

# 2-Norbornyl bromide

<b>Other names:</b>	2-Bromobicyclo[2.2.1]heptane 2-Bromo-norbornane Norbornane, 2-bromo-
<b>Inchi:</b>	InChI=1S/C7H11Br/c8-7-4-5-1-2-6(7)3-5/h5-7H,1-4H2
<b>InchiKey:</b>	QXYOAWHKJRWNID-UHFFFAOYSA-N
<b>Formula:</b>	C7H11Br
<b>SMILES:</b>	BrC1CC2CCC1C2
<b>Mol. weight [g/mol]:</b>	175.07
<b>CAS:</b>	29342-65-2

## Physical Properties

Property code	Value	Unit	Source
gf	124.07	kJ/mol	Joback Method
hf	-42.38	kJ/mol	Joback Method
hfus	14.41	kJ/mol	Joback Method
hvap	37.30	kJ/mol	Joback Method
log10ws	-2.60		Crippen Method
logp	2.570		Crippen Method
mcvol	105.270	ml/mol	McGowan Method
pc	4062.13	kPa	Joback Method
tb	438.80	K	Joback Method
tc	662.17	K	Joback Method
tf	256.57	K	Joback Method
vc	0.395	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	205.03	J/molxK	438.80	Joback Method
cpg	220.89	J/molxK	476.03	Joback Method
cpg	235.60	J/molxK	513.26	Joback Method
cpg	249.23	J/molxK	550.48	Joback Method
cpg	261.87	J/molxK	587.71	Joback Method
cpg	273.59	J/molxK	624.94	Joback Method

cpg	284.46	J/mol×K	662.17	Joback Method
dvisc	0.0011366	Paxs	256.57	Joback Method
dvisc	0.0010516	Paxs	286.94	Joback Method
dvisc	0.0009876	Paxs	317.31	Joback Method
dvisc	0.0009377	Paxs	347.69	Joback Method
dvisc	0.0008977	Paxs	378.06	Joback Method
dvisc	0.0008651	Paxs	408.43	Joback Method
dvisc	0.0008379	Paxs	438.80	Joback Method

## Sources

<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=C29342652&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C29342652&amp;Units=SI</a>
<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>

## 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>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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