## 9-Bromoanthracene

Other names: 9-Anthracenyl bromide

anthracene, 9-bromo-

InChl=1S/C14H9Br/c15-14-12-7-3-1-5-10(12)9-11-6-2-4-8-13(11)14/h1-9H

InchiKey: ZIRVQSRSPDUEOJ-UHFFFAOYSA-N

Formula: C14H9Br

SMILES: Brc1c2cccc2cc2cccc12

Mol. weight [g/mol]: 257.12 CAS: 1564-64-3

### **Physical Properties**

Property code	Value	Unit	Source
ea	0.61 ± 0.10	eV	NIST Webbook
gf	387.77	kJ/mol	Joback Method
hf	289.77	kJ/mol	Joback Method
hfus	24.60	kJ/mol	Joback Method
hvap	60.07	kJ/mol	Joback Method
ie	7.58	eV	NIST Webbook
ie	$7.48 \pm 0.03$	eV	NIST Webbook
log10ws	-6.24		Crippen Method
logp	4.755		Crippen Method
mcvol	162.940	ml/mol	McGowan Method
рс	3452.08	kPa	Joback Method
rinpol	2125.00		NIST Webbook
rinpol	2143.00		NIST Webbook
rinpol	2125.00		NIST Webbook
rinpol	2140.00		NIST Webbook
rinpol	2125.00		NIST Webbook
tb	660.48	K	Joback Method
tc	927.78	K	Joback Method
tf	374.15	К	Thermochemical and vapor pressure behavior of anthracene and brominated anthracene mixtures
VC	0.618	m3/kmol	Joback Method

# **Temperature Dependent Properties**

Property code	Value	Unit	Temperature [K]	Source
cpg	367.67	J/mol×K	660.48	Joback Method
cpg	380.45	J/mol×K	705.03	Joback Method
cpg	392.09	J/mol×K	749.58	Joback Method
cpg	402.76	J/mol×K	794.13	Joback Method
cpg	412.64	J/mol×K	838.68	Joback Method
cpg	421.90	J/mol×K	883.23	Joback Method
cpg	430.71	J/mol×K	927.78	Joback Method
dvisc	0.0014982	Paxs	424.20	Joback Method
dvisc	0.0011666	Paxs	463.58	Joback Method
dvisc	0.0009447	Paxs	502.96	Joback Method
dvisc	0.0007888	Paxs	542.34	Joback Method
dvisc	0.0006749	Paxs	581.72	Joback Method
dvisc	0.0005890	Paxs	621.10	Joback Method
dvisc	0.0005224	Paxs	660.48	Joback Method
hsubt	100.50 ± 1.80	kJ/mol	341.50	NIST Webbook
psub	1.87e-05	kPa	325.50	The effect of halogen hetero-atoms on the vapor pressures and thermodynamics of polycyclic aromatic compounds measured via the Knudsen effusion technique
psub	9.70e-06	kPa	319.00	The effect of halogen hetero-atoms on the vapor pressures and thermodynamics of polycyclic aromatic compounds measured via the Knudsen effusion technique

psub	1.16e-05	kPa	321.80	The effect of halogen hetero-atoms on the vapor pressures and thermodynamics of polycyclic aromatic compounds measured via the Knudsen effusion technique	
psub	1.57e-05	kPa	322.90	The effect of halogen hetero-atoms on the vapor pressures and thermodynamics of polycyclic aromatic compounds measured via the Knudsen effusion technique	
psub	6.00e-06	kPa	315.60	The effect of halogen hetero-atoms on the vapor pressures and thermodynamics of polycyclic aromatic compounds measured via the Knudsen effusion technique	
psub	2.95e-05	kPa	329.20	The effect of halogen hetero-atoms on the vapor pressures and thermodynamics of polycyclic aromatic compounds measured via the Knudsen effusion technique	
psub	3.40e-05	kPa	330.80	The effect of halogen hetero-atoms on the vapor pressures and thermodynamics of polycyclic aromatic compounds measured via the Knudsen effusion technique	

