

# Cyclopentane, bromo-

<b>Other names:</b>	Bromocyclopentane Cyclopentyl bromide
<b>Inchi:</b>	InChI=1S/C5H9Br/c6-5-3-1-2-4-5/h5H,1-4H2
<b>InchiKey:</b>	BRTFVKHPEHKBQF-UHFFFAOYSA-N
<b>Formula:</b>	C5H9Br
<b>SMILES:</b>	BrC1CCCC1
<b>Mol. weight [g/mol]:</b>	149.03
<b>CAS:</b>	137-43-9

## Physical Properties

Property code	Value	Unit	Source
gf	42.09	kJ/mol	Joback Method
hf	9.83	kJ/mol	NIST Webbook
hfus	7.93	kJ/mol	Joback Method
hvap	33.42	kJ/mol	Joback Method
ie	9.94 ± 0.02	eV	NIST Webbook
log10ws	-2.35		Crippen Method
logp	2.324		Crippen Method
mvol	87.950	ml/mol	McGowan Method
pc	4762.81	kPa	Joback Method
rmpol	905.00		NIST Webbook
rmpol	905.00		NIST Webbook
tb	410.70	K	NIST Webbook
tc	614.76	K	Joback Method
tf	216.81	K	Joback Method
vc	0.319	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	140.14	J/mol×K	395.24	Joback Method
cpg	152.64	J/mol×K	431.83	Joback Method
cpg	164.36	J/mol×K	468.41	Joback Method
cpg	175.35	J/mol×K	505.00	Joback Method

cpg	185.63	J/molxK	541.59	Joback Method
cpg	195.25	J/molxK	578.17	Joback Method
cpg	204.23	J/molxK	614.76	Joback Method
dvisc	0.0034445	Paxs	216.81	Joback Method
dvisc	0.0020003	Paxs	246.55	Joback Method
dvisc	0.0013058	Paxs	276.29	Joback Method
dvisc	0.0009261	Paxs	306.02	Joback Method
dvisc	0.0006980	Paxs	335.76	Joback Method
dvisc	0.0005509	Paxs	365.50	Joback Method
dvisc	0.0004505	Paxs	395.24	Joback Method

## Sources

<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci990307I">http://pubs.acs.org/doi/abs/10.1021/ci990307I</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=C137439&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C137439&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>ie:</b>	Ionization energy
<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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