

Bicyclo[2.2.1]hept-2-en-7-ol

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|-----------------------------|--|
| Inchi: | InChI=1S/C7H10O/c8-7-5-1-2-6(7)4-3-5/h1-2,5-8H,3-4H2 |
| InchiKey: | PSKWDFVNXXATCG-UHFFFAOYSA-N |
| Formula: | C7H10O |
| SMILES: | OC1C2C=CC1CC2 |
| Mol. weight [g/mol]: | 110.15 |
| CAS: | 53783-87-2 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | 2.89 | kJ/mol | Joback Method |
| hf | -163.16 | kJ/mol | Joback Method |
| hfus | 14.44 | kJ/mol | Joback Method |
| hvap | 47.84 | kJ/mol | Joback Method |
| log10ws | -1.29 | | Crippen Method |
| logp | 0.943 | | Crippen Method |
| mcvol | 89.340 | ml/mol | McGowan Method |
| pc | 4385.77 | kPa | Joback Method |
| rinpol | 1164.00 | | NIST Webbook |
| rinpol | 1164.00 | | NIST Webbook |
| ripol | 1709.00 | | NIST Webbook |
| ripol | 1709.00 | | NIST Webbook |
| tb | 463.98 | K | Joback Method |
| tc | 658.77 | K | Joback Method |
| tf | 258.35 | K | Joback Method |
| vc | 0.338 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 202.39 | J/molxK | 463.98 | Joback Method |
| cpg | 214.87 | J/molxK | 496.44 | Joback Method |
| cpg | 226.55 | J/molxK | 528.91 | Joback Method |
| cpg | 237.49 | J/molxK | 561.37 | Joback Method |
| cpg | 247.73 | J/molxK | 593.84 | Joback Method |

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|-------|-----------|---------|--------|---------------|
| cpg | 257.31 | J/mol×K | 626.30 | Joback Method |
| cpg | 266.29 | J/mol×K | 658.77 | Joback Method |
| dvisc | 0.0075063 | Paxs | 258.35 | Joback Method |
| dvisc | 0.0037559 | Paxs | 292.62 | Joback Method |
| dvisc | 0.0021730 | Paxs | 326.89 | Joback Method |
| dvisc | 0.0013948 | Paxs | 361.16 | Joback Method |
| dvisc | 0.0009668 | Paxs | 395.44 | Joback Method |
| dvisc | 0.0007104 | Paxs | 429.71 | Joback Method |
| dvisc | 0.0005464 | Paxs | 463.98 | Joback Method |

Sources

| | |
|------------------------|---|
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci9903071 |
| Crippen Method: | https://www.chemeo.com/doc/models/crippen_log10ws |
| Joback Method: | https://en.wikipedia.org/wiki/Joback_method |
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=C53783872&Units=SI |

Legend

| | |
|-----------------|---|
| cpg: | Ideal gas heat capacity |
| dvisc: | Dynamic viscosity |
| gf: | Standard Gibbs free energy of formation |
| hf: | 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 |
| mccvol: | McGowan's characteristic volume |
| pc: | Critical Pressure |
| rinpol: | Non-polar retention indices |
| ripol: | Polar retention indices |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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