

# Diethylmalonic acid, hexadecyl isobutyl ester

|                             |  |
|-----------------------------|--|
| <b>Inchi:</b>               | InChI=1S/C27H52O4/c1-6-9-10-11-12-13-14-15-16-17-18-19-20-21-22-30-25(28)27(7-2, |
| <b>InchiKey:</b>            | BBIUNWCCHOALN-UHFFFAOYSA-N   |
| <b>Formula:</b>             | C27H52O4   |
| <b>SMILES:</b>              | CCCCCCCCCCCCCCCCOC(=O)C(CC)(CC)C(=O)OCC(C)C                                      |
| <b>Mol. weight [g/mol]:</b> | 440.70   |

## Physical Properties

| Property code | Value    | Unit                 | Source         |
|---------------|----------|----------------------|----------------|
| gf            | -290.98  | kJ/mol               | Joback Method  |
| hf            | -1104.24 | kJ/mol               | Joback Method  |
| hfus          | 60.32    | kJ/mol               | Joback Method  |
| hvap          | 92.32    | kJ/mol               | Joback Method  |
| log10ws       | -8.37    |                      | Crippen Method |
| logp          | 8.017    |                      | Crippen Method |
| mvol          | 406.170  | ml/mol               | McGowan Method |
| pc            | 734.03   | kPa                  | Joback Method  |
| rinpol        | 2753.00  |                      | NIST Webbook   |
| tb            | 966.07   | K                    | Joback Method  |
| tc            | 1188.14  | K                    | Joback Method  |
| tf            | 525.79   | K                    | Joback Method  |
| vc            | 1.579    | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 1409.68   | J/molxK | 966.07          | Joback Method |
| cpg           | 1500.92   | J/molxK | 1151.12         | Joback Method |
| cpg           | 1485.63   | J/molxK | 1114.11         | Joback Method |
| cpg           | 1468.94   | J/molxK | 1077.10         | Joback Method |
| cpg           | 1450.77   | J/molxK | 1040.09         | Joback Method |
| cpg           | 1431.04   | J/molxK | 1003.08         | Joback Method |
| cpg           | 1514.88   | J/molxK | 1188.14         | Joback Method |
| dvisc         | 0.0000128 | Paxs    | 966.07          | Joback Method |
| dvisc         | 0.0000178 | Paxs    | 892.69          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0000262 | Paxs | 819.31 | Joback Method |
| dvisc | 0.0000418 | Paxs | 745.93 | Joback Method |
| dvisc | 0.0000735 | Paxs | 672.55 | Joback Method |
| dvisc | 0.0001487 | Paxs | 599.17 | Joback Method |
| dvisc | 0.0003661 | Paxs | 525.79 | Joback Method |

## Sources

|                        |   |
|------------------------|---|
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=U369434&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U369434&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>                         |
| <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>                     |

## 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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