

Benzene, 1-methyl-4-(dimethoxymethyl)-

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|-----------------------------|--|
| Other names: | p-(dimethoxymethyl)toluene |
| Inchi: | InChI=1S/C10H14O2/c1-8-4-6-9(7-5-8)10(11-2)12-3/h4-7,10H,1-3H3 |
| InchiKey: | MKZGLAOLSFUPRT-UHFFFAOYSA-N |
| Formula: | C10H14O2 |
| SMILES: | COC(OC)c1ccc(C)cc1 |
| Mol. weight [g/mol]: | 166.22 |
| CAS: | 3395-83-3 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -76.34 | kJ/mol | Joback Method |
| hf | -294.39 | kJ/mol | Joback Method |
| hfus | 14.16 | kJ/mol | Joback Method |
| hvap | 45.22 | kJ/mol | Joback Method |
| log10ws | -2.30 | | Crippen Method |
| logp | 2.286 | | Crippen Method |
| mcvol | 139.740 | ml/mol | McGowan Method |
| pc | 2805.41 | kPa | Joback Method |
| tb | 504.26 | K | Joback Method |
| tc | 711.53 | K | Joback Method |
| tf | 270.86 | K | Joback Method |
| vc | 0.517 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 306.21 | J/molxK | 504.26 | Joback Method |
| cpg | 320.60 | J/molxK | 538.81 | Joback Method |
| cpg | 334.35 | J/molxK | 573.35 | Joback Method |
| cpg | 347.46 | J/molxK | 607.90 | Joback Method |
| cpg | 359.93 | J/molxK | 642.44 | Joback Method |
| cpg | 371.75 | J/molxK | 676.99 | Joback Method |
| cpg | 382.94 | J/molxK | 711.53 | Joback Method |
| dvisc | 0.0019953 | Paxs | 270.86 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0009879 | Paxs | 309.76 | Joback Method |
| dvisc | 0.0005722 | Paxs | 348.66 | Joback Method |
| dvisc | 0.0003698 | Paxs | 387.56 | Joback Method |
| dvisc | 0.0002588 | Paxs | 426.46 | Joback Method |
| dvisc | 0.0001923 | Paxs | 465.36 | Joback Method |
| dvisc | 0.0001495 | Paxs | 504.26 | 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=C3395833&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 |
| 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/85-652-4/Benzene-1-methyl-4-dimethoxymethyl.pdf>

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