

Geraniol, picolinyloxydimethylsilyl ether

Inchi: InChI=1S/C18H29NO2Si/c1-16(2)8-6-9-17(3)11-13-20-22(4,5)21-15-18-10-7-12-19-14-16
InchiKey: PKFQQWWGVSLTBQ-GZTJUZNSA-N
Formula: C18H29NO2Si
SMILES: CC(C)=CCCC(C)=CCO[Si](C)(C)OCc1cccnc1
Mol. weight [g/mol]: 319.51

Physical Properties

Property code	Value	Unit	Source
log10ws	-3.65		Crippen Method
logp	5.009		Crippen Method
rinpol	2123.90		NIST Webbook
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Sources

NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=U334073&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/49-531-8/Geraniol-picolinyloxydimethylsilyl-ether.pdf>

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