

# 1-Propene, 1-(1-methoxy-1-methylethoxy)-, (Z)-

<b>Inchi:</b>	InChI=1S/C7H14O2/c1-5-6-9-7(2,3)8-4/h5-6H,1-4H3/b6-5-
<b>InchiKey:</b>	FEJFIWROECCXPD-WAYWQWQTSA-N
<b>Formula:</b>	C7H14O2
<b>SMILES:</b>	CC=COC(C)(C)OC
<b>Mol. weight [g/mol]:</b>	130.18
<b>CAS:</b>	62322-41-2

## Physical Properties

Property code	Value	Unit	Source
gf	-118.88	kJ/mol	Joback Method
hf	-343.78	kJ/mol	Joback Method
hfus	9.05	kJ/mol	Joback Method
hvap	34.66	kJ/mol	Joback Method
log10ws	-1.88		Crippen Method
logp	1.919		Crippen Method
mcvol	116.930	ml/mol	McGowan Method
pc	2912.39	kPa	Joback Method
tb	405.33	K	Joback Method
tc	589.82	K	Joback Method
tf	210.45	K	Joback Method
vc	0.432	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	229.13	J/molxK	405.33	Joback Method
cpg	241.52	J/molxK	436.08	Joback Method
cpg	253.38	J/molxK	466.83	Joback Method
cpg	264.71	J/molxK	497.58	Joback Method
cpg	275.54	J/molxK	528.33	Joback Method
cpg	285.88	J/molxK	559.07	Joback Method
cpg	295.73	J/molxK	589.82	Joback Method
dvisc	0.0046981	Paxs	210.45	Joback Method
dvisc	0.0018933	Paxs	242.93	Joback Method

dvisc	0.0009454	Paxs	275.41	Joback Method
dvisc	0.0005466	Paxs	307.89	Joback Method
dvisc	0.0003508	Paxs	340.37	Joback Method
dvisc	0.0002433	Paxs	372.85	Joback Method
dvisc	0.0001789	Paxs	405.33	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=C62322412&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C62322412&amp;Units=SI</a>
<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci990307l">http://pubs.acs.org/doi/abs/10.1021/ci990307l</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>mccvol:</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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