

5-Phenylvaleric acid, heptadecyl ester

| | |
|-----------------------------|-----------------------------------------------------------------------------------|
| Inchi: | InChI=1S/C28H48O2/c1-2-3-4-5-6-7-8-9-10-11-12-13-14-15-21-26-30-28(29)25-20-19-24 |
| InchiKey: | OKDGRPXPSAHHRS-UHFFFAOYSA-N |
| Formula: | C28H48O2 |
| SMILES: | CCCCCCCCCCCCCCCCOC(=O)CCCCc1ccccc1 |
| Mol. weight [g/mol]: | 416.68 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | 63.37 | kJ/mol | Joback Method |
| hf | -629.52 | kJ/mol | Joback Method |
| hfus | 65.10 | kJ/mol | Joback Method |
| hvap | 89.35 | kJ/mol | Joback Method |
| log10ws | -9.51 | | Crippen Method |
| logp | 8.814 | | Crippen Method |
| mvol | 389.060 | ml/mol | McGowan Method |
| pc | 802.97 | kPa | Joback Method |
| rinpol | 3130.00 | | NIST Webbook |
| rinpol | 3130.00 | | NIST Webbook |
| tb | 943.01 | K | Joback Method |
| tc | 1155.06 | K | Joback Method |
| tf | 503.90 | K | Joback Method |
| vc | 1.520 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 1320.09 | J/molxK | 943.01 | Joback Method |
| cpg | 1410.47 | J/molxK | 1119.72 | Joback Method |
| cpg | 1394.93 | J/molxK | 1084.38 | Joback Method |
| cpg | 1378.20 | J/molxK | 1049.03 | Joback Method |
| cpg | 1360.20 | J/molxK | 1013.69 | Joback Method |
| cpg | 1340.86 | J/molxK | 978.35 | Joback Method |
| cpg | 1424.89 | J/molxK | 1155.06 | Joback Method |
| dvisc | 0.0000237 | Paxs | 943.01 | Joback Method |

| | | | | |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0000319 | Paxs | 869.82 | Joback Method |
| dvisc | 0.0000454 | Paxs | 796.64 | Joback Method |
| dvisc | 0.0000695 | Paxs | 723.45 | Joback Method |
| dvisc | 0.0001170 | Paxs | 650.27 | Joback Method |
| dvisc | 0.0002249 | Paxs | 577.08 | Joback Method |
| dvisc | 0.0005223 | Paxs | 503.90 | Joback Method |

Sources

| | |
|------------------------|-------------------------------------------------------------------------------------------------------------------------------------------|
| 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 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U406086&Units=SI |

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/87-818-8/5-Phenylvaleric-acid-heptadecyl-ester.pdf>

Generated by Cheméo on 2024-04-26 17:39:31.225131082 +0000 UTC m=+16442420.145708402.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.