

# Diethylmalonic acid, 2-ethylphenyl undecyl ester

**Inchi:** InChI=1S/C26H42O4/c1-5-9-10-11-12-13-14-15-18-21-29-24(27)26(7-3,8-4)25(28)30-23

**InchiKey:** NZZILWTWPOZJDK-UHFFFAOYSA-N

**Formula:** C26H42O4

**SMILES:** CCCCCCCCCCOC(=O)C(CC)(CC)C(=O)Oc1ccccc1CC

**Mol. weight [g/mol]:** 418.61

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -194.18 | kJ/mol  | Joback Method  |
| hf            | -853.26 | kJ/mol  | Joback Method  |
| hfus          | 54.91   | kJ/mol  | Joback Method  |
| hvap          | 93.42   | kJ/mol  | Joback Method  |
| log10ws       | -7.91   |         | Crippen Method |
| logp          | 7.035   |         | Crippen Method |
| mcvol         | 368.320 | ml/mol  | McGowan Method |
| pc            | 925.55  | kPa     | Joback Method  |
| rinpol        | 2707.00 |         | NIST Webbook   |
| tb            | 975.29  | K       | Joback Method  |
| tc            | 1194.04 | K       | Joback Method  |
| tf            | 568.46  | K       | Joback Method  |
| vc            | 1.421   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 1243.60   | J/mol×K | 975.29          | Joback Method |
| cpg           | 1317.86   | J/mol×K | 1157.58         | Joback Method |
| cpg           | 1305.54   | J/mol×K | 1121.12         | Joback Method |
| cpg           | 1292.02   | J/mol×K | 1084.66         | Joback Method |
| cpg           | 1277.24   | J/mol×K | 1048.21         | Joback Method |
| cpg           | 1261.12   | J/mol×K | 1011.75         | Joback Method |
| cpg           | 1329.05   | J/mol×K | 1194.04         | Joback Method |
| dvisc         | 0.0000173 | Pa×s    | 975.29          | Joback Method |
| dvisc         | 0.0000230 | Pa×s    | 907.49          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0000320 | Pa×s | 839.68 | Joback Method |
| dvisc | 0.0000472 | Pa×s | 771.88 | Joback Method |
| dvisc | 0.0000751 | Pa×s | 704.07 | Joback Method |
| dvisc | 0.0001319 | Pa×s | 636.26 | Joback Method |
| dvisc | 0.0002647 | Pa×s | 568.46 | Joback Method |

## Sources

|                        |                                                                                                                                           |
|------------------------|-------------------------------------------------------------------------------------------------------------------------------------------|
| <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>                     |
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=U370250&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U370250&amp;Units=SI</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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