

# 3-oxy-7«alpha»-hydroxy-4-cholestenoate, methyl ester-trimethylsilyl ether

Inchi:	InChI=1S/C34H60O4Si2/c1-23(13-12-14-24(2)32(35)36-5)27-15-16-28-31-29(18-20-34(2
InchiKey:	UDKJVEQQSDKFME-NAUBTBHQSA-N
Formula:	C34H60O4Si2
SMILES:	COC(=O)C(C)CCCC(C)C1CCC2C3C(O[Si](C)(C)C)CC4=CC(O[Si](C)(C)C)=CCC4(C)C3
Mol. weight [g/mol]:	589.01

## Physical Properties

Property code	Value	Unit	Source
log10ws	-5.15		Crippen Method
logp	9.356		Crippen Method
rinsol	3535.00		NIST Webbook
rinsol	3535.00		NIST Webbook

## Sources

Crippen Method:	<a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</a>
Crippen Method:	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>
NIST Webbook:	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=R493879&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R493879&amp;Units=SI</a>

## Legend

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

Latest version available from:

<https://www.chemeo.com/cid/58-050-2/3-oxy-7-alpha-hydroxy-4-cholestenoate-methyl-ester-trimethylsilyl-ether.pdf>

Generated by Cheméo on 2024-08-06 13:55:29.314947871 +0000 UTC m=+1642398.562053233.

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