

4-Iodophenol, isoBOC

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| Inchi: | InChI=1S/C11H13IO3/c1-8(2)7-14-11(13)15-10-5-3-9(12)4-6-10/h3-6,8H,7H2,1-2H3 |
| InchiKey: | KZWDGEAZQPAOGO-UHFFFAOYSA-N |
| Formula: | C11H13IO3 |
| SMILES: | CC(C)COC(=O)Oc1ccc(I)cc1 |
| Mol. weight [g/mol]: | 320.12 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | -138.72 | kJ/mol | Joback Method |
| hf | -350.74 | kJ/mol | Joback Method |
| hfus | 22.76 | kJ/mol | Joback Method |
| hvap | 63.57 | kJ/mol | Joback Method |
| log10ws | -4.01 | | Crippen Method |
| logp | 3.463 | | Crippen Method |
| mcvol | 181.220 | ml/mol | McGowan Method |
| pc | 2648.83 | kPa | Joback Method |
| rinpol | 1775.00 | | NIST Webbook |
| rinpol | 1775.00 | | NIST Webbook |
| tb | 674.15 | K | Joback Method |
| tc | 912.43 | K | Joback Method |
| tf | 390.12 | K | Joback Method |
| vc | 0.667 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 420.27 | J/molxK | 674.15 | Joback Method |
| cpg | 476.49 | J/molxK | 872.71 | Joback Method |
| cpg | 467.04 | J/molxK | 833.00 | Joback Method |
| cpg | 456.72 | J/molxK | 793.29 | Joback Method |
| cpg | 445.49 | J/molxK | 753.58 | Joback Method |
| cpg | 433.35 | J/molxK | 713.86 | Joback Method |
| cpg | 485.07 | J/molxK | 912.43 | Joback Method |
| dvisc | 0.0001321 | Paxs | 674.15 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0001692 | Paxs | 626.81 | Joback Method |
| dvisc | 0.0002258 | Paxs | 579.47 | Joback Method |
| dvisc | 0.0003170 | Paxs | 532.13 | Joback Method |
| dvisc | 0.0004758 | Paxs | 484.80 | Joback Method |
| dvisc | 0.0007795 | Paxs | 437.46 | Joback Method |
| dvisc | 0.0014397 | Paxs | 390.12 | 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=R235202&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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