

Silane, dimethyl(pentafluorobenzyloxy)pentadecyloxy-

Inchi: InChI=1S/C24H39F5O2Si/c1-4-5-6-7-8-9-10-11-12-13-14-15-16-17-30-32(2,3)31-18-19-20
InchiKey: XHOKGJFGAOJNOY-UHFFFAOYSA-N
Formula: C24H39F5O2Si
SMILES: CCCCCCCCCCCCCCO[Si](C)(C)OCc1c(F)c(F)c(F)c(F)c1F
Mol. weight [g/mol]: 482.64

Physical Properties

Property code	Value	Unit	Source
log10ws	-7.93		Crippen Method
logp	8.708		Crippen Method
rinpol	2397.00		NIST Webbook
rinpol	2397.00		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=U347271&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/122-193-2/Silane-dimethyl-pentafluorobenzyloxy-pentadecyloxy.pdf>

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