

Formamide, n-(7-nitrofluoren-2-yl)-

| | |
|-----------------------------|--|
| Inchi: | InChI=1S/C14H10N2O3/c17-8-15-11-1-3-13-9(6-11)5-10-7-12(16(18)19)2-4-14(10)13/h1 |
| InchiKey: | NNVUDKLVHMGGOQ-UHFFFAOYSA-N |
| Formula: | C14H10N2O3 |
| SMILES: | O=CNc1ccc2c(c1)Cc1cc([N+](=O)[O-])ccc1-2 |
| Mol. weight [g/mol]: | 254.24 |
| CAS: | 6583-71-7 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | 371.38 | kJ/mol | Joback Method |
| hf | 157.48 | kJ/mol | Joback Method |
| hfus | 38.55 | kJ/mol | Joback Method |
| hvap | 83.58 | kJ/mol | Joback Method |
| log10ws | -4.86 | | Crippen Method |
| logp | 2.734 | | Crippen Method |
| mcvol | 178.710 | ml/mol | McGowan Method |
| pc | 3272.78 | kPa | Joback Method |
| tb | 846.54 | K | Joback Method |
| tc | 1105.79 | K | Joback Method |
| tf | 617.95 | K | Joback Method |
| vc | 0.712 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 501.05 | J/molxK | 846.54 | Joback Method |
| cpg | 511.50 | J/molxK | 889.75 | Joback Method |
| cpg | 521.35 | J/molxK | 932.96 | Joback Method |
| cpg | 530.75 | J/molxK | 976.17 | Joback Method |
| cpg | 539.87 | J/molxK | 1019.38 | Joback Method |
| cpg | 548.86 | J/molxK | 1062.59 | Joback Method |
| cpg | 557.88 | J/molxK | 1105.79 | Joback Method |

Sources

| | |
|------------------------|---|
| 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 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=C6583717&Units=SI |

Legend

| | |
|-----------------|---|
| cpg: | Ideal gas heat capacity |
| 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/124-100-2/Formamide-n-7-nitrofluoren-2-yl.pdf>

Generated by Cheméo on 2024-04-26 18:52:55.535947177 +0000 UTC m=+16446824.456524498.

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