



OPEN ACCESS

EDITED AND REVIEWED BY
Taishi Kobayashi,
Kyoto University, Japan

*CORRESPONDENCE
F. Bok,
✉ f.bok@hzdr.de

RECEIVED 11 September 2023
ACCEPTED 02 October 2023
PUBLISHED 17 October 2023

CITATION
Bok F, Moog HC and Brendler V (2023),
Corrigendum: The solubility of oxygen in
water and saline solutions.
Front. Nucl. Eng. 2:1292254.
doi: 10.3389/fnuen.2023.1292254

COPYRIGHT
© 2023 Bok, Moog and Brendler. This is
an open-access article distributed under
the terms of the [Creative Commons
Attribution License \(CC BY\)](#). The use,
distribution or reproduction in other
forums is permitted, provided the original
author(s) and the copyright owner(s) are
credited and that the original publication
in this journal is cited, in accordance with
accepted academic practice. No use,
distribution or reproduction is permitted
which does not comply with these terms.

Corrigendum: The solubility of oxygen in water and saline solutions

F. Bok^{1*}, H. C. Moog² and V. Brendler¹

¹Actinide Thermodynamics Department, Institute of Resource Ecology, Helmholtz-Zentrum Dresden-Rossendorf e.V., Dresden, Germany, ²Department Repository Research, Gesellschaft für Anlagen- und Reaktorsicherheit (GRS) gGmbH, Braunschweig, Germany

KEYWORDS

oxygen solubility, electrolyte solutions, water, Pitzer ion-interaction approach, E_H equation

A Corrigendum on The solubility of oxygen in water and saline solutions

by Bok F, Moog HC and Brendler V (2023). *Front. Nucl. Eng.* 2:1158109. doi: 10.3389/fnuen.2023.1158109

In the published article, there was an error in [Table 5](#). In line 3 [without table head, $O_2(aq)$] in the T_{min}/T_{max} column the stated temperature range is wrong. Since no temperature coefficients for $C_p^0(T)$ of $O_2(aq)$ can be given, this field must stay empty. The table's footnote has been adjusted accordingly. The corrected [Table 5](#) and its caption *Standard formation data for $H_2O(l)$, $O_2(g)$, $H_2(g)$, and $O_2(aq)$. Temperature parameters for $C_p^0(T)$ with regard to the general function ([Eq. 12](#)) were fitted to values from [Chase \(1998\)](#).* appears below.

In the published article, there was an error in [Table 9](#). In line 7 [without table head, $O_2(aq)-OH^-$] in the last column (X_2) it has to be 3.934 ± 0.882 , not 3.6934 ± 0.882 . In line 15 [without table head, $O_2(aq)-K^+-SO_4^{2-}$] in the second column (X_0) it has to be -0.1618 ± 0.078 not -0.1575 ± 0.0688 and in the third column (X_1) it has to be 164.7 ± 102 not 183.5 ± 70.3 . The corrected [Table 9](#) and its caption *Parameters for the temperature dependency function of the binary and ternary interaction coefficients (λ , ζ). The uncertainty information refers to one standard deviation.* appears below.

The authors apologize for these errors and state that this does not change the scientific conclusions of the article in any way. The original article has been updated.

TABLE 5 Standard formation data for H₂O(l), O₂(g), H₂(g), and O₂(aq). Temperature parameters for $C_p^0(T)$ with regard to the general function (Eq. 12) were fitted to values from Chase (1998).

Species	$\Delta_f H_{i,T=T_0}^0$	References	$S_{i,T=T_0}^0$	References	$\Delta_f G_{i,T=T_0}^0$	References	$C_p^0(T)$ (J/mol K)				T_{\min}/T_{\max}	$C_{p,T=T_0}^0$	References
	(J/mol)		(J/mol K)		(J/mol)		A_1	A_2	A_5	A_6	(K)	(J/mol K)	
H ₂ O(l)	-285,830 ± 40 (2)	Cox et al. (1989)/ Chase (1998)	69.95 ± 0.03 (2)	Cox et al. (1989)/ Chase (1998)	-237,140 ± 41	(1)	149 ± 11	-0.33 ± 0.04	-1,056,714 ± 250,728	0.00042 ± 0.00004	280/500	75.418	Chase (1998)
O ₂ (g)	0	(By definition)	205.152 ± 0.005 (2)	Cox et al. (1989)	0	(By definition)	24.64 ± 0.03	0.0121 ± 0.0001	100,230 ± 1,433	0	298.15/500	29.376	Chase (1998)
O ₂ (aq)	-12,411 ± 971	(This work)	108 ± 14	(This work)	16,593 ± 3,168	(This work)	(3)	(3)	(3)	(3)	(3)	249.90 ± 0.07	(This work)
H ₂ (g)	0	(By definition)	130.68 ± 0.003 (2)	Cox et al. (1989)/ Chase (1998)	0	(By definition)	33.6 ± 0.2	-0.012 ± 0.001	-174,946 ± 3,652	1.01 ± 0.05 × 10 ⁻⁵	298.15/500	28.836	Chase (1998)

