

Dimethylaminoacetaldehyde dimethyl acetal

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|-----------------------------|--|
| Other names: | Ethanamine, 2,2-dimethoxy-N,N-dimethyl- 2,2-Dimethoxy-N,N-dimethylethanamine 2,2-dimethoxyethyl(dimethyl)amine |
| Inchi: | InChI=1S/C6H15NO2/c1-7(2)5-6(8-3)9-4/h6H,5H2,1-4H3 |
| InchiKey: | HUYAEQCJNXODLQ-UHFFFAOYSA-N |
| Formula: | C6H15NO2 |
| SMILES: | COC(CN(C)C)OC |
| Mol. weight [g/mol]: | 133.19 |
| CAS: | 38711-20-5 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | -102.02 | kJ/mol | Joback Method |
| hf | -369.36 | kJ/mol | Joback Method |
| hfus | 13.17 | kJ/mol | Joback Method |
| hvap | 35.42 | kJ/mol | Joback Method |
| log10ws | 0.31 | | Crippen Method |
| logp | 0.167 | | Crippen Method |
| mcvol | 117.120 | ml/mol | McGowan Method |
| pc | 3012.33 | kPa | Joback Method |
| tb | 393.52 | K | Joback Method |
| tc | 561.37 | K | Joback Method |
| tf | 219.31 | K | Joback Method |
| vc | 0.419 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 232.06 | J/molxK | 393.52 | Joback Method |
| cpg | 243.55 | J/molxK | 421.49 | Joback Method |
| cpg | 254.72 | J/molxK | 449.47 | Joback Method |
| cpg | 265.55 | J/molxK | 477.44 | Joback Method |
| cpg | 276.05 | J/molxK | 505.42 | Joback Method |
| cpg | 286.20 | J/molxK | 533.39 | Joback Method |

cpg

296.02

J/mol×K

561.37

Joback Method

Pressure Dependent Properties

| Property code | Value | Unit | Pressure [kPa] | Source |
|---------------|--------|------|----------------|--------------|
| tbrp | 353.70 | K | 14.90 | NIST Webbook |

Sources

| | |
|------------------------|---|
| 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=C38711205&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307i |

Legend

| | |
|-----------------|---|
| cpg: | Ideal gas heat capacity |
| 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 |
| tbrp: | Boiling point at reduced pressure |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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<https://www.chemeo.com/cid/98-471-1/Dimethylaminoacetaldehyde-dimethyl-acetal.pdf>

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