

endo-5,6-Bis(hydroxymethyl)bicyclo[2.2.1]hept-2-

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|-----------------------------|--|
| Inchi: | InChI=1S/C9H14O2/c10-4-8-6-1-2-7(3-6)9(8)5-11/h1-2,6-11H,3-5H2 |
| InchiKey: | IGHHPVIMEQGKNE-UHFFFAOYSA-N |
| Formula: | C9H14O2 |
| SMILES: | OCC1C2C=CC(C2)C1CO |
| Mol. weight [g/mol]: | 154.21 |
| CAS: | 699-97-8 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------------|----------------------|----------------|
| gf | -124.80 | kJ/mol | Joback Method |
| hf | -377.01 | kJ/mol | Joback Method |
| hfus | 24.78 | kJ/mol | Joback Method |
| hvap | 68.66 | kJ/mol | Joback Method |
| log10ws | -0.79 | | Crippen Method |
| logp | 0.409 | | Crippen Method |
| mvol | 123.390 | ml/mol | McGowan Method |
| pc | 3704.46 | kPa | Joback Method |
| tb | 597.25 | K | Joback Method |
| tc | 777.84 | K | Joback Method |
| tf | 337.47 | K | Joback Method |
| tt | 353.00 ± 1.50 | K | NIST Webbook |
| vc | 0.468 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 347.39 | J/molxK | 597.25 | Joback Method |
| cpg | 359.28 | J/molxK | 627.35 | Joback Method |
| cpg | 370.51 | J/molxK | 657.45 | Joback Method |
| cpg | 381.13 | J/molxK | 687.54 | Joback Method |
| cpg | 391.17 | J/molxK | 717.64 | Joback Method |
| cpg | 400.67 | J/molxK | 747.74 | Joback Method |
| cpg | 409.67 | J/molxK | 777.84 | Joback Method |
| dvisc | 0.0119374 | Paxs | 337.47 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0040781 | Paxs | 380.77 | Joback Method |
| dvisc | 0.0017348 | Paxs | 424.06 | Joback Method |
| dvisc | 0.0008646 | Paxs | 467.36 | Joback Method |
| dvisc | 0.0004849 | Paxs | 510.66 | Joback Method |
| dvisc | 0.0002977 | Paxs | 553.95 | Joback Method |
| dvisc | 0.0001962 | Paxs | 597.25 | Joback Method |

Sources

| | |
|------------------------|---|
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=C699978&Units=SI |
| 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 |

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 |
| mcvol: | McGowan's characteristic volume |
| pc: | Critical Pressure |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| tt: | Triple Point Temperature |
| vc: | Critical Volume |

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