

2-(Methylsulphonyl)propane

Other names:	Isopropyl methyl sulphone
Inchi:	InChI=1S/C4H10O2S/c1-4(2)7(3,5)6/h4H,1-3H3
InchiKey:	VTWYQAQIXXAXOR-UHFFFAOYSA-N
Formula:	C4H10O2S
SMILES:	CC(C)S(C)(=O)=O
Mol. weight [g/mol]:	122.19
CAS:	4853-74-1

Physical Properties

Property code	Value	Unit	Source
chl	-3101.30 ± 0.63	kJ/mol	NIST Webbook
gf	-488.18	kJ/mol	Joback Method
hf	-433.00 ± 3.00	kJ/mol	NIST Webbook
hfl	-503.80 ± 1.30	kJ/mol	NIST Webbook
hfus	13.97	kJ/mol	Joback Method
hvap	70.00 ± 3.00	kJ/mol	NIST Webbook
log10ws	-0.44		Crippen Method
logp	0.439		Crippen Method
mcvol	95.310	ml/mol	McGowan Method
pc	4829.24	kPa	Joback Method
tb	338.26	K	Joback Method
tc	504.95	K	Joback Method
tf	289.00 ± 3.00	K	NIST Webbook
vc	0.380	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	157.92	J/mol×K	338.26	Joback Method
cpg	166.83	J/mol×K	366.04	Joback Method
cpg	175.49	J/mol×K	393.82	Joback Method
cpg	183.89	J/mol×K	421.61	Joback Method
cpg	192.04	J/mol×K	449.39	Joback Method
cpg	199.93	J/mol×K	477.17	Joback Method

Sources

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=C4853741&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws

Legend

chl:	Standard liquid enthalpy of combustion
cpg:	Ideal gas heat capacity
gf:	Standard Gibbs free energy of formation
hf:	Enthalpy of formation at standard conditions
hfl:	Liquid phase enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion 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
tf:	Normal melting (fusion) point
vc:	Critical Volume

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