

# cis-Cyclohex-4-en-1,2-dicarboxylic acid, 3-methylbutyl pentyl ester

|                      |   |
|----------------------|---|
| Inchi:               | InChI=1S/C18H30O4/c1-4-5-8-12-21-17(19)15-9-6-7-10-16(15)18(20)22-13-11-14(2)3/h6 |
| InchiKey:            | GADNAZPUYKALMC-UHFFFAOYSA-N   |
| Formula:             | C18H30O4  |
| SMILES:              | CCCCCOC(=O)C1CC=CCC1C(=O)OCCC(C)C   |
| Mol. weight [g/mol]: | 310.43  |

## Physical Properties

| Property code | Value   | Unit                 | Source         |
|---------------|---------|----------------------|----------------|
| gf            | -322.90 | kJ/mol               | Joback Method  |
| hf            | -817.97 | kJ/mol               | Joback Method  |
| hfus          | 38.55   | kJ/mol               | Joback Method  |
| hvap          | 74.00   | kJ/mol               | Joback Method  |
| log10ws       | -4.11   |                      | Crippen Method |
| logp          | 3.891   |                      | Crippen Method |
| mcvol         | 264.200 | ml/mol               | McGowan Method |
| pc            | 1436.98 | kPa                  | Joback Method  |
| rinpol        | 2053.00 |                      | NIST Webbook   |
| tb            | 777.42  | K                    | Joback Method  |
| tc            | 975.18  | K                    | Joback Method  |
| tf            | 425.84  | K                    | Joback Method  |
| vc            | 1.004   | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 822.66    | J/mol×K | 777.42          | Joback Method |
| cpg           | 841.13    | J/mol×K | 810.38          | Joback Method |
| cpg           | 858.38    | J/mol×K | 843.34          | Joback Method |
| cpg           | 874.42    | J/mol×K | 876.30          | Joback Method |
| cpg           | 889.25    | J/mol×K | 909.26          | Joback Method |
| cpg           | 902.89    | J/mol×K | 942.22          | Joback Method |
| cpg           | 915.36    | J/mol×K | 975.18          | Joback Method |
| dvisc         | 0.0013807 | Paxs    | 425.84          | Joback Method |
| dvisc         | 0.0006672 | Paxs    | 484.44          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0003772 | Paxs | 543.03 | Joback Method |
| dvisc | 0.0002383 | Paxs | 601.63 | Joback Method |
| dvisc | 0.0001633 | Paxs | 660.23 | Joback Method |
| dvisc | 0.0001191 | Paxs | 718.82 | Joback Method |
| dvisc | 0.0000910 | Paxs | 777.42 | 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=U382822&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U382822&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.cheméo.com/doc/models/crippen_log10ws">https://www.cheméo.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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