

2-Methoxybenzyl alcohol, 2-methylpropyl ether

| | |
|----------------------|---|
| Inchi: | InChI=1S/C12H18O2/c1-10(2)8-14-9-11-6-4-5-7-12(11)13-3/h4-7,10H,8-9H2,1-3H3 |
| InchiKey: | SAHQWLNZCJVIOC-UHFFFAOYSA-N |
| Formula: | C12H18O2 |
| SMILES: | COc1ccccc1COCC(C)C |
| Mol. weight [g/mol]: | 194.27 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -59.50 | kJ/mol | Joback Method |
| hf | -335.67 | kJ/mol | Joback Method |
| hfus | 19.34 | kJ/mol | Joback Method |
| hvap | 49.68 | kJ/mol | Joback Method |
| log10ws | -2.99 | | Crippen Method |
| logp | 2.868 | | Crippen Method |
| mcvol | 167.920 | ml/mol | McGowan Method |
| pc | 2311.39 | kPa | Joback Method |
| rinsol | 1430.00 | | NIST Webbook |
| tb | 550.02 | K | Joback Method |
| tc | 751.26 | K | Joback Method |
| tf | 293.40 | K | Joback Method |
| vc | 0.629 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 398.86 | J/molxK | 550.02 | Joback Method |
| cpg | 415.03 | J/molxK | 583.56 | Joback Method |
| cpg | 430.43 | J/molxK | 617.10 | Joback Method |
| cpg | 445.09 | J/molxK | 650.64 | Joback Method |
| cpg | 459.00 | J/molxK | 684.18 | Joback Method |
| cpg | 472.18 | J/molxK | 717.72 | Joback Method |
| cpg | 484.61 | J/molxK | 751.26 | Joback Method |
| dvisc | 0.0019359 | Paxs | 293.40 | Joback Method |
| dvisc | 0.0009244 | Paxs | 336.17 | Joback Method |

| | | | | |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0005216 | Paxs | 378.94 | Joback Method |
| dvisc | 0.0003305 | Paxs | 421.71 | Joback Method |
| dvisc | 0.0002278 | Paxs | 464.48 | Joback Method |
| dvisc | 0.0001672 | Paxs | 507.25 | Joback Method |
| dvisc | 0.0001287 | Paxs | 550.02 | 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=U378190&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 |
| mcvol: | McGowan's characteristic volume |
| pc: | Critical Pressure |
| rinpol: | Non-polar retention indices |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

Latest version available from:

<https://www.chemeo.com/cid/65-242-1/2-Methoxybenzyl-alcohol-2-methylpropyl-ether.pdf>

Generated by Cheméo on 2024-04-27 19:27:50.32491864 +0000 UTC m=+16535319.245495961.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.