

# 3-Bromobenzoic acid, undec-2-enyl ester

<b>Inchi:</b>	InChI=1S/C18H25BrO2/c1-2-3-4-5-6-7-8-9-10-14-21-18(20)16-12-11-13-17(19)15-16/h9-
<b>InchiKey:</b>	ZKCZARNVCCJLSL-MDZDMXLPSA-N
<b>Formula:</b>	C18H25BrO2
<b>SMILES:</b>	CCCCCCCC=CCOC(=O)c1cccc(Br)c1
<b>Mol. weight [g/mol]:</b>	353.29

## Physical Properties

Property code	Value	Unit	Source
gf	64.08	kJ/mol	Joback Method
hf	-291.04	kJ/mol	Joback Method
hfus	44.30	kJ/mol	Joback Method
hvap	74.15	kJ/mol	Joback Method
log10ws	-6.92		Crippen Method
logp	5.913		Crippen Method
mcvol	261.360	ml/mol	McGowan Method
pc	1648.43	kPa	Joback Method
rmpol	2330.00		NIST Webbook
tb	789.51	K	Joback Method
tc	998.63	K	Joback Method
tf	458.44	K	Joback Method
vc	1.002	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	731.35	J/molxK	789.51	Joback Method
cpg	800.58	J/molxK	963.78	Joback Method
cpg	788.44	J/molxK	928.92	Joback Method
cpg	775.51	J/molxK	894.07	Joback Method
cpg	761.72	J/molxK	859.22	Joback Method
cpg	747.02	J/molxK	824.36	Joback Method
cpg	811.97	J/molxK	998.63	Joback Method
dvisc	0.0000696	Paxs	789.51	Joback Method
dvisc	0.0000892	Paxs	734.33	Joback Method

dvisc	0.0001192	Paxs	679.15	Joback Method
dvisc	0.0001676	Paxs	623.97	Joback Method
dvisc	0.0002517	Paxs	568.80	Joback Method
dvisc	0.0004127	Paxs	513.62	Joback Method
dvisc	0.0007619	Paxs	458.44	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=U299147&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U299147&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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