

# Benzene, (1-chloropentyl)

<b>Inchi:</b>	InChI=1S/C11H15Cl/c1-2-3-9-11(12)10-7-5-4-6-8-10/h4-8,11H,2-3,9H2,1H3
<b>InchiKey:</b>	OSOLHARQJQEMBZ-UHFFFAOYSA-N
<b>Formula:</b>	C11H15Cl
<b>SMILES:</b>	CCCCC(Cl)c1ccccc1
<b>Mol. weight [g/mol]:</b>	182.69

## Physical Properties

Property code	Value	Unit	Source
gf	139.78	kJ/mol	Joback Method
hf	-54.86	kJ/mol	Joback Method
hfus	18.96	kJ/mol	Joback Method
hvap	46.35	kJ/mol	Joback Method
log10ws	-4.14		Crippen Method
logp	4.157		Crippen Method
mcvol	154.330	ml/mol	McGowan Method
pc	2571.50	kPa	Joback Method
rinpola	1322.00		NIST Webbook
tb	514.75	K	Joback Method
tc	726.81	K	Joback Method
tf	255.07	K	Joback Method
vc	0.587	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	330.19	J/molxK	514.75	Joback Method
cpg	399.52	J/molxK	691.46	Joback Method
cpg	387.37	J/molxK	656.12	Joback Method
cpg	374.40	J/molxK	620.78	Joback Method
cpg	360.58	J/molxK	585.44	Joback Method
cpg	345.85	J/molxK	550.09	Joback Method
cpg	410.90	J/molxK	726.81	Joback Method
dvisc	0.0002142	Paxs	514.75	Joback Method
dvisc	0.0002843	Paxs	471.47	Joback Method

dvisc	0.0003997	Paxs	428.19	Joback Method
dvisc	0.0006066	Paxs	384.91	Joback Method
dvisc	0.0010234	Paxs	341.63	Joback Method
dvisc	0.0020092	Paxs	298.35	Joback Method
dvisc	0.0049594	Paxs	255.07	Joback Method

## Sources

<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>
<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=R131863&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R131863&amp;Units=SI</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>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>rinpol:</b>	Non-polar retention indices
<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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