

Cyclohexanemethanol

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| Other names: | (hydroxymethyl)cyclohexane Benzyl alcohol, hexahydro- Cyclohexanecarbinol Cyclohexylmethyl alcohol Methanol, cyclohexyl- NSC 5288 USAF DO-49 cyclohexylcarbinol cyclohexylmethanol hexahydrobenzyl alcohol |
| Inchi: | InChI=1S/C7H14O/c8-6-7-4-2-1-3-5-7/h7-8H,1-6H2 |
| InchiKey: | VSSAZBXXNIABDN-UHFFFAOYSA-N |
| Formula: | C7H14O |
| SMILES: | OCC1CCCCC1 |
| Mol. weight [g/mol]: | 114.19 |
| CAS: | 100-49-2 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|----------|---------|----------------|
| affp | 802.10 | kJ/mol | NIST Webbook |
| basg | 771.70 | kJ/mol | NIST Webbook |
| chl | -4377.30 | kJ/mol | NIST Webbook |
| gf | -104.31 | kJ/mol | Joback Method |
| hf | -285.72 | kJ/mol | Joback Method |
| hfl | -378.00 | kJ/mol | NIST Webbook |
| hfus | 9.81 | kJ/mol | Joback Method |
| hvap | 48.28 | kJ/mol | Joback Method |
| log10ws | -1.67 | | Crippen Method |
| logp | 1.559 | | Crippen Method |
| mcvol | 104.500 | ml/mol | McGowan Method |
| pc | 3940.66 | kPa | Joback Method |
| tb | 456.20 | K | NIST Webbook |
| tb | 454.20 | K | NIST Webbook |
| tc | 664.09 | K | Joback Method |
| tf | 236.85 | K | Joback Method |
| vc | 0.380 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---|
| cpg | 270.29 | J/molxK | 567.69 | Joback Method |
| cpg | 304.24 | J/molxK | 664.09 | Joback Method |
| cpg | 293.54 | J/molxK | 631.95 | Joback Method |
| cpg | 282.23 | J/molxK | 599.82 | Joback Method |
| cpg | 257.70 | J/molxK | 535.56 | Joback Method |
| cpg | 244.45 | J/molxK | 503.42 | Joback Method |
| cpg | 230.52 | J/molxK | 471.29 | Joback Method |
| cpl | 230.56 | J/molxK | 293.83 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |
| cpl | 275.87 | J/molxK | 334.67 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |
| cpl | 276.54 | J/molxK | 334.67 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |
| cpl | 270.55 | J/molxK | 329.57 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |
| cpl | 271.38 | J/molxK | 329.57 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |
| cpl | 265.56 | J/molxK | 324.46 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |
| cpl | 265.98 | J/molxK | 324.46 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |
| cpl | 259.49 | J/molxK | 319.35 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |
| cpl | 260.24 | J/molxK | 319.35 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |
| cpl | 253.34 | J/molxK | 314.25 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |

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|-----|--------|---------|--------|--|
| cpl | 254.26 | J/molxK | 314.25 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |
| cpl | 281.78 | J/molxK | 339.77 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |
| cpl | 247.94 | J/molxK | 309.14 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |
| cpl | 241.62 | J/molxK | 304.04 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |
| cpl | 242.28 | J/molxK | 304.04 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |
| cpl | 235.55 | J/molxK | 298.93 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |
| cpl | 247.36 | J/molxK | 309.14 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |
| cpl | 229.98 | J/molxK | 293.83 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |
| cpl | 280.86 | J/molxK | 339.77 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |
| cpl | 287.35 | J/molxK | 344.88 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |
| cpl | 286.52 | J/molxK | 344.88 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |
| cpl | 236.46 | J/molxK | 298.93 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |
| cpl | 292.34 | J/molxK | 349.98 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |
| cpl | 291.51 | J/molxK | 349.98 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |
| cpl | 296.83 | J/molxK | 355.09 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |

