

(4-Iodo-phenyl)-hydrazine

Inchi:

InChI=1S/C6H7IN2/c7-5-1-3-6(9-8)4-2-5/h1-4,9H,8H2

InchiKey:

IQMLCMKMSBMMGR-UHFFFAOYSA-N

Formula:

C6H7IN2

SMILES:

NNc1ccc(I)cc1

Mol. weight [g/mol]:

234.04

Physical Properties

Property code	Value	Unit	Source
hf	240.80	kJ/mol	Cheméo Relay (1.0)
hfus	19.65	kJ/mol	Joback Method
hvap	76.77	kJ/mol	Cheméo Relay (1.0)
ie	7.64	eV	Cheméo Relay (1.0)
log10ws	-1.94		Cheméo Relay (1.0)
logp	1.577		Crippen Method
mcvol	117.420	ml/mol	McGowan Method
tb	585.00	K	Cheméo Relay (1.0)
tf	374.65	K	Cheméo Relay (1.0)

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	226.76	J/mol×K	584.18	Joback Method
cpg	236.24	J/mol×K	628.57	Joback Method
cpg	244.89	J/mol×K	672.96	Joback Method
cpg	252.78	J/mol×K	717.35	Joback Method
cpg	259.98	J/mol×K	761.74	Joback Method
cpg	266.53	J/mol×K	806.13	Joback Method
cpg	272.52	J/mol×K	850.53	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Cheméo Relay (1.0):	
Cheméo Relay (1.0, beta):	
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C13116273&Units=SI
Joback Method:	https://en.wikipedia.org/wiki/Joback_method

Legend

cpg:	Ideal gas heat capacity
hf:	Enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion at standard conditions
hvap:	Enthalpy of vaporization at standard conditions
ie:	Ionization energy
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
tb:	Normal Boiling Point Temperature
tf:	Normal melting (fusion) point

Latest version available from:

<https://www.chemeo.com/cid/61-074-2/4-Iodo-phenyl-hydrazine.pdf>

Generated by Cheméo on 2026-07-22 02:45:50.949249203 +0000 UTC m=+199446.791134593.

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