

# 5-Bromovaleric acid, octyl ester

<b>Inchi:</b>	InChI=1S/C13H25BrO2/c1-2-3-4-5-6-9-12-16-13(15)10-7-8-11-14/h2-12H2,1H3
<b>InchiKey:</b>	WJGKHCBEUVDHNP-UHFFFAOYSA-N
<b>Formula:</b>	C13H25BrO2
<b>SMILES:</b>	CCCCCCCCOC(=O)CCCCBr
<b>Mol. weight [g/mol]:</b>	293.24
<b>CAS:</b>	13931-44-7

## Physical Properties

Property code	Value	Unit	Source
gf	-161.02	kJ/mol	Joback Method
hf	-530.12	kJ/mol	Joback Method
hfus	37.50	kJ/mol	Joback Method
hvap	60.12	kJ/mol	Joback Method
log10ws	-4.56		Crippen Method
logp	4.455		Crippen Method
mvol	218.970	ml/mol	McGowan Method
pc	1809.23	kPa	Joback Method
rinpol	1823.00		NIST Webbook
tb	639.29	K	Joback Method
tc	819.82	K	Joback Method
tf	368.23	K	Joback Method
vc	0.850	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	564.45	J/molxK	639.29	Joback Method
cpg	634.77	J/molxK	789.74	Joback Method
cpg	622.07	J/molxK	759.65	Joback Method
cpg	608.71	J/molxK	729.56	Joback Method
cpg	594.66	J/molxK	699.47	Joback Method
cpg	579.92	J/molxK	669.38	Joback Method
cpg	646.83	J/molxK	819.82	Joback Method
dvisc	0.0001536	Paxs	639.29	Joback Method

dvisc	0.0001993	Paxs	594.11	Joback Method
dvisc	0.0002698	Paxs	548.94	Joback Method
dvisc	0.0003857	Paxs	503.76	Joback Method
dvisc	0.0005916	Paxs	458.58	Joback Method
dvisc	0.0009964	Paxs	413.41	Joback Method
dvisc	0.0019071	Paxs	368.23	Joback Method

## Sources

<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=C13931447&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C13931447&amp;Units=SI</a>
<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci990307l">http://pubs.acs.org/doi/abs/10.1021/ci990307l</a>
<b>Crippen Method:</b>	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</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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