

Phthalic acid, propyl 2-propylpentyl ester

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|-----------------------------|---|
| Inchi: | InChI=1S/C19H28O4/c1-4-9-15(10-5-2)14-23-19(21)17-12-8-7-11-16(17)18(20)22-13-6-3 |
| InchiKey: | UXNAJCYEFNCVQD-UHFFFAOYSA-N |
| Formula: | C19H28O4 |
| SMILES: | CCCOC(=O)c1ccccc1C(=O)OCC(CCC)CCC |
| Mol. weight [g/mol]: | 320.42 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -258.40 | kJ/mol | Joback Method |
| hf | -705.31 | kJ/mol | Joback Method |
| hfus | 40.67 | kJ/mol | Joback Method |
| hvap | 78.75 | kJ/mol | Joback Method |
| log10ws | -5.48 | | Crippen Method |
| logp | 4.627 | | Crippen Method |
| mvol | 269.690 | ml/mol | McGowan Method |
| pc | 1450.14 | kPa | Joback Method |
| rinpol | 2148.00 | | NIST Webbook |
| rinpol | 2148.00 | | NIST Webbook |
| tb | 817.92 | K | Joback Method |
| tc | 1019.49 | K | Joback Method |
| tf | 472.15 | K | Joback Method |
| vc | 1.034 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 819.59 | J/molxK | 817.92 | Joback Method |
| cpg | 835.66 | J/molxK | 851.51 | Joback Method |
| cpg | 850.61 | J/molxK | 885.11 | Joback Method |
| cpg | 864.46 | J/molxK | 918.70 | Joback Method |
| cpg | 877.24 | J/molxK | 952.30 | Joback Method |
| cpg | 888.96 | J/molxK | 985.89 | Joback Method |
| cpg | 899.64 | J/molxK | 1019.49 | Joback Method |
| dvisc | 0.0007448 | Paxs | 472.15 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0003857 | Paxs | 529.78 | Joback Method |
| dvisc | 0.0002273 | Paxs | 587.41 | Joback Method |
| dvisc | 0.0001472 | Paxs | 645.03 | Joback Method |
| dvisc | 0.0001024 | Paxs | 702.66 | Joback Method |
| dvisc | 0.0000752 | Paxs | 760.29 | Joback Method |
| dvisc | 0.0000577 | Paxs | 817.92 | 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=U377918&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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<https://www.chemeo.com/cid/25-041-8/Phthalic-acid-propyl-2-propylpentyl-ester.pdf>

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