

1,7-Di(2-chlorophenyl)-2,2,4,4,6,6-hexamethyl-1,3,

Inchi: InChI=1S/C18H26Cl2O4Si3/c1-25(2,21-17-13-9-7-11-15(17)19)23-27(5,6)24-26(3,4)22-1
InchiKey: ADBQHPZJKQIGMK-UHFFFAOYSA-N
Formula: C18H26Cl2O4Si3
SMILES: C[Si](C)(Oc1ccccc1Cl)O[Si](C)(C)O[Si](C)(C)Oc1ccccc1Cl
Mol. weight [g/mol]: 461.56

Physical Properties

Property code	Value	Unit	Source
log10ws	-0.48		Crippen Method
logp	6.590		Crippen Method
rinpol	2273.00		NIST Webbook
rinpol	2273.00		NIST Webbook

Sources

Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=U347096&Units=SI>
Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci990307I>

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/114-220-0/1-7-Di-2-chlorophenyl-2-2-4-4-6-6-hexamethyl-1-3-5-7-tetraoxa-2-4-6-trisilah>

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