

2-Naphthamide, N-2-ethylhexyl-

Inchi: InChI=1S/C19H25NO/c1-3-5-8-15(4-2)14-20-19(21)18-12-11-16-9-6-7-10-17(16)13-18/h
InchiKey: SMCUWTMUHGHVVMY-UHFFFAOYSA-N
Formula: C19H25NO
SMILES: CCCCC(CC)CN=C(O)c1ccc2ccccc2c1
Mol. weight [g/mol]: 283.41

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|--------|----------------|
| hf | -104.44 | kJ/mol | Joback Method |
| hvap | 82.15 | kJ/mol | Joback Method |
| log10ws | -5.85 | | Crippen Method |
| logp | 5.361 | | Crippen Method |
| mcvol | 246.900 | ml/mol | McGowan Method |
| pc | 1624.60 | kPa | Joback Method |
| rinsol | 2532.00 | | NIST Webbook |
| tb | 853.06 | K | Joback Method |
| tc | 1067.08 | K | Joback Method |

Sources

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=U407353&Units=SI>
Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci990307l>
Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws

Legend

hf: Enthalpy of formation 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 |
| rinpol: | Non-polar retention indices |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |

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