(1) Internally calculated. (2) Data uncertainty adopted from Cox et al. (1989). (3) With our selection for the temperature dependence of Henry's law constant (Eq. 8 and Table 2) no temperature function for the standard molar heat capacity of reaction $\Delta_r C_{p,m}^0(T)$ (6) can be derived. Hence, no temperature dependence for the standard molar heat capacity for O₂(aq) can be given.

TABLE 9 Parameters for the temperature dependency function of the binary and ternary interaction coefficients (λ , ζ). The uncertainty information refers to one standard deviation.

λ/ζ (Species: 1-2-3)	X_0	X_1	X_2
O ₂ (aq)-Cl ⁻	0 (by definition)	0 (by definition)	0 (by definition)
O ₂ (aq)-H ⁺	0.02598 ± 0.00171	4,941 ± 533	16.4 ± 1.8
O ₂ (aq)-Na ⁺	0.1315 ± 0.0093	2,897 ± 274	9.614 ± 0.947
O ₂ (aq)-K ⁺	0.135 ± 0.010	-1,004 ± 527	-3.513 ± 1.79
O ₂ (aq)-Mg ²⁺	0.2293 ± 0.0032	2,981 ± 860	9.9 ± 2.9
O ₂ (aq)-Ca ²⁺	0.2519 ± 0.0063	2,821 ± 651	9.449 ± 2.26
O ₂ (aq)-OH ⁻	0.06785 ± 0.00965	1,230 ± 280	3.934 ± 0.882
O ₂ (aq)-SO ₄ ²⁻	0.1334 ± 0.0082	-6,654 ± 1,150	-22.01 ± 3.38
O ₂ (aq)-HSO ₄ ⁻	0.03842 ± 0.00201	-45.38 ± 27.7	0 (fixed)
O ₂ (aq)-H ⁺ -HSO ₄ ⁻	-0.002472 ± 0.000302	0 (fixed)	0 (fixed)
O ₂ (aq)-Na ⁺ -Cl ⁻	-0.003767 ± 0.00579	-1,075 ± 128	-3.663 ± 0.442
O ₂ (aq)-Na ⁺ -OH ⁻	0 (fixed)	-40.7 ± 17.9	0 (fixed)
O ₂ (aq)-Na ⁺ -SO ₄ ²⁻	-0.0381 ± 0.0073	4,216 ± 1,820	13.49 ± 5.92
O ₂ (aq)-K ⁺ -Cl ⁻	-0.01711 ± 0.00647	747.7 ± 319	2.464 ± 1.07
O ₂ (aq)-K ⁺ -SO ₄ ²⁻	-0.1618 ± 0.078	164.7 ± 102	0 (fixed)
O ₂ (aq)-Mg ²⁺ -Cl ⁻	-0.006612 ± 0.00101	-875.5 ± 351	-2.974 ± 1.18
O ₂ (aq)-Mg ²⁺ -SO ₄ ²⁻	-0.05115 ± 0.00586	5,181 ± 550	16.86 ± 1.86
O ₂ (aq)-Ca ²⁺ -Cl ⁻	-0.01269 ± 0.00291	24.49 ± 1.81	0 (fixed)

Publisher's note

All claims expressed in this article are solely those of the authors and do not necessarily represent those of their affiliated

organizations, or those of the publisher, the editors and the reviewers. Any product that may be evaluated in this article, or claim that may be made by its manufacturer, is not guaranteed or endorsed by the publisher.