

bis[(4Z)-Dec-4-en-1-yloxy](dimethyl)silane

Inchi: InChI=1S/C22H44O2Si/c1-5-7-9-11-13-15-17-19-21-23-25(3,4)24-22-20-18-16-14-12-10
InchiKey: RTAYPUXYDPVRCM-VMNXYWKNSA-N
Formula: C22H44O2Si
SMILES: CCCCCC=CCCCO[Si](C)(C)OCCCC=CCCCC
Mol. weight [g/mol]: 368.67

Physical Properties

Property code	Value	Unit	Source
log10ws	-5.55		Crippen Method
logp	7.555		Crippen Method
rinpol	2183.50		NIST Webbook
rinpol	2183.50		NIST Webbook

Sources

NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=U334103&Units=SI>
Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci9903071>
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

Latest version available from:

<https://www.chemeo.com/cid/14-114-9/bis-4Z-Dec-4-en-1-yloxy-dimethyl-silane.pdf>

Generated by Cheméo on 2024-04-29 00:26:43.712095802 +0000 UTC m=+16639652.632673114.

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