

4-(2,3,4-trimethylphenyl)-2-butanone

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|-----------------------------|--|
| Inchi: | InChI=1S/C13H18O/c1-9-5-7-13(8-6-10(2)14)12(4)11(9)3/h5,7H,6,8H2,1-4H3 |
| InchiKey: | WOYNOSTWAOGOET-UHFFFAOYSA-N |
| Formula: | C13H18O |
| SMILES: | CC(=O)CCc1ccc(C)c(C)c1C |
| Mol. weight [g/mol]: | 190.28 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | 13.18 | kJ/mol | Joback Method |
| hf | -222.11 | kJ/mol | Joback Method |
| hfus | 23.90 | kJ/mol | Joback Method |
| hvap | 55.54 | kJ/mol | Joback Method |
| log10ws | -3.82 | | Crippen Method |
| logp | 3.133 | | Crippen Method |
| mcvol | 171.840 | ml/mol | McGowan Method |
| pc | 2231.30 | kPa | Joback Method |
| ripol | 2258.00 | | NIST Webbook |
| ripol | 2258.00 | | NIST Webbook |
| tb | 592.33 | K | Joback Method |
| tc | 800.18 | K | Joback Method |
| tf | 350.18 | K | Joback Method |
| vc | 0.661 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 416.32 | J/molxK | 592.33 | Joback Method |
| cpg | 431.75 | J/molxK | 626.97 | Joback Method |
| cpg | 446.37 | J/molxK | 661.61 | Joback Method |
| cpg | 460.21 | J/molxK | 696.26 | Joback Method |
| cpg | 473.29 | J/molxK | 730.90 | Joback Method |
| cpg | 485.63 | J/molxK | 765.54 | Joback Method |
| cpg | 497.26 | J/molxK | 800.18 | Joback Method |
| dvisc | 0.0013712 | Paxs | 350.18 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0008401 | Paxs | 390.54 | Joback Method |
| dvisc | 0.0005641 | Paxs | 430.90 | Joback Method |
| dvisc | 0.0004056 | Paxs | 471.25 | Joback Method |
| dvisc | 0.0003072 | Paxs | 511.61 | Joback Method |
| dvisc | 0.0002423 | Paxs | 551.97 | Joback Method |
| dvisc | 0.0001974 | Paxs | 592.33 | Joback Method |

Sources

| | |
|------------------------|---|
| 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 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=R494387&Units=SI |

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

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