

1H-Indole-3-acetamide, N,1-bis(trimethylsilyl)-

Other names:	Indole-3-acetamide, TMS 3-Indolylacetamide, 2tms derivative
Inchi:	InChI=1S/C16H26N2OSi2/c1-20(2,3)17-16(19)11-13-12-18(21(4,5)6)15-10-8-7-9-14(13)
InchiKey:	KJUSHXFSECJYLY-UHFFFAOYSA-N
Formula:	C16H26N2OSi2
SMILES:	C[Si](C)(C)NC(=O)Cc1cn([Si](C)(C)C)c2ccccc12
Mol. weight [g/mol]:	318.56
CAS:	72101-39-4

Physical Properties

Property code	Value	Unit	Source
log10ws	-0.69		Crippen Method
logp	3.818		Crippen Method
rinpol	2066.00		NIST Webbook
rinpol	2066.00		NIST Webbook

Sources

Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C72101394&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071

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/51-576-6/1H-Indole-3-acetamide-N-1-bis-trimethylsilyl.pdf>

Generated by Cheméo on 2024-04-28 10:12:46.790319987 +0000 UTC m=+16588415.710897299.

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