

# 1,2-Cyclohexanedicarboxylic acid, ethyl 3-pentyl ester

|                      |  |
|----------------------|--|
| Inchi:               | InChI=1S/C15H26O4/c1-4-11(5-2)19-15(17)13-10-8-7-9-12(13)14(16)18-6-3/h11-13H,4- |
| InchiKey:            | CBLYCSVNHRLRGM-UHFFFAOYSA-N  |
| Formula:             | C15H26O4   |
| SMILES:              | CCOC(=O)C1CCCCC1C(=O)OC(CC)CC  |
| Mol. weight [g/mol]: | 270.36   |

## Physical Properties

| Property code | Value   | Unit                 | Source         |
|---------------|---------|----------------------|----------------|
| gf            | -378.12 | kJ/mol               | Joback Method  |
| hf            | -813.83 | kJ/mol               | Joback Method  |
| hfus          | 29.56   | kJ/mol               | Joback Method  |
| hvap          | 67.03   | kJ/mol               | Joback Method  |
| log10ws       | -3.35   |                      | Crippen Method |
| logp          | 3.088   |                      | Crippen Method |
| mcvol         | 226.230 | ml/mol               | McGowan Method |
| pc            | 1756.54 | kPa                  | Joback Method  |
| rinpol        | 1758.00 |                      | NIST Webbook   |
| tb            | 709.62  | K                    | Joback Method  |
| tc            | 910.33  | K                    | Joback Method  |
| tf            | 391.27  | K                    | Joback Method  |
| vc            | 0.850   | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 674.70    | J/molxK | 709.62          | Joback Method |
| cpg           | 693.58    | J/molxK | 743.07          | Joback Method |
| cpg           | 711.32    | J/molxK | 776.52          | Joback Method |
| cpg           | 727.90    | J/molxK | 809.98          | Joback Method |
| cpg           | 743.34    | J/molxK | 843.43          | Joback Method |
| cpg           | 757.63    | J/molxK | 876.88          | Joback Method |
| cpg           | 770.79    | J/molxK | 910.33          | Joback Method |
| dvisc         | 0.0019030 | Paxs    | 391.27          | Joback Method |
| dvisc         | 0.0009222 | Paxs    | 444.33          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0005216 | Paxs | 497.39 | Joback Method |
| dvisc | 0.0003292 | Paxs | 550.44 | Joback Method |
| dvisc | 0.0002254 | Paxs | 603.50 | Joback Method |
| dvisc | 0.0001640 | Paxs | 656.56 | Joback Method |
| dvisc | 0.0001251 | Paxs | 709.62 | 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=U339502&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U339502&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>                                 |

## 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/93-422-0/1-2-Cyclohexanedicarboxylic-acid-ethyl-3-pentyl-ester.pdf>

Generated by Cheméo on 2024-04-20 10:26:45.717529106 +0000 UTC m=+15898054.638106427.

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