

2-Ethoxycarbonylphenyl isothiocyanate

Inchi: InChI=1S/C10H9NO2S/c1-2-13-10(12)8-5-3-4-6-9(8)11-7-14/h3-6H,2H2,1H3
InchiKey: TYEPWLJHQLYWSR-UHFFFAOYSA-N
Formula: C10H9NO2S
SMILES: CCOC(=O)c1ccccc1N=C=S
Mol. weight [g/mol]: 207.25
CAS: 99960-09-5

Physical Properties

Property code	Value	Unit	Source
hf	14.60	kJ/mol	Joback Method
hvap	60.39	kJ/mol	Joback Method
log10ws	-3.05		Crippen Method
logp	2.598		Crippen Method
mcvol	153.170	ml/mol	McGowan Method
pc	3107.10	kPa	Joback Method
tb	682.10	K	Joback Method
tc	933.38	K	Joback Method

Sources

Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci9903071>
Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
Joback Method: https://en.wikipedia.org/wiki/Joback_method
McGowan Method: <http://link.springer.com/article/10.1007/BF02311772>
NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=C99960095&Units=SI>

Legend

hf: Enthalpy of formation at standard conditions
hvap: Enthalpy of vaporization at standard conditions
log10ws: Log10 of Water solubility in mol/l

logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
pc:	Critical Pressure
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature

Latest version available from:

<https://www.cheméo.com/cid/83-500-4/2-Ethoxycarbonylphenyl-isothiocyanate.pdf>

Generated by Cheméo on 2024-04-28 23:18:29.994156952 +0000 UTC m=+16635558.914734267.

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