

1,7-Di(2,5-dichlorophenyl)-2,2,4,4,6,6-hexamethyl-

Inchi: InChI=1S/C18H24Cl4O4Si3/c1-27(2,23-17-11-13(19)7-9-15(17)21)25-29(5,6)26-28(3,4)2
InchiKey: KGTFRJOXCOZHFA-UHFFFAOYSA-N
Formula: C18H24Cl4O4Si3
SMILES: C[Si](C)(Oc1cc(Cl)ccc1Cl)O[Si](C)(C)O[Si](C)(C)Oc1cc(Cl)ccc1Cl
Mol. weight [g/mol]: 530.45

Physical Properties

Property code	Value	Unit	Source
log10ws	-1.85		Crippen Method
logp	7.897		Crippen Method
rinpol	2540.00		NIST Webbook
rinpol	2540.00		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=U347379&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/118-556-4/1-7-Di-2-5-dichlorophenyl-2-2-4-4-6-6-hexamethyl-1-3-5-7-tetraoxa-2-4-6-tris>

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