

Aniline, 2-chloro-n-isopropyl-

| | |
|-----------------------------|---|
| Inchi: | InChI=1S/C9H12ClN/c1-7(2)11-9-6-4-3-5-8(9)10/h3-7,11H,1-2H3 |
| InchiKey: | JCRMYLUIMGTZSR-UHFFFAOYSA-N |
| Formula: | C9H12ClN |
| SMILES: | CC(C)Nc1ccccc1Cl |
| Mol. weight [g/mol]: | 169.65 |
| CAS: | 78235-07-1 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | 202.70 | kJ/mol | Joback Method |
| hf | 28.42 | kJ/mol | Joback Method |
| hfus | 18.49 | kJ/mol | Joback Method |
| hvap | 49.00 | kJ/mol | Joback Method |
| log10ws | -3.11 | | Crippen Method |
| logp | 3.160 | | Crippen Method |
| mcvol | 136.130 | ml/mol | McGowan Method |
| pc | 3149.09 | kPa | Joback Method |
| tb | 524.14 | K | Joback Method |
| tc | 744.97 | K | Joback Method |
| tf | 297.71 | K | Joback Method |
| vc | 0.509 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 287.98 | J/mol×K | 524.14 | Joback Method |
| cpg | 301.47 | J/mol×K | 560.94 | Joback Method |
| cpg | 314.13 | J/mol×K | 597.75 | Joback Method |
| cpg | 325.98 | J/mol×K | 634.55 | Joback Method |
| cpg | 337.07 | J/mol×K | 671.36 | Joback Method |
| cpg | 347.43 | J/mol×K | 708.16 | Joback Method |
| cpg | 357.10 | J/mol×K | 744.97 | Joback Method |

Sources

| | |
|------------------------|---|
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=C78235071&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 |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

Latest version available from:

<https://www.chemeo.com/cid/49-054-8/Aniline-2-chloro-n-isopropyl.pdf>

Generated by Cheméo on 2024-04-24 20:21:01.129075629 +0000 UTC m=+16279310.049652944.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.