

Glutaric acid, hexyl 3-phenylpropyl ester

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|-----------------------------|---|
| Inchi: | InChI=1S/C20H30O4/c1-2-3-4-8-16-23-19(21)14-9-15-20(22)24-17-10-13-18-11-6-5-7-12 |
| InchiKey: | PWUCASNGNDJJOM-UHFFFAOYSA-N |
| Formula: | C20H30O4 |
| SMILES: | CCCCCOC(=O)CCCC(=O)OCCc1ccccc1 |
| Mol. weight [g/mol]: | 334.45 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -237.91 | kJ/mol | Joback Method |
| hf | -709.20 | kJ/mol | Joback Method |
| hfus | 47.17 | kJ/mol | Joback Method |
| hvap | 80.70 | kJ/mol | Joback Method |
| log10ws | -5.02 | | Crippen Method |
| logp | 4.456 | | Crippen Method |
| mvol | 283.780 | ml/mol | McGowan Method |
| pc | 1356.63 | kPa | Joback Method |
| rinpol | 2544.00 | | NIST Webbook |
| rinpol | 2544.00 | | NIST Webbook |
| tb | 836.26 | K | Joback Method |
| tc | 1035.40 | K | Joback Method |
| tf | 485.90 | K | Joback Method |
| vc | 1.095 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 878.33 | J/molxK | 836.26 | Joback Method |
| cpg | 894.55 | J/molxK | 869.45 | Joback Method |
| cpg | 909.64 | J/molxK | 902.64 | Joback Method |
| cpg | 923.63 | J/molxK | 935.83 | Joback Method |
| cpg | 936.55 | J/molxK | 969.02 | Joback Method |
| cpg | 948.42 | J/molxK | 1002.21 | Joback Method |
| cpg | 959.28 | J/molxK | 1035.40 | Joback Method |
| dvisc | 0.0006973 | Paxs | 485.90 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0003625 | Paxs | 544.29 | Joback Method |
| dvisc | 0.0002139 | Paxs | 602.69 | Joback Method |
| dvisc | 0.0001385 | Paxs | 661.08 | Joback Method |
| dvisc | 0.0000963 | Paxs | 719.47 | Joback Method |
| dvisc | 0.0000707 | Paxs | 777.87 | Joback Method |
| dvisc | 0.0000542 | Paxs | 836.26 | Joback Method |

Sources

| | |
|------------------------|---|
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U360135&Units=SI |
| 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 |

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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