

# Silane, methylvinyl(3-methylbutoxy)octyloxy-

<b>Inchi:</b>	InChI=1S/C16H34O2Si/c1-6-8-9-10-11-12-14-17-19(5,7-2)18-15-13-16(3)4/h7,16H,2,6,8
<b>InchiKey:</b>	RAZBHAUDIXTNDZ-UHFFFAOYSA-N
<b>Formula:</b>	C16H34O2Si
<b>SMILES:</b>	C=C[Si](C)(OCCCCCCCC)OCCC(C)C
<b>Mol. weight [g/mol]:</b>	286.53

## Physical Properties

Property code	Value	Unit	Source
log10ws	-3.03		Crippen Method
logp	5.223		Crippen Method
rinpol	1692.00		NIST Webbook
rinpol	1692.00		NIST Webbook

## Sources

<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=U416314&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U416314&amp;Units=SI</a>
<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci990307l">http://pubs.acs.org/doi/abs/10.1021/ci990307l</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

Latest version available from:

<https://www.chemeo.com/cid/123-903-2/Silane-methylvinyl-3-methylbutoxy-octyloxy.pdf>

Generated by Cheméo on 2024-05-24 10:44:25.243734626 +0000 UTC m=+18836714.164311937.

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