

2,6,7-trimethyl-dibenzothiophene

Inchi:	InChI=1S/C15H14S/c1-9-4-7-14-13(8-9)12-6-5-10(2)11(3)15(12)16-14/h4-8H,1-3H3
InchiKey:	DESPWWKBIWQGNF-UHFFFAOYSA-N
Formula:	C15H14S
SMILES:	Cc1ccc2sc3c(C)c(C)ccc3c2c1
Mol. weight [g/mol]:	226.34

Physical Properties

Property code	Value	Unit	Source
log10ws	-6.27		Crippen Method
logp	4.980		Crippen Method
mcvol	180.180	ml/mol	McGowan Method
rinpol	352.91		NIST Webbook

Sources

Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=R436183&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071

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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