

# Benzenesulfonamide, N-butyl-

<b>Other names:</b>	Benzenesulfonic acid butyl amide N-Butylbenzenesulfonamide Plastomoll BMB N-(n-Butyl)benzenesulfonamide Dellatol BBS Plasthall BSA Uniplex 214 BM 4 (sulfonamide) Cetamol BMB NSC 3536 N-butylbenzenesulphonamide
<b>Inchi:</b>	InChI=1S/C10H15NO2S/c1-2-3-9-11-14(12,13)10-7-5-4-6-8-10/h4-8,11H,2-3,9H2,1H3
<b>InchiKey:</b>	IPRJXAGUEGOFGG-UHFFFAOYSA-N
<b>Formula:</b>	C10H15NO2S
<b>SMILES:</b>	CCCCNS(=O)(=O)c1ccccc1
<b>Mol. weight [g/mol]:</b>	213.30
<b>CAS:</b>	3622-84-2

## Physical Properties

Property code	Value	Unit	Source
gf	-233.42	kJ/mol	Joback Method
hf	-413.08	kJ/mol	Joback Method
hfus	32.17	kJ/mol	Joback Method
hvap	65.20	kJ/mol	Joback Method
log10ws	-2.60		Crippen Method
logp	1.765		Crippen Method
mcvol	166.070	ml/mol	McGowan Method
pc	3547.31	kPa	Joback Method
rinpol	1794.00		NIST Webbook
rinpol	1797.00		NIST Webbook
ripol	2956.00		NIST Webbook
tb	552.83	K	Joback Method
tc	752.37	K	Joback Method
tf	320.10	K	Joback Method
vc	0.648	m <sup>3</sup> /kmol	Joback Method

# Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	386.59	J/mol×K	552.83	Joback Method
cpg	401.71	J/mol×K	586.09	Joback Method
cpg	415.96	J/mol×K	619.34	Joback Method
cpg	429.35	J/mol×K	652.60	Joback Method
cpg	441.92	J/mol×K	685.86	Joback Method
cpg	453.66	J/mol×K	719.12	Joback Method
cpg	464.61	J/mol×K	752.37	Joback Method

## Sources

Joback Method:	<a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>
McGowan Method:	<a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>
NIST Webbook:	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C3622842&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C3622842&amp;Units=SI</a>
Crippen Method:	<a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</a>
Crippen Method:	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>

## 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>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>ripol:</b>	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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