

1-Octadecanol, picolinyloxydimethylsilyl ether

Inchi: InChI=1S/C26H49NO2Si/c1-4-5-6-7-8-9-10-11-12-13-14-15-16-17-18-19-23-28-30(2,3)2
InchiKey: GLZCCIQRBXHIKK-UHFFFAOYSA-N
Formula: C₂₆H₄₉NO₂Si
SMILES: CCCCCCCCCCCCCCCCCCO[Si](C)(C)OCc1cccnc1
Mol. weight [g/mol]: 435.76

Physical Properties

Property code	Value	Unit	Source
log10ws	-7.30		Crippen Method
logp	8.578		Crippen Method
rinpol	2930.00		NIST Webbook
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Sources

Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci990307l>
Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=U334129&Units=SI>

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/61-557-6/1-Octadecanol-picolinyloxydimethylsilyl-ether.pdf>

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