

# Benzene, 4-chloro-3-(chloromethyl)-1-methyl

|                             |  |
|-----------------------------|--|
| <b>Inchi:</b>               | InChI=1S/C8H8Cl2/c1-6-2-3-8(10)7(4-6)5-9/h2-4H,5H2,1H3 |
| <b>InchiKey:</b>            | PYEZGJNZDIRIIM-UHFFFAOYSA-N                            |
| <b>Formula:</b>             | C8H8Cl2  |
| <b>SMILES:</b>              | Cc1ccc(Cl)c(CCl)c1                                     |
| <b>Mol. weight [g/mol]:</b> | 175.06   |

## Physical Properties

| Property code | Value   | Unit                 | Source         |
|---------------|---------|----------------------|----------------|
| gf            | 85.77   | kJ/mol               | Joback Method  |
| hf            | -26.34  | kJ/mol               | Joback Method  |
| hfus          | 18.13   | kJ/mol               | Joback Method  |
| hvap          | 45.77   | kJ/mol               | Joback Method  |
| log10ws       | -3.66   |                      | Crippen Method |
| logp          | 3.387   |                      | Crippen Method |
| mcvol         | 124.300 | ml/mol               | McGowan Method |
| pc            | 3231.98 | kPa                  | Joback Method  |
| rinpola       | 1238.00 |                      | NIST Webbook   |
| tb            | 493.94  | K                    | Joback Method  |
| tc            | 719.52  | K                    | Joback Method  |
| tf            | 291.22  | K                    | Joback Method  |
| vc            | 0.473   | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 226.48    | J/molxK | 493.94          | Joback Method |
| cpg           | 237.10    | J/molxK | 531.54          | Joback Method |
| cpg           | 247.09    | J/molxK | 569.13          | Joback Method |
| cpg           | 256.47    | J/molxK | 606.73          | Joback Method |
| cpg           | 265.25    | J/molxK | 644.33          | Joback Method |
| cpg           | 273.48    | J/molxK | 681.92          | Joback Method |
| cpg           | 281.16    | J/molxK | 719.52          | Joback Method |
| dvisc         | 0.0017277 | Paxs    | 291.22          | Joback Method |
| dvisc         | 0.0010719 | Paxs    | 325.01          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0007276 | Paxs | 358.79 | Joback Method |
| dvisc | 0.0005279 | Paxs | 392.58 | Joback Method |
| dvisc | 0.0004030 | Paxs | 426.37 | Joback Method |
| dvisc | 0.0003201 | Paxs | 460.15 | Joback Method |
| dvisc | 0.0002624 | Paxs | 493.94 | Joback Method |

## Sources

|                        |   |
|------------------------|---|
| <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>                     |
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=R132188&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R132188&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>                                 |

## Legend

|                 |   |
|-----------------|---|
| <b>cpg:</b>     | Ideal gas heat capacity                         |
| <b>dvisc:</b>   | Dynamic viscosity                               |
| <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>rinpol:</b>  | Non-polar retention indices                     |
| <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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