

3,4-Dichlorofluorobenzene

Other names:	3,4-Dichlorofluorobenzene Benzene, 1,2-dichloro-4-fluoro-
Inchi:	InChI=1S/C6H3Cl2F/c7-5-2-1-4(9)3-6(5)8/h1-3H
InchiKey:	QSDKXMGVRLVIQV-UHFFFAOYSA-N
Formula:	C6H3Cl2F
SMILES:	Fc1ccc(Cl)c(Cl)c1
Mol. weight [g/mol]:	164.99

Physical Properties

Property code	Value	Unit	Source
af	0.3319		Cheméo Relay (1.0)
gf	-125.88	kJ/mol	Joback Method
hf	-165.20	kJ/mol	Cheméo Relay (1.0)
hfus	16.03	kJ/mol	Joback Method
hvap	50.38	kJ/mol	Cheméo Relay (1.0)
ie	9.16 ± 0.02	eV	NIST Webbook
log10ws	-3.10		Cheméo Relay (1.0)
logp	3.132		Crippen Method
mcvol	97.890	ml/mol	McGowan Method
pc	3848.31	kPa	Joback Method
tb	445.20	K	NIST Webbook
tc	680.19	K	Cheméo Relay (1.0)
tf	278.72	K	Cheméo Relay (1.0)

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	157.45	J/mol×K	447.45	Joback Method
cpg	164.82	J/mol×K	484.15	Joback Method
cpg	171.73	J/mol×K	520.85	Joback Method
cpg	178.19	J/mol×K	557.55	Joback Method
cpg	184.23	J/mol×K	594.25	Joback Method
cpg	189.87	J/mol×K	630.94	Joback Method
cpg	195.11	J/mol×K	667.64	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=C1435490&Units=SI
Joback Method:	https://en.wikipedia.org/wiki/Joback_method

Legend

af:	Acentric Factor
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
ie:	Ionization energy
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

Latest version available from:

<https://www.chemeo.com/cid/96-789-1/3-4-Dichlorofluorobenzene.pdf>

Generated by Cheméo on 2026-07-21 17:38:50.679974324 +0000 UTC m=+166626.521859717.

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