may de	4 700 05	I-De	222.40		
psub	4.79e-05	kPa	333.10	The effect of halogen hetero-atoms on the vapor pressures and thermodynamics of polycyclic aromatic compounds measured via the Knudsen effusion technique	
psub	6.28e-05	kPa	336.20	The effect of halogen hetero-atoms on the vapor pressures and thermodynamics of polycyclic aromatic compounds measured via the Knudsen effusion technique	
psub	9.05e-05	kPa	339.20	The effect of halogen hetero-atoms on the vapor pressures and thermodynamics of polycyclic aromatic compounds measured via the Knudsen effusion technique	
psub	9.04e-05	kPa	339.90	The effect of halogen hetero-atoms on the vapor pressures and thermodynamics of polycyclic aromatic compounds measured via the Knudsen effusion technique	
psub	9.19e-05	kPa	340.00	The effect of halogen hetero-atoms on the vapor pressures and thermodynamics of polycyclic aromatic compounds measured via the Knudsen effusion technique	

psub	1.02e-04	kPa	342.10	The effect of halogen hetero-atoms on the vapor pressures and thermodynamics of polycyclic aromatic compounds measured via the Knudsen effusion technique	
psub	1.16e-04	kPa	342.30	The effect of halogen hetero-atoms on the vapor pressures and thermodynamics of polycyclic aromatic compounds measured via the Knudsen effusion technique	
psub	1.36e-04	kPa	344.10	The effect of halogen hetero-atoms on the vapor pressures and thermodynamics of polycyclic aromatic compounds measured via the Knudsen effusion technique	
psub	1.41e-04	kPa	344.30	The effect of halogen hetero-atoms on the vapor pressures and thermodynamics of polycyclic aromatic compounds measured via the Knudsen effusion technique	
psub	2.08e-04	kPa	347.10	The effect of halogen hetero-atoms on the vapor pressures and thermodynamics of polycyclic aromatic compounds measured via the Knudsen effusion technique	

psub	2.29e-04	kPa	348.20	The effect of halogen hetero-atoms on the vapor pressures and thermodynamics of polycyclic aromatic compounds measured via the Knudsen effusion technique	
psub	3.53e-04	kPa	352.10	The effect of halogen hetero-atoms on the vapor pressures and thermodynamics of polycyclic aromatic compounds measured via the Knudsen effusion technique	
psub	4.50e-04	kPa	355.70	The effect of halogen hetero-atoms on the vapor pressures and thermodynamics of polycyclic aromatic compounds measured via the Knudsen effusion technique	
psub	8.30e-04	kPa	362.80	The effect of halogen hetero-atoms on the vapor pressures and thermodynamics of polycyclic aromatic compounds measured via the Knudsen effusion technique	
psub	9.66e-04	kPa	364.00	The effect of halogen hetero-atoms on the vapor pressures and thermodynamics of polycyclic aromatic compounds measured via the Knudsen effusion technique	

1.48e-03 kPa 367.80 psub

The effect of halogen hetero-atoms on the vapor pressures and thermodynamics of polycyclic aromatic compounds measured via the Knudsen effusion technique

#### Sources

McGowan Method: http://link.springer.com/article/10.1007/BF02311772

**NIST Webbook:** http://webbook.nist.gov/cgi/cbook.cgi?ID=C1564643&Units=SI

**Crippen Method:** http://pubs.acs.org/doi/abs/10.1021/ci990307l

**Crippen Method:** https://www.chemeo.com/doc/models/crippen\_log10ws

Thermochemical and vapor pressure behavior of anthracene and brominated Thankeeher halages: hetero-atoms on the vapor pressures and the vapor pressures are the vapor pressures and the vapor pressures and the vapor pressures and the vapor pressures and the vapor pressures are the va

https://www.doi.org/10.1016/j.jct.2007.09.006

https://www.doi.org/10.1016/j.fluid.2012.12.036

compounds measured via the Knudsen

effusion technique:

### Legend

cpg: Ideal gas heat capacity

dvisc: Dynamic viscosity ea: Electron affinity

Standard Gibbs free energy of formation gf: hf: Enthalpy of formation at standard conditions hfus: Enthalpy of fusion at standard conditions

hsubt: Enthalpy of sublimation at a given temperature hvap: Enthalpy of vaporization at standard conditions

ie: Ionization energy

Log10 of Water solubility in mol/l log10ws: Octanol/Water partition coefficient logp: McGowan's characteristic volume mcvol:

Critical Pressure pc:

Sublimation pressure psub:

rinpol: Non-polar retention indices

Normal Boiling Point Temperature tb:

tc: Critical Temperature tf: Normal melting (fusion) point

vc: Critical Volume

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