

19-Hydroxy PGF2 TMS

Inchi: InChI=1S/C32H64O6Si4/c1-26(35-39(2,3)4)19-18-20-27(33)23-24-29-28(21-16-14-15-17)
InchiKey: YBVFPNBVTWJBFJ-JCSKAYPWSA-N
Formula: C32H64O6Si4
SMILES: CC(CCCC(=O)C=CC1C(O[Si](C)(C)C)CC(O[Si](C)(C)C)C1CC=CCCCC(=O)O[Si](C)(C)C
Mol. weight [g/mol]: 657.19

Physical Properties

Property code	Value	Unit	Source
log10ws	-0.28		Crippen Method
logp	9.093		Crippen Method
rinpol	2892.00		NIST Webbook

Sources

Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=R16812&Units=SI>
Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci9903071>

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/61-373-0/19-Hydroxy-PGF2-TMS.pdf>

Generated by Cheméo on 2024-04-19 14:10:42.403113901 +0000 UTC m=+15825091.323691222.

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