

2,4-Dimethylpentan-3-ol, dimethylpentafluorophenylsilyl ether

Inchi: InChI=1S/C15H21F5OSi/c1-7(2)14(8(3)4)21-22(5,6)15-12(19)10(17)9(16)11(18)13(15)20
InchiKey: YGDYVQRCZZPKQH-UHFFFAOYSA-N
Formula: C15H21F5OSi
SMILES: CC(C)C(O[Si](C)(C)c1c(F)c(F)c(F)c1F)C(C)C
Mol. weight [g/mol]: 340.40

Physical Properties

Property code	Value	Unit	Source
log10ws	-7.61		Crippen Method
logp	4.491		Crippen Method
rinpol	1441.00		NIST Webbook
rinpol	1441.00		NIST Webbook

Sources

Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci9903071>
Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=U368264&Units=SI>

Legend

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

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<https://www.chemeo.com/cid/69-139-2/2-4-Dimethylpentan-3-ol-dimethylpentafluorophenylsilyl-ether.pdf>

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