

# Perfluorooctyl bromide

<b>Other names:</b>	1-Bromoperfluorooctane 1-bromoheptadecafluorooctane
<b>Inchi:</b>	InChI=1S/C8BrF17/c9-7(22,23)5(18,19)3(14,15)1(10,11)2(12,13)4(16,17)6(20,21)8(24,25)
<b>InchiKey:</b>	WTWWXOGTJWMJHI-UHFFFAOYSA-N
<b>Formula:</b>	C8BrF17
<b>SMILES:</b>	FC(F)(F)C(F)(F)C(F)(F)C(F)(F)C(F)(F)C(F)(F)C(F)(F)C(F)(F)C(F)(F)Br
<b>Mol. weight [g/mol]:</b>	498.96
<b>CAS:</b>	423-55-2

## Physical Properties

Property code	Value	Unit	Source
gf	-3258.25	kJ/mol	Joback Method
hf	-3585.99	kJ/mol	Joback Method
hfus	14.81	kJ/mol	Joback Method
hvap	45.60 ± 0.40	kJ/mol	NIST Webbook
log10ws	-6.96		Crippen Method
logp	6.348		Crippen Method
mcvol	171.170	ml/mol	McGowan Method
pc	1562.28	kPa	Joback Method
tb	410.35	K	Joback Method
tc	536.18	K	Joback Method
tf	269.11	K	Joback Method
vc	0.763	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	443.99	J/molxK	431.32	Joback Method
cpg	496.22	J/molxK	536.18	Joback Method
cpg	487.48	J/molxK	515.21	Joback Method
cpg	477.92	J/molxK	494.24	Joback Method
cpg	467.52	J/molxK	473.26	Joback Method
cpg	456.22	J/molxK	452.29	Joback Method
cpg	430.77	J/molxK	410.35	Joback Method

hfust	12.13	kJ/mol	278.90	NIST Webbook
hfust	1.60	kJ/mol	146.40	NIST Webbook
hfust	12.13	kJ/mol	278.90	NIST Webbook
kvisc	0.0000011	m <sup>2</sup> /s	298.15	Viscosities of Liquid Fluorocompounds
kvisc	0.0000010	m <sup>2</sup> /s	303.15	Viscosities of Liquid Fluorocompounds
kvisc	0.0000009	m <sup>2</sup> /s	308.15	Viscosities of Liquid Fluorocompounds
kvisc	0.0000008	m <sup>2</sup> /s	313.15	Viscosities of Liquid Fluorocompounds
kvisc	0.0000008	m <sup>2</sup> /s	318.15	Viscosities of Liquid Fluorocompounds
pvap	0.54	kPa	293.20	Densities and Vapor Pressures of Highly Fluorinated Compounds
pvap	0.98	kPa	303.26	Densities and Vapor Pressures of Highly Fluorinated Compounds
pvap	1.30	kPa	308.22	Densities and Vapor Pressures of Highly Fluorinated Compounds
pvap	1.67	kPa	313.18	Densities and Vapor Pressures of Highly Fluorinated Compounds
pvap	2.24	kPa	318.09	Densities and Vapor Pressures of Highly Fluorinated Compounds
pvap	2.89	kPa	322.92	Densities and Vapor Pressures of Highly Fluorinated Compounds
pvap	3.69	kPa	327.88	Densities and Vapor Pressures of Highly Fluorinated Compounds
pvap	0.38	kPa	288.22	Densities and Vapor Pressures of Highly Fluorinated Compounds

pvap	4.70	kPa	332.78	Densities and Vapor Pressures of Highly Fluorinated Compounds
pvap	0.69	kPa	298.08	Densities and Vapor Pressures of Highly Fluorinated Compounds
rhoI	1917.00	kg/m3	292.57	Low pressure solubility and thermodynamics of solvation of oxygen, carbon dioxide, and carbon monoxide in fluorinated liquids
rhoI	1906.56	kg/m3	298.78	Low pressure solubility and thermodynamics of solvation of oxygen, carbon dioxide, and carbon monoxide in fluorinated liquids
rhoI	1892.19	kg/m3	302.71	Low pressure solubility and thermodynamics of solvation of oxygen, carbon dioxide, and carbon monoxide in fluorinated liquids
rhoI	1881.37	kg/m3	307.01	Low pressure solubility and thermodynamics of solvation of oxygen, carbon dioxide, and carbon monoxide in fluorinated liquids
rhoI	1869.53	kg/m3	311.72	Low pressure solubility and thermodynamics of solvation of oxygen, carbon dioxide, and carbon monoxide in fluorinated liquids

rho1	1926.65	kg/m <sup>3</sup>	288.63	Low pressure solubility and thermodynamics of solvation of oxygen, carbon dioxide, and carbon monoxide in fluorinated liquids
sfust	10.93	J/molxK	146.40	NIST Webbook
sfust	43.49	J/molxK	278.90	NIST Webbook
srf	0.02	N/m	318.55	Surface Tension of Liquid Fluorocompounds
srf	0.02	N/m	313.45	Surface Tension of Liquid Fluorocompounds
srf	0.02	N/m	308.45	Surface Tension of Liquid Fluorocompounds
srf	0.02	N/m	303.35	Surface Tension of Liquid Fluorocompounds
srf	0.02	N/m	298.35	Surface Tension of Liquid Fluorocompounds
srf	0.02	N/m	292.75	Surface Tension of Liquid Fluorocompounds
srf	0.02	N/m	287.15	Surface Tension of Liquid Fluorocompounds

## Sources

### Crippen Method:

[https://www.chemeo.com/doc/models/crippen\\_log10ws](https://www.chemeo.com/doc/models/crippen_log10ws)

### Solubility of oxygen in substituted

#### perfluorocarbons:

<https://www.doi.org/10.1016/j.fluid.2005.09.011>

#### Low pressure solubility and

#### thermodynamics of solvation of

#### oxygen, carbon dioxide, and carbon

#### monoxide in fluorinated liquids:

#### NIST Webbook:

<https://www.doi.org/10.1016/j.jct.2006.11.012>

<https://www.doi.org/10.1021/je060199g>

<http://webbook.nist.gov/cgi/cbook.cgi?ID=C423552&Units=SI>

### McGowan Method:

<http://link.springer.com/article/10.1007/BF02311772>

### Joback Method:

[https://en.wikipedia.org/wiki/Joback\\_method](https://en.wikipedia.org/wiki/Joback_method)

### Crippen Method:

<http://pubs.acs.org/doi/abs/10.1021/ci990307l>

### liquid-liquid equilibrium of substituted

### perfluoro-n-octane + n-octane systems:

<https://www.doi.org/10.1016/j.fluid.2008.04.011>

### Viscosities of Liquid

### Fluorocompounds:

<https://www.doi.org/10.1021/je700632z>

### Densities and Vapor Pressures of

### Highly Fluorinated Compounds:

<https://www.doi.org/10.1021/je050056e>

# Legend

<b>cpg:</b>	Ideal gas heat capacity
<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>hfust:</b>	Enthalpy of fusion at a given temperature
<b>hvac:</b>	Enthalpy of vaporization at standard conditions
<b>kvisc:</b>	Kinematic viscosity
<b>log10ws:</b>	Log10 of Water solubility in mol/l
<b>logp:</b>	Octanol/Water partition coefficient
<b>mccvol:</b>	McGowan's characteristic volume
<b>pc:</b>	Critical Pressure
<b>pvap:</b>	Vapor pressure
<b>rho:</b>	Liquid Density
<b>sfust:</b>	Entropy of fusion at a given temperature
<b>srf:</b>	Surface Tension
<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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