

Terephthalic acid, heptyl 3-hexyl ester

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| Inchi: | InChI=1S/C21H32O4/c1-4-7-8-9-10-16-24-20(22)17-12-14-18(15-13-17)21(23)25-19(6-3 |
| InchiKey: | QXHNZVSDXMXHQV-UHFFFAOYSA-N |
| Formula: | C21H32O4 |
| SMILES: | CCCCCCCOC(=O)c1ccc(C(=O)OC(CC)CCC)cc1 |
| Mol. weight [g/mol]: | 348.48 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -241.56 | kJ/mol | Joback Method |
| hf | -746.59 | kJ/mol | Joback Method |
| hfus | 45.85 | kJ/mol | Joback Method |
| hvap | 83.20 | kJ/mol | Joback Method |
| log10ws | -6.67 | | Crippen Method |
| logp | 5.549 | | Crippen Method |
| mcvol | 297.870 | ml/mol | McGowan Method |
| pc | 1259.27 | kPa | Joback Method |
| rinpola | 2472.00 | | NIST Webbook |
| tb | 863.68 | K | Joback Method |
| tc | 1066.32 | K | Joback Method |
| tf | 494.69 | K | Joback Method |
| vc | 1.145 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 937.65 | J/molxK | 863.68 | Joback Method |
| cpg | 954.09 | J/molxK | 897.45 | Joback Method |
| cpg | 969.33 | J/molxK | 931.23 | Joback Method |
| cpg | 983.39 | J/molxK | 965.00 | Joback Method |
| cpg | 996.30 | J/molxK | 998.77 | Joback Method |
| cpg | 1008.08 | J/molxK | 1032.54 | Joback Method |
| cpg | 1018.76 | J/molxK | 1066.32 | Joback Method |
| dvisc | 0.0006043 | Paxs | 494.69 | Joback Method |
| dvisc | 0.0003060 | Paxs | 556.19 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0001775 | Paxs | 617.69 | Joback Method |
| dvisc | 0.0001136 | Paxs | 679.18 | Joback Method |
| dvisc | 0.0000783 | Paxs | 740.68 | Joback Method |
| dvisc | 0.0000571 | Paxs | 802.18 | Joback Method |
| dvisc | 0.0000436 | Paxs | 863.68 | Joback Method |

Sources

| | |
|------------------------|---|
| 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 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U356230&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307l |

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 |

Latest version available from:

<https://www.chemeo.com/cid/62-590-8/Terephthalic-acid-heptyl-3-hexyl-ester.pdf>

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