

# 2-Chlorobenzoic acid, heptadecyl ester

|                             |  |
|-----------------------------|--|
| <b>Inchi:</b>               | InChI=1S/C24H39ClO2/c1-2-3-4-5-6-7-8-9-10-11-12-13-14-15-18-21-27-24(26)22-19-16 |
| <b>InchiKey:</b>            | BODQEIRZCIRWTL-UHFFFAOYSA-N  |
| <b>Formula:</b>             | C24H39ClO2   |
| <b>SMILES:</b>              | CCCCCCCCCCCCCCCCOC(=O)c1ccccc1Cl   |
| <b>Mol. weight [g/mol]:</b> | 395.02   |

## Physical Properties

| Property code | Value   | Unit                 | Source         |
|---------------|---------|----------------------|----------------|
| gf            | 8.13    | kJ/mol               | Joback Method  |
| hf            | -574.17 | kJ/mol               | Joback Method  |
| hfus          | 58.55   | kJ/mol               | Joback Method  |
| hvap          | 85.50   | kJ/mol               | Joback Method  |
| log10ws       | -9.10   |                      | Crippen Method |
| logp          | 8.368   |                      | Crippen Method |
| mvol          | 344.940 | ml/mol               | McGowan Method |
| pc            | 978.40  | kPa                  | Joback Method  |
| rinpol        | 2674.30 |                      | NIST Webbook   |
| rinpol        | 2674.30 |                      | NIST Webbook   |
| tb            | 893.90  | K                    | Joback Method  |
| tc            | 1096.46 | K                    | Joback Method  |
| tf            | 501.26  | K                    | Joback Method  |
| vc            | 1.345   | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 1098.56   | J/molxK | 893.90          | Joback Method |
| cpg           | 1116.74   | J/molxK | 927.66          | Joback Method |
| cpg           | 1133.72   | J/molxK | 961.42          | Joback Method |
| cpg           | 1149.56   | J/molxK | 995.18          | Joback Method |
| cpg           | 1164.31   | J/molxK | 1028.94         | Joback Method |
| cpg           | 1178.02   | J/molxK | 1062.70         | Joback Method |
| cpg           | 1190.73   | J/molxK | 1096.46         | Joback Method |
| dvisc         | 0.0005578 | Paxs    | 501.26          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0002736 | Paxs | 566.70 | Joback Method |
| dvisc | 0.0001555 | Paxs | 632.14 | Joback Method |
| dvisc | 0.0000983 | Paxs | 697.58 | Joback Method |
| dvisc | 0.0000672 | Paxs | 763.02 | Joback Method |
| dvisc | 0.0000488 | Paxs | 828.46 | Joback Method |
| dvisc | 0.0000371 | Paxs | 893.90 | Joback Method |

## Sources

|                        |   |
|------------------------|---|
| <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>                                 |
| <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=U292283&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U292283&amp;Units=SI</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                                 |

Latest version available from:

<https://www.chemeo.com/cid/95-610-9/2-Chlorobenzoic-acid-heptadecyl-ester.pdf>

Generated by Cheméo on 2024-04-25 16:17:20.323801882 +0000 UTC m=+16351089.244379192.

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