

# Fumaric acid, ethyl 2,3,5-trichlorophenyl ester

|                             |  |
|-----------------------------|--|
| <b>Inchi:</b>               | InChI=1S/C12H9Cl3O4/c1-2-18-10(16)3-4-11(17)19-9-6-7(13)5-8(14)12(9)15/h3-6H,2H2 |
| <b>InchiKey:</b>            | RWERCTSWXBEGQR-ONEGZZNKSA-N  |
| <b>Formula:</b>             | C12H9Cl3O4   |
| <b>SMILES:</b>              | CCOC(=O)C=CC(=O)Oc1cc(Cl)cc(Cl)c1Cl  |
| <b>Mol. weight [g/mol]:</b> | 323.56   |

## Physical Properties

| Property code | Value   | Unit                 | Source         |
|---------------|---------|----------------------|----------------|
| gf            | -289.73 | kJ/mol               | Joback Method  |
| hf            | -508.49 | kJ/mol               | Joback Method  |
| hfus          | 38.08   | kJ/mol               | Joback Method  |
| hvap          | 77.99   | kJ/mol               | Joback Method  |
| log10ws       | -4.23   |                      | Crippen Method |
| logp          | 3.671   |                      | Crippen Method |
| mcvol         | 203.480 | ml/mol               | McGowan Method |
| pc            | 2377.22 | kPa                  | Joback Method  |
| rinsol        | 2114.00 |                      | NIST Webbook   |
| tb            | 784.61  | K                    | Joback Method  |
| tc            | 1014.85 | K                    | Joback Method  |
| tf            | 517.98  | K                    | Joback Method  |
| vc            | 0.774   | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 477.14    | J/molxK | 784.61          | Joback Method |
| cpg           | 516.55    | J/molxK | 976.48          | Joback Method |
| cpg           | 510.23    | J/molxK | 938.11          | Joback Method |
| cpg           | 503.14    | J/molxK | 899.73          | Joback Method |
| cpg           | 495.27    | J/molxK | 861.36          | Joback Method |
| cpg           | 486.61    | J/molxK | 822.98          | Joback Method |
| cpg           | 522.10    | J/molxK | 1014.85         | Joback Method |
| dvisc         | 0.0000962 | Paxs    | 784.61          | Joback Method |
| dvisc         | 0.0001169 | Paxs    | 740.17          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0001457 | Paxs | 695.73 | Joback Method |
| dvisc | 0.0001870 | Paxs | 651.30 | Joback Method |
| dvisc | 0.0002490 | Paxs | 606.86 | Joback Method |
| dvisc | 0.0003469 | Paxs | 562.42 | Joback Method |
| dvisc | 0.0005116 | Paxs | 517.98 | 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=U348139&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U348139&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                                 |

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