2-Aminobenzoic acid, N-heptafluorobutyryl-, N,O-bis(tert.-butyldimethylsilyl)-

Inchi: InChI=1S/C23H34F7NO3Si2/c1-19(2,3)35(7,8)31(18(33)21(24,25)22(26,27)23(28,29)30)
InchiKey: ANLVKEWFVROJOC-UHFFFAOYSA-N
Formula: C23H34F7NO3Si2
SMILES: CC(C)(C)[Si](C)(C)OC(=O)c1ccccc1N(C(=O)C(F)(F)C(F)(F)C(F)(F)F)[Si](C)(C)C(C)(C)C
Mol. weight [g/mol]: 561.68

Physical Properties

Property code	Value	Unit	Source
log10ws	-4.23		Crippen Method
logp	8.020		Crippen Method
rinpol	1956.00		NIST Webbook
rinpol	1956.00		NIST Webbook

Sources

Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
NIST Webbook: http://webbook.nist.gov/cgi/cbook.cgi?ID=U375101&Units=SI

Legend

log10ws: Log10 of Water solubility in mol/l
logp: Octanol/Water partition coefficient
rinpol: Non-polar retention indices