

1,4-Benzenediol, 2,5-dichloro-

Other names:	Hydroquinone, 2,5-dichloro- 2,5-Dichloro-p-benzohydroquinone 2,5-Dichloro-p-hydroquinone 2,5-Dichloro-1,4-hydroquinone 2,5-Dichlorohydroquinone 2,5-Dichloro-1,4-benzenediol 2,5-Dichloro-1,4-dihydroxybenzene
Inchi:	InChI=1S/C6H4Cl2O2/c7-3-1-5(9)4(8)2-6(3)10/h1-2,9-10H
InchiKey:	AYNPIRVEWMUJDE-UHFFFAOYSA-N
Formula:	C6H4Cl2O2
SMILES:	Oc1cc(Cl)c(O)cc1Cl
Mol. weight [g/mol]:	179.00
CAS:	824-69-1

Physical Properties

Property code	Value	Unit	Source
chs	-2547.00	kJ/mol	NIST Webbook
chs	-2552.70 ± 8.40	kJ/mol	NIST Webbook
gf	-230.68	kJ/mol	Joback Method
hf	-328.21	kJ/mol	Joback Method
hfs	-437.60	kJ/mol	NIST Webbook
hfus	24.91	kJ/mol	Joback Method
hvap	66.69	kJ/mol	Joback Method
log10ws	-1.94		Crippen Method
logp	2.405		Crippen Method
mcvol	107.860	ml/mol	McGowan Method
pc	6359.24	kPa	Joback Method
tb	604.44	K	Joback Method
tc	861.92	K	Joback Method
tf	479.60	K	Joback Method
vc	0.293	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	220.56	J/mol×K	604.44	Joback Method
cpg	226.43	J/mol×K	647.35	Joback Method
cpg	231.76	J/mol×K	690.27	Joback Method
cpg	236.70	J/mol×K	733.18	Joback Method
cpg	241.40	J/mol×K	776.09	Joback Method
cpg	246.05	J/mol×K	819.00	Joback Method
cpg	250.80	J/mol×K	861.92	Joback Method
dvisc	0.0001179	Pa×s	479.60	Joback Method
dvisc	0.0000687	Pa×s	500.41	Joback Method
dvisc	0.0000418	Pa×s	521.21	Joback Method
dvisc	0.0000264	Pa×s	542.02	Joback Method
dvisc	0.0000173	Pa×s	562.83	Joback Method
dvisc	0.0000117	Pa×s	583.63	Joback Method
dvisc	0.0000081	Pa×s	604.44	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C824691&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Joback Method:	https://en.wikipedia.org/wiki/Joback_method

Legend

chs:	Standard solid enthalpy of combustion
cpg:	Ideal gas heat capacity
dvisc:	Dynamic viscosity
gf:	Standard Gibbs free energy of formation
hf:	Enthalpy of formation at standard conditions
hfs:	Solid phase 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

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