

Diethylmalonic acid, dodecyl ethyl ester

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| Inchi: | InChI=1S/C21H40O4/c1-5-9-10-11-12-13-14-15-16-17-18-25-20(23)21(6-2,7-3)19(22)24 |
| InchiKey: | JBVFCWPIUMQZRI-UHFFFAOYSA-N |
| Formula: | C21H40O4 |
| SMILES: | CCCCCCCCCCCCOC(=O)C(CC)(CC)C(=O)OCC |
| Mol. weight [g/mol]: | 356.54 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | -339.06 | kJ/mol | Joback Method |
| hf | -975.12 | kJ/mol | Joback Method |
| hfus | 48.31 | kJ/mol | Joback Method |
| hvap | 79.36 | kJ/mol | Joback Method |
| log10ws | -6.10 | | Crippen Method |
| logp | 5.820 | | Crippen Method |
| mcvol | 321.630 | ml/mol | McGowan Method |
| pc | 1025.31 | kPa | Joback Method |
| rinpol | 2202.00 | | NIST Webbook |
| tb | 829.23 | K | Joback Method |
| tc | 1017.76 | K | Joback Method |
| tf | 473.17 | K | Joback Method |
| vc | 1.248 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 1031.00 | J/molxK | 829.23 | Joback Method |
| cpg | 1113.29 | J/molxK | 986.33 | Joback Method |
| cpg | 1098.90 | J/molxK | 954.91 | Joback Method |
| cpg | 1083.51 | J/molxK | 923.49 | Joback Method |
| cpg | 1067.09 | J/molxK | 892.07 | Joback Method |
| cpg | 1049.60 | J/molxK | 860.65 | Joback Method |
| cpg | 1126.71 | J/molxK | 1017.76 | Joback Method |
| dvisc | 0.0000366 | Paxs | 829.23 | Joback Method |
| dvisc | 0.0000496 | Paxs | 769.89 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0000707 | Paxs | 710.54 | Joback Method |
| dvisc | 0.0001074 | Paxs | 651.20 | Joback Method |
| dvisc | 0.0001776 | Paxs | 591.86 | Joback Method |
| dvisc | 0.0003284 | Paxs | 532.51 | Joback Method |
| dvisc | 0.0007084 | Paxs | 473.17 | Joback Method |

Sources

| | |
|------------------------|---|
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U369712&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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