Methyl 5-«beta»-cholan-12-«alpha»-ol-3-one-24-oate, oxime, TMS #1

Inchi: InChI=1S/C31H57NO4Si2/c1-21(11-16-29(33)34-4)25-14-15-26-24-13-12-22-19-23(32-3
InchiKey: QIWSBVFPWMELCB-NAALBDNSSA-N
Formula: C31H57NO4Si2
SMILES: COC(=O)CCC(C)C1CCC2C3CCC4CC(=NO[Si](C)(C)C)CCC4(C)C3CC(O[Si](C)(C)C)C1
Mol. weight [g/mol]: 563.96

Physical Properties

Property code	Value	Unit	Source
log10ws	-6.44		Cheméo Relay (1.0)
logp	8.272		Crippen Method
rinpol	3301.00		NIST Webbook
rinpol	3302.00		NIST Webbook
rinpol	3301.00		NIST Webbook

Sources

Cheméo Relay (1.0):
Cheméo Relay (1.0, beta):
NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=R215773&Units=SI
Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws

Legend

log10ws: Log10 of Water solubility in mol/l
logp: Octanol/Water partition coefficient
rinpol: Non-polar retention indices

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