6«alpha»-Hydroxy-Methandienone, tris-TMS

Inchi: InChI=1S/C29H52O3Si3/c1-27-16-13-21(30-33(4,5)6)19-25(27)26(31-34(7,8)9)20-22-23(
InchiKey: CBUDFJGEOLVXDB-GCDQPWEESA-N
Formula: C29H52O3Si3
SMILES: CC12C=CC(O[Si](C)(C)C)=CC1=C(O[Si](C)(C)C)CC1C2CCC2(C)C1CCC2(C)O[Si](C)(C
Mol. weight [g/mol]: 532.98

Physical Properties

Property code	Value	Unit	Source
logp	8.860		Crippen Method
rinpol	2888.00		NIST Webbook
rinpol	2888.00		NIST Webbook
rinpol	2888.00		NIST Webbook

Sources

Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
Cheméo Relay (1.0):
Cheméo Relay (1.0, beta):
NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=R495944&Units=SI
Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l

Legend

logp: Octanol/Water partition coefficient
rinpol: Non-polar retention indices