

1-Mercaptopentyl-2-acetate

| | |
|-----------------------------|--|
| Inchi: | InChI=1S/C7H14O2S/c1-3-4-7(5-10)9-6(2)8/h7,10H,3-5H2,1-2H3 |
| InchiKey: | AYVKTHMRZAYNPJ-UHFFFAOYSA-N |
| Formula: | C7H14O2S |
| SMILES: | CCCC(CS)OC(C)=O |
| Mol. weight [g/mol]: | 162.25 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | -198.91 | kJ/mol | Joback Method |
| hf | -399.41 | kJ/mol | Joback Method |
| hfus | 17.19 | kJ/mol | Joback Method |
| hvap | 46.68 | kJ/mol | Joback Method |
| log10ws | -1.80 | | Crippen Method |
| logp | 1.648 | | Crippen Method |
| mcvol | 133.280 | ml/mol | McGowan Method |
| pc | 3184.73 | kPa | Joback Method |
| tb | 498.27 | K | Joback Method |
| tc | 699.54 | K | Joback Method |
| tf | 262.27 | K | Joback Method |
| vc | 0.499 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 285.82 | J/molxK | 498.27 | Joback Method |
| cpg | 297.93 | J/molxK | 531.81 | Joback Method |
| cpg | 309.52 | J/molxK | 565.36 | Joback Method |
| cpg | 320.59 | J/molxK | 598.90 | Joback Method |
| cpg | 331.14 | J/molxK | 632.45 | Joback Method |
| cpg | 341.18 | J/molxK | 665.99 | Joback Method |
| cpg | 350.72 | J/molxK | 699.54 | Joback Method |

Sources

| | |
|------------------------|---|
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=R611646&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307I |
| Crippen Method: | https://www.chemeo.com/doc/models/crippen_log10ws |
| Joback Method: | https://en.wikipedia.org/wiki/Joback_method |

Legend

| | |
|-----------------|---|
| cpg: | Ideal gas heat capacity |
| 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 |
| 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/96-942-0/1-Mercaptopentyl-2-acetate.pdf>

Generated by Cheméo on 2024-05-01 10:12:34.252382135 +0000 UTC m=+16847603.172959448.

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