

# Cyclopropyl carbinol

<b>Other names:</b>	Cyclopropanemethanol Cyclopropanemethyl alcohol Cyclopropylcarbinyl alcohol Cyclopropylmethanol Cyclopropylmethyl alcohol Hydroxymethylcyclopropane
<b>Inchi:</b>	InChI=1S/C4H8O/c5-3-4-1-2-4/h4-5H,1-3H2
<b>InchiKey:</b>	GUDMZGLFZNLYEY-UHFFFAOYSA-N
<b>Formula:</b>	C4H8O
<b>SMILES:</b>	OCC1CC1
<b>Mol. weight [g/mol]:</b>	72.11
<b>CAS:</b>	2516-33-8

## Physical Properties

Property code	Value	Unit	Source
gf	-93.27	kJ/mol	Joback Method
hf	-205.32	kJ/mol	Joback Method
hfus	8.34	kJ/mol	Joback Method
hvap	41.09	kJ/mol	Joback Method
log10ws	-0.41		Crippen Method
logp	0.389		Crippen Method
mcvol	62.230	ml/mol	McGowan Method
pc	5251.00	kPa	Joback Method
rinpol	648.00		NIST Webbook
rinpol	656.00		NIST Webbook
rinpol	648.00		NIST Webbook
tb	389.84	K	Joback Method
tc	565.69	K	Joback Method
tf	213.60	K	Joback Method
vc	0.235	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
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cpg	160.14	J/molxK	565.69	Joback Method
cpg	125.31	J/molxK	419.15	Joback Method
cpg	133.09	J/molxK	448.46	Joback Method
cpg	140.43	J/molxK	477.76	Joback Method
cpg	147.38	J/molxK	507.07	Joback Method
cpg	153.94	J/molxK	536.38	Joback Method
cpg	117.09	J/molxK	389.84	Joback Method
dvisc	0.0006896	Paxs	360.47	Joback Method
dvisc	0.0010983	Paxs	331.09	Joback Method
dvisc	0.0019150	Paxs	301.72	Joback Method
dvisc	0.0037646	Paxs	272.35	Joback Method
dvisc	0.0087142	Paxs	242.97	Joback Method
dvisc	0.0004645	Paxs	389.84	Joback Method
dvisc	0.0254091	Paxs	213.60	Joback Method
rho1	883.84	kg/m3	328.15	Densities and Viscosities for the Ternary System of Cyclopropanemethanol (1) + n-Dodecane (2) + Butylcyclohexane (3) and Corresponding Binaries at T = 293.15 343.15 K
rho1	879.49	kg/m3	333.15	Densities and Viscosities for the Ternary System of Cyclopropanemethanol (1) + n-Dodecane (2) + Butylcyclohexane (3) and Corresponding Binaries at T = 293.15 343.15 K
rho1	875.07	kg/m3	338.15	Densities and Viscosities for the Ternary System of Cyclopropanemethanol (1) + n-Dodecane (2) + Butylcyclohexane (3) and Corresponding Binaries at T = 293.15 343.15 K

rho1	870.62	kg/m3	343.15	Densities and Viscosities for the Ternary System of Cyclopropanemethanol (1) + n-Dodecane (2) + Butylcyclohexane (3) and Corresponding Binaries at T = 293.15 343.15 K
rho1	888.13	kg/m3	323.15	Densities and Viscosities for the Ternary System of Cyclopropanemethanol (1) + n-Dodecane (2) + Butylcyclohexane (3) and Corresponding Binaries at T = 293.15 343.15 K
rho1	908.96	kg/m3	298.15	Densities and Viscosities for the Ternary System of 1,2,3,4-Tetrahydronaphthalene + Isopropylcyclohexane + Cyclopropanemethanol and Corresponding Binaries at T = (293.15 to 343.15) K
rho1	904.87	kg/m3	303.15	Densities and Viscosities for the Ternary System of 1,2,3,4-Tetrahydronaphthalene + Isopropylcyclohexane + Cyclopropanemethanol and Corresponding Binaries at T = (293.15 to 343.15) K

