

N,N-Dimethyldodecanamide

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| Other names: | Dodecanamide, N,N-dimethyl- Hallcomid M 12 Lauryl N,N-dimethylamide N,N-Dimethyldodecamide N,N-Dimethylauramide 1-Dodecanamide, N,N-dimethyl- |
| Inchi: | InChI=1S/C14H29NO/c1-4-5-6-7-8-9-10-11-12-13-14(16)15(2)3/h4-13H2,1-3H3 |
| InchiKey: | BDYUSDIJIDGWCY-UHFFFAOYSA-N |
| Formula: | C14H29NO |
| SMILES: | CCCCCCCCCCCC(=O)N(C)C |
| Mol. weight [g/mol]: | 227.39 |
| CAS: | 3007-53-2 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | 48.86 | kJ/mol | Joback Method |
| hf | -377.34 | kJ/mol | Joback Method |
| hfus | 36.64 | kJ/mol | Joback Method |
| hvap | 55.55 | kJ/mol | Joback Method |
| log10ws | -4.03 | | Crippen Method |
| logp | 3.996 | | Crippen Method |
| mvol | 219.670 | ml/mol | McGowan Method |
| pc | 1598.72 | kPa | Joback Method |
| rinpol | 1504.00 | | NIST Webbook |
| tb | 586.03 | K | Joback Method |
| tc | 752.99 | K | Joback Method |
| tf | 329.94 | K | Joback Method |
| vc | 0.844 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 576.06 | J/mol×K | 586.03 | Joback Method |
| cpg | 593.60 | J/mol×K | 613.86 | Joback Method |

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|-----|--------|---------|--------|---------------|
| cpg | 610.38 | J/mol×K | 641.68 | Joback Method |
| cpg | 626.41 | J/mol×K | 669.51 | Joback Method |
| cpg | 641.73 | J/mol×K | 697.34 | Joback Method |
| cpg | 656.35 | J/mol×K | 725.16 | Joback Method |
| cpg | 670.31 | J/mol×K | 752.99 | Joback Method |

Pressure Dependent Properties

| Property code | Value | Unit | Pressure [kPa] | Source |
|---------------|---------------|------|----------------|--------------|
| tbrp | 453.50 ± 0.50 | K | 2.00 | NIST Webbook |

Sources

| | |
|------------------------|---|
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=C3007532&Units=SI |
| 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 |

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 |
| rinpol: | Non-polar retention indices |
| 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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