

# 3,5-Dimethoxybenzoyl chloride

|                             |  |
|-----------------------------|--|
| <b>Other names:</b>         | Benzoyl chloride, 3,5-dimethoxy-                               |
| <b>Inchi:</b>               | InChI=1S/C9H9ClO3/c1-12-7-3-6(9(10)11)4-8(5-7)13-2/h3-5H,1-2H3 |
| <b>InchiKey:</b>            | FTHPLWDYWAKYCY-UHFFFAOYSA-N                                    |
| <b>Formula:</b>             | C9H9ClO3   |
| <b>SMILES:</b>              | COc1cc(OC)cc(C(=O)Cl)c1  |
| <b>Mol. weight [g/mol]:</b> | 200.62   |
| <b>CAS:</b>                 | 17213-57-9   |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -232.80 | kJ/mol  | Joback Method  |
| hf            | -408.26 | kJ/mol  | Joback Method  |
| hfus          | 20.50   | kJ/mol  | Joback Method  |
| hvap          | 55.18   | kJ/mol  | Joback Method  |
| log10ws       | -2.60   |         | Crippen Method |
| logp          | 2.083   |         | Crippen Method |
| mcvol         | 139.460 | ml/mol  | McGowan Method |
| pc            | 3124.49 | kPa     | Joback Method  |
| tb            | 578.10  | K       | Joback Method  |
| tc            | 798.52  | K       | Joback Method  |
| tf            | 366.96  | K       | Joback Method  |
| vc            | 0.522   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 302.40    | J/molxK | 578.10          | Joback Method |
| cpg           | 313.52    | J/molxK | 614.84          | Joback Method |
| cpg           | 324.06    | J/molxK | 651.57          | Joback Method |
| cpg           | 334.01    | J/molxK | 688.31          | Joback Method |
| cpg           | 343.37    | J/molxK | 725.05          | Joback Method |
| cpg           | 352.12    | J/molxK | 761.78          | Joback Method |
| cpg           | 360.25    | J/molxK | 798.52          | Joback Method |
| dvisc         | 0.0010339 | Paxs    | 366.96          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0006811 | Paxs | 402.15 | Joback Method |
| dvisc | 0.0004798 | Paxs | 437.34 | Joback Method |
| dvisc | 0.0003562 | Paxs | 472.53 | Joback Method |
| dvisc | 0.0002755 | Paxs | 507.72 | Joback Method |
| dvisc | 0.0002203 | Paxs | 542.91 | Joback Method |
| dvisc | 0.0001811 | Paxs | 578.10 | Joback Method |

## Pressure Dependent Properties

| Property code | Value  | Unit | Pressure [kPa] | Source       |
|---------------|--------|------|----------------|--------------|
| tbrp          | 430.70 | K    | 2.10           | NIST Webbook |

## 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=C17213579&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C17213579&amp;Units=SI</a> |
| <b>Crippen Method:</b> | <a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</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>tb:</b>      | Normal Boiling Point Temperature                |
| <b>tbrp:</b>    | Boiling point at reduced pressure               |
| <b>tc:</b>      | Critical Temperature                            |

**tf:** Normal melting (fusion) point

**vc:** Critical Volume

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