

# Diethylmalonic acid, decyl 1-naphthyl ester

|                             |   |
|-----------------------------|---|
| <b>Inchi:</b>               | InChI=1S/C27H38O4/c1-4-7-8-9-10-11-12-15-21-30-25(28)27(5-2,6-3)26(29)31-24-20-16 |
| <b>InchiKey:</b>            | KTAFMWSPZDQDMW-UHFFFAOYSA-N   |
| <b>Formula:</b>             | C27H38O4  |
| <b>SMILES:</b>              | CCCCCCCCCOC(=O)C(CC)(CC)C(=O)Oc1cccc2ccccc12                                      |
| <b>Mol. weight [g/mol]:</b> | 426.59  |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -79.11  | kJ/mol  | Joback Method  |
| hf            | -682.83 | kJ/mol  | Joback Method  |
| hfus          | 54.52   | kJ/mol  | Joback Method  |
| hvap          | 97.29   | kJ/mol  | Joback Method  |
| log10ws       | -8.49   |         | Crippen Method |
| logp          | 7.235   |         | Crippen Method |
| mcvol         | 362.950 | ml/mol  | McGowan Method |
| pc            | 1015.53 | kPa     | Joback Method  |
| rinpol        | 3014.00 |         | NIST Webbook   |
| tb            | 1017.15 | K       | Joback Method  |
| tc            | 1245.65 | K       | Joback Method  |
| tf            | 612.43  | K       | Joback Method  |
| vc            | 1.399   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 1224.23   | J/molxK | 1017.15         | Joback Method |
| cpg           | 1240.62   | J/molxK | 1055.23         | Joback Method |
| cpg           | 1255.87   | J/molxK | 1093.32         | Joback Method |
| cpg           | 1270.09   | J/molxK | 1131.40         | Joback Method |
| cpg           | 1283.39   | J/molxK | 1169.49         | Joback Method |
| cpg           | 1295.87   | J/molxK | 1207.57         | Joback Method |
| cpg           | 1307.66   | J/molxK | 1245.65         | Joback Method |
| dvisc         | 0.0003046 | Paxs    | 612.43          | Joback Method |
| dvisc         | 0.0001724 | Paxs    | 679.88          | Joback Method |

|       |           |      |         |               |
|-------|-----------|------|---------|---------------|
| dvisc | 0.0001081 | Paxs | 747.34  | Joback Method |
| dvisc | 0.0000732 | Paxs | 814.79  | Joback Method |
| dvisc | 0.0000527 | Paxs | 882.24  | Joback Method |
| dvisc | 0.0000397 | Paxs | 949.70  | Joback Method |
| dvisc | 0.0000310 | Paxs | 1017.15 | 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=U369874&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U369874&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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