O2Si/c1-16-14-17-15-18(25)8-11-22(17,2)19-9-12-23(3)20(21(16)19)1
InChIKey: WOJBEUHZDDQLML-PDBPDUHWSA-N
Formula: C24H42O2Si
SMILES: CC1CC2CC(=O)CCC2(C)C2CCC3(C)C(CCC3(C)O[Si](C)(C)C)C12
Mol. weight [g/mol]: 390.67

Physical Properties

Property code	Value	Unit	Source
log10ws	-4.29		Crippen Method
logp	6.454		Crippen Method
rinpol	2641.00		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=R257188&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/31-951-1/17-alpha-Hydroxy-7-alpha-17-alpha-dimethyl-5-beta-androstan-3-one-17-TMS>

Generated by Cheméo on 2025-12-05 14:42:33.932372569 +0000 UTC m=+4693951.462413241.

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