

Terephthalic acid, ethyl 2-heptyl ester

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|-----------------------------|---|
| Inchi: | InChI=1S/C17H24O4/c1-4-6-7-8-13(3)21-17(19)15-11-9-14(10-12-15)16(18)20-5-2/h9-13 |
| InchiKey: | NDYLCUKKNUTEQI-UHFFFAOYSA-N |
| Formula: | C17H24O4 |
| SMILES: | CCCCC(C)OC(=O)c1ccc(C(=O)OCC)cc1 |
| Mol. weight [g/mol]: | 292.37 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -275.24 | kJ/mol | Joback Method |
| hf | -664.03 | kJ/mol | Joback Method |
| hfus | 35.49 | kJ/mol | Joback Method |
| hvap | 74.30 | kJ/mol | Joback Method |
| log10ws | -5.00 | | Crippen Method |
| logp | 3.989 | | Crippen Method |
| mvol | 241.510 | ml/mol | McGowan Method |
| pc | 1687.95 | kPa | Joback Method |
| rinpol | 2073.00 | | NIST Webbook |
| tb | 772.16 | K | Joback Method |
| tc | 975.26 | K | Joback Method |
| tf | 449.61 | K | Joback Method |
| vc | 0.921 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 704.89 | J/molxK | 772.16 | Joback Method |
| cpg | 720.52 | J/molxK | 806.01 | Joback Method |
| cpg | 735.10 | J/molxK | 839.86 | Joback Method |
| cpg | 748.65 | J/molxK | 873.71 | Joback Method |
| cpg | 761.19 | J/molxK | 907.56 | Joback Method |
| cpg | 772.73 | J/molxK | 941.41 | Joback Method |
| cpg | 783.29 | J/molxK | 975.26 | Joback Method |
| dvisc | 0.0009014 | Paxs | 449.61 | Joback Method |
| dvisc | 0.0004783 | Paxs | 503.37 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0002868 | Paxs | 557.13 | Joback Method |
| dvisc | 0.0001882 | Paxs | 610.88 | Joback Method |
| dvisc | 0.0001322 | Paxs | 664.64 | Joback Method |
| dvisc | 0.0000979 | Paxs | 718.40 | Joback Method |
| dvisc | 0.0000756 | Paxs | 772.16 | Joback Method |

Sources

| | |
|------------------------|---|
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U356285&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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<https://www.chemeo.com/cid/37-610-3/Terephthalic-acid-ethyl-2-heptyl-ester.pdf>

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