

# Cyclobutane, butyl-

<b>Other names:</b>	Butylcyclobutane
<b>Inchi:</b>	InChI=1S/C8H16/c1-2-3-5-8-6-4-7-8/h8H,2-7H2,1H3
<b>InchiKey:</b>	RFLJRIWHBUSIOM-UHFFFAOYSA-N
<b>Formula:</b>	C8H16
<b>SMILES:</b>	CCCCC1CCC1
<b>Mol. weight [g/mol]:</b>	112.21
<b>CAS:</b>	13152-44-8

## Physical Properties

Property code	Value	Unit	Source
gf	65.13	kJ/mol	Joback Method
hf	-141.81	kJ/mol	Joback Method
hfus	12.51	kJ/mol	Joback Method
hvap	33.49	kJ/mol	Joback Method
log10ws	-2.82		Crippen Method
logp	2.977		Crippen Method
mcvol	112.720	ml/mol	McGowan Method
pc	2982.79	kPa	Joback Method
tb	393.45	K	Joback Method
tc	578.37	K	Joback Method
tf	194.34	K	Joback Method
vc	0.432	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	215.12	J/mol×K	393.45	Joback Method
cpg	230.64	J/mol×K	424.27	Joback Method
cpg	245.42	J/mol×K	455.09	Joback Method
cpg	259.49	J/mol×K	485.91	Joback Method
cpg	272.87	J/mol×K	516.73	Joback Method
cpg	285.58	J/mol×K	547.55	Joback Method
cpg	297.67	J/mol×K	578.37	Joback Method
dvisc	0.0024213	Paxs	194.34	Joback Method

dvisc	0.0013675	Paxs	227.53	Joback Method
dvisc	0.0008932	Paxs	260.71	Joback Method
dvisc	0.0006424	Paxs	293.89	Joback Method
dvisc	0.0004939	Paxs	327.08	Joback Method
dvisc	0.0003986	Paxs	360.26	Joback Method
dvisc	0.0003335	Paxs	393.45	Joback Method

## Sources

<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C13152448&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C13152448&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>
<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>

## 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

Latest version available from:

<https://www.chemeo.com/cid/35-533-1/Cyclobutane-butyl.pdf>

Generated by Cheméo on 2024-04-20 16:03:31.788552986 +0000 UTC m=+15918260.709130298.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.