

Benzamide, 4-bromo-N-ethyl-N-hexyl-

Inchi:	InChI=1S/C15H22BrNO/c1-3-5-6-7-12-17(4-2)15(18)13-8-10-14(16)11-9-13/h8-11H,3-7,
InchiKey:	QSFXPXMLQVLUJR-UHFFFAOYSA-N
Formula:	C15H22BrNO
SMILES:	CCCCCN(CC)C(=O)c1ccc(Br)cc1
Mol. weight [g/mol]:	312.25

Physical Properties

Property code	Value	Unit	Source
gf	174.38	kJ/mol	Joback Method
hf	-146.59	kJ/mol	Joback Method
hfus	38.16	kJ/mol	Joback Method
hvap	67.15	kJ/mol	Joback Method
log10ws	-5.29		Crippen Method
logp	4.492		Crippen Method
mvol	227.500	ml/mol	McGowan Method
pc	2043.76	kPa	Joback Method
rmpol	2250.00		NIST Webbook
rmpol	2250.00		NIST Webbook
tb	706.73	K	Joback Method
tc	915.27	K	Joback Method
tf	439.95	K	Joback Method
vc	0.854	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	602.25	J/mol×K	706.73	Joback Method
cpg	617.96	J/mol×K	741.49	Joback Method
cpg	632.69	J/mol×K	776.24	Joback Method
cpg	646.49	J/mol×K	811.00	Joback Method
cpg	659.41	J/mol×K	845.75	Joback Method
cpg	671.52	J/mol×K	880.51	Joback Method
cpg	682.88	J/mol×K	915.27	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U415455&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Joback Method:	https://en.wikipedia.org/wiki/Joback_method
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772

Legend

cpg:	Ideal gas heat capacity
gf:	Standard Gibbs free energy of formation
hf:	Enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion at standard conditions
hvp:	Enthalpy of vaporization at standard conditions
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
pc:	Critical Pressure
rinp:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

Latest version available from:

<https://www.chemeo.com/cid/124-109-3/Benzamide-4-bromo-N-ethyl-N-hexyl.pdf>

Generated by Cheméo on 2024-04-28 18:38:30.438273284 +0000 UTC m=+16618759.358850597.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.