

Carbonic acid, (1R)-(-)-menthyl dodecyl ester

Inchi: InChI=1S/C23H44O3/c1-5-6-7-8-9-10-11-12-13-14-17-25-23(24)26-22-18-21(19(2)3)16-
InchiKey: WAESBELRAYYZQT-UHFFFAOYSA-N
Formula: C23H44O3
SMILES: CCCCCCCCCCOC(=O)OC1CC(C(C)C)CCC1C
Mol. weight [g/mol]: 368.59

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -189.55 | kJ/mol | Joback Method |
| hf | -886.71 | kJ/mol | Joback Method |
| hfus | 49.75 | kJ/mol | Joback Method |
| hvap | 77.78 | kJ/mol | Joback Method |
| log10ws | -7.66 | | Crippen Method |
| logp | 7.521 | | Crippen Method |
| mvol | 337.380 | ml/mol | McGowan Method |
| pc | 946.75 | kPa | Joback Method |
| rinpol | 2500.00 | | NIST Webbook |
| rinpol | 2500.00 | | NIST Webbook |
| tb | 834.12 | K | Joback Method |
| tc | 1026.28 | K | Joback Method |
| tf | 427.26 | K | Joback Method |
| vc | 1.290 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 1132.53 | J/molxK | 834.12 | Joback Method |
| cpg | 1154.52 | J/molxK | 866.15 | Joback Method |
| cpg | 1175.05 | J/molxK | 898.17 | Joback Method |
| cpg | 1194.14 | J/molxK | 930.20 | Joback Method |
| cpg | 1211.82 | J/molxK | 962.23 | Joback Method |
| cpg | 1228.10 | J/molxK | 994.25 | Joback Method |
| cpg | 1243.00 | J/molxK | 1026.28 | Joback Method |
| dvisc | 0.0011474 | Paxs | 427.26 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0004933 | Paxs | 495.07 | Joback Method |
| dvisc | 0.0002599 | Paxs | 562.88 | Joback Method |
| dvisc | 0.0001572 | Paxs | 630.69 | Joback Method |
| dvisc | 0.0001048 | Paxs | 698.50 | Joback Method |
| dvisc | 0.0000751 | Paxs | 766.31 | Joback Method |
| dvisc | 0.0000568 | Paxs | 834.12 | Joback Method |

Sources

| | |
|------------------------|---|
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U392441&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307l |
| Crippen Method: | https://www.chemeo.com/doc/models/crippen_log10ws |
| Joback Method: | https://en.wikipedia.org/wiki/Joback_method |

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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