

5-Decanone, 6-hydroxy-

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|-----------------------------|--|
| Other names: | 4-hydroxydecan-5-one |
| Inchi: | InChI=1S/C10H20O2/c1-3-5-7-9(11)10(12)8-6-4-2/h9,11H,3-8H2,1-2H3 |
| InchiKey: | ORLDHCIQVZTBQQ-UHFFFAOYSA-N |
| Formula: | C10H20O2 |
| SMILES: | CCCCC(=O)C(O)CCCC |
| Mol. weight [g/mol]: | 172.26 |
| CAS: | 6540-98-3 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | -234.86 | kJ/mol | Joback Method |
| hf | -519.82 | kJ/mol | Joback Method |
| hfus | 23.82 | kJ/mol | Joback Method |
| hvap | 60.89 | kJ/mol | Joback Method |
| log10ws | -2.66 | | Crippen Method |
| logp | 2.297 | | Crippen Method |
| mcvol | 159.200 | ml/mol | McGowan Method |
| pc | 2462.92 | kPa | Joback Method |
| tb | 573.81 | K | Joback Method |
| tc | 743.57 | K | Joback Method |
| tf | 298.21 | K | Joback Method |
| vc | 0.615 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 403.15 | J/molxK | 573.81 | Joback Method |
| cpg | 415.84 | J/molxK | 602.10 | Joback Method |
| cpg | 427.99 | J/molxK | 630.40 | Joback Method |
| cpg | 439.61 | J/molxK | 658.69 | Joback Method |
| cpg | 450.71 | J/molxK | 686.99 | Joback Method |
| cpg | 461.31 | J/molxK | 715.28 | Joback Method |
| cpg | 471.43 | J/molxK | 743.57 | Joback Method |
| dvisc | 0.0185428 | Paxs | 298.21 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0043186 | Paxs | 344.14 | Joback Method |
| dvisc | 0.0014176 | Paxs | 390.08 | Joback Method |
| dvisc | 0.0005884 | Paxs | 436.01 | Joback Method |
| dvisc | 0.0002888 | Paxs | 481.94 | Joback Method |
| dvisc | 0.0001605 | Paxs | 527.88 | Joback Method |
| dvisc | 0.0000979 | Paxs | 573.81 | Joback Method |

Sources

| | |
|------------------------|---|
| 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=C6540983&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307l |
| Crippen Method: | https://www.chemeo.com/doc/models/crippen_log10ws |

Legend

| | |
|----------------------------|---|
| cp_g: | Ideal gas heat capacity |
| dvisc: | Dynamic viscosity |
| g_f: | Standard Gibbs free energy of formation |
| h_f: | Enthalpy of formation at standard conditions |
| h_{fus}: | Enthalpy of fusion at standard conditions |
| h_{vap}: | Enthalpy of vaporization at standard conditions |
| log₁₀ws: | Log ₁₀ of Water solubility in mol/l |
| log_p: | 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 |
| vc: | Critical Volume |

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