

Dibenzothiophene, 2,4-dimethyl

Other names:	2,4-dimethyl-dibenzothiophene
Inchi:	InChI=1S/C14H12S/c1-9-7-10(2)14-12(8-9)11-5-3-4-6-13(11)15-14/h3-8H,1-2H3
InchiKey:	BNOHAWZYMVMJNG-UHFFFAOYSA-N
Formula:	C14H12S
SMILES:	<chem>Cc1cc(C)c2sc3ccccc3c2c1</chem>
Mol. weight [g/mol]:	212.31

Physical Properties

Property code	Value	Unit	Source
log10ws	-5.79		Crippen Method
logp	4.671		Crippen Method
mcvol	166.090	ml/mol	McGowan Method
rinpol	328.05		NIST Webbook
rinpol	330.72		NIST Webbook
rinpol	329.95		NIST Webbook
rinpol	329.92		NIST Webbook
rinpol	329.92		NIST Webbook
rinpol	329.95		NIST Webbook
rinpol	2050.00		NIST Webbook
rinpol	2050.00		NIST Webbook
rinpol	2050.00		NIST Webbook
rinpol	2050.00		NIST Webbook

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=R67275&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307I
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772

Legend

log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
rinpol:	Non-polar retention indices

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