

exo-1H-Indene, 3a,4,7,7a-tetrahydro-4,7-ethano

| | |
|-----------------------------|---|
| Inchi: | InChI=1S/C11H14/c1-2-10-8-4-6-9(7-5-8)11(10)3-1/h1-2,4,6,8-11H,3,5,7H2/t8-,9+,10-,11+ |
| InchiKey: | MOKDIRQGOJTXJX-YTWAJWBKSA-N |
| Formula: | C11H14 |
| SMILES: | C1=CC2C3C=CC(CC3)C2C1 |
| Mol. weight [g/mol]: | 146.23 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | 252.00 | kJ/mol | Joback Method |
| hf | 30.93 | kJ/mol | Joback Method |
| hfus | 17.97 | kJ/mol | Joback Method |
| hvap | 40.44 | kJ/mol | Joback Method |
| log10ws | -2.85 | | Crippen Method |
| logp | 2.775 | | Crippen Method |
| mcvol | 124.670 | ml/mol | McGowan Method |
| pc | 3103.64 | kPa | Joback Method |
| rinpol | 1142.00 | | NIST Webbook |
| tb | 473.49 | K | Joback Method |
| tc | 697.09 | K | Joback Method |
| tf | 257.79 | K | Joback Method |
| vc | 0.477 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 287.58 | J/molxK | 473.49 | Joback Method |
| cpg | 307.88 | J/molxK | 510.76 | Joback Method |
| cpg | 326.67 | J/molxK | 548.02 | Joback Method |
| cpg | 344.05 | J/molxK | 585.29 | Joback Method |
| cpg | 360.13 | J/molxK | 622.56 | Joback Method |
| cpg | 375.02 | J/molxK | 659.82 | Joback Method |
| cpg | 388.81 | J/molxK | 697.09 | Joback Method |
| dvisc | 0.0007144 | Paxs | 257.79 | Joback Method |
| dvisc | 0.0007907 | Paxs | 293.74 | Joback Method |

| | | | | |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0008560 | Paxs | 329.69 | Joback Method |
| dvisc | 0.0009123 | Paxs | 365.64 | Joback Method |
| dvisc | 0.0009614 | Paxs | 401.59 | Joback Method |
| dvisc | 0.0010043 | Paxs | 437.54 | Joback Method |
| dvisc | 0.0010423 | Paxs | 473.49 | 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=R128060&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307l |

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 |
| rinpol: | Non-polar retention indices |
| 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/37-725-6/exo-1H-Indene-3a-4-7-7a-tetrahydro-4-7-ethano.pdf>

Generated by Cheméo on 2024-04-18 20:33:15.169084425 +0000 UTC m=+15761644.089661740.

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