

# Dimethylmalonic acid, heptyl nonyl ester

|                             |  |
|-----------------------------|--|
| <b>Inchi:</b>               | InChI=1S/C21H40O4/c1-5-7-9-11-12-14-16-18-25-20(23)21(3,4)19(22)24-17-15-13-10-8 |
| <b>InchiKey:</b>            | IUSWVICQWUHOSN-UHFFFAOYSA-N  |
| <b>Formula:</b>             | C21H40O4   |
| <b>SMILES:</b>              | CCCCCCCCCOC(=O)C(C)(C)C(=O)OCCCCCCC  |
| <b>Mol. weight [g/mol]:</b> | 356.54   |

## Physical Properties

| Property code | Value   | Unit                 | Source         |
|---------------|---------|----------------------|----------------|
| gf            | -339.06 | kJ/mol               | Joback Method  |
| hf            | -975.12 | kJ/mol               | Joback Method  |
| hfus          | 48.31   | kJ/mol               | Joback Method  |
| hvap          | 79.36   | kJ/mol               | Joback Method  |
| log10ws       | -6.10   |                      | Crippen Method |
| logp          | 5.820   |                      | Crippen Method |
| mvol          | 321.630 | ml/mol               | McGowan Method |
| pc            | 1025.31 | kPa                  | Joback Method  |
| rinpol        | 2209.00 |                      | NIST Webbook   |
| rinpol        | 2209.00 |                      | NIST Webbook   |
| tb            | 829.23  | K                    | Joback Method  |
| tc            | 1017.76 | K                    | Joback Method  |
| tf            | 473.17  | K                    | Joback Method  |
| vc            | 1.248   | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 1031.00   | J/molxK | 829.23          | Joback Method |
| cpg           | 1049.60   | J/molxK | 860.65          | Joback Method |
| cpg           | 1067.09   | J/molxK | 892.07          | Joback Method |
| cpg           | 1083.51   | J/molxK | 923.49          | Joback Method |
| cpg           | 1098.90   | J/molxK | 954.91          | Joback Method |
| cpg           | 1113.29   | J/molxK | 986.33          | Joback Method |
| cpg           | 1126.71   | J/molxK | 1017.76         | Joback Method |
| dvisc         | 0.0007084 | Paxs    | 473.17          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0003284 | Paxs | 532.51 | Joback Method |
| dvisc | 0.0001776 | Paxs | 591.86 | Joback Method |
| dvisc | 0.0001074 | Paxs | 651.20 | Joback Method |
| dvisc | 0.0000707 | Paxs | 710.54 | Joback Method |
| dvisc | 0.0000496 | Paxs | 769.89 | Joback Method |
| dvisc | 0.0000366 | Paxs | 829.23 | 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=U361706&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U361706&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>                                 |

## 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/18-984-0/Dimethylmalonic-acid-heptyl-nonyl-ester.pdf>

Generated by Cheméo on 2024-04-28 03:24:19.812432131 +0000 UTC m=+16563908.733009446.

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