

Furazabol, 17-TMS

Other names:	Furazabole mono-TMS
Inchi:	InChI=1S/C23H38N2O2Si/c1-21-14-20-19(24-27-25-20)13-15(21)7-8-16-17(21)9-11-22(21)23
InchiKey:	TZFIHXONFDFXBC-UFOSUTSGSA-N
Formula:	C23H38N2O2Si
SMILES:	CC12Cc3nonc3CC1CCC1C2CCC2(C)C1CCC2(C)O[Si](C)(C)C
Mol. weight [g/mol]:	402.65

Physical Properties

Property code	Value	Unit	Source
log10ws	-8.84		Crippen Method
logp	5.637		Crippen Method
rinpol	2994.00		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=R172973&Units=SI

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/11-332-0/Furazabol-17-TMS.pdf>

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