

manganese

Other names:	manganese element
Inchi:	InChI=1S/Mn
InchiKey:	PWHULOQIROXLJO-UHFFFAOYSA-N
Formula:	Mn
SMILES:	[Mn]
Mol. weight [g/mol]:	54.94
CAS:	7439-96-5

Physical Properties

Property code	Value	Unit	Source
affp	797.30	kJ/mol	NIST Webbook
basg	774.40	kJ/mol	NIST Webbook
ie	7.43 ± 0.00	eV	NIST Webbook
ie	7.43 ± 0.00	eV	NIST Webbook
ie	7.43	eV	NIST Webbook
ie	7.43 ± 0.00	eV	NIST Webbook
ie	7.43	eV	NIST Webbook
ie	7.43	eV	NIST Webbook
ie	7.43	eV	NIST Webbook

Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/(T + C)$
Coeff. A	1.52661e+01
Coeff. B	-2.26890e+04
Coeff. C	-2.04920e+02
Temperature range (K), min.	1228.15
Temperature range (K), max.	2333.15

Sources

Crystal structure and thermochemical properties of a novel coordination compound of manganese(II) with Mn_2O_4 (less than x less than 0.1) spinel phase: calorimetric method for determining the thermochemical energy storage capacities of redox metal oxides:
NIST Webbook:

<https://www.doi.org/10.1016/j.jct.2014.03.009>

The Yaws Handbook of Vapor Pressure: High temperature calorimetric examination of enthalpies of mixing in liquid (gadolinium germanium manganese) alloys:

<https://www.doi.org/10.1016/j.tca.2014.11.003>

<https://www.doi.org/10.1016/j.tca.2019.01.008>

<http://webbook.nist.gov/cgi/cbook.cgi?ID=C7439965&Units=SI>

<https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure>

<https://www.doi.org/10.1016/j.jct.2005.09.002>

Legend

affp:	Proton affinity
basg:	Gas basicity
ie:	Ionization energy
pvap:	Vapor pressure

Latest version available from:

<https://www.cheméo.com/cid/29-517-6/manganese.pdf>

Generated by Cheméo on 2024-05-15 07:47:49.779822928 +0000 UTC m=+18048518.700400240.

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