2-Propylpentanol, dimethylpentafluorophenylsilyl ether

Inchi: InChI=1S/C16H23F5OSi/c1-5-7-10(8-6-2)9-22-23(3,4)16-14(20)12(18)11(17)13(19)15(16

InchiKey: VXUXLZOPPMNTKB-UHFFFAOYSA-N

Formula: C16H23F5OSi

SMILES: CCCC(CCC)CO[Si](C)(C)c1c(F)c(F)c(F)c(F)c1F

Mol. weight [g/mol]: 354.43

Physical Properties

Property code	Value	Unit	Source
log10ws	-8.16		Crippen Method
logp	5.027		Crippen Method
rinpol	1546.00		NIST Webbook

Sources

Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws

NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=U368750&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