

1,2-Cyclohexanedicarboxylic acid, hexyl isopropyl ester

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|----------------------|---|
| Inchi: | InChI=1S/C17H30O4/c1-4-5-6-9-12-20-16(18)14-10-7-8-11-15(14)17(19)21-13(2)3/h13-1 |
| InchiKey: | DGXFWVNYSUOIBQ-UHFFFAOYSA-N |
| Formula: | C17H30O4 |
| SMILES: | CCCCCOC(=O)C1CCCCC1C(=O)OC(C)C |
| Mol. weight [g/mol]: | 298.42 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -361.28 | kJ/mol | Joback Method |
| hf | -855.11 | kJ/mol | Joback Method |
| hfus | 34.74 | kJ/mol | Joback Method |
| hvap | 71.48 | kJ/mol | Joback Method |
| log10ws | -4.19 | | Crippen Method |
| logp | 3.868 | | Crippen Method |
| mvol | 254.410 | ml/mol | McGowan Method |
| pc | 1504.65 | kPa | Joback Method |
| rinpol | 2150.00 | | NIST Webbook |
| rinpol | 2150.00 | | NIST Webbook |
| tb | 755.38 | K | Joback Method |
| tc | 952.93 | K | Joback Method |
| tf | 413.81 | K | Joback Method |
| vc | 0.962 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 790.12 | J/molxK | 755.38 | Joback Method |
| cpg | 809.29 | J/molxK | 788.31 | Joback Method |
| cpg | 827.24 | J/molxK | 821.23 | Joback Method |
| cpg | 843.97 | J/molxK | 854.16 | Joback Method |
| cpg | 859.49 | J/molxK | 887.08 | Joback Method |
| cpg | 873.82 | J/molxK | 920.01 | Joback Method |
| cpg | 886.95 | J/molxK | 952.93 | Joback Method |
| dvisc | 0.0015981 | Paxs | 413.81 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0007558 | Paxs | 470.74 | Joback Method |
| dvisc | 0.0004201 | Paxs | 527.67 | Joback Method |
| dvisc | 0.0002618 | Paxs | 584.60 | Joback Method |
| dvisc | 0.0001775 | Paxs | 641.52 | Joback Method |
| dvisc | 0.0001282 | Paxs | 698.45 | Joback Method |
| dvisc | 0.0000972 | Paxs | 755.38 | Joback Method |

Sources

| | |
|------------------------|---|
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U339640&Units=SI |
| 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 |

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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