

Propane, 2,2-bis(3,4-xylyl)-

| | |
|-----------------------------|--|
| Inchi: | InChI=1S/C19H24/c1-13-7-9-17(11-15(13)3)19(5,6)18-10-8-14(2)16(4)12-18/h7-12H,1-6H |
| InchiKey: | BGTZYQSCNQIZER-UHFFFAOYSA-N |
| Formula: | C19H24 |
| SMILES: | <chem>Cc1ccc(C(C)(C)c2ccc(C)c(C)c2)cc1C</chem> |
| Mol. weight [g/mol]: | 252.39 |
| CAS: | 6970-01-0 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | 298.24 | kJ/mol | Joback Method |
| hf | -17.06 | kJ/mol | Joback Method |
| hfus | 24.08 | kJ/mol | Joback Method |
| hvap | 63.79 | kJ/mol | Joback Method |
| log10ws | -5.99 | | Crippen Method |
| logp | 5.246 | | Crippen Method |
| mcvol | 231.050 | ml/mol | McGowan Method |
| pc | 1706.12 | kPa | Joback Method |
| tb | 704.17 | K | Joback Method |
| tc | 937.43 | K | Joback Method |
| tf | 409.23 | K | Joback Method |
| vc | 0.873 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 633.49 | J/molxK | 704.17 | Joback Method |
| cpg | 718.25 | J/molxK | 898.56 | Joback Method |
| cpg | 703.65 | J/molxK | 859.68 | Joback Method |
| cpg | 687.96 | J/molxK | 820.80 | Joback Method |
| cpg | 671.09 | J/molxK | 781.92 | Joback Method |
| cpg | 652.96 | J/molxK | 743.05 | Joback Method |
| cpg | 731.84 | J/molxK | 937.43 | Joback Method |
| dvisc | 0.0000922 | Paxs | 704.17 | Joback Method |
| dvisc | 0.0001166 | Paxs | 655.01 | Joback Method |

| | | | | |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0001532 | Paxs | 605.86 | Joback Method |
| dvisc | 0.0002112 | Paxs | 556.70 | Joback Method |
| dvisc | 0.0003097 | Paxs | 507.54 | Joback Method |
| dvisc | 0.0004932 | Paxs | 458.39 | Joback Method |
| dvisc | 0.0008782 | Paxs | 409.23 | 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=C6970010&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci9903071 |

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 |

Latest version available from:

<https://www.chemeo.com/cid/19-425-9/Propane-2-2-bis-3-4-xylyl.pdf>

Generated by Cheméo on 2024-04-25 21:17:06.577938421 +0000 UTC m=+16369075.498515733.

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