

# 2-Propenoic acid, 2-methyl-3,3-bis[(trimethylsilyl)oxy]-, trimethylsilyl ester

<b>Inchi:</b>	InChI=1S/C13H30O4Si3/c1-11(12(14)15-18(2,3)4)13(16-19(5,6)7)17-20(8,9)10/h1-10H3
<b>InchiKey:</b>	NZIQCKNREBFDLB-UHFFFAOYSA-N
<b>Formula:</b>	C13H30O4Si3
<b>SMILES:</b>	CC(C(=O)O[Si](C)(C)C)=C(O[Si](C)(C)C)O[Si](C)(C)C
<b>Mol. weight [g/mol]:</b>	334.63
<b>CAS:</b>	40333-08-2

## Physical Properties

Property code	Value	Unit	Source
log10ws	2.70		Crippen Method
logp	4.299		Crippen Method
rinpol	1424.00		NIST Webbook
rinpol	1424.00		NIST Webbook

## Sources

<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</a>
<b>Crippen Method:</b>	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>
<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C40333082&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C40333082&amp;Units=SI</a>

## Legend

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

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