

# Diethylmalonic acid, heptyl octyl ester

|                             |  |
|-----------------------------|--|
| <b>Inchi:</b>               | InChI=1S/C22H42O4/c1-5-9-11-13-15-17-19-26-21(24)22(7-3,8-4)20(23)25-18-16-14-12 |
| <b>InchiKey:</b>            | HKTLWTIKAXGXFV-UHFFFAOYSA-N  |
| <b>Formula:</b>             | C22H42O4   |
| <b>SMILES:</b>              | CCCCCCCCOC(=O)C(CC)(CC)C(=O)OCCCCCCC   |
| <b>Mol. weight [g/mol]:</b> | 370.57   |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -330.64 | kJ/mol  | Joback Method  |
| hf            | -995.76 | kJ/mol  | Joback Method  |
| hfus          | 50.90   | kJ/mol  | Joback Method  |
| hvap          | 81.58   | kJ/mol  | Joback Method  |
| log10ws       | -6.51   |         | Crippen Method |
| logp          | 6.210   |         | Crippen Method |
| mcvol         | 335.720 | ml/mol  | McGowan Method |
| pc            | 965.07  | kPa     | Joback Method  |
| rinpol        | 2257.00 |         | NIST Webbook   |
| tb            | 852.11  | K       | Joback Method  |
| tc            | 1044.11 | K       | Joback Method  |
| tf            | 484.44  | K       | Joback Method  |
| vc            | 1.304   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 1092.56   | J/molxK | 852.11          | Joback Method |
| cpg           | 1176.22   | J/molxK | 1012.11         | Joback Method |
| cpg           | 1161.66   | J/molxK | 980.11          | Joback Method |
| cpg           | 1146.06   | J/molxK | 948.11          | Joback Method |
| cpg           | 1129.36   | J/molxK | 916.11          | Joback Method |
| cpg           | 1111.54   | J/molxK | 884.11          | Joback Method |
| cpg           | 1189.77   | J/molxK | 1044.11         | Joback Method |
| dvisc         | 0.0000313 | Paxs    | 852.11          | Joback Method |
| dvisc         | 0.0000425 | Paxs    | 790.83          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0000608 | Paxs | 729.55 | Joback Method |
| dvisc | 0.0000927 | Paxs | 668.27 | Joback Method |
| dvisc | 0.0001539 | Paxs | 607.00 | Joback Method |
| dvisc | 0.0002865 | Paxs | 545.72 | Joback Method |
| dvisc | 0.0006242 | Paxs | 484.44 | 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=U369517&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U369517&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                                 |

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