

# H CBr radical

**Inchi:** InChI=1S/CHBr/c1-2/h1H  
**InchiKey:** PRCFIJGWTCBHPP-UHFFFAOYSA-N  
**Formula:** CHBr  
**SMILES:** [CH]Br  
**Mol. weight [g/mol]:** 92.92  
**CAS:** 17141-28-5

## Physical Properties

Property code	Value	Unit	Source
ea	1.45 ± 0.01	eV	NIST Webbook
ea	1.56 ± 0.01	eV	NIST Webbook
log10ws	-0.73		Crippen Method
logp	1.050		Crippen Method
mcvol	38.150	ml/mol	McGowan Method

## Sources

**McGowan Method:** <http://link.springer.com/article/10.1007/BF02311772>  
**NIST Webbook:** <http://webbook.nist.gov/cgi/cbook.cgi?ID=C17141285&Units=SI>  
**Crippen Method:** <http://pubs.acs.org/doi/abs/10.1021/ci9903071>  
**Crippen Method:** [https://www.chemeo.com/doc/models/crippen\\_log10ws](https://www.chemeo.com/doc/models/crippen_log10ws)

## Legend

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

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<https://www.chemeo.com/cid/64-887-7/H CBr-radical.pdf>

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