

Terephthalic acid, 2-heptyl tridecyl ester

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| Inchi: | InChI=1S/C28H46O4/c1-4-6-8-9-10-11-12-13-14-15-17-23-31-27(29)25-19-21-26(22-20- |
| InchiKey: | CPBIRCPTZBDDSK-UHFFFAOYSA-N |
| Formula: | C28H46O4 |
| SMILES: | CCCCCCCCCCCCOC(=O)c1ccc(C(=O)OC(C)CCCC)cc1 |
| Mol. weight [g/mol]: | 446.66 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | -182.62 | kJ/mol | Joback Method |
| hf | -891.07 | kJ/mol | Joback Method |
| hfus | 63.98 | kJ/mol | Joback Method |
| hvap | 98.78 | kJ/mol | Joback Method |
| log10ws | -9.60 | | Crippen Method |
| logp | 8.280 | | Crippen Method |
| mcvol | 396.500 | ml/mol | McGowan Method |
| pc | 821.01 | kPa | Joback Method |
| rinpol | 3200.00 | | NIST Webbook |
| tb | 1023.84 | K | Joback Method |
| tc | 1257.25 | K | Joback Method |
| tf | 573.58 | K | Joback Method |
| vc | 1.538 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 1369.87 | J/molxK | 1023.84 | Joback Method |
| cpg | 1443.15 | J/molxK | 1218.35 | Joback Method |
| cpg | 1431.78 | J/molxK | 1179.45 | Joback Method |
| cpg | 1418.83 | J/molxK | 1140.55 | Joback Method |
| cpg | 1404.24 | J/molxK | 1101.64 | Joback Method |
| cpg | 1387.95 | J/molxK | 1062.74 | Joback Method |
| cpg | 1453.02 | J/molxK | 1257.25 | Joback Method |
| dvisc | 0.0000152 | Paxs | 1023.84 | Joback Method |
| dvisc | 0.0000203 | Paxs | 948.80 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0000283 | Paxs | 873.75 | Joback Method |
| dvisc | 0.0000422 | Paxs | 798.71 | Joback Method |
| dvisc | 0.0000682 | Paxs | 723.67 | Joback Method |
| dvisc | 0.0001234 | Paxs | 648.62 | Joback Method |
| dvisc | 0.0002604 | Paxs | 573.58 | Joback Method |

Sources

| | |
|------------------------|---|
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U356297&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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<https://www.chemeo.com/cid/18-154-1/Terephthalic-acid-2-heptyl-tridecyl-ester.pdf>

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