

1,2-Benzenediamine, 4-chloro-

Other names:	o-Phenylenediamine, 4-chloro- C.I. 76015 Ursol Olive 6G 1,2-Diamino-4-chlorobenzene 2-Amino-4-chloroaniline 3,4-Diaminobenzene 4-Chloro-o-phenylenediamine 4-Chloro-1,2-benzenediamine 4-Chloro-1,2-diaminobenzene 4-Chloro-1,2-phenylenediamine 4-Chloro-ortho-phenylenediamine NCI-C03292 p-Chloro-o-phenylenediamine 4-Cl-o-PD 3,4-Diamino-1-chlorobenzene 4-Chloro-2-aminoaniline NSC 6157
Inchi:	InChI=1S/C6H7ClN2/c7-4-1-2-5(8)6(9)3-4/h1-3H,8-9H2
InchiKey:	BXIXXXYDDJVHDL-UHFFFAOYSA-N
Formula:	C6H7ClN2
SMILES:	<chem>Nc1ccc(Cl)cc1N</chem>
Mol. weight [g/mol]:	142.59
CAS:	95-83-0

Physical Properties

Property code	Value	Unit	Source
gf	213.76	kJ/mol	Joback Method
hf	98.26	kJ/mol	Joback Method
hfus	19.15	kJ/mol	Joback Method
hvap	58.22	kJ/mol	Joback Method
log10ws	-1.42		Crippen Method
logp	1.504		Crippen Method
mcpvol	103.840	ml/mol	McGowan Method
pc	4987.39	kPa	Joback Method
tb	555.81	K	Joback Method
tc	805.36	K	Joback Method
tf	405.28	K	Joback Method

vc

0.370

m³/kmol

Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	216.19	J/mol×K	555.81	Joback Method
cpg	225.31	J/mol×K	597.40	Joback Method
cpg	233.79	J/mol×K	638.99	Joback Method
cpg	241.64	J/mol×K	680.58	Joback Method
cpg	248.91	J/mol×K	722.17	Joback Method
cpg	255.62	J/mol×K	763.77	Joback Method
cpg	261.79	J/mol×K	805.36	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=C95830&Units=SI>

Crippen Method:

<http://pubs.acs.org/doi/abs/10.1021/ci990307l>

Crippen Method:

https://www.chemeo.com/doc/models/crippen_log10ws

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
mcvol:	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/44-620-4/1-2-Benzenediamine-4-chloro.pdf>

Generated by Cheméo on 2024-04-25 01:33:54.930580044 +0000 UTC m=+16298083.851157359.

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