

Glutaric acid, 3-bromobenzyl butyl ester

| | |
|-----------------------------|---|
| Inchi: | InChI=1S/C16H21BrO4/c1-2-3-10-20-15(18)8-5-9-16(19)21-12-13-6-4-7-14(17)11-13/h4, |
| InchiKey: | YHYNOBQASBYBNQ-UHFFFAOYSA-N |
| Formula: | C16H21BrO4 |
| SMILES: | CCCCOC(=O)CCCC(=O)OCc1cccc(Br)c1 |
| Mol. weight [g/mol]: | 357.24 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -266.90 | kJ/mol | Joback Method |
| hf | -611.78 | kJ/mol | Joback Method |
| hfus | 41.71 | kJ/mol | Joback Method |
| hvap | 78.90 | kJ/mol | Joback Method |
| log10ws | -5.01 | | Crippen Method |
| logp | 4.006 | | Crippen Method |
| mcvol | 244.920 | ml/mol | McGowan Method |
| pc | 1928.74 | kPa | Joback Method |
| rinpola | 2628.00 | | NIST Webbook |
| rinpola | 2628.00 | | NIST Webbook |
| tb | 815.88 | K | Joback Method |
| tc | 1027.86 | K | Joback Method |
| tf | 513.14 | K | Joback Method |
| vc | 0.933 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 687.61 | J/molxK | 815.88 | Joback Method |
| cpg | 701.14 | J/molxK | 851.21 | Joback Method |
| cpg | 713.65 | J/molxK | 886.54 | Joback Method |
| cpg | 725.19 | J/molxK | 921.87 | Joback Method |
| cpg | 735.77 | J/molxK | 957.20 | Joback Method |
| cpg | 745.41 | J/molxK | 992.53 | Joback Method |
| cpg | 754.13 | J/molxK | 1027.86 | Joback Method |
| dvisc | 0.0005914 | Paxs | 513.14 | Joback Method |

| | | | | |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0003615 | Paxs | 563.60 | Joback Method |
| dvisc | 0.0002396 | Paxs | 614.05 | Joback Method |
| dvisc | 0.0001691 | Paxs | 664.51 | Joback Method |
| dvisc | 0.0001253 | Paxs | 714.97 | Joback Method |
| dvisc | 0.0000966 | Paxs | 765.42 | Joback Method |
| dvisc | 0.0000769 | Paxs | 815.88 | 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=U378001&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci9903071 |
| Crippen Method: | https://www.chemeo.com/doc/models/crippen_log10ws |

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 |

Latest version available from:

<https://www.chemeo.com/cid/114-755-7/Glutaric-acid-3-bromobenzyl-butyl-ester.pdf>

Generated by Cheméo on 2024-05-02 22:07:18.822927831 +0000 UTC m=+16976887.743505143.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.