

4-Bromophenyl isocyanate

Other names:	p-Bromophenylisocyanate Isocyanic acid p-bromophenyl ester Benzene, 1-bromo-4-isocyanato- p-Bromocarbanil p-Bromophenyl carbonimide
Inchi:	InChI=1S/C7H4BrNO/c8-6-1-3-7(4-2-6)9-5-10/h1-4H
InchiKey:	CZQIJQFTRGDODI-UHFFFAOYSA-N
Formula:	C7H4BrNO
SMILES:	O=C=Nc1ccc(Br)cc1
Mol. weight [g/mol]:	198.02
CAS:	2493-02-9

Physical Properties

Property code	Value	Unit	Source
hf	58.17	kJ/mol	Joback Method
hvap	50.08	kJ/mol	Joback Method
log10ws	-7.12		Crippen Method
logp	2.416		Crippen Method
mcvol	110.480	ml/mol	McGowan Method
pc	4883.38	kPa	Joback Method
tb	524.05	K	Joback Method
tc	766.95	K	Joback Method

Sources

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307I
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=C2493029&Units=SI

Legend

hf:	Enthalpy of formation at standard conditions
h_{vap}:	Enthalpy of vaporization at standard conditions
log₁₀ws:	Log ₁₀ of Water solubility in mol/l
log_p:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
pc:	Critical Pressure
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature

Latest version available from:

<https://www.cheméo.com/cid/59-966-5/4-Bromophenyl-isocyanate.pdf>

Generated by Cheméo on 2024-04-19 20:17:43.263355834 +0000 UTC m=+15847112.183933156.

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