

# Chlorodifluoroamine

|                      |                                                                                                                                           |
|----------------------|-------------------------------------------------------------------------------------------------------------------------------------------|
| Other names:         | Nitrogen chloride fluoride (nclf2)<br>Difluorochloroamine<br>Chlorodifluoroamine<br>Chlorodifluoroammonia<br>Nitrogen chloride difluoride |
| Inchi:               | InChI=1S/ClF2N/c1-4(2)3                                                                                                                   |
| InchiKey:            | ZIOUHCMXEAFYSA-UHFFFAOYSA-N                                                                                                               |
| Formula:             | ClF2N                                                                                                                                     |
| SMILES:              | FN(F)Cl                                                                                                                                   |
| Mol. weight [g/mol]: | 87.46                                                                                                                                     |
| CAS:                 | 13637-87-1                                                                                                                                |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -341.65 | kJ/mol  | Joback Method  |
| hf            | -383.76 | kJ/mol  | Joback Method  |
| hfus          | 9.13    | kJ/mol  | Joback Method  |
| hvap          | 20.39   | kJ/mol  | Joback Method  |
| log10ws       | -1.22   |         | Crippen Method |
| logp          | 1.211   |         | Crippen Method |
| mcvol         | 36.620  | ml/mol  | McGowan Method |
| pc            | 5511.44 | kPa     | Joback Method  |
| tb            | 247.81  | K       | Joback Method  |
| tc            | 393.72  | K       | Joback Method  |
| tf            | 153.33  | K       | Joback Method  |
| vc            | 0.139   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value | Unit    | Temperature [K] | Source        |
|---------------|-------|---------|-----------------|---------------|
| cpg           | 48.43 | J/mol×K | 247.81          | Joback Method |
| cpg           | 50.71 | J/mol×K | 272.13          | Joback Method |
| cpg           | 52.87 | J/mol×K | 296.45          | Joback Method |
| cpg           | 54.90 | J/mol×K | 320.77          | Joback Method |

|     |       |         |        |               |
|-----|-------|---------|--------|---------------|
| cpg | 56.82 | J/mol×K | 345.08 | Joback Method |
| cpg | 58.63 | J/mol×K | 369.40 | Joback Method |
| cpg | 60.34 | J/mol×K | 393.72 | Joback Method |

## Sources

|                        |                                                                                                                                               |
|------------------------|-----------------------------------------------------------------------------------------------------------------------------------------------|
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C13637871&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C13637871&amp;Units=SI</a> |
| <b>Crippen Method:</b> | <a href="http://pubs.acs.org/doi/abs/10.1021/ci990307l">http://pubs.acs.org/doi/abs/10.1021/ci990307l</a>                                     |
| <b>Crippen Method:</b> | <a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>                             |
| <b>Joback Method:</b>  | <a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>                                         |
| <b>McGowan Method:</b> | <a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>                         |

## Legend

|                 |                                                 |
|-----------------|-------------------------------------------------|
| <b>cpg:</b>     | Ideal gas heat capacity                         |
| <b>gf:</b>      | Standard Gibbs free energy of formation         |
| <b>hf:</b>      | Enthalpy of formation at standard conditions    |
| <b>hfus:</b>    | Enthalpy of fusion at standard conditions       |
| <b>hvap:</b>    | Enthalpy of vaporization at standard conditions |
| <b>log10ws:</b> | Log10 of Water solubility in mol/l              |
| <b>logp:</b>    | Octanol/Water partition coefficient             |
| <b>mcvol:</b>   | McGowan's characteristic volume                 |
| <b>pc:</b>      | Critical Pressure                               |
| <b>tb:</b>      | Normal Boiling Point Temperature                |
| <b>tc:</b>      | Critical Temperature                            |
| <b>tf:</b>      | Normal melting (fusion) point                   |
| <b>vc:</b>      | Critical Volume                                 |

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