1-phenyl,3,10-dimethylsilatrane, b,c

Inchi: InChI=1S/C14H21NO3Si/c1-12-10-15-8-9-16-19(17-12,18-13(2)11-15)14-6-4-3-5-7-14/h3
InchiKey: HOBATWHQQHWNCD-UHFFFAOYSA-N
Formula: C14H21NO3Si
SMILES: CC1CN2CCO[Si](c3ccccc3)(O1)OC(C)C2
Mol. weight [g/mol]: 279.41

Physical Properties

Property code	Value	Unit	Source
log10ws	-3.82		Crippen Method
logp	0.989		Crippen Method
rinpol	2067.00		NIST Webbook
rinpol	2067.00		NIST Webbook

Sources

Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=R145901&Units=SI
Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l

Legend

log10ws: Log10 of Water solubility in mol/l
logp: Octanol/Water partition coefficient
rinpol: Non-polar retention indices