

6-Chloroanthranilonitrile

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| Other names: | 2-Amino-6-chlorobenzonitrile Benzonitrile, 2-amino-6-chloro- 6-Amino-2-chlorobenzonitrile |
| Inchi: | InChI=1S/C7H5ClN2/c8-6-2-1-3-7(10)5(6)4-9/h1-3H,10H2 |
| InchiKey: | MEJVTQKBWPYBFG-UHFFFAOYSA-N |
| Formula: | C7H5ClN2 |
| SMILES: | N#Cc1c(N)cccc1Cl |
| Mol. weight [g/mol]: | 152.58 |
| CAS: | 6575-11-7 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | 288.91 | kJ/mol | Joback Method |
| hf | 208.71 | kJ/mol | Joback Method |
| hfus | 18.05 | kJ/mol | Joback Method |
| hvap | 60.28 | kJ/mol | Joback Method |
| log10ws | -2.14 | | Crippen Method |
| logp | 1.794 | | Crippen Method |
| mcvol | 109.330 | ml/mol | McGowan Method |
| pc | 3935.71 | kPa | Joback Method |
| tb | 608.24 | K | Joback Method |
| tc | 859.69 | K | Joback Method |
| tf | 398.28 | K | Joback Method |
| vc | 0.423 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 223.84 | J/molxK | 608.24 | Joback Method |
| cpg | 231.65 | J/molxK | 650.15 | Joback Method |
| cpg | 238.88 | J/molxK | 692.06 | Joback Method |
| cpg | 245.54 | J/molxK | 733.97 | Joback Method |
| cpg | 251.68 | J/molxK | 775.87 | Joback Method |
| cpg | 257.31 | J/molxK | 817.78 | Joback Method |

Sources

| | |
|------------------------|---|
| 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=C6575117&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci9903071 |
| Crippen Method: | https://www.chemeo.com/doc/models/crippen_log10ws |

Legend

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|-----------------|---|
| 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 |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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