

# 1,3-bis([(2Z)-3,7-Dimethylocta-2,6-dien-1-yl]oxy)-1

<b>Other names:</b>	1,3-bis{[(2Z)-3,7-Dimethylocta-2,6-dien-1-yl]oxy}-1,1,3,3-tetramethyldisiloxane
<b>Inchi:</b>	InChI=1S/C24H46O3Si2/c1-21(2)13-11-15-23(5)17-19-25-28(7,8)27-29(9,10)26-20-18-2
<b>InchiKey:</b>	BBMZRMYYZZWZAH-BOYKQRMESA-N
<b>Formula:</b>	C24H46O3Si2
<b>SMILES:</b>	CC(C)=CCCC(C)=CCO[Si](C)(C)O[Si](C)(C)OCC=C(C)CCC=C(C)C
<b>Mol. weight [g/mol]:</b>	438.79

## Physical Properties

Property code	Value	Unit	Source
log10ws	-3.81		Crippen Method
logp	7.825		Crippen Method
rinsol	2219.40		NIST Webbook
rinsol	2219.40		NIST Webbook

## Sources

<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=U352642&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U352642&amp;Units=SI</a>
<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>

## Legend

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

Latest version available from:

<https://www.chemeo.com/cid/119-141-3/1-3-bis-2Z-3-7-Dimethylocta-2-6-dien-1-yl-oxy-1-1-3-3-tetramethyldisiloxane>

Generated by Cheméo on 2024-05-04 14:23:35.978708403 +0000 UTC m=+17121864.899285718.

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