

2-Bromo-4-methylphenyl isothiocyanate

Inchi: InChI=1S/C8H6BrNS/c1-6-2-3-8(10-5-11)7(9)4-6/h2-4H,1H3
InchiKey: DQXSCOWMTPPFCJ-UHFFFAOYSA-N
Formula: C8H6BrNS
SMILES: Cc1ccc(N=C=S)c(Br)c1
Mol. weight [g/mol]: 228.11
CAS: 19241-39-5

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|--------|----------------|
| hf | 315.54 | kJ/mol | Joback Method |
| hvap | 53.88 | kJ/mol | Joback Method |
| log10ws | -4.03 | | Crippen Method |
| logp | 3.492 | | Crippen Method |
| mcvol | 135.050 | ml/mol | McGowan Method |
| pc | 3935.71 | kPa | Joback Method |
| tb | 631.19 | K | Joback Method |
| tc | 906.91 | 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=C19241395&Units=SI>
Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci9903071>
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 |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |

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