

# (E)-2-Butenylcyclopropane

<b>Other names:</b>	(2E)-2-Butenylcyclopropane trans-2-Butenyl-cyclopropane
<b>Inchi:</b>	InChI=1S/C7H12/c1-2-3-4-7-5-6-7/h2-3,7H,4-6H2,1H3/b3-2+
<b>InchiKey:</b>	HFWBELFXCMGIGS-NSCUHMNNSA-N
<b>Formula:</b>	C7H12
<b>SMILES:</b>	CC=CCC1CC1
<b>Mol. weight [g/mol]:</b>	96.17
<b>CAS:</b>	76588-98-2

## Physical Properties

Property code	Value	Unit	Source
gf	149.03	kJ/mol	Joback Method
hf	2.21	kJ/mol	Joback Method
hfus	12.22	kJ/mol	Joback Method
hvap	31.05	kJ/mol	Joback Method
log10ws	-2.26		Crippen Method
logp	2.363		Crippen Method
mvol	94.330	ml/mol	McGowan Method
pc	3427.87	kPa	Joback Method
rinpol	715.40		NIST Webbook
rinpol	715.00		NIST Webbook
tb	370.46	K	Joback Method
tc	558.31	K	Joback Method
tf	181.51	K	Joback Method
vc	0.364	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	163.59	J/molxK	370.46	Joback Method
cpg	176.97	J/molxK	401.77	Joback Method
cpg	189.56	J/molxK	433.08	Joback Method
cpg	201.42	J/molxK	464.39	Joback Method
cpg	212.58	J/molxK	495.70	Joback Method

cpg	223.08	J/molxK	527.00	Joback Method
cpg	232.96	J/molxK	558.31	Joback Method
dvisc	0.0008992	Paxs	181.51	Joback Method
dvisc	0.0006335	Paxs	213.00	Joback Method
dvisc	0.0004884	Paxs	244.49	Joback Method
dvisc	0.0003996	Paxs	275.99	Joback Method
dvisc	0.0003407	Paxs	307.48	Joback Method
dvisc	0.0002992	Paxs	338.97	Joback Method
dvisc	0.0002686	Paxs	370.46	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=C76588982&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C76588982&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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