

# Methane sulfonamide

<b>Other names:</b>	Methylsulfonamide methanesulphonamide
<b>Inchi:</b>	InChI=1S/CH5NO2S/c1-5(2,3)4/h1H3,(H2,2,3,4)
<b>InchiKey:</b>	HNQIVZYLMDVSB-UHFFFAOYSA-N
<b>Formula:</b>	CH5NO2S
<b>SMILES:</b>	CS(N)(=O)=O
<b>Mol. weight [g/mol]:</b>	95.12
<b>CAS:</b>	3144-09-0

## Physical Properties

Property code	Value	Unit	Source
gf	-444.55	kJ/mol	Joback Method
hf	-483.53	kJ/mol	Joback Method
hfus	14.92	kJ/mol	Joback Method
hvap	47.10	kJ/mol	Joback Method
log10ws	0.50		Crippen Method
logp	-1.095		Crippen Method
mvol	63.020	ml/mol	McGowan Method
pc	8234.49	kPa	Joback Method
tb	342.59	K	Joback Method
tc	523.91	K	Joback Method
tf	222.85	K	Joback Method
vc	0.246	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	105.22	J/mol×K	342.59	Joback Method
cpg	110.31	J/mol×K	372.81	Joback Method
cpg	115.29	J/mol×K	403.03	Joback Method
cpg	120.13	J/mol×K	433.25	Joback Method
cpg	124.84	J/mol×K	463.47	Joback Method
cpg	129.39	J/mol×K	493.69	Joback Method
cpg	133.80	J/mol×K	523.91	Joback Method

# Sources

<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci990307l">http://pubs.acs.org/doi/abs/10.1021/ci990307l</a>
<b>Crippen Method:</b>	<a href="https://www.cheméo.com/doc/models/crippen_log10ws">https://www.cheméo.com/doc/models/crippen_log10ws</a>
<b>Joback Method:</b>	<a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>
<b>McGowan Method:</b>	<a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>
<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C3144090&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C3144090&amp;Units=SI</a>

# Legend

<b>cpg:</b>	Ideal gas heat capacity
<b>gf:</b>	Standard Gibbs free energy of formation
<b>hf:</b>	Enthalpy of formation at standard conditions
<b>hfus:</b>	Enthalpy of fusion at standard conditions
<b>hvap:</b>	Enthalpy of vaporization at standard conditions
<b>log10ws:</b>	Log10 of Water solubility in mol/l
<b>logp:</b>	Octanol/Water partition coefficient
<b>mcvol:</b>	McGowan's characteristic volume
<b>pc:</b>	Critical Pressure
<b>tb:</b>	Normal Boiling Point Temperature
<b>tc:</b>	Critical Temperature
<b>tf:</b>	Normal melting (fusion) point
<b>vc:</b>	Critical Volume

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