## National Institute of Advanced Industrial Science and Technology

# National Metrology Institute of Japan



Reference Material Certificate

NMIJ CRM 7405-b No. +++



Trace Elements and Arsenic Compounds in Seaweed (Hijiki)

This certified reference material (CRM) is produced in accordance with the NMIJ's management system and is in compliance with ISO 17034 and ISO/IEC 17025. This CRM is intended for the calibration of instruments, and validation of analytical methods and instruments used for the quantification of arsenic compounds and trace elements in hijiki and other seaweeds.

## **Certified Values**

The certified values for arsenic compounds, including arsenate and arsenosugar compounds, and five elements (As, Cd, Cu, Mn and Zn) in this CRM, are given in the tables below. The values are expressed as a mass fraction, based on a dry mass (the drying procedure is given in this certificate). The uncertainty of the certified value is the expanded uncertainty obtained by multiplying the combined standard uncertainty by a coverage factor (k) of 2, and it is the half-width of an interval estimated to have a level of confidence of approximately 95 %.

Compound	Certified value,	Expanded uncertainty,	Analytical methods
Compound	Mass fraction (mg/kg, as As)	Mass fraction (mg/kg, as As)	(see below)
Arsenic acid	24.4	0.7	1), 2), 3), 4)
[As(V)]	24.4	0.7	
Arsenosugar-408*			1), 5), 6), 7), 8)
(Arsenosugar-	1 4 1	0.04	
SO <sub>4</sub> , Molecular	1.41	0.04	
weight 408)			
Arsenosugar-328*			1), 5), 6), 7), 9)
(Arsenosugar-OH,		0.03	
Molecular weight	0.44	0.02	
328)			

\*IUPAC and common names and structural formulas of the arsenosugar compound are described in the Technical Information below.

Analytical methods: High-performance liquid chromatography-inductively coupled plasma mass spectrometry (HPLC-ICP-MS)

Measurement condition: [Extraction method] / [Extraction solvent] / [Extraction temperature] / [Column] / [Mobile phase] / [Calibrant]

- 1) Ultrasonication / Water / Room temperature / Reversed-phase  $C_{18}$  / 10 mmol L<sup>-1</sup> sodium 1-butanesulfonate-4 mmol L<sup>-1</sup> malonic acid-4 mmol L<sup>-1</sup> tetramethylammonium hydroxide-0.05 % methanol (pH 3.0) / As(V)
- 2) Ultrasonication / Water / Room temperature / Reversed-phase C<sub>30</sub> / 10 mmol L<sup>-1</sup> sodium 1-butanesulfonate-4 mmol L<sup>-1</sup> malonic acid-4 mmol L<sup>-1</sup> tetramethylammonium hydroxide-2 mmol L<sup>-1</sup> ammonium dihydrogen phosphate-0.05 % methanol (pH 2.7) / As(V)
- 3) Dry block bath / 0.07 mol L<sup>-1</sup> HCl+ 0.1 g L<sup>-1</sup> pepsin / 37 °C / Reversed-phase  $C_{28}$  / 10 mmol L<sup>-1</sup> sodium 1-butanesulfonate-4 mmol L<sup>-1</sup> malonic acid-4 mmol L<sup>-1</sup> tetramethylammonium hydroxide-0.05 % methanol (pH 3.0)/As(V)
- Ultrasonication / Water / Room temperature / Anion-exchange / 10 mmol L<sup>-1</sup> ammonium dihydrogen phosphate- -0.05 % methanol (pH 8.25) / As(V)