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|-----|--------|---------|--------|--|
| cpl | 199.13 | J/molxK | 258.09 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |
| cpl | 198.88 | J/molxK | 258.09 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |
| cpl | 202.87 | J/molxK | 263.20 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |
| cpl | 202.54 | J/molxK | 263.20 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |
| cpl | 207.03 | J/molxK | 268.30 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |
| cpl | 206.61 | J/molxK | 268.30 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |
| cpl | 211.52 | J/molxK | 273.41 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |
| cpl | 210.94 | J/molxK | 273.41 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |
| cpl | 215.68 | J/molxK | 278.51 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |
| cpl | 215.18 | J/molxK | 278.51 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |
| cpl | 220.83 | J/molxK | 283.62 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |
| cpl | 219.83 | J/molxK | 283.62 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |
| cpl | 225.32 | J/molxK | 288.72 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |
| cpl | 224.74 | J/molxK | 288.72 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |

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|-------|-----------|---------|--------|---|
| cpl | 296.33 | J/molxK | 355.09 | Heat Capacities of 2-Propenol and Selected Cyclohexylalcohols |
| dvisc | 0.0757236 | Paxs | 236.85 | Joback Method |
| dvisc | 0.0140832 | Paxs | 275.92 | Joback Method |
| dvisc | 0.0039756 | Paxs | 315.00 | Joback Method |
| dvisc | 0.0014837 | Paxs | 354.07 | Joback Method |
| dvisc | 0.0006736 | Paxs | 393.14 | Joback Method |
| dvisc | 0.0003527 | Paxs | 432.22 | Joback Method |
| dvisc | 0.0002056 | Paxs | 471.29 | Joback Method |
| rho1 | 900.67 | kg/m3 | 333.15 | Density, Viscosity, Refractive Index, and Surface Tension for Six Binary Systems of Adamantane Derivatives with 1-Heptanol and Cyclohexylmethanol |
| rho1 | 929.76 | kg/m3 | 293.15 | Density, Viscosity, Refractive Index, and Surface Tension for Six Binary Systems of Adamantane Derivatives with 1-Heptanol and Cyclohexylmethanol |
| rho1 | 900.78 | kg/m3 | 333.15 | Density, Viscosity, and Freezing Point for Four Binary Systems of n-Dodecane or Methylcyclohexane Mixed with 1-Heptanol or Cyclohexylmethanol |
| rho1 | 904.57 | kg/m3 | 328.15 | Density, Viscosity, and Freezing Point for Four Binary Systems of n-Dodecane or Methylcyclohexane Mixed with 1-Heptanol or Cyclohexylmethanol |
| rho1 | 908.31 | kg/m3 | 323.15 | Density, Viscosity, and Freezing Point for Four Binary Systems of n-Dodecane or Methylcyclohexane Mixed with 1-Heptanol or Cyclohexylmethanol |

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|------|--------|-------|--------|---|
| rhoI | 912.00 | kg/m3 | 318.15 | Density, Viscosity, and Freezing Point for Four Binary Systems of n-Dodecane or Methylcyclohexane Mixed with 1-Heptanol or Cyclohexylmethanol |
| rhoI | 915.66 | kg/m3 | 313.15 | Density, Viscosity, and Freezing Point for Four Binary Systems of n-Dodecane or Methylcyclohexane Mixed with 1-Heptanol or Cyclohexylmethanol |
| rhoI | 919.28 | kg/m3 | 308.15 | Density, Viscosity, and Freezing Point for Four Binary Systems of n-Dodecane or Methylcyclohexane Mixed with 1-Heptanol or Cyclohexylmethanol |
| rhoI | 926.25 | kg/m3 | 298.15 | Density, Viscosity, Refractive Index, and Surface Tension for Six Binary Systems of Adamantane Derivatives with 1-Heptanol and Cyclohexylmethanol |
| rhoI | 926.43 | kg/m3 | 298.15 | Density, Viscosity, and Freezing Point for Four Binary Systems of n-Dodecane or Methylcyclohexane Mixed with 1-Heptanol or Cyclohexylmethanol |
| rhoI | 929.96 | kg/m3 | 293.15 | Density, Viscosity, and Freezing Point for Four Binary Systems of n-Dodecane or Methylcyclohexane Mixed with 1-Heptanol or Cyclohexylmethanol |

