

Bis(tert-butyldimethylsilyl) 2-[(tert-butyldimethylsilyl)amino]ethylphosphonate

Other names:	Phosphonic acid, [2-[(tert-butyldimethylsilyl)amino]ethyl]-, bis(tert-butyldimethylsilyl) ester 2-Bis(tert-butyldimethylsilyloxy)phosphoryl-N-tert-butyldimethylsilyl-ethanamine
Inchi:	InChI=1S/C20H50NO3PSi3/c1-18(2,3)26(10,11)21-16-17-25(22,23-27(12,13)19(4,5)6)24
InchiKey:	MWAOCBMMJSWPBR-UHFFFAOYSA-N
Formula:	C20H50NO3PSi3
SMILES:	CC(C)(C)[Si](C)(C)NCCP(=O)(O[Si](C)(C)C(C)(C)C)O[Si](C)(C)C(C)(C)C
Mol. weight [g/mol]:	467.85

Physical Properties

Property code	Value	Unit	Source
log10ws	-1.98		Crippen Method
logp	7.818		Crippen Method
rinpol	2137.00		NIST Webbook

Sources

Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U373307&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/46-332-2/Bis-tert-butyldimethylsilyl-2-tert-butyldimethylsilyl-amino-ethylphosphonate.pdf>

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