

N,N-Diethylpropionamide

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| Other names: | Diethylpropionamide N,N-Diethylpropanamide Propanamide, N,N-diethyl- Propionamide, N,N-diethyl- Diethylamide of propionic acid |
| Inchi: | InChI=1S/C7H15NO/c1-4-7(9)8(5-2)6-3/h4-6H2,1-3H3 |
| InchiKey: | YKOQQFDCCBKROY-UHFFFAOYSA-N |
| Formula: | C7H15NO |
| SMILES: | CCC(=O)N(CC)CC |
| Mol. weight [g/mol]: | 129.20 |
| CAS: | 1114-51-8 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | -10.08 | kJ/mol | Joback Method |
| hf | -232.86 | kJ/mol | Joback Method |
| hfus | 18.51 | kJ/mol | Joback Method |
| hvap | 39.97 | kJ/mol | Joback Method |
| log10ws | -1.10 | | Crippen Method |
| logp | 1.265 | | Crippen Method |
| mcvol | 121.040 | ml/mol | McGowan Method |
| pc | 2989.32 | kPa | Joback Method |
| tb | 425.87 | K | Joback Method |
| tc | 600.81 | K | Joback Method |
| tf | 251.05 | K | Joback Method |
| vc | 0.452 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 244.86 | J/molxK | 425.87 | Joback Method |
| cpg | 257.29 | J/molxK | 455.03 | Joback Method |
| cpg | 269.17 | J/molxK | 484.18 | Joback Method |
| cpg | 280.54 | J/molxK | 513.34 | Joback Method |

| | | | | |
|-----|--------|---------|--------|---------------|
| cpg | 291.41 | J/mol×K | 542.50 | Joback Method |
| cpg | 301.79 | J/mol×K | 571.65 | Joback Method |
| cpg | 311.69 | J/mol×K | 600.81 | Joback Method |

Pressure Dependent Properties

| Property code | Value | Unit | Pressure [kPa] | Source |
|---------------|--------|------|----------------|--------------|
| tbrp | 350.20 | K | 1.60 | 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=C1114518&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307l |

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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