

Dibenzothiophene, 4,6-dimethyl-

Other names:	4,6-dimethyldibenzo[b,d]thiophene 4,6-dimethyldibenzothiophene
Inchi:	InChI=1S/C14H12S/c1-9-5-3-7-11-12-8-4-6-10(2)14(12)15-13(9)11/h3-8H,1-2H3
InchiKey:	MYAQZIAVOLKEGW-UHFFFAOYSA-N
Formula:	C14H12S
SMILES:	<chem>Cc1cccc2c1sc1c(C)cccc12</chem>
Mol. weight [g/mol]:	212.31
CAS:	1207-12-1

Physical Properties

Property code	Value	Unit	Source
log10ws	-5.79		Crippen Method
logp	4.671		Crippen Method
mccvol	166.090	ml/mol	McGowan Method
rinpol	329.68		NIST Webbook
rinpol	2029.00		NIST Webbook
rinpol	2029.00		NIST Webbook
rinpol	328.80		NIST Webbook
rinpol	327.55		NIST Webbook
rinpol	328.32		NIST Webbook
rinpol	329.46		NIST Webbook
rinpol	2033.00		NIST Webbook
rinpol	329.20		NIST Webbook
rinpol	333.60		NIST Webbook
rinpol	329.17		NIST Webbook
rinpol	327.77		NIST Webbook
rinpol	325.75		NIST Webbook
rinpol	329.17		NIST Webbook
rinpol	329.60		NIST Webbook
rinpol	329.20		NIST Webbook
rinpol	329.20		NIST Webbook

Sources

Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
Liquid-Liquid Equilibria for <https://www.doi.org/10.1021/je100991y>
(1-Ethyl-3-methylimidazolium
Diethylphosphate) and Partial Molar Volumes at <https://www.doi.org/10.1021/je200327s>
in Ethyl Diphosphonate and Solid-Liquid
Equilibria of Dodecane + <http://link.springer.com/article/10.1007/BF02311772>
Diethylphosphate, Dodecane +
Alkane Systems, Dodecane + <http://webbook.nist.gov/cgi/cbook.cgi?ID=C1207121&Units=SI>
Dodecane Systems at 298.2 K and
213.2 K <http://pubs.acs.org/doi/abs/10.1021/ci990307i>
Crippen Method:

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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