

Pimelic acid, heptyl octyl ester

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|-----------------------------|--|
| Inchi: | InChI=1S/C22H42O4/c1-3-5-7-9-11-16-20-26-22(24)18-14-12-13-17-21(23)25-19-15-10- |
| InchiKey: | RHEBDLAIZPJQKQ-UHFFFAOYSA-N |
| Formula: | C22H42O4 |
| SMILES: | CCCCCCCCOC(=O)CCCCC(=O)OCCCCCCC |
| Mol. weight [g/mol]: | 370.57 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -333.48 | kJ/mol | Joback Method |
| hf | -987.01 | kJ/mol | Joback Method |
| hfus | 58.31 | kJ/mol | Joback Method |
| hvap | 82.88 | kJ/mol | Joback Method |
| log10ws | -6.76 | | Crippen Method |
| logp | 6.354 | | Crippen Method |
| mvol | 335.720 | ml/mol | McGowan Method |
| pc | 953.78 | kPa | Joback Method |
| rinpol | 2565.00 | | NIST Webbook |
| rinpol | 2565.00 | | NIST Webbook |
| tb | 855.34 | K | Joback Method |
| tc | 1047.20 | K | Joback Method |
| tf | 482.02 | K | Joback Method |
| vc | 1.315 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 1092.06 | J/molxK | 855.34 | Joback Method |
| cpg | 1176.14 | J/molxK | 1015.23 | Joback Method |
| cpg | 1161.63 | J/molxK | 983.25 | Joback Method |
| cpg | 1145.99 | J/molxK | 951.27 | Joback Method |
| cpg | 1129.19 | J/molxK | 919.29 | Joback Method |
| cpg | 1111.23 | J/molxK | 887.32 | Joback Method |
| cpg | 1189.55 | J/molxK | 1047.20 | Joback Method |
| dvisc | 0.0000397 | Paxs | 855.34 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0000529 | Paxs | 793.12 | Joback Method |
| dvisc | 0.0000739 | Paxs | 730.90 | Joback Method |
| dvisc | 0.0001100 | Paxs | 668.68 | Joback Method |
| dvisc | 0.0001775 | Paxs | 606.46 | Joback Method |
| dvisc | 0.0003195 | Paxs | 544.24 | Joback Method |
| dvisc | 0.0006693 | Paxs | 482.02 | 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=U393873&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307l |
| Crippen Method: | https://www.chemeo.com/doc/models/crippen_log10ws |

Legend

| | |
|----------------------------|---|
| cp_g: | 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: | Log ₁₀ of Water solubility in mol/l |
| logp: | Octanol/Water partition coefficient |
| m_{cvol}: | 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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