

Diethylmalonic acid, 4-methoxyphenyl nonyl ester

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|----------------------|---|
| Inchi: | InChI=1S/C23H36O5/c1-5-8-9-10-11-12-13-18-27-21(24)23(6-2,7-3)22(25)28-20-16-14-1 |
| InchiKey: | VBPXFOGDMOWLGX-UHFFFAOYSA-N |
| Formula: | C23H36O5 |
| SMILES: | CCCCCCCCCOC(=O)C(CC)(CC)C(=O)Oc1ccc(OC)cc1 |
| Mol. weight [g/mol]: | 392.53 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -324.44 | kJ/mol | Joback Method |
| hf | -923.56 | kJ/mol | Joback Method |
| hfus | 48.33 | kJ/mol | Joback Method |
| hvap | 89.16 | kJ/mol | Joback Method |
| log10ws | -6.39 | | Crippen Method |
| logp | 5.701 | | Crippen Method |
| mcvol | 331.920 | ml/mol | McGowan Method |
| pc | 1099.35 | kPa | Joback Method |
| rinsol | 2649.00 | | NIST Webbook |
| tb | 929.07 | K | Joback Method |
| tc | 1139.98 | K | Joback Method |
| tf | 556.88 | K | Joback Method |
| vc | 1.270 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 1088.08 | J/molxK | 929.07 | Joback Method |
| cpg | 1104.27 | J/molxK | 964.22 | Joback Method |
| cpg | 1119.11 | J/molxK | 999.37 | Joback Method |
| cpg | 1132.62 | J/molxK | 1034.52 | Joback Method |
| cpg | 1144.84 | J/molxK | 1069.67 | Joback Method |
| cpg | 1155.83 | J/molxK | 1104.82 | Joback Method |
| cpg | 1165.60 | J/molxK | 1139.98 | Joback Method |
| dvisc | 0.0002633 | Paxs | 556.88 | Joback Method |
| dvisc | 0.0001388 | Paxs | 618.91 | Joback Method |

| | | | | |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0000822 | Paxs | 680.94 | Joback Method |
| dvisc | 0.0000532 | Paxs | 742.98 | Joback Method |
| dvisc | 0.0000368 | Paxs | 805.01 | Joback Method |
| dvisc | 0.0000268 | Paxs | 867.04 | Joback Method |
| dvisc | 0.0000204 | Paxs | 929.07 | 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=U369844&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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