

1,3-Difluoro-5-trihexylsilyloxybenzene

Inchi: InChI=1S/C24H42F2OSi/c1-4-7-10-13-16-28(17-14-11-8-5-2,18-15-12-9-6-3)27-24-20-22
InchiKey: PJOKLIQUPQTJQK-UHFFFAOYSA-N
Formula: C24H42F2OSi
SMILES: CCCCCC[Si](CCCCC)(CCCCC)Oc1cc(F)cc(F)c1
Mol. weight [g/mol]: 412.67

Physical Properties

Property code	Value	Unit	Source
log10ws	-7.42		Crippen Method
logp	9.030		Crippen Method
rinpol	2257.00		NIST Webbook

Sources

Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci990307l>
Crippen Method: https://www.cheméo.com/doc/models/crippen_log10ws
NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=U299091&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.cheméo.com/cid/65-943-3/1-3-Difluoro-5-trihexylsilyloxybenzene.pdf>

Generated by Cheméo on 2024-04-23 07:32:25.190694942 +0000 UTC m=+16146794.111272254.

Cheméo (<https://www.cheméo.com>) is the biggest free database of chemical and physical data for the process industry.