

# Benzene, (2-bromopropyl)-

<b>Other names:</b>	(2-Bromopropyl)benzene ./-.-2-Bromo-1-phenylpropane 1-Phenyl-2-bromopropane 2-Bromo-1-phenylpropane
<b>Inchi:</b>	InChI=1S/C9H11Br/c1-8(10)7-9-5-3-2-4-6-9/h2-6,8H,7H2,1H3
<b>InchiKey:</b>	NVYOCAOZCSNIHR-UHFFFAOYSA-N
<b>Formula:</b>	C9H11Br
<b>SMILES:</b>	CC(Br)Cc1ccccc1
<b>Mol. weight [g/mol]:</b>	199.09
<b>CAS:</b>	2114-39-8

## Physical Properties

Property code	Value	Unit	Source
gf	149.19	kJ/mol	Joback Method
hf	28.49	kJ/mol	Joback Method
hfus	14.87	kJ/mol	Joback Method
hvap	43.95	kJ/mol	Joback Method
log10ws	-3.24		Crippen Method
logp	3.013		Crippen Method
mcvol	131.410	ml/mol	McGowan Method
pc	3568.53	kPa	Joback Method
tb	497.72	K	Joback Method
tc	727.72	K	Joback Method
tf	262.41	K	Joback Method
vc	0.487	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	254.95	J/molxK	497.72	Joback Method
cpg	314.51	J/molxK	689.39	Joback Method
cpg	304.28	J/molxK	651.05	Joback Method
cpg	293.27	J/molxK	612.72	Joback Method
cpg	281.41	J/molxK	574.39	Joback Method

cpg	268.65	J/molxK	536.05	Joback Method
cpg	323.99	J/molxK	727.72	Joback Method
dvisc	0.0002656	Paxs	497.72	Joback Method
dvisc	0.0003445	Paxs	458.50	Joback Method
dvisc	0.0004692	Paxs	419.28	Joback Method
dvisc	0.0006810	Paxs	380.07	Joback Method
dvisc	0.0010769	Paxs	340.85	Joback Method
dvisc	0.0019186	Paxs	301.63	Joback Method
dvisc	0.0040622	Paxs	262.41	Joback Method

## Pressure Dependent Properties

Property code	Value	Unit	Pressure [kPa]	Source
tbrp	381.20	K	2.10	NIST Webbook

## Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/(T + C)$
Coeff. A	1.37293e+01
Coeff. B	-3.83321e+03
Coeff. C	-7.81910e+01
Temperature range (K), min.	363.36
Temperature range (K), max.	533.56

## 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=C2114398&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C2114398&amp;Units=SI</a>
<b>The Yaws Handbook of Vapor Pressure:</b>	<a href="https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure">https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure</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>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>pvap:</b>	Vapor pressure
<b>tb:</b>	Normal Boiling Point Temperature
<b>tbrp:</b>	Boiling point at reduced pressure
<b>tc:</b>	Critical Temperature
<b>tf:</b>	Normal melting (fusion) point
<b>vc:</b>	Critical Volume

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