

Phenanthro(1,2-b)thiophene

Inchi: InChI=1S/C16H10S/c1-2-4-13-11(3-1)5-8-15-14(13)7-6-12-9-10-17-16(12)15/h1-10H
InchiKey: FYSWUOGCANSBCW-UHFFFAOYSA-N
Formula: C16H10S
SMILES: c1ccc2c(c1)ccc1c2ccc2ccsc21
Mol. weight [g/mol]: 234.32
CAS: 58426-99-6

Physical Properties

Property code	Value	Unit	Source
log10ws	-6.64		Crippen Method
logp	5.208		Crippen Method
mcvol	174.810	ml/mol	McGowan Method
rinpol	395.52		NIST Webbook
rinpol	395.79		NIST Webbook
rinpol	396.01		NIST Webbook
rinpol	396.29		NIST Webbook
rinpol	396.29		NIST Webbook

Sources

NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=C58426996&Units=SI>
Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci9903071>
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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<https://www.chemeo.com/cid/78-374-1/Phenanthro-1-2-b-thiophene.pdf>

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