

Homoserine, ethoxycarbonylated, TBDMS

Inchi: InChI=1S/C19H41NO5Si2/c1-12-23-17(22)20-15(13-14-24-26(8,9)18(2,3)4)16(21)25-27(28)
InchiKey: CACLSQDKDKSIFH-UHFFFAOYSA-N
Formula: C19H41NO5Si2
SMILES: CCOC(=O)NC(CCO[Si](C)(C)C(C)(C)C)C(=O)O[Si](C)(C)C(C)(C)C
Mol. weight [g/mol]: 419.70

Physical Properties

Property code	Value	Unit	Source
log10ws	-0.97		Crippen Method
logp	5.061		Crippen Method
rinpol	2119.00		NIST Webbook

Sources

Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=R564446&Units=SI>
Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci990307l>

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/59-548-9/Homoserine-ethoxycarbonylated-TBDMS.pdf>

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