

N,N,2,4-Tetramethylbenzenamine

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|-----------------------------|---|
| Other names: | N,N-Dimethyl-2,4-xylidine |
| Inchi: | InChI=1S/C10H15N/c1-8-5-6-10(11(3)4)9(2)7-8/h5-7H,1-4H3 |
| InchiKey: | DZXAIYQRCQALGE-UHFFFAOYSA-N |
| Formula: | C10H15N |
| SMILES: | <chem>Cc1ccc(N(C)C)c(C)c1</chem> |
| Mol. weight [g/mol]: | 149.23 |
| CAS: | 769-53-9 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | 237.25 | kJ/mol | Joback Method |
| hf | 31.39 | kJ/mol | Joback Method |
| hfus | 17.94 | kJ/mol | Joback Method |
| hvap | 43.50 | kJ/mol | Joback Method |
| ie | 7.17 | eV | NIST Webbook |
| ie | 7.79 | eV | NIST Webbook |
| log10ws | -2.33 | | Crippen Method |
| logp | 2.369 | | Crippen Method |
| mcvol | 137.980 | ml/mol | McGowan Method |
| pc | 2832.35 | kPa | Joback Method |
| rinpol | 1149.60 | | NIST Webbook |
| rinpol | 1149.60 | | NIST Webbook |
| ripol | 1798.10 | | NIST Webbook |
| tb | 477.28 | K | Joback Method |
| tc | 681.78 | K | Joback Method |
| tf | 286.39 | K | Joback Method |
| vc | 0.505 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 290.18 | J/mol×K | 477.28 | Joback Method |
| cpg | 305.33 | J/mol×K | 511.36 | Joback Method |
| cpg | 319.69 | J/mol×K | 545.45 | Joback Method |

| | | | | |
|-----|--------|---------|--------|---------------|
| cpg | 333.27 | J/mol×K | 579.53 | Joback Method |
| cpg | 346.11 | J/mol×K | 613.61 | Joback Method |
| cpg | 358.24 | J/mol×K | 647.70 | Joback Method |
| cpg | 369.68 | J/mol×K | 681.78 | Joback Method |

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=C769539&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci9903071 |

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 |
| ie: | Ionization energy |
| log10ws: | Log10 of Water solubility in mol/l |
| logp: | Octanol/Water partition coefficient |
| mcvol: | McGowan's characteristic volume |
| pc: | Critical Pressure |
| rinpola: | Non-polar retention indices |
| ripola: | Polar retention indices |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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