

Octane, 2,3,7-trimethyl-

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|----------------------|---|
| Inchi: | InChI=1S/C11H24/c1-9(2)7-6-8-11(5)10(3)4/h9-11H,6-8H2,1-5H3 |
| InchiKey: | XJKKSYAVEVAGFX-UHFFFAOYSA-N |
| Formula: | C11H24 |
| SMILES: | CC(C)CCCC(C)C(C)C |
| Mol. weight [g/mol]: | 156.31 |
| CAS: | 62016-34-6 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | 34.42 | kJ/mol | Joback Method |
| hf | -286.21 | kJ/mol | Joback Method |
| hfus | 13.68 | kJ/mol | Joback Method |
| hvap | 38.92 | kJ/mol | Joback Method |
| log10ws | -3.70 | | Crippen Method |
| logp | 4.105 | | Crippen Method |
| mcvol | 165.850 | ml/mol | McGowan Method |
| pc | 1975.31 | kPa | Joback Method |
| tb | 449.76 | K | Joback Method |
| tc | 621.98 | K | Joback Method |
| tf | 168.73 | K | Joback Method |
| vc | 0.633 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 360.14 | J/molxK | 449.76 | Joback Method |
| cpg | 439.69 | J/molxK | 593.28 | Joback Method |
| cpg | 425.08 | J/molxK | 564.57 | Joback Method |
| cpg | 409.83 | J/molxK | 535.87 | Joback Method |
| cpg | 393.94 | J/molxK | 507.17 | Joback Method |
| cpg | 377.38 | J/molxK | 478.46 | Joback Method |
| cpg | 453.69 | J/molxK | 621.98 | Joback Method |
| dvisc | 0.0001963 | Paxs | 449.76 | Joback Method |
| dvisc | 0.0002863 | Paxs | 402.92 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0004611 | Paxs | 356.08 | Joback Method |
| dvisc | 0.0008579 | Paxs | 309.25 | Joback Method |
| dvisc | 0.0019921 | Paxs | 262.41 | Joback Method |
| dvisc | 0.0066714 | Paxs | 215.57 | Joback Method |
| dvisc | 0.0437057 | Paxs | 168.73 | Joback Method |

Sources

| | |
|------------------------|---|
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=C62016346&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307l |
| 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 |
| 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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