

Naphtho[2,1-c]picene

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|-----------------------------|-----------------------------------------------------------------------------------|
| Inchi: | InChI=1S/C30H18/c1-3-7-22-19(5-1)11-13-25-24(22)15-16-27-26(25)17-18-29-28(27)14- |
| InchiKey: | RFYQCYCUNLXVDB-UHFFFAOYSA-N |
| Formula: | C30H18 |
| SMILES: | <chem>c1ccc2c(c1)ccc1c2ccc2c1ccc1c2ccc2ccc3ccccc3c21</chem> |
| Mol. weight [g/mol]: | 378.46 |
| CAS: | 115747-43-8 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | 905.88 | kJ/mol | Joback Method |
| hf | 663.07 | kJ/mol | Joback Method |
| hfus | 47.67 | kJ/mol | Joback Method |
| hvap | 97.80 | kJ/mol | Joback Method |
| log10ws | -12.30 | | Crippen Method |
| logp | 8.606 | | Crippen Method |
| mvol | 293.040 | ml/mol | McGowan Method |
| pc | 1777.34 | kPa | Joback Method |
| tb | 1051.26 | K | Joback Method |
| tc | 1330.47 | K | Joback Method |
| tf | 713.08 | K | Joback Method |
| vc | 1.139 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 909.76 | J/molxK | 1051.26 | Joback Method |
| cpg | 929.78 | J/molxK | 1097.80 | Joback Method |
| cpg | 951.06 | J/molxK | 1144.33 | Joback Method |
| cpg | 974.05 | J/molxK | 1190.87 | Joback Method |
| cpg | 999.18 | J/molxK | 1237.40 | Joback Method |
| cpg | 1026.90 | J/molxK | 1283.94 | Joback Method |
| cpg | 1057.66 | J/molxK | 1330.47 | Joback Method |
| dvisc | 0.0046199 | Paxs | 713.08 | Joback Method |
| dvisc | 0.0041148 | Paxs | 769.44 | Joback Method |

| | | | | |
|-------|-----------|------|---------|---------------|
| dvisc | 0.0037234 | Paxs | 825.81 | Joback Method |
| dvisc | 0.0034125 | Paxs | 882.17 | Joback Method |
| dvisc | 0.0031604 | Paxs | 938.53 | Joback Method |
| dvisc | 0.0029526 | Paxs | 994.90 | Joback Method |
| dvisc | 0.0027786 | Paxs | 1051.26 | Joback Method |

Sources

| | |
|------------------------|-------------------------------------------------------------------------------------------------------------------------------------------------|
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=C115747438&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 |
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |

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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