

Glutaric acid, (2-methylcyclohex-1-enyl)methyl hept-2-yl

Inchi:
ester

InChI=1S/C20H34O4/c1-4-5-6-11-17(3)24-20(22)14-9-13-19(21)23-15-18-12-8-7-10-16(

InchiKey:

CWMNUSXCTVHAGX-UHFFFAOYSA-N

Formula:

C20H34O4

SMILES:

CCCCC(C)OC(=O)CCCC(=O)OCC1=C(C)CCCC1

Mol. weight [g/mol]:

338.48

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -309.90 | kJ/mol | Joback Method |
| hf | -841.51 | kJ/mol | Joback Method |
| hfus | 40.82 | kJ/mol | Joback Method |
| hvap | 80.39 | kJ/mol | Joback Method |
| log10ws | -5.78 | | Crippen Method |
| logp | 5.102 | | Crippen Method |
| mcvol | 292.380 | ml/mol | McGowan Method |
| pc | 1280.99 | kPa | Joback Method |
| rinqol | 2297.00 | | NIST Webbook |
| tb | 842.48 | K | Joback Method |
| tc | 1042.81 | K | Joback Method |
| tf | 481.90 | K | Joback Method |
| vc | 1.117 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 934.80 | J/molxK | 842.48 | Joback Method |
| cpg | 952.54 | J/molxK | 875.87 | Joback Method |
| cpg | 969.04 | J/molxK | 909.26 | Joback Method |
| cpg | 984.32 | J/molxK | 942.64 | Joback Method |
| cpg | 998.39 | J/molxK | 976.03 | Joback Method |
| cpg | 1011.27 | J/molxK | 1009.42 | Joback Method |
| cpg | 1022.99 | J/molxK | 1042.81 | Joback Method |
| dvisc | 0.0007030 | Paxs | 481.90 | Joback Method |
| dvisc | 0.0003412 | Paxs | 542.00 | Joback Method |

| | | | | |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0001913 | Paxs | 602.09 | Joback Method |
| dvisc | 0.0001192 | Paxs | 662.19 | Joback Method |
| dvisc | 0.0000803 | Paxs | 722.29 | Joback Method |
| dvisc | 0.0000575 | Paxs | 782.38 | Joback Method |
| dvisc | 0.0000432 | Paxs | 842.48 | 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=U405504&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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