

Glutaric acid, 2-ethylhexyl 2,6-dimethylnon-1-en-3-yn-5-yl ester

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|----------------------|--|
| Inchi: | InChI=1S/C24H40O4/c1-7-10-13-21(9-3)18-27-23(25)14-11-15-24(26)28-22(17-16-19(4) |
| InchiKey: | TYHBDRKGLMFQSE-UHFFFAOYSA-N |
| Formula: | C24H40O4 |
| SMILES: | C=C(C)C#CC(OC(=O)CCCC(=O)OCC(CC)CCCC)C(C)CCC |
| Mol. weight [g/mol]: | 392.57 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -41.87 | kJ/mol | Joback Method |
| hf | -656.19 | kJ/mol | Joback Method |
| hfus | 53.45 | kJ/mol | Joback Method |
| hvap | 87.73 | kJ/mol | Joback Method |
| log10ws | -6.87 | | Crippen Method |
| logp | 5.844 | | Crippen Method |
| mcvol | 351.000 | ml/mol | McGowan Method |
| pc | 977.17 | kPa | Joback Method |
| rinpol | 2398.00 | | NIST Webbook |
| rinpol | 2398.00 | | NIST Webbook |
| tb | 905.34 | K | Joback Method |
| tc | 1110.27 | K | Joback Method |
| tf | 549.94 | K | Joback Method |
| vc | 1.353 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|---------|---------|-----------------|---------------|
| cpg | 1136.81 | J/mol×K | 905.34 | Joback Method |
| cpg | 1155.08 | J/mol×K | 939.50 | Joback Method |
| cpg | 1172.05 | J/mol×K | 973.65 | Joback Method |
| cpg | 1187.74 | J/mol×K | 1007.81 | Joback Method |
| cpg | 1202.19 | J/mol×K | 1041.96 | Joback Method |
| cpg | 1215.43 | J/mol×K | 1076.12 | Joback Method |
| cpg | 1227.50 | J/mol×K | 1110.27 | Joback Method |

Sources

| | |
|------------------------|---|
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U393966&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 |
| gf: | Standard Gibbs free energy of formation |
| hf: | Enthalpy of formation at standard conditions |
| hfus: | Enthalpy of fusion at standard conditions |
| hvpap: | 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 |
| rinppl: | Non-polar retention indices |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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