

N-(5-Chloro-2-methylphenyl)-N-(trifluoroacetyl)-2,

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|-----------------------------|--|
| Other names: | 5-Chloro-2-methyl-aniline, N,N-bis-trifluoroacetyl- |
| Inchi: | InChI=1S/C11H6ClF6NO2/c1-5-2-3-6(12)4-7(5)19(8(20)10(13,14)15)9(21)11(16,17)18/h |
| InchiKey: | QKMNNYAERWCXHO-UHFFFAOYSA-N |
| Formula: | C11H6ClF6NO2 |
| SMILES: | <chem>Cc1ccc(Cl)cc1N(C(=O)C(F)(F)F)C(=O)C(F)(F)F</chem> |
| Mol. weight [g/mol]: | 333.61 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|----------|----------------------|----------------|
| gf | -1187.28 | kJ/mol | Joback Method |
| hf | -1424.31 | kJ/mol | Joback Method |
| hfus | 31.58 | kJ/mol | Joback Method |
| hvap | 56.11 | kJ/mol | Joback Method |
| log10ws | -4.25 | | Crippen Method |
| logp | 3.633 | | Crippen Method |
| mcvol | 178.070 | ml/mol | McGowan Method |
| pc | 2229.20 | kPa | Joback Method |
| rinpol | 1179.00 | | NIST Webbook |
| rinpol | 1179.00 | | NIST Webbook |
| tb | 634.49 | K | Joback Method |
| tc | 824.55 | K | Joback Method |
| tf | 435.82 | K | Joback Method |
| vc | 0.709 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 454.54 | J/mol×K | 634.49 | Joback Method |
| cpg | 464.60 | J/mol×K | 666.17 | Joback Method |
| cpg | 473.83 | J/mol×K | 697.84 | Joback Method |
| cpg | 482.31 | J/mol×K | 729.52 | Joback Method |
| cpg | 490.08 | J/mol×K | 761.20 | Joback Method |
| cpg | 497.20 | J/mol×K | 792.87 | Joback Method |
| cpg | 503.74 | J/mol×K | 824.55 | Joback Method |

Sources

| | |
|------------------------|---|
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U373427&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 |
| hvpap: | 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 |
| rinppl: | Non-polar retention indices |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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