

# Phenol, 4-bromo-2,5-dichloro-

<b>Other names:</b>	Phosvel phenol 2,5-Dichloro-4-bromophenol 4-Bromo-2,5-dichlorophenol Leptophos phenol
<b>Inchi:</b>	InChI=1S/C6H3BrCl2O/c7-3-1-5(9)6(10)2-4(3)8/h1-2,10H
<b>InchiKey:</b>	HWWKEEKUMAZJLL-UHFFFAOYSA-N
<b>Formula:</b>	C6H3BrCl2O
<b>SMILES:</b>	Oc1cc(Cl)c(Br)cc1Cl
<b>Mol. weight [g/mol]:</b>	241.90
<b>CAS:</b>	1940-42-7

## Physical Properties

Property code	Value	Unit	Source
gf	-71.37	kJ/mol	Joback Method
hf	-136.04	kJ/mol	Joback Method
hfus	24.02	kJ/mol	Joback Method
hvap	60.77	kJ/mol	Joback Method
log10ws	-3.55		Crippen Method
logp	3.462		Crippen Method
mcvol	119.490	ml/mol	McGowan Method
pc	5495.11	kPa	Joback Method
tb	594.96	K	Joback Method
tc	857.43	K	Joback Method
tf	343.97 ± 0.20	K	NIST Webbook
vc	0.390	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	209.03	J/molxK	594.96	Joback Method
cpg	214.88	J/molxK	638.71	Joback Method
cpg	220.18	J/molxK	682.45	Joback Method
cpg	225.03	J/molxK	726.20	Joback Method
cpg	229.55	J/molxK	769.94	Joback Method

cpg	233.85	J/mol×K	813.69	Joback Method
cpg	238.05	J/mol×K	857.43	Joback Method
dvisc	0.0003421	Paxs	465.99	Joback Method
dvisc	0.0005619	Paxs	440.20	Joback Method
dvisc	0.0002194	Paxs	491.79	Joback Method
dvisc	0.0001471	Paxs	517.58	Joback Method
dvisc	0.0001024	Paxs	543.37	Joback Method
dvisc	0.0000737	Paxs	569.17	Joback Method
dvisc	0.0000546	Paxs	594.96	Joback Method
hfust	22.11	kJ/mol	343.40	NIST Webbook
hfust	22.11	kJ/mol	343.40	NIST Webbook

## Sources

<b>Crippen Method:</b>	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>
<b>Joback Method:</b>	<a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>
<b>McGowan Method:</b>	<a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>
<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C1940427&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C1940427&amp;Units=SI</a>
<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</a>

## Legend

<b>cpg:</b>	Ideal gas heat capacity
<b>dvisc:</b>	Dynamic viscosity
<b>gf:</b>	Standard Gibbs free energy of formation
<b>hf:</b>	Enthalpy of formation at standard conditions
<b>hfus:</b>	Enthalpy of fusion at standard conditions
<b>hfust:</b>	Enthalpy of fusion at a given temperature
<b>hvap:</b>	Enthalpy of vaporization at standard conditions
<b>log10ws:</b>	Log10 of Water solubility in mol/l
<b>logp:</b>	Octanol/Water partition coefficient
<b>mcvol:</b>	McGowan's characteristic volume
<b>pc:</b>	Critical Pressure
<b>tb:</b>	Normal Boiling Point Temperature
<b>tc:</b>	Critical Temperature
<b>tf:</b>	Normal melting (fusion) point
<b>vc:</b>	Critical Volume

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