

# Dimethylmalonic acid, butyl tetradecyl ester

|                             |  |
|-----------------------------|--|
| <b>Inchi:</b>               | InChI=1S/C23H44O4/c1-5-7-9-10-11-12-13-14-15-16-17-18-20-27-22(25)23(3,4)21(24)2 |
| <b>InchiKey:</b>            | VKJXLZZWCYMFDS-UHFFFAOYSA-N  |
| <b>Formula:</b>             | C23H44O4   |
| <b>SMILES:</b>              | CCCCCCCCCCCCCOC(=O)C(C)(C)C(=O)OCCCC   |
| <b>Mol. weight [g/mol]:</b> | 384.59   |

## Physical Properties

| Property code | Value    | Unit    | Source         |
|---------------|----------|---------|----------------|
| gf            | -322.22  | kJ/mol  | Joback Method  |
| hf            | -1016.40 | kJ/mol  | Joback Method  |
| hfus          | 53.49    | kJ/mol  | Joback Method  |
| hvap          | 83.81    | kJ/mol  | Joback Method  |
| log10ws       | -6.93    |         | Crippen Method |
| logp          | 6.600    |         | Crippen Method |
| mcvol         | 349.810  | ml/mol  | McGowan Method |
| pc            | 909.98   | kPa     | Joback Method  |
| rinpol        | 2434.00  |         | NIST Webbook   |
| rinpol        | 2434.00  |         | NIST Webbook   |
| tb            | 874.99   | K       | Joback Method  |
| tc            | 1071.31  | K       | Joback Method  |
| tf            | 495.71   | K       | Joback Method  |
| vc            | 1.361    | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 1154.79   | J/molxK | 874.99          | Joback Method |
| cpg           | 1239.91   | J/molxK | 1038.59         | Joback Method |
| cpg           | 1225.18   | J/molxK | 1005.87         | Joback Method |
| cpg           | 1209.35   | J/molxK | 973.15          | Joback Method |
| cpg           | 1192.37   | J/molxK | 940.43          | Joback Method |
| cpg           | 1174.20   | J/molxK | 907.71          | Joback Method |
| cpg           | 1253.57   | J/molxK | 1071.31         | Joback Method |
| dvisc         | 0.0000268 | Paxs    | 874.99          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0000364 | Paxs | 811.78 | Joback Method |
| dvisc | 0.0000521 | Paxs | 748.56 | Joback Method |
| dvisc | 0.0000798 | Paxs | 685.35 | Joback Method |
| dvisc | 0.0001331 | Paxs | 622.14 | Joback Method |
| dvisc | 0.0002494 | Paxs | 558.92 | Joback Method |
| dvisc | 0.0005484 | Paxs | 495.71 | Joback Method |

## Sources

|                        |   |
|------------------------|---|
| <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=U361761&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U361761&amp;Units=SI</a> |
| <b>Crippen Method:</b> | <a href="http://pubs.acs.org/doi/abs/10.1021/ci990307l">http://pubs.acs.org/doi/abs/10.1021/ci990307l</a>                                 |
| <b>Crippen Method:</b> | <a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</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/16-131-8/Dimethylmalonic-acid-butyl-tetradecyl-ester.pdf>

Generated by Cheméo on 2024-04-19 01:50:17.925404006 +0000 UTC m=+15780666.845981324.

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