- 5) Dry block bath / 0.07 mol L<sup>-1</sup> HCl+ 0.1 g L<sup>-1</sup> pepsin / 37 °C / Reversed-phase C<sub>28</sub> / 10 mmol L<sup>-1</sup> sodium 1butanesulfonate-4 mmol L<sup>-1</sup> malonic acid-4 mmol L<sup>-1</sup> tetramethylammonium hydroxide-2 mmol L<sup>-1</sup> ammonium dihydrogen phosphate-0.05 % methanol (pH 2.7) / As(V)
- $\label{eq:constraint} \begin{array}{l} \mbox{ 6) } & \mbox{ Ultrasonication / Water / Room temperature / Reversed-phase $C_{18}$ / 10 mmol $L^{-1}$ ammonium dihydrogen phosphate $0.05 \%$ methanol (pH 8.25) / As(V) \\ \end{array}$
- 7) Dry block bath / 0.07 mol L<sup>-1</sup> HCl+ 0.1 g L<sup>-1</sup> pepsin / 37 °C / Reversed-phase C<sub>28</sub> / 10 mmol L<sup>-1</sup> sodium 1butanesulfonate-4 mmol L<sup>-1</sup> malonic acid-4 mmol L<sup>-1</sup> tetramethylammonium hydroxide-2 mmol L<sup>-1</sup> ammonium dihydrogen phosphate-0.05 % methanol (pH 2.7) / Arsenosugar-408 and -328
- 8) Rotating wheel / Water / Room temperature / Anion-exchange / 5 mmol L<sup>-1</sup> formic acid (pH 4.0) / DMA
- 9) Rotating wheel / Water / Room temperature / Reversed-phase  $C_{18}$  / 5 mmol L<sup>-1</sup> acetic acid (pH 5.0) / DMA

Element	Certified value,	Expanded uncertainty,	Analytical methods
	Mass fraction (mg/kg)	Mass fraction (mg/kg)	(see bel <mark>ow</mark> )
As	49.5	1.0	1, 3, 4, 5
Cd	1.25	0.04	1, 2, 4, 5
Cu	4.48	0.12	1, 2, 4, 5
Mn	22.6	0.5	1, 3, 4, 5
Zn	13.6	0.5	1, 2, 4, 5

Analytical methods:

- 1 Inductively coupled plasma mass spectrometry (ICP-MS)
- 2 Isotope dilution-inductively coupled plasma mass spectrometry (ID-ICP-MS)
- 3 High resolution inductively coupled plasma mass spectrometry
- 4 Inductively coupled plasma optical emission spectrometry (ICP-OES)
- 5 Graphite furnace atomic absorption spectrometry

The sample digestion method for 2 was microwave digestion with nitric acid, hydrofluoric acid, and perchloric acid. The digestion method for all others was microwave digestion with nitric acid, hydrofluoric acid, and hydrogen peroxide.

#### Analysis

These certified values are the weighted means of the results from two or more analytical methods conducted at NMIJ. Quantitative analysis of arsenic compounds was made after extraction by HPLC-ICP-MS. The different analytical methods were used, with combinations of different heating methods, extracts, and types of HPLC columns. The quantitative analysis of elements was made by the aforementioned analytical methods of 1 to 5, and combinations of these are based on: (1) a single primary method (ID-ICP-MS) with one or more reference methods or (2) three or more reference methods.

The expanded uncertainty in each certified value is equal to  $U = ku_c$ , where  $u_c$  is the combined standard uncertainty derived from: (a) the analytical results, (b) the method-to-method variance, (c) the dry mass correction, (d) the concentration of a standard solution, and (e) the sample homogeneity.

## **Metrological Traceability**

The certified values were determined by isotope dilution mass spectrometry, or other accurate methods, with NMIJ CRM 7912a Arsenate [As(V)] Solution and JCSS (Japan Calibration Service System) standard solutions. All sample preparation was carried out by a gravimetric method, using a balance calibrated by JCSS. Therefore, the certified values are traceable to the International System of Units (SI).

## Mutual Recognition Arrangement under Meter Convention

The certified values of elements and As(V) of this CRM are recognized for international equivalence based on the Mutual Recognition Arrangement under the Metre Convention (CIPM MRA). The calibration measurement capabilities (CMC) of NMIJ related to this CRM are registered in the Key Comparison Database (KCDB) (see https://www.bipm.org/kcdb/) of the International Bureau of Weights and Measures (BIPM).

#### **Expiration of Certification**

This certificate is valid for one year from the date of shipment, provided that the material is stored in accordance with the instructions given in this certificate.

