

11-keto-methylprednisolone, diMO-diTMS (1)

Inchi: InChI=1S/C30H50N2O5Si2/c1-20-16-22-23-13-15-30(37-39(9,10)11,26(32-35-5)19-36-3
InchiKey: SFOCNCWFMBQHIL-NKLIHJSYSA-N
Formula: C30H50N2O5Si2
SMILES: CON=C1CC2(C)C(CCC2(O[Si](C)(C)C)C(CO[Si](C)(C)C)=NOC)C2CC(C)C3=CC(=O)C=C
Mol. weight [g/mol]: 574.90

Physical Properties

Property code	Value	Unit	Source
log10ws	-2.37		Crippen Method
logp	6.597		Crippen Method
rinpol	3195.00		NIST Webbook
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Sources

Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=R251982&Units=SI>
Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci990307l>

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/124-886-1/11-keto-methylprednisolone-diMO-diTMS-1.pdf>

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