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|------|--------|-------|--------|--|
| rhoI | 922.71 | kg/m3 | 303.15 | Density, Viscosity, Refractive Index, and Surface Tension for Six Binary Systems of Adamantane Derivatives with 1-Heptanol and Cyclohexylmethanol |
| rhoI | 919.14 | kg/m3 | 308.15 | Density, Viscosity, Refractive Index, and Surface Tension for Six Binary Systems of Adamantane Derivatives with 1-Heptanol and Cyclohexylmethanol |
| rhoI | 915.53 | kg/m3 | 313.15 | Density, Viscosity, Refractive Index, and Surface Tension for Six Binary Systems of Adamantane Derivatives with 1-Heptanol and Cyclohexylmethanol |
| rhoI | 911.88 | kg/m3 | 318.15 | Density, Viscosity, Refractive Index, and Surface Tension for Six Binary Systems of Adamantane Derivatives with 1-Heptanol and Cyclohexylmethanol |
| rhoI | 908.19 | kg/m3 | 323.15 | Density, Viscosity, Refractive Index, and Surface Tension for Six Binary Systems of Adamantane Derivatives with 1-Heptanol and Cyclohexylmethanol |
| rhoI | 904.45 | kg/m3 | 328.15 | Density, Viscosity, Refractive Index, and Surface Tension for Six Binary Systems of Adamantane Derivatives with 1-Heptanol and Cyclohexylmethanol |

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|------|--------|-------|--------|--|
| rhoI | 922.87 | kg/m3 | 303.15 | Density, Viscosity, and Freezing Point for Four Binary Systems of n-Dodecane or Methylcyclohexane Mixed with 1-Heptanol or Cyclohexylmethanol |
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Sources

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| Crippen Method: | https://www.chemeo.com/doc/models/crippen_log10ws |
| Heat Capacities of 2-Propanol and Selected Cyclohexylalcohols: | https://www.doi.org/10.1016/j.tca.2014.03.043 |
| Density, Viscosity, and Freezing Point for Four Binary Systems of n-Dodecane or Methylcyclohexane Mixed with 1-Heptanol or Cyclohexylmethanol: | https://www.doi.org/10.1021/acs.jced.6b00688 |
| Density, Viscosity, Refractive Index, and Surface Tension for Six Binary Systems of Adamantane Derivatives with 1-Heptanol and Cyclohexylmethanol: | https://www.doi.org/10.1021/je500381c |
| Joback Method: | https://en.wikipedia.org/wiki/Joback_method |
| NIST Webbook: | http://link.springer.com/article/10.1007/BF02311772 |
| Crippen Method: | http://webbook.nist.gov/cgi/cbook.cgi?ID=C100492&Units=SI |
| | http://pubs.acs.org/doi/abs/10.1021/ci990307I |

Legend

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|-----------------|---|
| affp: | Proton affinity |
| basg: | Gas basicity |
| chl: | Standard liquid enthalpy of combustion |
| cpg: | Ideal gas heat capacity |
| cpl: | Liquid phase heat capacity |
| dvisc: | Dynamic viscosity |
| gf: | Standard Gibbs free energy of formation |
| hf: | Enthalpy of formation at standard conditions |
| hfl: | Liquid phase enthalpy of formation at standard conditions |
| hfus: | Enthalpy of fusion at standard conditions |
| hvap: | Enthalpy of vaporization at standard conditions |
| log10ws: | Log10 of Water solubility in mol/l |
| logp: | Octanol/Water partition coefficient |
| mcvol: | McGowan's characteristic volume |
| pc: | Critical Pressure |
| rhoI: | Liquid Density |
| tb: | Normal Boiling Point Temperature |

tc: Critical Temperature
tf: Normal melting (fusion) point
vc: Critical Volume

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