

Diethylmalonic acid, 2,3-dichlorophenyl pentyl ester

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|-----------------------------|---|
| Inchi: | InChI=1S/C18H24Cl2O4/c1-4-7-8-12-23-16(21)18(5-2,6-3)17(22)24-14-11-9-10-13(19)15 |
| InchiKey: | RARHNJHWDWBCSM-UHFFFAOYSA-N |
| Formula: | C18H24Cl2O4 |
| SMILES: | CCCCCOC(=O)C(CC)(CC)C(=O)Oc1cccc(Cl)c1Cl |
| Mol. weight [g/mol]: | 375.29 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | -295.03 | kJ/mol | Joback Method |
| hf | -731.09 | kJ/mol | Joback Method |
| hfus | 42.19 | kJ/mol | Joback Method |
| hvap | 85.05 | kJ/mol | Joback Method |
| log10ws | -5.96 | | Crippen Method |
| logp | 5.439 | | Crippen Method |
| mcvol | 280.080 | ml/mol | McGowan Method |
| pc | 1470.23 | kPa | Joback Method |
| rinsol | 2324.00 | | NIST Webbook |
| tb | 872.09 | K | Joback Method |
| tc | 1087.50 | K | Joback Method |
| tf | 550.66 | K | Joback Method |
| vc | 1.071 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 813.48 | J/molxK | 872.09 | Joback Method |
| cpg | 870.07 | J/molxK | 1051.60 | Joback Method |
| cpg | 860.77 | J/molxK | 1015.70 | Joback Method |
| cpg | 850.50 | J/molxK | 979.79 | Joback Method |
| cpg | 839.23 | J/molxK | 943.89 | Joback Method |
| cpg | 826.90 | J/molxK | 907.99 | Joback Method |
| cpg | 878.44 | J/molxK | 1087.50 | Joback Method |
| dvisc | 0.0000430 | Paxs | 872.09 | Joback Method |
| dvisc | 0.0000549 | Paxs | 818.52 | Joback Method |

| | | | | |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0000725 | Paxs | 764.95 | Joback Method |
| dvisc | 0.0000999 | Paxs | 711.38 | Joback Method |
| dvisc | 0.0001450 | Paxs | 657.80 | Joback Method |
| dvisc | 0.0002249 | Paxs | 604.23 | Joback Method |
| dvisc | 0.0003800 | Paxs | 550.66 | Joback Method |

Sources

| | |
|------------------------|---|
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U370033&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 |
| 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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