

DL-3-Methyl-2-butanol, tert-butyldimethylsilyl ether

Other names:	3-Methyl-2-butanol, tbdms derivative
Inchi:	InChI=1S/C11H26OSi/c1-9(2)10(3)12-13(7,8)11(4,5)6/h9-10H,1-8H3
InchiKey:	WGFVINWONVITGE-UHFFFAOYSA-N
Formula:	C11H26OSi
SMILES:	CC(C)C(C)O[Si](C)(C)C(C)(C)C
Mol. weight [g/mol]:	202.41

Physical Properties

Property code	Value	Unit	Source
logp	4.053		Crippen Method
rinpol	1031.40		NIST Webbook

Sources

Cheméo Relay (1.0, beta):

NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=U333062&Units=SI>

Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci990307l>

Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws

Cheméo Relay (1.0):

Legend

logp: Octanol/Water partition coefficient

rinpol: Non-polar retention indices

Latest version available from:

<https://www.chemeo.com/cid/12-425-6/DL-3-Methyl-2-butanol-tert-butyldimethylsilyl-ether.pdf>

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