

# 1,4-Dioxepane-2-carboxylic acid, 3-phenyl, ethyl ester, cis

|                      |   |
|----------------------|---|
| Inchi:               | InChI=1S/C14H18O4/c1-2-16-14(15)13-12(17-9-6-10-18-13)11-7-4-3-5-8-11/h3-5,7-8,12 |
| InchiKey:            | JCVYDFPRMXWSHP-CHWSQXEVSAN  |
| Formula:             | C14H18O4  |
| SMILES:              | CCOC(=O)C1OCCOC1c1ccccc1  |
| Mol. weight [g/mol]: | 250.29  |

## Physical Properties

| Property code | Value   | Unit                 | Source         |
|---------------|---------|----------------------|----------------|
| gf            | -222.11 | kJ/mol               | Joback Method  |
| hf            | -576.74 | kJ/mol               | Joback Method  |
| hfus          | 35.61   | kJ/mol               | Joback Method  |
| hvap          | 67.50   | kJ/mol               | Joback Method  |
| log10ws       | -2.29   |                      | Crippen Method |
| logp          | 2.096   |                      | Crippen Method |
| mcvol         | 192.680 | ml/mol               | McGowan Method |
| pc            | 2485.07 | kPa                  | Joback Method  |
| rinpol        | 1776.00 |                      | NIST Webbook   |
| tb            | 695.74  | K                    | Joback Method  |
| tc            | 932.17  | K                    | Joback Method  |
| tf            | 398.88  | K                    | Joback Method  |
| vc            | 0.702   | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 555.58    | J/molxK | 695.74          | Joback Method |
| cpg           | 634.92    | J/molxK | 892.76          | Joback Method |
| cpg           | 621.94    | J/molxK | 853.36          | Joback Method |
| cpg           | 607.54    | J/molxK | 813.95          | Joback Method |
| cpg           | 591.69    | J/molxK | 774.55          | Joback Method |
| cpg           | 574.37    | J/molxK | 735.14          | Joback Method |
| cpg           | 646.48    | J/molxK | 932.17          | Joback Method |
| dvisc         | 0.0001198 | Paxs    | 695.74          | Joback Method |
| dvisc         | 0.0001590 | Paxs    | 646.26          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0002213 | Paxs | 596.79 | Joback Method |
| dvisc | 0.0003268 | Paxs | 547.31 | Joback Method |
| dvisc | 0.0005215 | Paxs | 497.83 | Joback Method |
| dvisc | 0.0009227 | Paxs | 448.36 | Joback Method |
| dvisc | 0.0018807 | Paxs | 398.88 | Joback Method |

## Sources

|                        |   |
|------------------------|---|
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=R329436&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R329436&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>                                     |
| <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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