

# 1-bromo-dibenzofuran

<b>Inchi:</b>	InChI=1S/C12H7BrO/c13-9-5-3-7-11-12(9)8-4-1-2-6-10(8)14-11/h1-7H
<b>InchiKey:</b>	WUYYYVOWEBMOELQ-UHFFFAOYSA-N
<b>Formula:</b>	C12H7BrO
<b>SMILES:</b>	BrC1CCCC2OC3CCCCC3C12
<b>Mol. weight [g/mol]:</b>	247.09

## Physical Properties

Property code	Value	Unit	Source
log10ws	-9.98		Crippen Method
logp	4.348		Crippen Method
mcvol	144.930	ml/mol	McGowan Method
rinsol	1837.00		NIST Webbook

## Sources

<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=R171574&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R171574&amp;Units=SI</a>
<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci990307I">http://pubs.acs.org/doi/abs/10.1021/ci990307I</a>
<b>Crippen Method:</b>	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>

## Legend

<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>rinsol:</b>	Non-polar retention indices

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