

2-Methoxybenzyl isothiocyanate

Inchi:	InChI=1S/C9H9NOS/c1-11-9-5-3-2-4-8(9)6-10-7-12/h2-5H,6H2,1H3
InchiKey:	HUYOVUYNRGOCLZ-UHFFFAOYSA-N
Formula:	C9H9NOS
SMILES:	COc1ccccc1CN=C=S
Mol. weight [g/mol]:	179.24
CAS:	17608-09-2

Physical Properties

Property code	Value	Unit	Source
hf	147.82	kJ/mol	Joback Method
hvap	51.42	kJ/mol	Joback Method
log10ws	-2.82		Crippen Method
logp	2.298		Crippen Method
mcvol	137.510	ml/mol	McGowan Method
pc	3224.64	kPa	Joback Method
tb	605.35	K	Joback Method
tc	856.17	K	Joback Method

Sources

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=C17608092&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307I

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/60-378-6/2-Methoxybenzyl-isothiocyanate.pdf>

Generated by Cheméo on 2025-12-23 00:41:16.758110775 +0000 UTC m=+6198674.288151439.

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