

Glutaric acid, 2-ethylphenyl nonyl ester

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|-----------------------------|--|
| Inchi: | InChI=1S/C22H34O4/c1-3-5-6-7-8-9-12-18-25-21(23)16-13-17-22(24)26-20-15-11-10-14 |
| InchiKey: | VVNBKHUXWSOGNDN-UHFFFAOYSA-N |
| Formula: | C22H34O4 |
| SMILES: | CCCCCCCCCOC(=O)CCCC(=O)Oc1ccccc1CC |
| Mol. weight [g/mol]: | 362.50 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -230.70 | kJ/mol | Joback Method |
| hf | -761.95 | kJ/mol | Joback Method |
| hfus | 51.96 | kJ/mol | Joback Method |
| hvap | 85.82 | kJ/mol | Joback Method |
| log10ws | -6.47 | | Crippen Method |
| logp | 5.619 | | Crippen Method |
| mcvol | 311.960 | ml/mol | McGowan Method |
| pc | 1171.22 | kPa | Joback Method |
| rinpol | 2676.00 | | NIST Webbook |
| tb | 887.00 | K | Joback Method |
| tc | 1090.20 | K | Joback Method |
| tf | 520.96 | K | Joback Method |
| vc | 1.208 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 997.29 | J/molxK | 887.00 | Joback Method |
| cpg | 1013.85 | J/molxK | 920.87 | Joback Method |
| cpg | 1029.18 | J/molxK | 954.73 | Joback Method |
| cpg | 1043.31 | J/molxK | 988.60 | Joback Method |
| cpg | 1056.27 | J/molxK | 1022.47 | Joback Method |
| cpg | 1068.08 | J/molxK | 1056.33 | Joback Method |
| cpg | 1078.77 | J/molxK | 1090.20 | Joback Method |
| dvisc | 0.0004813 | Paxs | 520.96 | Joback Method |
| dvisc | 0.0002579 | Paxs | 581.97 | Joback Method |

| | | | | |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0001556 | Paxs | 642.97 | Joback Method |
| dvisc | 0.0001024 | Paxs | 703.98 | Joback Method |
| dvisc | 0.0000721 | Paxs | 764.99 | Joback Method |
| dvisc | 0.0000534 | Paxs | 825.99 | Joback Method |
| dvisc | 0.0000413 | Paxs | 887.00 | Joback Method |

Sources

| | |
|------------------------|---|
| 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=U358509&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307l |

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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