

(2,5-Dimethoxyphenyl)acetone

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| Other names: | 2-Propanone, (2,5-dimethoxy)phenyl- 2-Propanone, 1-(2,5-dimethoxyphenyl)- |
| Inchi: | InChI=1S/C11H14O3/c1-8(12)6-9-7-10(13-2)4-5-11(9)14-3/h4-5,7H,6H2,1-3H3 |
| InchiKey: | RUCPISZJMCQMPY-UHFFFAOYSA-N |
| Formula: | C11H14O3 |
| SMILES: | COc1ccc(OC)c(CC(C)=O)c1 |
| Mol. weight [g/mol]: | 194.23 |
| CAS: | 14293-24-4 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -204.03 | kJ/mol | Joback Method |
| hf | -433.80 | kJ/mol | Joback Method |
| hfus | 21.48 | kJ/mol | Joback Method |
| hvap | 55.25 | kJ/mol | Joback Method |
| log10ws | -2.21 | | Crippen Method |
| logp | 1.835 | | Crippen Method |
| mvol | 155.400 | ml/mol | McGowan Method |
| pc | 2646.11 | kPa | Joback Method |
| tb | 586.43 | K | Joback Method |
| tc | 795.52 | K | Joback Method |
| tf | 359.58 | K | Joback Method |
| vc | 0.586 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 370.29 | J/molxK | 586.43 | Joback Method |
| cpg | 383.95 | J/molxK | 621.28 | Joback Method |
| cpg | 396.95 | J/molxK | 656.13 | Joback Method |
| cpg | 409.28 | J/molxK | 690.97 | Joback Method |
| cpg | 420.93 | J/molxK | 725.82 | Joback Method |
| cpg | 431.89 | J/molxK | 760.67 | Joback Method |
| cpg | 442.17 | J/molxK | 795.52 | Joback Method |

| | | | | |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0010431 | Paxs | 359.58 | Joback Method |
| dvisc | 0.0006535 | Paxs | 397.39 | Joback Method |
| dvisc | 0.0004440 | Paxs | 435.20 | Joback Method |
| dvisc | 0.0003209 | Paxs | 473.00 | Joback Method |
| dvisc | 0.0002434 | Paxs | 510.81 | Joback Method |
| dvisc | 0.0001918 | Paxs | 548.62 | Joback Method |
| dvisc | 0.0001558 | Paxs | 586.43 | Joback Method |

Sources

| | |
|------------------------|---|
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=C14293244&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 |

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 |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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