

# holmium

Inchi:	InChI=1S/Ho
InchiKey:	KJZYNXUDTRRSPN-UHFFFAOYSA-N
Formula:	Ho
SMILES:	[Ho]
Mol. weight [g/mol]:	164.93
CAS:	7440-60-0

## Physical Properties

Property code	Value	Unit	Source
ie	6.02 ± 0.00	eV	NIST Webbook
ie	6.02	eV	NIST Webbook
ie	6.02 ± 0.00	eV	NIST Webbook
ie	6.02 ± 0.00	eV	NIST Webbook
ie	6.02 ± 0.00	eV	NIST Webbook
ie	6.02 ± 0.00	eV	NIST Webbook
ie	6.10 ± 0.60	eV	NIST Webbook
ie	6.00 ± 0.10	eV	NIST Webbook
ie	5.80 ± 0.20	eV	NIST Webbook
ie	5.89 ± 0.03	eV	NIST Webbook
ie	5.85 ± 0.10	eV	NIST Webbook
ie	6.02 ± 0.02	eV	NIST Webbook
ie	6.08 ± 0.09	eV	NIST Webbook
ie	6.19 ± 0.02	eV	NIST Webbook

## Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/(T + C)$
Coeff. A	1.33238e+01
Coeff. B	-2.35472e+04
Coeff. C	-2.68270e+02
Temperature range (K), min.	1432.15
Temperature range (K), max.	2973.15

# Sources

Thermodynamic stability of RNi<sub>2</sub> Laves phases: <https://www.doi.org/10.1016/j.jct.2013.05.044>  
Investigation in the variation of Gibbs energy of formation of RE<sub>6</sub>UO<sub>12</sub> (RE = La, Pr, Sm, Eu, Gd, Tb, Dy, Ho, Er, Y) <https://www.doi.org/10.1016/j.jct.2019.06.030>  
Thermodynamic stability of Ln<sub>2</sub>Co<sub>3</sub> (Ln = Dy, Ho, Er, Y) <https://www.doi.org/10.1016/j.tca.2008.09.023>  
(Ln = Dy, Ho, Er) by self-state electrochemical cells: <http://webbook.nist.gov/cgi/cbook.cgi?ID=C7440600&Units=SI>  
The Yaws Handbook of Vapor Pressure: <https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure>

# Legend

ie: Ionization energy  
pvap: Vapor pressure

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