

Hydrazine, (1-methylethyl)-

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|----------------------|---|
| Inchi: | InChI=1S/C3H10N2/c1-3(2)5-4/h3,5H,4H2,1-2H3 |
| InchiKey: | KJAQRHMKLVGSCG-UHFFFAOYSA-N |
| Formula: | C3H10N2 |
| SMILES: | CC(C)NN |
| Mol. weight [g/mol]: | 74.12 |
| CAS: | 2257-52-5 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|-------------|---------|----------------|
| gf | 127.78 | kJ/mol | Joback Method |
| hf | -23.27 | kJ/mol | Joback Method |
| hfus | 10.30 | kJ/mol | Joback Method |
| hvap | 38.96 | kJ/mol | Joback Method |
| ie | 8.42 ± 0.05 | eV | NIST Webbook |
| ie | 9.05 | eV | NIST Webbook |
| log10ws | -0.80 | | Crippen Method |
| logp | -0.142 | | Crippen Method |
| mcvol | 73.090 | ml/mol | McGowan Method |
| pc | 4903.92 | kPa | Joback Method |
| tb | 390.30 | K | Joback Method |
| tc | 583.17 | K | Joback Method |
| tf | 244.49 | K | Joback Method |
| vc | 0.262 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 139.67 | J/molxK | 390.30 | Joback Method |
| cpg | 148.17 | J/molxK | 422.44 | Joback Method |
| cpg | 156.32 | J/molxK | 454.59 | Joback Method |
| cpg | 164.11 | J/molxK | 486.73 | Joback Method |
| cpg | 171.55 | J/molxK | 518.88 | Joback Method |
| cpg | 178.65 | J/molxK | 551.02 | Joback Method |
| cpg | 185.42 | J/molxK | 583.17 | Joback Method |

Sources

| | |
|------------------------|---|
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=C2257525&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 |
| hvp: | Enthalpy of vaporization at standard conditions |
| ie: | Ionization energy |
| log10ws: | Log10 of Water solubility in mol/l |
| logp: | Octanol/Water partition coefficient |
| mvol: | McGowan's characteristic volume |
| pc: | Critical Pressure |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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