

# Propane, 1-bromo-2-methoxy

<b>Inchi:</b>	InChI=1S/C4H9BrO/c1-4(3-5)6-2/h4H,3H2,1-2H3
<b>InchiKey:</b>	ULHBRIULGNSYDP-UHFFFAOYSA-N
<b>Formula:</b>	C4H9BrO
<b>SMILES:</b>	COC(C)CBr
<b>Mol. weight [g/mol]:</b>	153.02

## Physical Properties

Property code	Value	Unit	Source
gf	-110.32	kJ/mol	Joback Method
hf	-237.06	kJ/mol	Joback Method
hfus	9.07	kJ/mol	Joback Method
hvap	32.95	kJ/mol	Joback Method
log10ws	-1.13		Crippen Method
logp	1.416		Crippen Method
mvol	90.590	ml/mol	McGowan Method
pc	4205.63	kPa	Joback Method
rinpol	754.00		NIST Webbook
rinpol	754.00		NIST Webbook
tb	379.06	K	Joback Method
tc	569.53	K	Joback Method
tf	201.87	K	Joback Method
vc	0.334	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	147.73	J/molxK	379.06	Joback Method
cpg	184.26	J/molxK	537.79	Joback Method
cpg	177.50	J/molxK	506.04	Joback Method
cpg	170.47	J/molxK	474.30	Joback Method
cpg	163.17	J/molxK	442.55	Joback Method
cpg	155.59	J/molxK	410.81	Joback Method
cpg	190.74	J/molxK	569.53	Joback Method
dvisc	0.0003160	Paxs	379.06	Joback Method

dvisc	0.0004067	Paxs	349.53	Joback Method
dvisc	0.0005483	Paxs	320.00	Joback Method
dvisc	0.0007856	Paxs	290.47	Joback Method
dvisc	0.0012210	Paxs	260.93	Joback Method
dvisc	0.0021239	Paxs	231.40	Joback Method
dvisc	0.0043438	Paxs	201.87	Joback Method

## Sources

<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</a>
<b>Crippen Method:</b>	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.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=R12042&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R12042&amp;Units=SI</a>

## Legend

<b>cpg:</b>	Ideal gas heat capacity
<b>dvisc:</b>	Dynamic viscosity
<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>rinpol:</b>	Non-polar retention indices
<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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