

2-(2-isopropoxyethoxy)ethanol

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| Inchi: | InChI=1S/C7H16O3/c1-7(2)10-6-5-9-4-3-8/h7-8H,3-6H2,1-2H3 |
| InchiKey: | HRWADRITRNUCIY-UHFFFAOYSA-N |
| Formula: | C7H16O3 |
| SMILES: | CC(C)OCCOCCO |
| Mol. weight [g/mol]: | 148.20 |
| CAS: | 5412-01-1 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | -341.20 | kJ/mol | Joback Method |
| hf | -609.76 | kJ/mol | Joback Method |
| hfus | 16.83 | kJ/mol | Joback Method |
| hvap | 52.29 | kJ/mol | Joback Method |
| log10ws | -0.30 | | Crippen Method |
| logp | 0.420 | | Crippen Method |
| mcvol | 127.100 | ml/mol | McGowan Method |
| pc | 3005.73 | kPa | Joback Method |
| tb | 496.14 | K | Joback Method |
| tc | 659.77 | K | Joback Method |
| tf | 258.93 | K | Joback Method |
| vc | 0.476 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 295.11 | J/molxK | 496.14 | Joback Method |
| cpg | 305.70 | J/molxK | 523.41 | Joback Method |
| cpg | 315.98 | J/molxK | 550.68 | Joback Method |
| cpg | 325.94 | J/molxK | 577.96 | Joback Method |
| cpg | 335.57 | J/molxK | 605.23 | Joback Method |
| cpg | 344.88 | J/molxK | 632.50 | Joback Method |
| cpg | 353.86 | J/molxK | 659.77 | Joback Method |
| dvisc | 0.0271956 | Paxs | 258.93 | Joback Method |
| dvisc | 0.0059483 | Paxs | 298.47 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0018565 | Paxs | 338.00 | Joback Method |
| dvisc | 0.0007395 | Paxs | 377.53 | Joback Method |
| dvisc | 0.0003507 | Paxs | 417.07 | Joback Method |
| dvisc | 0.0001893 | Paxs | 456.61 | Joback Method |
| dvisc | 0.0001127 | Paxs | 496.14 | 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=C5412011&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 |
| mccvol: | McGowan's characteristic volume |
| pc: | Critical Pressure |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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