

# Fumaric acid, 2-chloropropyl decyl ester

|                             |                                                                                  |
|-----------------------------|----------------------------------------------------------------------------------|
| <b>Inchi:</b>               | InChI=1S/C17H29ClO4/c1-3-4-5-6-7-8-9-10-13-21-16(19)11-12-17(20)22-14-15(2)18/h1 |
| <b>InchiKey:</b>            | JKNGOSKOOJMAKL-VAWYXSNFSA-N                                                      |
| <b>Formula:</b>             | C17H29ClO4                                                                       |
| <b>SMILES:</b>              | CCCCCCCCCOC(=O)C=CC(=O)OCC(C)Cl                                                  |
| <b>Mol. weight [g/mol]:</b> | 332.86                                                                           |

## Physical Properties

| Property code | Value   | Unit                 | Source         |
|---------------|---------|----------------------|----------------|
| gf            | -309.73 | kJ/mol               | Joback Method  |
| hf            | -787.61 | kJ/mol               | Joback Method  |
| hfus          | 46.24   | kJ/mol               | Joback Method  |
| hvap          | 75.70   | kJ/mol               | Joback Method  |
| log10ws       | -4.78   |                      | Crippen Method |
| logp          | 4.397   |                      | Crippen Method |
| mvol          | 273.210 | ml/mol               | McGowan Method |
| pc            | 1336.86 | kPa                  | Joback Method  |
| rinpol        | 2272.00 |                      | NIST Webbook   |
| rinpol        | 2272.00 |                      | NIST Webbook   |
| tb            | 782.09  | K                    | Joback Method  |
| tc            | 970.10  | K                    | Joback Method  |
| tf            | 435.51  | K                    | Joback Method  |
| vc            | 1.058   | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 798.12    | J/molxK | 782.09          | Joback Method |
| cpg           | 867.99    | J/molxK | 938.76          | Joback Method |
| cpg           | 855.73    | J/molxK | 907.43          | Joback Method |
| cpg           | 842.63    | J/molxK | 876.09          | Joback Method |
| cpg           | 828.68    | J/molxK | 844.76          | Joback Method |
| cpg           | 813.85    | J/molxK | 813.42          | Joback Method |
| cpg           | 879.45    | J/molxK | 970.10          | Joback Method |
| dvisc         | 0.0000564 | Paxs    | 782.09          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0000753 | Paxs | 724.33 | Joback Method |
| dvisc | 0.0001057 | Paxs | 666.56 | Joback Method |
| dvisc | 0.0001582 | Paxs | 608.80 | Joback Method |
| dvisc | 0.0002577 | Paxs | 551.04 | Joback Method |
| dvisc | 0.0004705 | Paxs | 493.27 | Joback Method |
| dvisc | 0.0010078 | Paxs | 435.51 | 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=U348568&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U348568&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>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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