

Benzenamine, 4-cyclohexyl-

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|-----------------------------|---|
| Other names: | Aniline, p-cyclohexyl- p-Cyclohexylaniline 4-Cyclohexylaniline |
| Inchi: | InChI=1S/C12H17N/c13-12-8-6-11(7-9-12)10-4-2-1-3-5-10/h6-10H,1-5,13H2 |
| InchiKey: | JLNMBIKJQAKQBH-UHFFFAOYSA-N |
| Formula: | C12H17N |
| SMILES: | <chem>Nc1ccc(C2CCCCC2)cc1</chem> |
| Mol. weight [g/mol]: | 175.27 |
| CAS: | 6373-50-8 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | 243.84 | kJ/mol | Joback Method |
| hf | 22.16 | kJ/mol | Joback Method |
| hfus | 17.52 | kJ/mol | Joback Method |
| hvap | 56.31 | kJ/mol | Joback Method |
| log10ws | -3.44 | | Crippen Method |
| logp | 3.316 | | Crippen Method |
| mcvol | 155.300 | ml/mol | McGowan Method |
| pc | 3114.04 | kPa | Joback Method |
| tb | 597.70 | K | Joback Method |
| tc | 847.98 | K | Joback Method |
| tf | 354.58 | K | Joback Method |
| vc | 0.561 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 400.45 | J/molxK | 597.70 | Joback Method |
| cpg | 420.35 | J/molxK | 639.41 | Joback Method |
| cpg | 438.78 | J/molxK | 681.13 | Joback Method |
| cpg | 455.79 | J/molxK | 722.84 | Joback Method |
| cpg | 471.46 | J/molxK | 764.56 | Joback Method |
| cpg | 485.86 | J/molxK | 806.27 | Joback Method |

cpg

499.04

J/mol×K

847.98

Joback Method

Pressure Dependent Properties

| Property code | Value | Unit | Pressure [kPa] | Source |
|---------------|--------|------|----------------|--------------|
| tbrp | 439.20 | K | 1.70 | NIST Webbook |

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=C6373508&Units=SI |

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/75-455-4/Benzenamine-4-cyclohexyl.pdf>

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