

# Diethylmalonic acid, butyl 2-chlorophenyl ester

|                      |   |
|----------------------|---|
| Inchi:               | InChI=1S/C17H23ClO4/c1-4-7-12-21-15(19)17(5-2,6-3)16(20)22-14-11-9-8-10-13(14)18/ |
| InchiKey:            | XNHPRBYQSCHXED-UHFFFAOYSA-N   |
| Formula:             | C17H23ClO4  |
| SMILES:              | CCCCOC(=O)C(CC)(CC)C(=O)Oc1ccccc1Cl   |
| Mol. weight [g/mol]: | 326.81  |

## Physical Properties

| Property code | Value   | Unit                 | Source         |
|---------------|---------|----------------------|----------------|
| gf            | -281.89 | kJ/mol               | Joback Method  |
| hf            | -683.24 | kJ/mol               | Joback Method  |
| hfus          | 35.79   | kJ/mol               | Joback Method  |
| hvap          | 77.78   | kJ/mol               | Joback Method  |
| log10ws       | -4.86   |                      | Crippen Method |
| logp          | 4.395   |                      | Crippen Method |
| mcvol         | 253.750 | ml/mol               | McGowan Method |
| pc            | 1652.46 | kPa                  | Joback Method  |
| rinsol        | 2064.00 |                      | NIST Webbook   |
| tb            | 806.80  | K                    | Joback Method  |
| tc            | 1018.69 | K                    | Joback Method  |
| tf            | 496.95  | K                    | Joback Method  |
| vc            | 0.966   | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 732.92    | J/molxK | 806.80          | Joback Method |
| cpg           | 747.41    | J/molxK | 842.11          | Joback Method |
| cpg           | 760.83    | J/molxK | 877.43          | Joback Method |
| cpg           | 773.21    | J/molxK | 912.74          | Joback Method |
| cpg           | 784.59    | J/molxK | 948.06          | Joback Method |
| cpg           | 795.00    | J/molxK | 983.37          | Joback Method |
| cpg           | 804.50    | J/molxK | 1018.69         | Joback Method |
| dvisc         | 0.0006071 | Paxs    | 496.95          | Joback Method |
| dvisc         | 0.0003412 | Paxs    | 548.59          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0002118 | Paxs | 600.23 | Joback Method |
| dvisc | 0.0001417 | Paxs | 651.88 | Joback Method |
| dvisc | 0.0001006 | Paxs | 703.52 | Joback Method |
| dvisc | 0.0000749 | Paxs | 755.16 | Joback Method |
| dvisc | 0.0000579 | Paxs | 806.80 | Joback Method |

## Sources

|                        |   |
|------------------------|---|
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=U369615&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U369615&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>                                 |
| <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>                     |

## 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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