

2-Methylbutyl isothiocyanate

Inchi:	InChI=1S/C6H11NS/c1-3-6(2)4-7-5-8/h6H,3-4H2,1-2H3
InchiKey:	MKWQGOSIKKPMLW-UHFFFAOYSA-N
Formula:	C6H11NS
SMILES:	CCC(C)CN=C=S
Mol. weight [g/mol]:	129.22
CAS:	4404-51-7

Physical Properties

Property code	Value	Unit	Source
hf	111.62	kJ/mol	Joback Method
hvac	39.00	kJ/mol	Joback Method
log10ws	-2.03		Crippen Method
logp	2.135		Crippen Method
mccvol	113.130	ml/mol	McGowan Method
pc	3213.68	kPa	Joback Method
ripol	1412.00		NIST Webbook
ripol	1412.00		NIST Webbook
tb	482.19	K	Joback Method
tc	701.74	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=C4404517&Units=SI

Legend

hf: Enthalpy of formation at standard conditions

h_{vap}:	Enthalpy of vaporization at standard conditions
log₁₀w_s:	Log ₁₀ of Water solubility in mol/l
log_p:	Octanol/Water partition coefficient
mc_{vol}:	McGowan's characteristic volume
p_c:	Critical Pressure
ri_{pol}:	Polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature

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