1-Hexanamine, mono-TMS

Inchi: InChI=1S/C9H23NSi/c1-5-6-7-8-9-10-11(2,3)4/h10H,5-9H2,1-4H3
InchiKey: GYOPIQDCFVVUEH-UHFFFAOYSA-N
Formula: C9H23NSi
SMILES: CCCCCCN[Si](C)(C)C
Mol. weight [g/mol]: 173.37

Physical Properties

Property code	Value	Unit	Source
log10ws	-0.83		Crippen Method
logp	2.991		Crippen Method
rinpol	1055.00		NIST Webbook

Sources

NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=R64984&Units=SI
Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l
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