

Benzenemethanol, 2,5-dimethyl-

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| Other names: | 2,5-Dimethylbenzyl alcohol |
| Inchi: | InChI=1S/C9H12O/c1-7-3-4-8(2)9(5-7)6-10/h3-5,10H,6H2,1-2H3 |
| InchiKey: | LEBQTCCCNMTXSF-UHFFFAOYSA-N |
| Formula: | C9H12O |
| SMILES: | Cc1ccc(C)c(CO)c1 |
| Mol. weight [g/mol]: | 136.19 |
| CAS: | 53957-33-8 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -18.77 | kJ/mol | Joback Method |
| hf | -167.73 | kJ/mol | Joback Method |
| hfus | 16.42 | kJ/mol | Joback Method |
| hvap | 55.91 | kJ/mol | Joback Method |
| log10ws | -2.57 | | Crippen Method |
| logp | 1.796 | | Crippen Method |
| mcvol | 119.780 | ml/mol | McGowan Method |
| pc | 3505.43 | kPa | Joback Method |
| tb | 506.20 | K | NIST Webbook |
| tc | 731.21 | K | Joback Method |
| tf | 303.47 | K | Joback Method |
| vc | 0.451 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 266.07 | J/molxK | 534.14 | Joback Method |
| cpg | 277.11 | J/molxK | 566.99 | Joback Method |
| cpg | 287.59 | J/molxK | 599.83 | Joback Method |
| cpg | 297.54 | J/molxK | 632.68 | Joback Method |
| cpg | 306.97 | J/molxK | 665.52 | Joback Method |
| cpg | 315.91 | J/molxK | 698.37 | Joback Method |
| cpg | 324.37 | J/molxK | 731.21 | Joback Method |
| dvisc | 0.0061106 | Paxs | 303.47 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0021587 | Paxs | 341.92 | Joback Method |
| dvisc | 0.0009411 | Paxs | 380.36 | Joback Method |
| dvisc | 0.0004778 | Paxs | 418.81 | Joback Method |
| dvisc | 0.0002719 | Paxs | 457.25 | Joback Method |
| dvisc | 0.0001689 | Paxs | 495.69 | Joback Method |
| dvisc | 0.0001123 | Paxs | 534.14 | Joback Method |

Sources

| | |
|------------------------|---|
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=C53957338&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307l |
| Crippen Method: | https://www.chemeo.com/doc/models/crippen_log10ws |
| Joback Method: | https://en.wikipedia.org/wiki/Joback_method |

Legend

| | |
|----------------------------|---|
| cp_g: | 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: | Log ₁₀ 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 |
| vc: | Critical Volume |

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<https://www.chemeo.com/cid/50-822-3/Benzenemethanol-2-5-dimethyl.pdf>

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