

# 2-Propenoic acid, 2-methyl-, 2-propenyl ester

Other names:	2-Methyl-2-propenyl 2-propenoate 2-Propenoic acid, 2-methyl-, 2-propen-1-yl ester 2-methylpropenoic acid, allyl ester 2-propenyl 2-methylpropenoate AGEFLEX AMA Allyl 2-methylacrylate Allyl ester of methacrylic acid Allyl methacrylate Allylester kyseliny methakrylove Methacrylic acid, allyl ester NSC 18597 allyl 2-methylpropenoate
Inchi:	InChI=1S/C7H10O2/c1-4-5-9-7(8)6(2)3/h4H,1-2,5H2,3H3
InchiKey:	FBCQUCJYYPMKRO-UHFFFAOYSA-N
Formula:	C7H10O2
SMILES:	C=CCOC(=O)C(=C)C
Mol. weight [g/mol]:	126.15
CAS:	96-05-9

## Physical Properties

Property code	Value	Unit	Source
gf	-58.73	kJ/mol	Joback Method
hf	-191.54	kJ/mol	Joback Method
hfus	12.80	kJ/mol	Joback Method
hvap	39.07	kJ/mol	Joback Method
log10ws	-1.32		Crippen Method
logp	1.292		Crippen Method
mcvol	108.330	ml/mol	McGowan Method
pc	3246.73	kPa	Joback Method
rinpol	849.00		NIST Webbook
ripol	1188.00		NIST Webbook
tb	429.09	K	Joback Method
tc	616.46	K	Joback Method
tf	223.33	K	Joback Method
vc	0.414	m3/kmol	Joback Method

# Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	207.19	J/molxK	429.09	Joback Method
cpg	217.16	J/molxK	460.32	Joback Method
cpg	226.71	J/molxK	491.55	Joback Method
cpg	235.86	J/molxK	522.78	Joback Method
cpg	244.60	J/molxK	554.01	Joback Method
cpg	252.95	J/molxK	585.24	Joback Method
cpg	260.92	J/molxK	616.46	Joback Method
rhoI	928.00	kg/m3	298.15	Densities and derived thermodynamic properties of the binary systems of 1,1-dimethylethyl methyl ether with allyl methacrylate, butyl methacrylate, methacrylic acid, and vinyl acetate at T = (298.15 and 308.15) K
rhoI	917.42	kg/m3	308.15	Densities and derived thermodynamic properties of the binary systems of 1,1-dimethylethyl methyl ether with allyl methacrylate, butyl methacrylate, methacrylic acid, and vinyl acetate at T = (298.15 and 308.15) K
rhoI	932.97	kg/m3	293.15	Density, excess volume, and excess coefficient of thermal expansion of the binary systems of dimethyl carbonate with butyl methacrylate, allyl methacrylate, styrene, and vinyl acetate at T = (293.15, 303.15, and 313.15) K

rhoI	922.52	kg/m3	303.15	Density, excess volume, and excess coefficient of thermal expansion of the binary systems of dimethyl carbonate with butyl methacrylate, allyl methacrylate, styrene, and vinyl acetate at T = (293.15, 303.15, and 313.15) K
rhoI	912.07	kg/m3	313.15	Density, excess volume, and excess coefficient of thermal expansion of the binary systems of dimethyl carbonate with butyl methacrylate, allyl methacrylate, styrene, and vinyl acetate at T = (293.15, 303.15, and 313.15) K
rhoI	928.00	kg/m3	298.15	Densities, isobaric thermal compressibilities and derived thermodynamic properties of the binary systems of cyclohexane with allyl methacrylate, butyl methacrylate, methacrylic acid, and vinyl acetate at t = (298.15 and 308.15)K

rho	917.42	kg/m <sup>3</sup>	308.15	Densities, isobaric thermal compressibilities and derived thermodynamic properties of the binary systems of cyclohexane with allyl methacrylate, butyl methacrylate, methacrylic acid, and vinyl acetate at t = (298.15 and 308.15)K
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## Pressure Dependent Properties

Property code	Value	Unit	Pressure [kPa]	Source
tbrp	333.20	K	5.70	NIST Webbook

## Sources

NIST Webbook:	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C96059&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C96059&amp;Units=SI</a>
Crippen Method:	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>
Crippen Method:	<a href="http://pubs.acs.org/doi/abs/10.1021/ci990307I">http://pubs.acs.org/doi/abs/10.1021/ci990307I</a>
KDB:	<a href="https://www.thermo.com/files/research/kdb/mol/mol1179.mol">https://www.thermo.com/files/research/kdb/mol/mol1179.mol</a>
Densities and derived thermodynamic properties of the binary systems of cyclohexane with allyl methacrylate, butyl methacrylate, methacrylic acid, and vinyl acetate at t = (298.15 and 308.15) K:	<a href="https://www.doi.org/10.1016/j.jct.2004.11.012">https://www.doi.org/10.1016/j.jct.2004.11.012</a>
Densities and derived thermodynamic properties of the binary systems of cyclohexane with allyl methacrylate, butyl methacrylate, methacrylic acid, and vinyl acetate at t = (298.15 and 308.15) K:	<a href="https://www.doi.org/10.1016/j.jct.2008.06.017">https://www.doi.org/10.1016/j.jct.2008.06.017</a>
Densities and derived thermodynamic properties of the binary systems of cyclohexane with allyl methacrylate, butyl methacrylate, methacrylic acid, and vinyl acetate at t = (298.15 and 308.15) K:	<a href="https://www.doi.org/10.1016/j.tca.2005.06.007">https://www.doi.org/10.1016/j.tca.2005.06.007</a>
Densities and derived thermodynamic properties of the binary systems of cyclohexane with allyl methacrylate, butyl methacrylate, methacrylic acid, and vinyl acetate at t = (298.15 and 308.15) K:	<a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>
Densities and derived thermodynamic properties of the binary systems of cyclohexane with allyl methacrylate, butyl methacrylate, methacrylic acid, and vinyl acetate at t = (298.15 and 308.15) K:	<a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>

## Legend

cp <sub>g</sub> :	Ideal gas heat capacity
gf:	Standard Gibbs free energy of formation
hf:	Enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion at standard conditions
h <sub>vap</sub> :	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>rho:</b>	Liquid Density
<b>rinpol:</b>	Non-polar retention indices
<b>ripol:</b>	Polar retention indices
<b>tb:</b>	Normal Boiling Point Temperature
<b>tbrp:</b>	Boiling point at reduced pressure
<b>tc:</b>	Critical Temperature
<b>tf:</b>	Normal melting (fusion) point
<b>vc:</b>	Critical Volume

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