

# Dimethylmalonic acid, butyl isohexyl ester

|                             |  |
|-----------------------------|--|
| <b>Inchi:</b>               | InChI=1S/C15H28O4/c1-6-7-10-18-13(16)15(4,5)14(17)19-11-8-9-12(2)3/h12H,6-11H2,1 |
| <b>InchiKey:</b>            | WAGBEBUVCKCTQK-UHFFFAOYSA-N  |
| <b>Formula:</b>             | C15H28O4   |
| <b>SMILES:</b>              | CCCCOC(=O)C(C)(C)C(=O)OCCCC(C)C  |
| <b>Mol. weight [g/mol]:</b> | 272.38   |

## Physical Properties

| Property code | Value   | Unit                 | Source         |
|---------------|---------|----------------------|----------------|
| gf            | -392.02 | kJ/mol               | Joback Method  |
| hf            | -856.56 | kJ/mol               | Joback Method  |
| hfus          | 29.24   | kJ/mol               | Joback Method  |
| hvap          | 65.61   | kJ/mol               | Joback Method  |
| log10ws       | -3.34   |                      | Crippen Method |
| logp          | 3.335   |                      | Crippen Method |
| mcvol         | 237.090 | ml/mol               | McGowan Method |
| pc            | 1551.22 | kPa                  | Joback Method  |
| rinpol        | 1608.00 |                      | NIST Webbook   |
| tb            | 691.51  | K                    | Joback Method  |
| tc            | 876.12  | K                    | Joback Method  |
| tf            | 390.55  | K                    | Joback Method  |
| vc            | 0.906   | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 680.11    | J/molxK | 691.51          | Joback Method |
| cpg           | 696.84    | J/molxK | 722.28          | Joback Method |
| cpg           | 712.70    | J/molxK | 753.05          | Joback Method |
| cpg           | 727.68    | J/molxK | 783.82          | Joback Method |
| cpg           | 741.83    | J/molxK | 814.59          | Joback Method |
| cpg           | 755.14    | J/molxK | 845.35          | Joback Method |
| cpg           | 767.65    | J/molxK | 876.12          | Joback Method |
| dvisc         | 0.0017437 | Paxs    | 390.55          | Joback Method |
| dvisc         | 0.0007852 | Paxs    | 440.71          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0004162 | Paxs | 490.87 | Joback Method |
| dvisc | 0.0002481 | Paxs | 541.03 | Joback Method |
| dvisc | 0.0001615 | Paxs | 591.19 | Joback Method |
| dvisc | 0.0001124 | Paxs | 641.35 | Joback Method |
| dvisc | 0.0000825 | Paxs | 691.51 | Joback Method |

## Sources

|                        |   |
|------------------------|---|
| <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=U361716&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U361716&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>                                     |

## 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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