

6.beta.-Hydroxy-oral turinabol, tri-trimethylsilyl

Inchi: InChI=1S/C29H51ClO3Si3/c1-27-16-15-23(31-34(4,5)6)26(30)25(27)24(32-35(7,8)9)19-2
InchiKey: RJRDWQNBTOPZNE-UHFFFAOYSA-N
Formula: C29H51ClO3Si3
SMILES: CC12C=CC(O[Si](C)(C)C)=C(Cl)C1=C(O[Si](C)(C)C)CC1C2CCC2(C)C1CCC2(C)O[Si](C)(C)C
Mol. weight [g/mol]: 567.44

Physical Properties

Property code	Value	Unit	Source
logp	9.426		Crippen Method

Sources

NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=U412281&Units=SI>
Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci990307l>
Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
Cheméo Relay (1.0):
Cheméo Relay (1.0, beta):

Legend

logp: Octanol/Water partition coefficient

Latest version available from:

<https://www.chemeo.com/cid/74-643-6/6-beta-Hydroxy-oral-turinabol-tri-trimethylsilyl.pdf>

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