rhoI	900.74	kg/m <sup>3</sup>	308.15	Densities and Viscosities for the Ternary System of 1,2,3,4-Tetrahydronaphthalene + Isopropylcyclohexane + Cyclopropanemethanol and Corresponding Binaries at T = (293.15 to 343.15) K
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rhoI	896.59	kg/m <sup>3</sup>	313.15	Densities and Viscosities for the Ternary System of 1,2,3,4-Tetrahydronaphthalene + Isopropylcyclohexane + Cyclopropanemethanol and Corresponding Binaries at T = (293.15 to 343.15) K
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rhoI	892.38	kg/m <sup>3</sup>	318.15	Densities and Viscosities for the Ternary System of 1,2,3,4-Tetrahydronaphthalene + Isopropylcyclohexane + Cyclopropanemethanol and Corresponding Binaries at T = (293.15 to 343.15) K
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rhoI	888.17	kg/m <sup>3</sup>	323.15	Densities and Viscosities for the Ternary System of 1,2,3,4-Tetrahydronaphthalene + Isopropylcyclohexane + Cyclopropanemethanol and Corresponding Binaries at T = (293.15 to 343.15) K
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rho1	883.86	kg/m3	328.15	Densities and Viscosities for the Ternary System of 1,2,3,4-Tetrahydronaphthalene + Isopropylcyclohexane + Cyclopropanemethanol and Corresponding Binaries at T = (293.15 to 343.15) K
rho1	879.52	kg/m3	333.15	Densities and Viscosities for the Ternary System of 1,2,3,4-Tetrahydronaphthalene + Isopropylcyclohexane + Cyclopropanemethanol and Corresponding Binaries at T = (293.15 to 343.15) K
rho1	875.10	kg/m3	338.15	Densities and Viscosities for the Ternary System of 1,2,3,4-Tetrahydronaphthalene + Isopropylcyclohexane + Cyclopropanemethanol and Corresponding Binaries at T = (293.15 to 343.15) K
rho1	870.62	kg/m3	343.15	Densities and Viscosities for the Ternary System of 1,2,3,4-Tetrahydronaphthalene + Isopropylcyclohexane + Cyclopropanemethanol and Corresponding Binaries at T = (293.15 to 343.15) K

rho1	892.38	kg/m3	318.15	Densities and Viscosities for the Ternary System of Cyclopropanemethanol (1) + n-Dodecane (2) + Butylcyclohexane (3) and Corresponding Binaries at T = 293.15 343.15 K
rho1	896.58	kg/m3	313.15	Densities and Viscosities for the Ternary System of Cyclopropanemethanol (1) + n-Dodecane (2) + Butylcyclohexane (3) and Corresponding Binaries at T = 293.15 343.15 K
rho1	900.74	kg/m3	308.15	Densities and Viscosities for the Ternary System of Cyclopropanemethanol (1) + n-Dodecane (2) + Butylcyclohexane (3) and Corresponding Binaries at T = 293.15 343.15 K
rho1	904.85	kg/m3	303.15	Densities and Viscosities for the Ternary System of Cyclopropanemethanol (1) + n-Dodecane (2) + Butylcyclohexane (3) and Corresponding Binaries at T = 293.15 343.15 K
rho1	908.94	kg/m3	298.15	Densities and Viscosities for the Ternary System of Cyclopropanemethanol (1) + n-Dodecane (2) + Butylcyclohexane (3) and Corresponding Binaries at T = 293.15 343.15 K

rho1	912.99	kg/m3	293.15	Densities and Viscosities for the Ternary System of Cyclopropanemethanol (1) + n-Dodecane (2) + Butylcyclohexane (3) and Corresponding Binaries at T = 293.15 343.15 K
rho1	913.02	kg/m3	293.15	Densities and Viscosities for the Ternary System of 1,2,3,4-Tetrahydronaphthalene + Isopropylcyclohexane + Cyclopropanemethanol and Corresponding Binaries at T = (293.15 to 343.15) K

## Pressure Dependent Properties

Property code	Value	Unit	Pressure [kPa]	Source
tbrp	396.70	K	98.40	NIST Webbook

## Sources

Joback Method:	<a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>
McGowan Method:	<a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>
NIST Webbook:	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C2516338&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C2516338&amp;Units=SI</a>
Crippen Method:	<a href="http://pubs.acs.org/doi/abs/10.1021/ci990307I">http://pubs.acs.org/doi/abs/10.1021/ci990307I</a>
Crippen Method:	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>
Densities and Viscosities for the Ternary System of Cyclopropanemethanol (1) + n-Dodecane (2) + Butylcyclohexane (3) and Corresponding Binaries at T = 293.15 343.15 K	<a href="https://www.doi.org/10.1021/acs.jced.7b00201">https://www.doi.org/10.1021/acs.jced.7b00201</a>
Densities and Viscosities for the Ternary System of 1,2,3,4-Tetrahydronaphthalene + Isopropylcyclohexane + Cyclopropanemethanol and Corresponding Binaries at T = (293.15 to 343.15) K	<a href="https://www.doi.org/10.1021/acs.jced.8b00662">https://www.doi.org/10.1021/acs.jced.8b00662</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>rho:</b>	Liquid Density
<b>rinpol:</b>	Non-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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