

Naphtho[2,1-b]thiophene, 9-methyl

Other names: Naphtho[2.1-b]thiophene, 9-methyl
Inchi: InChI=1S/C13H10S/c1-9-3-2-4-10-5-6-12-11(13(9)10)7-8-14-12/h2-8H,1H3
InchiKey: CFPORZKFHCXNOA-UHFFFAOYSA-N
Formula: C13H10S
SMILES: Cc1cccc2ccc3sccc3c12
Mol. weight [g/mol]: 198.28

Physical Properties

Property code	Value	Unit	Source
log10ws	-5.31		Crippen Method
logp	4.363		Crippen Method
mcvol	152.000	ml/mol	McGowan Method
rinpol	322.83		NIST Webbook
rinpol	325.25		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=R21649&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

Latest version available from:

<https://www.chemeo.com/cid/80-019-2/Naphtho-2-1-b-thiophene-9-methyl.pdf>

Generated by Cheméo on 2025-12-05 13:29:34.372206803 +0000 UTC m=+4689571.902247469.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.