#### **Description of the Material**

This CRM was prepared from hijiki seaweed that was powdered by freeze-pulverization. The CRM is in the form of brown powder, which is placed into amber glass bottles (approx 20 g in each).

#### Homogeneity

The homogeneity of the CRM was determined by analyzing 10 bottles hierarchical-randomly sampled from 640 bottles. Each arsenic compound was determined by HPLC-ICP-MS after extraction procedure. Each trace element was determined by ICP-MS after microwave acid digestion with nitric acid, hydrofluoric acid, and perchloric acid. The inhomogeneity of the analytes, which was evaluated by ANOVA, is not significant and is reflected in the uncertainty of the certified value. This material is homogeneous within the range of the uncertainty of the certified value.

## **Instructions for Storage**

This CRM should be stored at a temperature between 5 °C and 35 °C and shielded from light.

#### Instructions for Use

(1) This CRM should be opened and used up as soon as possible after opening to prevent contamination. When the bottle is stored after opening, it should be sealed with tape and kept in a desiccator with silica gel to limit its absorption of moisture as much as possible.

(2) Dry mass correction is required when the CRM is analyzed, as each certified value is expressed as a mass fraction based on a dry mass. The correction factor should be obtained by the following procedure. Do not use the sample that is used for the correction for analysis.

- 1) Take 1.0 g of the CRM in a weighing glass vessel.
- 2) Dry the CRM in the vessel at 105 °C for 5 h in a drying oven.
- 3) Weigh the CRM with the vessel after cooling in a desiccator with silica gel for 30 min.
- 4) The difference in the masses before and after drying is assumed to be the moisture content.

At April 2019, the dry mass correction factor of the certified values was *ca*. 7.6 % (mass fraction).

- Care should be taken to address the following points when the CRM is weighed.
- 1) Do not weigh in conditions of high humidity (over 60 %).
- 2) Weighing needs to be performed as quickly as possible.
- 3) Do not leave the bottle open when not in use, in order to keep the time the CRM is weathered to a minimum.
- 4) Weighing for dry mass correction has to be done in parallel with weighing for analysis.
- (4) From the viewpoint of homogeneity, more than 0.5 g of CRM should be used for each analysis.

## **Precautions for Handling**

Wear a mask, gloves, and other protective equipment during handling. Refer to the safety data sheet (SDS) on this CRM before use.

## Preparation

(3)

Approximately 22 kg of hijiki seaweed was obtained from the sea near Japan and used for preparation of the CRM after polishing. The hijiki seaweed was dried at 60 °C, and then freeze pulverized. The hijiki seaweed was again dried at 60 °C, and then placed into amber glass bottles (about 20 g in each) using a splitting method. The bottles were individually vacuum sealed into seal bags (Lamizip Aluminum), and were sterilized with <sup>60</sup>Co  $\gamma$ -ray irradiation (about 20 kGy). The preparation of the candidate material and  $\gamma$ -ray irradiation were performed by KANSO TECHNOS CO., LTD. and Radiation Application Development Association, respectively.

#### **Technical Information**

The CRM provides information, mass fractions for 3 kinds of arsenic compounds and 13 kinds of elements obtained in the certification process (June 2019). Dimethylarsinic acid (DMA), arsenosugar-482 and -392 were determined by HPLC-ICP-MS based on NMIJ CRM 7912-a As (V). The element concentrations were determined by calibration curve method using JCSS standard solution. Ba, Co, Cr, Ni, P, Pb and Sr were quantified by ICP-MS, and Al, Ca, Fe, K, Mg and Na were quantified by ICP-OES.

Compound		Mass fraction (mg/kg) (as As)			
DMA			0.24		
Arsenosugar-482				0.20	
(Arsenosugar-PO <sub>4</sub> , Molecular weight 482)		)	0.20		
Arsenosugar-392				0.16	
(Arsenosugar-SO <sub>3</sub> , Molecular weight 392)		)		0.10	
Element	Mass fraction (mg/kg)	El	ement	Mass fraction (mg/kg)	
Al	310		Mg	8000	
Ba	17		Na	9000	
Ca	18000		Ni	3	
Со	1.9		Р	780	
Cr	5.5		Pb	0.2	
Fe	210		Sr	1600	
K	36000				

The name, chemical structure and the molecular weight of arsenosugar compounds determined in the CRM are as follows.

