

3-Methoxyphenyl isothiocyanate

Inchi: InChI=1S/C8H7NOS/c1-10-8-4-2-3-7(5-8)9-6-11/h2-5H,1H3
InchiKey: WHBYCPUKGYEYFU-UHFFFAOYSA-N
Formula: C8H7NOS
SMILES: COc1cccc(N=C=S)c1
Mol. weight [g/mol]: 165.21
CAS: 3125-64-2

Physical Properties

Property code	Value	Unit	Source
hf	168.46	kJ/mol	Joback Method
hvap	49.19	kJ/mol	Joback Method
log10ws	-2.51		Crippen Method
logp	2.429		Crippen Method
mcvol	123.420	ml/mol	McGowan Method
pc	3607.21	kPa	Joback Method
tb	582.47	K	Joback Method
tc	839.09	K	Joback Method

Sources

Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci990307l>
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=C3125642&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

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