

cis-Cyclohex-4-en-1,2-dicarboxylic acid, phenethyl tridecyl ester

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|----------------------|---|
| Inchi: | InChI=1S/C29H44O4/c1-2-3-4-5-6-7-8-9-10-11-17-23-32-28(30)26-20-15-16-21-27(26)29 |
| InchiKey: | VYSWASZPZUVTBG-UHFFFAOYSA-N |
| Formula: | C29H44O4 |
| SMILES: | CCCCCCCCCCCCOC(=O)C1CC=CCC1C(=O)OCCc1ccccc1 |
| Mol. weight [g/mol]: | 456.66 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -115.43 | kJ/mol | Joback Method |
| hf | -803.20 | kJ/mol | Joback Method |
| hfus | 64.61 | kJ/mol | Joback Method |
| hvap | 101.15 | kJ/mol | Joback Method |
| log10ws | -8.05 | | Crippen Method |
| logp | 7.209 | | Crippen Method |
| mvol | 395.430 | ml/mol | McGowan Method |
| pc | 880.00 | kPa | Joback Method |
| rinpol | 3328.00 | | NIST Webbook |
| rinpol | 3328.00 | | NIST Webbook |
| tb | 1056.22 | K | Joback Method |
| tc | 1293.87 | K | Joback Method |
| tf | 591.23 | K | Joback Method |
| vc | 1.518 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 1399.07 | J/molxK | 1056.22 | Joback Method |
| cpg | 1415.38 | J/molxK | 1095.83 | Joback Method |
| cpg | 1429.72 | J/molxK | 1135.44 | Joback Method |
| cpg | 1442.19 | J/molxK | 1175.04 | Joback Method |
| cpg | 1452.86 | J/molxK | 1214.65 | Joback Method |
| cpg | 1461.82 | J/molxK | 1254.26 | Joback Method |
| cpg | 1469.14 | J/molxK | 1293.87 | Joback Method |
| dvisc | 0.0003110 | Paxs | 591.23 | Joback Method |

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|-------|-----------|------|---------|---------------|
| dvisc | 0.0001538 | Paxs | 668.73 | Joback Method |
| dvisc | 0.0000880 | Paxs | 746.23 | Joback Method |
| dvisc | 0.0000560 | Paxs | 823.73 | Joback Method |
| dvisc | 0.0000385 | Paxs | 901.22 | Joback Method |
| dvisc | 0.0000281 | Paxs | 978.72 | Joback Method |
| dvisc | 0.0000214 | Paxs | 1056.22 | Joback Method |

Sources

| | |
|------------------------|---|
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci9903071 |
| Crippen Method: | https://www.chemeo.com/doc/models/crippen_log10ws |
| Joback Method: | https://en.wikipedia.org/wiki/Joback_method |
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U382800&Units=SI |

Legend

| | |
|-----------------|---|
| cpg: | Ideal gas heat capacity |
| dvisc: | Dynamic viscosity |
| gf: | Standard Gibbs free energy of formation |
| hf: | Enthalpy of formation at standard conditions |
| hfus: | Enthalpy of fusion at standard conditions |
| hvap: | Enthalpy of vaporization at standard conditions |
| log10ws: | Log10 of Water solubility in mol/l |
| logp: | Octanol/Water partition coefficient |
| mcvol: | McGowan's characteristic volume |
| pc: | Critical Pressure |
| rinpol: | Non-polar retention indices |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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