Names of Arsenosugars in NMIJ 7405-b					
Abbreviation		Name			
	IUPAC	2-(2,3-dihydroxypropoxy)-5-(dimethylarsorylmethyl)oxolane- 3,4-diol			
Arsenosugar-328 Arsenosugar-OH	Common IUPAC	3-[5'-deoxy-5'-(dimethylarsinoyl)-β-			
		ribofuranosyloxy]propylene glyco			
		3-[5-(dimethylarsorylmethyl)-3,4-dihdroxyoxolan-2-yl]oxy-2-			
A man a man 202 A man and then SO		hydroxypropane-1-sulfonic acid			
Aischosugar-392 Aischosugar-303	J <sub>3</sub>	3-[5'-deoxy-5'-(dimethylarsinoyl)-β-ribofuranosyloxy]-2-			
	Common	hydroxypropanesulfonic acid			
	IUPAC	[3-[5-(dimethylarsorylmethyl)-3,4-dihydroxyoxolan-2-yl]oxy-			
Arcenosugar 408 Arcenosugar SO		2-hydroxypropyl] hydrogen sulfate			
Alschosugar-106	Common	3-[5'-deoxy-5'-(dimethylarsinoyl)-β-ribofuranosyloxy]-2-			
		hydroxypropyl hydrogen sulfate			
	IUPAC	2,3-dihydroxypropyl [3-[5-(dimethylarsorylmethyl)-3,4-			
		dihydroxyoxolan-2-yl]oxy-2-hydroxypropyl]hydrogen			
Arsenosugar-482 Arsenosugar-PO <sub>4</sub>		phosphate			
	Common	3-[5'-deoxy-5'-(dimethylarsinoyl)-β-ribofuranosyloxy]-2-			
		hydroxypropyl 2,3-dihydroxypropyl hydrogen phosphate			



Arsenosugar-328 (As-Sugar-328): R=OH Arsenosugar-392 (As-Sugar-392): R=SO<sub>3</sub>H Arsenosugar-408 (As-Sugar-408): R=OSO<sub>3</sub>H Arsenosugar-482 (As-Sugar-482): R=OPO<sub>2</sub>(OH)CH<sub>2</sub>CH(OH)CH<sub>2</sub>OH

#### Structural formula of arsenosugars in NMIJ 7405-b

R

#### **NMIJ Analysts**

The technical manager is INAGAKI K., the production manager is NARUKAWA T., and the analysts are ZHU Y., MIYASHITA S., ARIGA T., KOGUCHI M. and KUDO I.

#### Information

If substantive technical changes occur that affect the certification before the expiration of this certificate, NMIJ will notify the registered customer. Customer registration on the NMIJ Website (given below) will facilitate notification. Technical reports regarding this CRM can be obtained from the contact details given below.

#### **Reproduction of Certificate**

In reproducing this certificate, it should be clearly indicated that the document is a copy.

#### Note

For the value assignment of the mass fractions of arsenosugar compounds in this CRM, the inter-laboratory study based on NMIJ-University of Graz Joint Research, "Analysis of arsenosugar compounds in Hijiki seaweed powder (2017-2018)", was performed together with the analysis by NMIJ. The certification of arsenosugar-328 and -408 were calculated from the analytical results performed by Dr. Georg Raber of Graz University and NMIJ.

April 1, 2020

ISHIMURA Kazuhiko President National Institute of Advanced Industrial Science and Technology

If you have any questions about this CRM, please contact: National Institute of Advanced Industrial Science and Technology, National Metrology Institute of Japan, Center for Quality Management of Metrology, Reference Materials Office, 1-1-1 Umezono, Tsukuba, Ibaraki 305-8563, Japan Phone: +81-29-861-4059, https://unit.aist.go.jp/nmij/english/refmate/

#### Revision history

November 15, 2022: Mass fraction of Mg in "Technical Information" had been revised.