

N-Trimethylsilylaniline

Other names:	1,1,1-Trimethyl-N-phenylsilamine 1,1,1-trimethyl-N-phenylsilanamine Aniline, tms derivative Anilino(trimethyl)silane N-(trimethylsilyl)aniline N-Phenyltrimethylsilylamine Silanamine, 1,1,1-trimethyl-N-phenyl- Silylamine, 1,1,1-trimethyl-N-phenyl- phenyl(trimethylsilyl)amine trimethyl(phenylamino)silane trimethyl-N-phenylsilylamine
Inchi:	InChI=1S/C9H15NSi/c1-11(2,3)10-9-7-5-4-6-8-9/h4-8,10H,1-3H3
InchiKey:	IPJPAQIHUIKFLV-UHFFFAOYSA-N
Formula:	C9H15NSi
SMILES:	<chem>C[Si](C)(C)Nc1ccccc1</chem>
Mol. weight [g/mol]:	165.31
CAS:	3768-55-6

Physical Properties

Property code	Value	Unit	Source
log10ws	-0.37		Crippen Method
logp	2.933		Crippen Method
rinpol	1186.00		NIST Webbook
rinpol	1186.00		NIST Webbook

Pressure Dependent Properties

Property code	Value	Unit	Pressure [kPa]	Source
tbp	329.15	K	0.40	Synthesis and characterization of organosilicon compounds as novel precursors for CVD processes

Sources

Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Synthesis and characterization of organosilicon compounds as novel	https://www.doi.org/10.1016/j.tca.2015.02.004
Module for CVD processes:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C3768556&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
tbp:	Boiling point at given pressure

Latest version available from:

<https://www.chemeo.com/cid/62-229-9/N-Trimethylsilylaniline.pdf>

Generated by Cheméo on 2024-04-18 09:53:59.69015292 +0000 UTC m=+15723288.610730253.

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