

Glutaric acid, 4-bromobenzyl pentyl ester

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|-----------------------------|----------------------------------------------------------------------------------|
| Inchi: | InChI=1S/C17H23BrO4/c1-2-3-4-12-21-16(19)6-5-7-17(20)22-13-14-8-10-15(18)11-9-14 |
| InchiKey: | SBDVRDKJPWAYEK-UHFFFAOYSA-N |
| Formula: | C17H23BrO4 |
| SMILES: | CCCCCOC(=O)CCCC(=O)OCc1ccc(Br)cc1 |
| Mol. weight [g/mol]: | 371.27 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -258.48 | kJ/mol | Joback Method |
| hf | -632.42 | kJ/mol | Joback Method |
| hfus | 44.30 | kJ/mol | Joback Method |
| hvap | 81.12 | kJ/mol | Joback Method |
| log10ws | -5.42 | | Crippen Method |
| logp | 4.396 | | Crippen Method |
| mvol | 259.010 | ml/mol | McGowan Method |
| pc | 1775.84 | kPa | Joback Method |
| rinpol | 2684.00 | | NIST Webbook |
| rinpol | 2684.00 | | NIST Webbook |
| tb | 838.76 | K | Joback Method |
| tc | 1049.57 | K | Joback Method |
| tf | 524.41 | K | Joback Method |
| vc | 0.990 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 744.01 | J/molxK | 838.76 | Joback Method |
| cpg | 802.83 | J/molxK | 1014.44 | Joback Method |
| cpg | 793.04 | J/molxK | 979.30 | Joback Method |
| cpg | 782.28 | J/molxK | 944.17 | Joback Method |
| cpg | 770.55 | J/molxK | 909.03 | Joback Method |
| cpg | 757.80 | J/molxK | 873.90 | Joback Method |
| cpg | 811.68 | J/molxK | 1049.57 | Joback Method |
| dvisc | 0.0000667 | Paxs | 838.76 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0000841 | Paxs | 786.37 | Joback Method |
| dvisc | 0.0001096 | Paxs | 733.98 | Joback Method |
| dvisc | 0.0001487 | Paxs | 681.59 | Joback Method |
| dvisc | 0.0002123 | Paxs | 629.19 | Joback Method |
| dvisc | 0.0003233 | Paxs | 576.80 | Joback Method |
| dvisc | 0.0005356 | Paxs | 524.41 | Joback Method |

Sources

| | |
|------------------------|-------------------------------------------------------------------------------------------------------------------------------------------|
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U377532&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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