

Dimethylmalonic acid, decyl isohexyl ester

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| Inchi: | InChI=1S/C21H40O4/c1-6-7-8-9-10-11-12-13-16-24-19(22)21(4,5)20(23)25-17-14-15-18 |
| InchiKey: | INJNAIKNMFBCCH-UHFFFAOYSA-N |
| Formula: | C21H40O4 |
| SMILES: | CCCCCCCCCOC(=O)C(C)(C)C(=O)OCCCC(C)C |
| Mol. weight [g/mol]: | 356.54 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -341.50 | kJ/mol | Joback Method |
| hf | -980.40 | kJ/mol | Joback Method |
| hfus | 44.78 | kJ/mol | Joback Method |
| hvap | 78.97 | kJ/mol | Joback Method |
| log10ws | -5.85 | | Crippen Method |
| logp | 5.676 | | Crippen Method |
| mvol | 321.630 | ml/mol | McGowan Method |
| pc | 1030.59 | kPa | Joback Method |
| rinpol | 2179.00 | | NIST Webbook |
| rinpol | 2179.00 | | NIST Webbook |
| tb | 828.79 | K | Joback Method |
| tc | 1018.11 | K | Joback Method |
| tf | 458.17 | K | Joback Method |
| vc | 1.242 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 1031.48 | J/molxK | 828.79 | Joback Method |
| cpg | 1050.15 | J/molxK | 860.34 | Joback Method |
| cpg | 1067.69 | J/molxK | 891.90 | Joback Method |
| cpg | 1084.15 | J/molxK | 923.45 | Joback Method |
| cpg | 1099.56 | J/molxK | 955.00 | Joback Method |
| cpg | 1113.95 | J/molxK | 986.55 | Joback Method |
| cpg | 1127.36 | J/molxK | 1018.11 | Joback Method |
| dvisc | 0.0008434 | Paxs | 458.17 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0003582 | Paxs | 519.94 | Joback Method |
| dvisc | 0.0001825 | Paxs | 581.71 | Joback Method |
| dvisc | 0.0001058 | Paxs | 643.48 | Joback Method |
| dvisc | 0.0000675 | Paxs | 705.25 | Joback Method |
| dvisc | 0.0000463 | Paxs | 767.02 | Joback Method |
| dvisc | 0.0000336 | Paxs | 828.79 | Joback Method |

Sources

| | |
|------------------------|---|
| 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=U361720&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307l |
| Crippen Method: | https://www.chemeo.com/doc/models/crippen_log10ws |

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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